

Package ‘tfprobability’

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Title Interface to 'TensorFlow Probability'

Version 0.15.0

Description Interface to 'TensorFlow Probability', a 'Python' library built on 'TensorFlow' that makes it easy to combine probabilistic models and deep learning on modern hardware ('TPU', 'GPU').
'TensorFlow Probability' includes a wide selection of probability distributions and bijectors, probabilistic layers, variational inference, Markov chain Monte Carlo, and optimizers such as Nelder-Mead, BFGS, and SGLD.

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URL <https://github.com/rstudio/tfprobability>

BugReports <https://github.com/rstudio/tfprobability/issues>

SystemRequirements TensorFlow Probability
(<https://www.tensorflow.org/probability>)

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glm_families	<i>GLM families</i>
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Description

A list of models that can be used as the `model` argument in `glm_fit()`:

Details

- Bernoulli: `Bernoulli(probs=mean)` where `mean = sigmoid(matmul(X, weights))`
- BernoulliNormalCDF: `Bernoulli(probs=mean)` where `mean = Normal(0, 1).cdf(matmul(X, weights))`
- GammaExp: `Gamma(concentration=1, rate=1 / mean)` where `mean = exp(matmul(X, weights))`
- GammaSoftplus: `Gamma(concentration=1, rate=1 / mean)` where `mean = softplus(matmul(X, weights))`
- LogNormal: `LogNormal(loc=log(mean) - log(2) / 2, scale=sqrt(log(2)))` where `mean = exp(matmul(X, weights))`.
- LogNormalSoftplus: `LogNormal(loc=log(mean) - log(2) / 2, scale=sqrt(log(2)))` where `mean = softplus(matmul(X, weights))`
- Normal: `Normal(loc=mean, scale=1)` where `mean = matmul(X, weights)`.
- NormalReciprocal: `Normal(loc=mean, scale=1)` where `mean = 1 / matmul(X, weights)`
- Poisson: `Poisson(rate=mean)` where `mean = exp(matmul(X, weights))`.
- PoissonSoftplus: `Poisson(rate=mean)` where `mean = softplus(matmul(X, weights))`.

Value

list of models that can be used as the `model` argument in `glm_fit()`

See Also

Other `glm_fit`: `glm_fit.tensorflow.tensor()`, `glm_fit_one_step.tensorflow.tensor()`

glm_fit	<i>Runs multiple Fisher scoring steps</i>
---------	---

Description

Runs multiple Fisher scoring steps

Usage

```
glm_fit(x, ...)
```

Arguments

x	float-like, matrix-shaped Tensor where each row represents a sample's features.
...	other arguments passed to specific methods.

Value

A glm_fit object with parameter estimates, number of iterations, etc.

See Also

[glm_fit.tensorflow.tensor\(\)](#)

glm_fit.tensorflow.tensor	<i>Runs multiple Fisher scoring steps</i>
---------------------------	---

Description

Runs multiple Fisher scoring steps

Usage

```
## S3 method for class 'tensorflow.tensor'  
glm_fit(  
  x,  
  response,  
  model,  
  model_coefficients_start = NULL,  
  predicted_linear_response_start = NULL,  
  l2_regularizer = NULL,  
  dispersion = NULL,  
  offset = NULL,  
  convergence_criteria_fn = NULL,  
)
```

```

    learning_rate = NULL,
    fast_unsafe_numerics = TRUE,
    maximum_iterations = NULL,
    name = NULL,
    ...
)

```

Arguments

<code>x</code>	float-like, matrix-shaped Tensor where each row represents a sample's features.
<code>response</code>	vector-shaped Tensor where each element represents a sample's observed response (to the corresponding row of features). Must have same dtype as <code>x</code> .
<code>model</code>	a string naming the model (see glm_families) or a <code>tfp\$glm\$ExponentialFamily</code> -like instance which implicitly characterizes a negative log-likelihood loss by specifying the distribution's mean, <code>gradient_mean</code> , and variance.
<code>model_coefficients_start</code>	Optional (batch of) vector-shaped Tensor representing the initial model coefficients, one for each column in <code>x</code> . Must have same dtype as <code>model_matrix</code> . Default value: Zeros.
<code>predicted_linear_response_start</code>	Optional Tensor with shape, dtype matching <code>response</code> ; represents offset shifted initial linear predictions based on <code>model_coefficients_start</code> . Default value: offset if <code>model_coefficients</code> is NULL, and <code>tf\$linalg\$matvec(x, model_coefficients_start) + offset</code> otherwise.
<code>l2_regularizer</code>	Optional scalar Tensor representing L2 regularization penalty. Default: NULL ie. no regularization.
<code>dispersion</code>	Optional (batch of) Tensor representing response dispersion.
<code>offset</code>	Optional Tensor representing constant shift applied to <code>predicted_linear_response</code> .
<code>convergence_criteria_fn</code>	callable taking: <code>is_converged_previous</code> , <code>iter_</code> , <code>model_coefficients_previous</code> , <code>predicted_linear_response_previous</code> , <code>model_coefficients_next</code> , <code>predicted_linear_response</code> , <code>response</code> , <code>model</code> , <code>dispersion</code> and returning a logical Tensor indicating that Fisher scoring has converged.
<code>learning_rate</code>	Optional (batch of) scalar Tensor used to dampen iterative progress. Typically only needed if optimization diverges, should be no larger than 1 and typically very close to 1. Default value: NULL (i.e., 1).
<code>fast_unsafe_numerics</code>	Optional Python bool indicating if faster, less numerically accurate methods can be employed for computing the weighted least-squares solution. Default value: TRUE (i.e., "fast but possibly diminished accuracy").
<code>maximum_iterations</code>	Optional maximum number of iterations of Fisher scoring to run; "and-ed" with result of <code>convergence_criteria_fn</code> . Default value: NULL (i.e., infinity).
<code>name</code>	used as name prefix to ops created by this function. Default value: "fit".
<code>...</code>	other arguments passed to specific methods.

Value

A glm_fit object with parameter estimates, and number of required steps.

See Also

Other glm_fit: [glm_families](#), [glm_fit_one_step.tensorflow.tensor\(\)](#)

glm_fit_one_step *Runs one Fisher scoring step*

Description

Runs one Fisher scoring step

Usage

```
glm_fit_one_step(x, ...)
```

Arguments

x float-like, matrix-shaped Tensor where each row represents a sample's features.
... other arguments passed to specific methods.

Value

A glm_fit object with parameter estimates, number of iterations, etc.

See Also

[glm_fit_one_step.tensorflow.tensor\(\)](#)

glm_fit_one_step.tensorflow.tensor
Runs one Fisher Scoring step

Description

Runs one Fisher Scoring step

Usage

```

## S3 method for class 'tensorflow.tensor'
glm_fit_one_step(
  x,
  response,
  model,
  model_coefficients_start = NULL,
  predicted_linear_response_start = NULL,
  l2_regularizer = NULL,
  dispersion = NULL,
  offset = NULL,
  learning_rate = NULL,
  fast_unsafe_numerics = TRUE,
  name = NULL,
  ...
)

```

Arguments

<code>x</code>	float-like, matrix-shaped Tensor where each row represents a sample's features.
<code>response</code>	vector-shaped Tensor where each element represents a sample's observed response (to the corresponding row of features). Must have same dtype as <code>x</code> .
<code>model</code>	a string naming the model (see glm_families) or a <code>tfp\$glm\$ExponentialFamily</code> -like instance which implicitly characterizes a negative log-likelihood loss by specifying the distribution's mean, <code>gradient_mean</code> , and variance.
<code>model_coefficients_start</code>	Optional (batch of) vector-shaped Tensor representing the initial model coefficients, one for each column in <code>x</code> . Must have same dtype as <code>model_matrix</code> . Default value: Zeros.
<code>predicted_linear_response_start</code>	Optional Tensor with shape, dtype matching <code>response</code> ; represents offset shifted initial linear predictions based on <code>model_coefficients_start</code> . Default value: offset if <code>model_coefficients</code> is NULL, and <code>tf\$linalg\$matvec(x, model_coefficients_start) + offset</code> otherwise.
<code>l2_regularizer</code>	Optional scalar Tensor representing L2 regularization penalty. Default: NULL ie. no regularization.
<code>dispersion</code>	Optional (batch of) Tensor representing response dispersion.
<code>offset</code>	Optional Tensor representing constant shift applied to <code>predicted_linear_response</code> .
<code>learning_rate</code>	Optional (batch of) scalar Tensor used to dampen iterative progress. Typically only needed if optimization diverges, should be no larger than 1 and typically very close to 1. Default value: NULL (i.e., 1).
<code>fast_unsafe_numerics</code>	Optional Python bool indicating if faster, less numerically accurate methods can be employed for computing the weighted least-squares solution. Default value: TRUE (i.e., "fast but possibly diminished accuracy").
<code>name</code>	used as name prefix to ops created by this function. Default value: "fit".
<code>...</code>	other arguments passed to specific methods.

Value

A `glm_fit` object with parameter estimates, and number of required steps.

See Also

Other `glm_fit`: [glm_families](#), [glm_fit.tensorflow.tensor\(\)](#)

`initializer_blockwise` *Blockwise Initializer*

Description

Initializer which concatenates other initializers

Usage

```
initializer_blockwise(initializers, sizes, validate_args = FALSE)
```

Arguments

<code>initializers</code>	list of Keras initializers, eg: <code>keras::initializer_glorot_uniform()</code> or <code>initializer_constant()</code> .
<code>sizes</code>	list of integers/scalars representing the number of elements associated with each initializer in <code>initializers</code> .
<code>validate_args</code>	bool indicating we should do (possibly expensive) graph-time assertions, if necessary. @return Initializer which concatenates other initializers

`install_tfprobability` *Installs TensorFlow Probability*

Description

Installs TensorFlow Probability

Usage

```
install_tfprobability(
  method = c("auto", "virtualenv", "conda"),
  conda = "auto",
  version = "default",
  tensorflow = "default",
  extra_packages = NULL,
  ...,
  pip_ignore_installed = TRUE
)
```

Arguments

method	Installation method. By default, "auto" automatically finds a method that will work in the local environment. Change the default to force a specific installation method. Note that the "virtualenv" method is not available on Windows.
conda	The path to a conda executable. Use "auto" to allow reticulate to automatically find an appropriate conda binary. See conda_binary() for more details on how reticulate tries to resolve the conda executable.
version	TensorFlow version to install. Valid values include: <ul style="list-style-type: none"> • "default" installs 2.7 • "release" installs the latest release version of tensorflow (which may be incompatible with the current version of the R package) • A version specification like "2.4" or "2.4.0". Note that if the patch version is not supplied, the latest patch release is installed (e.g., "2.4" today installs version "2.4.2") • nightly for the latest available nightly build. • To any specification, you can append "-cpu" to install the cpu version only of the package (e.g., "2.4-cpu") • The full URL or path to an installer binary or python *.whl file.
tensorflow	Synonym for version. Maintained for backwards.
extra_packages	Additional Python packages to install along with TensorFlow.
...	other arguments passed to reticulate::conda_install() or reticulate::virtualenv_install() , depending on the method used.
pip_ignore_installed	Whether pip should ignore installed python packages and reinstall all already installed python packages. This defaults to TRUE, to ensure that TensorFlow dependencies like NumPy are compatible with the prebuilt TensorFlow binaries.

Value

invisible

layer_autoregressive *Masked Autoencoder for Distribution Estimation*

Description

layer_autoregressive takes as input a Tensor of shape $[..., \text{event_size}]$ and returns a Tensor of shape $[..., \text{event_size}, \text{params}]$. The output satisfies the autoregressive property. That is, the layer is configured with some permutation ord of $\{0, \dots, \text{event_size}-1\}$ (i.e., an ordering of the input dimensions), and the output $\text{output}[\text{batch_idx}, i, \dots]$ for input dimension i depends only on inputs $x[\text{batch_idx}, j]$ where $\text{ord}(j) < \text{ord}(i)$.

Usage

```

layer_autoregressive(
    object,
    params,
    event_shape = NULL,
    hidden_units = NULL,
    input_order = "left-to-right",
    hidden_degrees = "equal",
    activation = NULL,
    use_bias = TRUE,
    kernel_initializer = "glorot_uniform",
    validate_args = FALSE,
    ...
)

```

Arguments

object	What to compose the new Layer instance with. Typically a Sequential model or a Tensor (e.g., as returned by <code>layer_input()</code>). The return value depends on object. If object is: <ul style="list-style-type: none"> • missing or NULL, the Layer instance is returned. • a Sequential model, the model with an additional layer is returned. • a Tensor, the output tensor from <code>layer_instance(object)</code> is returned.
params	integer specifying the number of parameters to output per input.
event_shape	list-like of positive integers (or a single int), specifying the shape of the input to this layer, which is also the <code>event_shape</code> of the distribution parameterized by this layer. Currently only rank-1 shapes are supported. That is, <code>event_shape</code> must be a single integer. If not specified, the event shape is inferred when this layer is first called or built.
hidden_units	list-like of non-negative integers, specifying the number of units in each hidden layer.
input_order	Order of degrees to the input units: 'random', 'left-to-right', 'right-to-left', or an array of an explicit order. For example, 'left-to-right' builds an autoregressive model: $p(x) = p(x_1) p(x_2 x_1) \dots p(x_D x_{<D})$. Default: 'left-to-right'.
hidden_degrees	Method for assigning degrees to the hidden units: 'equal', 'random'. If 'equal', hidden units in each layer are allocated equally (up to a remainder term) to each degree. Default: 'equal'.
activation	An activation function. See <code>keras::layer_dense</code> . Default: NULL.
use_bias	Whether or not the dense layers constructed in this layer should have a bias term. See <code>keras::layer_dense</code> . Default: TRUE.
kernel_initializer	Initializer for the kernel weights matrix. Default: 'glorot_uniform'.
validate_args	logical, default FALSE. When TRUE, layer parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs.

... Additional keyword arguments passed to the keras::layer_dense constructed by this layer.

Details

The autoregressive property allows us to use `output[batch_idx, i]` to parameterize conditional distributions: $p(x[\text{batch_idx}, i] \mid x[\text{batch_idx}, j] \text{ for } \text{ord}(j) < \text{ord}(i))$ which give us a tractable distribution over input `x[batch_idx]`:

$$p(x[\text{batch_idx}]) = \prod_i p(x[\text{batch_idx}, \text{ord}(i)] \mid x[\text{batch_idx}, \text{ord}(0:i)])$$

For example, when `params` is 2, the output of the layer can parameterize the location and log-scale of an autoregressive Gaussian distribution.

Value

a Keras layer

See Also

Other layers: [layer_conv_1d_flipout\(\)](#), [layer_conv_1d_reparameterization\(\)](#), [layer_conv_2d_flipout\(\)](#), [layer_conv_2d_reparameterization\(\)](#), [layer_conv_3d_flipout\(\)](#), [layer_conv_3d_reparameterization\(\)](#), [layer_dense_flipout\(\)](#), [layer_dense_local_reparameterization\(\)](#), [layer_dense_reparameterization\(\)](#), [layer_dense_variational\(\)](#), [layer_variable\(\)](#)

layer_autoregressive_transform

An autoregressive normalizing flow layer, given a layer_autoregressive.

Description

Following [Papamakarios et al. \(2017\)](#), given an autoregressive model $p(x)$ with conditional distributions in the location-scale family, we can construct a normalizing flow for $p(x)$.

Usage

```
layer_autoregressive_transform(object, made, ...)
```

Arguments

<code>object</code>	What to compose the new Layer instance with. Typically a Sequential model or a Tensor (e.g., as returned by <code>layer_input()</code>). The return value depends on <code>object</code> . If <code>object</code> is: <ul style="list-style-type: none"> missing or NULL, the Layer instance is returned. a Sequential model, the model with an additional layer is returned. a Tensor, the output tensor from <code>layer_instance(object)</code> is returned.
<code>made</code>	A Made layer, which must output two parameters for each input.
...	Additional parameters passed to Keras Layer.

Details

Specifically, suppose `made` is a `[layer_autoregressive()]` – a layer implementing a Masked Autoencoder for Distribution Estimation (MADE) – that computes location and log-scale parameters $made(x)[i]$ for each input $x[i]$. Then we can represent the autoregressive model $p(x)$ as $x = f(u)$ where u is drawn from some base distribution and where f is an invertible and differentiable function (i.e., a Bijector) and $f^{-1}(x)$ is defined by:

```
library(tensorflow)
library(zeallot)
f_inverse <- function(x) {
  c(shift, log_scale) %<-% tf$unstack(made(x), 2, axis = -1L)
  (x - shift) * tf$math$exp(-log_scale)
}
```

Given a `layer_autoregressive()` `made`, a `layer_autoregressive_transform()` transforms an input $tfd_* p(u)$ to an output $tfd_* p(x)$ where $x = f(u)$.

Value

a Keras layer

References

[Papamakarios et al. \(2017\)](#)

See Also

[tfb_masked_autoregressive_flow\(\)](#) and [layer_autoregressive\(\)](#)

layer_categorical_mixture_of_one_hot_categorical

*A OneHotCategorical mixture Keras layer from $k * (1 + d)$ params.*

Description

`k` (i.e., `num_components`) represents the number of component `OneHotCategorical` distributions and `d` (i.e., `event_size`) represents the number of categories within each `OneHotCategorical` distribution.

Usage

```
layer_categorical_mixture_of_one_hot_categorical(
  object,
  event_size,
  num_components,
  convert_to_tensor_fn = tfp$distributions$Distribution$sample,
  sample_dtype = NULL,
```

```

    validate_args = FALSE,
    ...
)

```

Arguments

<code>object</code>	What to compose the new <code>Layer</code> instance with. Typically a <code>Sequential</code> model or a <code>Tensor</code> (e.g., as returned by <code>layer_input()</code>). The return value depends on <code>object</code> . If <code>object</code> is: <ul style="list-style-type: none"> • missing or <code>NULL</code>, the <code>Layer</code> instance is returned. • a <code>Sequential</code> model, the model with an additional layer is returned. • a <code>Tensor</code>, the output tensor from <code>layer_instance(object)</code> is returned.
<code>event_size</code>	Scalar integer representing the size of single draw from this distribution.
<code>num_components</code>	Scalar integer representing the number of mixture components. Must be at least 1. (If <code>num_components=1</code> , it's more efficient to use the <code>OneHotCategorical</code> layer.)
<code>convert_to_tensor_fn</code>	A callable that takes a <code>tfd\$Distribution</code> instance and returns a <code>tf\$Tensor</code> -like object. Default value: <code>tfd\$distributions\$Distribution\$sample</code> .
<code>sample_dtype</code>	<code>dtype</code> of samples produced by this distribution. Default value: <code>NULL</code> (i.e., previous layer's <code>dtype</code>).
<code>validate_args</code>	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
<code>...</code>	Additional arguments passed to <code>args</code> of <code>keras::create_layer</code> .

Details

Typical choices for `convert_to_tensor_fn` include:

- `tfp$distributions$Distribution$sample`
- `tfp$distributions$Distribution$mean`
- `tfp$distributions$Distribution$mode`

Value

a Keras layer

See Also

For an example how to use in a Keras model, see [layer_independent_normal\(\)](#).

Other `distribution_layers`: [layer_distribution_lambda\(\)](#), [layer_independent_bernoulli\(\)](#), [layer_independent_logistic\(\)](#), [layer_independent_normal\(\)](#), [layer_independent_poisson\(\)](#), [layer_kl_divergence_add_loss\(\)](#), [layer_kl_divergence_regularizer\(\)](#), [layer_mixture_logistic\(\)](#), [layer_mixture_normal\(\)](#), [layer_mixture_same_family\(\)](#), [layer_multivariate_normal_tri_l\(\)](#), [layer_one_hot_categorical\(\)](#)

layer_conv_1d_flipout *1D convolution layer (e.g. temporal convolution) with Flipout*

Description

This layer creates a convolution kernel that is convolved (actually cross-correlated) with the layer input to produce a tensor of outputs. It may also include a bias addition and activation function on the outputs. It assumes the kernel and/or bias are drawn from distributions.

Usage

```
layer_conv_1d_flipout(
    object,
    filters,
    kernel_size,
    strides = 1,
    padding = "valid",
    data_format = "channels_last",
    dilation_rate = 1,
    activation = NULL,
    activity_regularizer = NULL,
    trainable = TRUE,
    kernel_posterior_fn = tfp$layers$util$default_mean_field_normal_fn(),
    kernel_posterior_tensor_fn = function(d) d %>% tfd_sample(),
    kernel_prior_fn = tfp$layers$util$default_multivariate_normal_fn,
    kernel_divergence_fn = function(q, p, ignore) tfd_kl_divergence(q, p),
    bias_posterior_fn = tfp$layers$util$default_mean_field_normal_fn(is_singular = TRUE),
    bias_posterior_tensor_fn = function(d) d %>% tfd_sample(),
    bias_prior_fn = NULL,
    bias_divergence_fn = function(q, p, ignore) tfd_kl_divergence(q, p),
    ...
)
```

Arguments

object	What to compose the new Layer instance with. Typically a Sequential model or a Tensor (e.g., as returned by layer_input()). The return value depends on object. If object is: <ul style="list-style-type: none"> • missing or NULL, the Layer instance is returned. • a Sequential model, the model with an additional layer is returned. • a Tensor, the output tensor from layer_instance(object) is returned.
filters	Integer, the dimensionality of the output space (i.e. the number of filters in the convolution).
kernel_size	An integer or list of a single integer, specifying the length of the 1D convolution window.

strides	An integer or list of a single integer, specifying the stride length of the convolution. Specifying any stride value $\neq 1$ is incompatible with specifying any dilation_rate value $\neq 1$.
padding	One of "valid" or "same" (case-insensitive).
data_format	A string, one of channels_last (default) or channels_first. The ordering of the dimensions in the inputs. channels_last corresponds to inputs with shape (batch, length, channels) while channels_first corresponds to inputs with shape (batch, channels, length).
dilation_rate	An integer or tuple/list of a single integer, specifying the dilation rate to use for dilated convolution. Currently, specifying any dilation_rate value $\neq 1$ is incompatible with specifying any strides value $\neq 1$.
activation	Activation function. Set it to None to maintain a linear activation.
activity_regularizer	Regularizer function for the output.
trainable	Whether the layer weights will be updated during training.
kernel_posterior_fn	Function which creates tfd\$Distribution instance representing the surrogate posterior of the kernel parameter. Default value: default_mean_field_normal_fn().
kernel_posterior_tensor_fn	Function which takes a tfd\$Distribution instance and returns a representative value. Default value: function(d) d %>% tfd_sample().
kernel_prior_fn	Function which creates tfd\$Distribution instance. See default_mean_field_normal_fn docstring for required parameter signature. Default value: tfd_normal(loc = 0, scale = 1).
kernel_divergence_fn	Function which takes the surrogate posterior distribution, prior distribution and random variate sample(s) from the surrogate posterior and computes or approximates the KL divergence. The distributions are tfd\$Distribution-like instances and the sample is a Tensor.
bias_posterior_fn	Function which creates a tfd\$Distribution instance representing the surrogate posterior of the bias parameter. Default value: default_mean_field_normal_fn(is_singular = TRUE) (which creates an instance of tfd_deterministic).
bias_posterior_tensor_fn	Function which takes a tfd\$Distribution instance and returns a representative value. Default value: function(d) d %>% tfd_sample().
bias_prior_fn	Function which creates tfd instance. See default_mean_field_normal_fn docstring for required parameter signature. Default value: NULL (no prior, no variational inference)
bias_divergence_fn	Function which takes the surrogate posterior distribution, prior distribution and random variate sample(s) from the surrogate posterior and computes or approximates the KL divergence. The distributions are tfd\$Distribution-like instances and the sample is a Tensor.
...	Additional keyword arguments passed to the keras::layer_dense constructed by this layer.

Details

This layer implements the Bayesian variational inference analogue to a dense layer by assuming the kernel and/or the bias are drawn from distributions.

By default, the layer implements a stochastic forward pass via sampling from the kernel and bias posteriors,

```
outputs = f(inputs; kernel, bias), kernel, bias ~ posterior
```

where f denotes the layer's calculation. It uses the Flipout estimator (Wen et al., 2018), which performs a Monte Carlo approximation of the distribution integrating over the kernel and bias. Flipout uses roughly twice as many floating point operations as the reparameterization estimator but has the advantage of significantly lower variance.

The arguments permit separate specification of the surrogate posterior ($q(W|x)$), prior ($p(W)$), and divergence for both the kernel and bias distributions.

Upon being built, this layer adds losses (accessible via the `losses` property) representing the divergences of kernel and/or bias surrogate posteriors and their respective priors. When doing minibatch stochastic optimization, make sure to scale this loss such that it is applied just once per epoch (e.g. if `kl` is the sum of losses for each element of the batch, you should pass `kl / num_examples_per_epoch` to your optimizer). You can access the kernel and/or bias posterior and prior distributions after the layer is built via the `kernel_posterior`, `kernel_prior`, `bias_posterior` and `bias_prior` properties.

Value

a Keras layer

References

- Yeming Wen, Paul Vicol, Jimmy Ba, Dustin Tran, and Roger Grosse. Flipout: Efficient Pseudo-Independent Weight Perturbations on Mini-Batches. In *International Conference on Learning Representations*, 2018.

See Also

Other layers: [layer_autoregressive\(\)](#), [layer_conv_1d_reparameterization\(\)](#), [layer_conv_2d_flipout\(\)](#), [layer_conv_2d_reparameterization\(\)](#), [layer_conv_3d_flipout\(\)](#), [layer_conv_3d_reparameterization\(\)](#), [layer_dense_flipout\(\)](#), [layer_dense_local_reparameterization\(\)](#), [layer_dense_reparameterization\(\)](#), [layer_dense_variational\(\)](#), [layer_variable\(\)](#)

layer_conv_1d_reparameterization

1D convolution layer (e.g. temporal convolution).

Description

This layer creates a convolution kernel that is convolved (actually cross-correlated) with the layer input to produce a tensor of outputs. It may also include a bias addition and activation function on the outputs. It assumes the kernel and/or bias are drawn from distributions.

Usage

```
layer_conv_1d_reparameterization(
  object,
  filters,
  kernel_size,
  strides = 1,
  padding = "valid",
  data_format = "channels_last",
  dilation_rate = 1,
  activation = NULL,
  activity_regularizer = NULL,
  trainable = TRUE,
  kernel_posterior_fn = tfp$layers$util$default_mean_field_normal_fn(),
  kernel_posterior_tensor_fn = function(d) d %>% tfd_sample(),
  kernel_prior_fn = tfp$layers$util$default_multivariate_normal_fn,
  kernel_divergence_fn = function(q, p, ignore) tfd_kl_divergence(q, p),
  bias_posterior_fn = tfp$layers$util$default_mean_field_normal_fn(is_singular = TRUE),
  bias_posterior_tensor_fn = function(d) d %>% tfd_sample(),
  bias_prior_fn = NULL,
  bias_divergence_fn = function(q, p, ignore) tfd_kl_divergence(q, p),
  ...
)
```

Arguments

object	What to compose the new Layer instance with. Typically a Sequential model or a Tensor (e.g., as returned by <code>layer_input()</code>). The return value depends on object. If object is: <ul style="list-style-type: none"> • missing or NULL, the Layer instance is returned. • a Sequential model, the model with an additional layer is returned. • a Tensor, the output tensor from <code>layer_instance(object)</code> is returned.
filters	Integer, the dimensionality of the output space (i.e. the number of filters in the convolution).
kernel_size	An integer or list of a single integer, specifying the length of the 1D convolution window.
strides	An integer or list of a single integer, specifying the stride length of the convolution. Specifying any stride value $\neq 1$ is incompatible with specifying any <code>dilation_rate</code> value $\neq 1$.
padding	One of "valid" or "same" (case-insensitive).

<code>data_format</code>	A string, one of <code>channels_last</code> (default) or <code>channels_first</code> . The ordering of the dimensions in the inputs. <code>channels_last</code> corresponds to inputs with shape (batch, length, channels) while <code>channels_first</code> corresponds to inputs with shape (batch, channels, length).
<code>dilation_rate</code>	An integer or tuple/list of a single integer, specifying the dilation rate to use for dilated convolution. Currently, specifying any <code>dilation_rate</code> value $\neq 1$ is incompatible with specifying any <code>strides</code> value $\neq 1$.
<code>activation</code>	Activation function. Set it to <code>None</code> to maintain a linear activation.
<code>activity_regularizer</code>	Regularizer function for the output.
<code>trainable</code>	Whether the layer weights will be updated during training.
<code>kernel_posterior_fn</code>	Function which creates <code>tfd\$Distribution</code> instance representing the surrogate posterior of the kernel parameter. Default value: <code>default_mean_field_normal_fn()</code> .
<code>kernel_posterior_tensor_fn</code>	Function which takes a <code>tfd\$Distribution</code> instance and returns a representative value. Default value: <code>function(d) d %>% tfd_sample()</code> .
<code>kernel_prior_fn</code>	Function which creates <code>tfd\$Distribution</code> instance. See <code>default_mean_field_normal_fn</code> docstring for required parameter signature. Default value: <code>tfd_normal(loc = 0, scale = 1)</code> .
<code>kernel_divergence_fn</code>	Function which takes the surrogate posterior distribution, prior distribution and random variate sample(s) from the surrogate posterior and computes or approximates the KL divergence. The distributions are <code>tfd\$Distribution</code> -like instances and the sample is a <code>Tensor</code> .
<code>bias_posterior_fn</code>	Function which creates a <code>tfd\$Distribution</code> instance representing the surrogate posterior of the bias parameter. Default value: <code>default_mean_field_normal_fn(is_singular = TRUE)</code> (which creates an instance of <code>tfd_deterministic</code>).
<code>bias_posterior_tensor_fn</code>	Function which takes a <code>tfd\$Distribution</code> instance and returns a representative value. Default value: <code>function(d) d %>% tfd_sample()</code> .
<code>bias_prior_fn</code>	Function which creates <code>tfd</code> instance. See <code>default_mean_field_normal_fn</code> docstring for required parameter signature. Default value: <code>NULL</code> (no prior, no variational inference)
<code>bias_divergence_fn</code>	Function which takes the surrogate posterior distribution, prior distribution and random variate sample(s) from the surrogate posterior and computes or approximates the KL divergence. The distributions are <code>tfd\$Distribution</code> -like instances and the sample is a <code>Tensor</code> .
<code>...</code>	Additional keyword arguments passed to the <code>keras::layer_dense</code> constructed by this layer.

Details

This layer implements the Bayesian variational inference analogue to a dense layer by assuming the kernel and/or the bias are drawn from distributions.

By default, the layer implements a stochastic forward pass via sampling from the kernel and bias posteriors,

```
outputs = f(inputs; kernel, bias), kernel, bias ~ posterior
```

where f denotes the layer's calculation. It uses the reparameterization estimator (Kingma and Welling, 2014), which performs a Monte Carlo approximation of the distribution integrating over the kernel and bias.

The arguments permit separate specification of the surrogate posterior ($q(W|x)$), prior ($p(W)$), and divergence for both the kernel and bias distributions.

Upon being built, this layer adds losses (accessible via the `losses` property) representing the divergences of kernel and/or bias surrogate posteriors and their respective priors. When doing minibatch stochastic optimization, make sure to scale this loss such that it is applied just once per epoch (e.g. if `kl` is the sum of losses for each element of the batch, you should pass `kl / num_examples_per_epoch` to your optimizer). You can access the kernel and/or bias posterior and prior distributions after the layer is built via the `kernel_posterior`, `kernel_prior`, `bias_posterior` and `bias_prior` properties.

Value

a Keras layer

References

- Diederik Kingma and Max Welling. Auto-Encoding Variational Bayes. In *International Conference on Learning Representations*, 2014.

See Also

Other layers: `layer_autoregressive()`, `layer_conv_1d_flipout()`, `layer_conv_2d_flipout()`, `layer_conv_2d_reparameterization()`, `layer_conv_3d_flipout()`, `layer_conv_3d_reparameterization()`, `layer_dense_flipout()`, `layer_dense_local_reparameterization()`, `layer_dense_reparameterization()`, `layer_dense_variational()`, `layer_variable()`

layer_conv_2d_flipout *2D convolution layer (e.g. spatial convolution over images) with Flipout*

Description

This layer creates a convolution kernel that is convolved (actually cross-correlated) with the layer input to produce a tensor of outputs. It may also include a bias addition and activation function on the outputs. It assumes the kernel and/or bias are drawn from distributions.

Usage

```

layer_conv_2d_flipout(
  object,
  filters,
  kernel_size,
  strides = 1,
  padding = "valid",
  data_format = "channels_last",
  dilation_rate = 1,
  activation = NULL,
  activity_regularizer = NULL,
  trainable = TRUE,
  kernel_posterior_fn = tfp$layers$util$default_mean_field_normal_fn(),
  kernel_posterior_tensor_fn = function(d) d %>% tfd_sample(),
  kernel_prior_fn = tfp$layers$util$default_multivariate_normal_fn(),
  kernel_divergence_fn = function(q, p, ignore) tfd_kl_divergence(q, p),
  bias_posterior_fn = tfp$layers$util$default_mean_field_normal_fn(is_singular = TRUE),
  bias_posterior_tensor_fn = function(d) d %>% tfd_sample(),
  bias_prior_fn = NULL,
  bias_divergence_fn = function(q, p, ignore) tfd_kl_divergence(q, p),
  ...
)

```

Arguments

object	What to compose the new Layer instance with. Typically a Sequential model or a Tensor (e.g., as returned by <code>layer_input()</code>). The return value depends on object. If object is: <ul style="list-style-type: none"> • missing or NULL, the Layer instance is returned. • a Sequential model, the model with an additional layer is returned. • a Tensor, the output tensor from <code>layer_instance(object)</code> is returned.
filters	Integer, the dimensionality of the output space (i.e. the number of filters in the convolution).
kernel_size	An integer or list of a single integer, specifying the length of the 1D convolution window.
strides	An integer or list of a single integer, specifying the stride length of the convolution. Specifying any stride value $\neq 1$ is incompatible with specifying any <code>dilation_rate</code> value $\neq 1$.
padding	One of "valid" or "same" (case-insensitive).
data_format	A string, one of <code>channels_last</code> (default) or <code>channels_first</code> . The ordering of the dimensions in the inputs. <code>channels_last</code> corresponds to inputs with shape (batch, length, channels) while <code>channels_first</code> corresponds to inputs with shape (batch, channels, length).
dilation_rate	An integer or tuple/list of a single integer, specifying the dilation rate to use for dilated convolution. Currently, specifying any <code>dilation_rate</code> value $\neq 1$ is incompatible with specifying any <code>strides</code> value $\neq 1$.

activation	Activation function. Set it to None to maintain a linear activation.
activity_regularizer	Regularizer function for the output.
trainable	Whether the layer weights will be updated during training.
kernel_posterior_fn	Function which creates tfd\$Distribution instance representing the surrogate posterior of the kernel parameter. Default value: default_mean_field_normal_fn().
kernel_posterior_tensor_fn	Function which takes a tfd\$Distribution instance and returns a representative value. Default value: function(d) d %>% tfd_sample().
kernel_prior_fn	Function which creates tfd\$Distribution instance. See default_mean_field_normal_fn docstring for required parameter signature. Default value: tfd_normal(loc = 0, scale = 1).
kernel_divergence_fn	Function which takes the surrogate posterior distribution, prior distribution and random variate sample(s) from the surrogate posterior and computes or approximates the KL divergence. The distributions are tfd\$Distribution-like instances and the sample is a Tensor.
bias_posterior_fn	Function which creates a tfd\$Distribution instance representing the surrogate posterior of the bias parameter. Default value: default_mean_field_normal_fn(is_singular = TRUE) (which creates an instance of tfd_deterministic).
bias_posterior_tensor_fn	Function which takes a tfd\$Distribution instance and returns a representative value. Default value: function(d) d %>% tfd_sample().
bias_prior_fn	Function which creates tfd instance. See default_mean_field_normal_fn docstring for required parameter signature. Default value: NULL (no prior, no variational inference)
bias_divergence_fn	Function which takes the surrogate posterior distribution, prior distribution and random variate sample(s) from the surrogate posterior and computes or approximates the KL divergence. The distributions are tfd\$Distribution-like instances and the sample is a Tensor.
...	Additional keyword arguments passed to the keras::layer_dense constructed by this layer.

Details

This layer implements the Bayesian variational inference analogue to a dense layer by assuming the kernel and/or the bias are drawn from distributions.

By default, the layer implements a stochastic forward pass via sampling from the kernel and bias posteriors,

```
outputs = f(inputs; kernel, bias), kernel, bias ~ posterior
```

where f denotes the layer's calculation. It uses the Flipout estimator (Wen et al., 2018), which performs a Monte Carlo approximation of the distribution integrating over the kernel and bias. Flipout uses roughly twice as many floating point operations as the reparameterization estimator but has the advantage of significantly lower variance.

The arguments permit separate specification of the surrogate posterior ($q(W|x)$), prior ($p(W)$), and divergence for both the kernel and bias distributions.

Upon being built, this layer adds losses (accessible via the `losses` property) representing the divergences of kernel and/or bias surrogate posteriors and their respective priors. When doing minibatch stochastic optimization, make sure to scale this loss such that it is applied just once per epoch (e.g. if `kl` is the sum of losses for each element of the batch, you should pass `kl / num_examples_per_epoch` to your optimizer). You can access the kernel and/or bias posterior and prior distributions after the layer is built via the `kernel_posterior`, `kernel_prior`, `bias_posterior` and `bias_prior` properties.

Value

a Keras layer

References

- Yeming Wen, Paul Vicol, Jimmy Ba, Dustin Tran, and Roger Grosse. Flipout: Efficient Pseudo-Independent Weight Perturbations on Mini-Batches. In *International Conference on Learning Representations*, 2018.

See Also

Other layers: [layer_autoregressive\(\)](#), [layer_conv_1d_flipout\(\)](#), [layer_conv_1d_reparameterization\(\)](#), [layer_conv_2d_reparameterization\(\)](#), [layer_conv_3d_flipout\(\)](#), [layer_conv_3d_reparameterization\(\)](#), [layer_dense_flipout\(\)](#), [layer_dense_local_reparameterization\(\)](#), [layer_dense_reparameterization\(\)](#), [layer_dense_variational\(\)](#), [layer_variable\(\)](#)

layer_conv_2d_reparameterization

2D convolution layer (e.g. spatial convolution over images)

Description

This layer creates a convolution kernel that is convolved (actually cross-correlated) with the layer input to produce a tensor of outputs. It may also include a bias addition and activation function on the outputs. It assumes the kernel and/or bias are drawn from distributions.

Usage

```
layer_conv_2d_reparameterization(
    object,
    filters,
    kernel_size,
```

```

    strides = 1,
    padding = "valid",
    data_format = "channels_last",
    dilation_rate = 1,
    activation = NULL,
    activity_regularizer = NULL,
    trainable = TRUE,
    kernel_posterior_fn = tfp$layers$util$default_mean_field_normal_fn(),
    kernel_posterior_tensor_fn = function(d) d %>% tfd_sample(),
    kernel_prior_fn = tfp$layers$util$default_multivariate_normal_fn,
    kernel_divergence_fn = function(q, p, ignore) tfd_kl_divergence(q, p),
    bias_posterior_fn = tfp$layers$util$default_mean_field_normal_fn(is_singular = TRUE),
    bias_posterior_tensor_fn = function(d) d %>% tfd_sample(),
    bias_prior_fn = NULL,
    bias_divergence_fn = function(q, p, ignore) tfd_kl_divergence(q, p),
    ...
)

```

Arguments

object	What to compose the new Layer instance with. Typically a Sequential model or a Tensor (e.g., as returned by <code>layer_input()</code>). The return value depends on object. If object is: <ul style="list-style-type: none"> • missing or NULL, the Layer instance is returned. • a Sequential model, the model with an additional layer is returned. • a Tensor, the output tensor from <code>layer_instance(object)</code> is returned.
filters	Integer, the dimensionality of the output space (i.e. the number of filters in the convolution).
kernel_size	An integer or list of a single integer, specifying the length of the 1D convolution window.
strides	An integer or list of a single integer, specifying the stride length of the convolution. Specifying any stride value $\neq 1$ is incompatible with specifying any <code>dilation_rate</code> value $\neq 1$.
padding	One of "valid" or "same" (case-insensitive).
data_format	A string, one of <code>channels_last</code> (default) or <code>channels_first</code> . The ordering of the dimensions in the inputs. <code>channels_last</code> corresponds to inputs with shape (batch, length, channels) while <code>channels_first</code> corresponds to inputs with shape (batch, channels, length).
dilation_rate	An integer or tuple/list of a single integer, specifying the dilation rate to use for dilated convolution. Currently, specifying any <code>dilation_rate</code> value $\neq 1$ is incompatible with specifying any <code>strides</code> value $\neq 1$.
activation	Activation function. Set it to None to maintain a linear activation.
activity_regularizer	Regularizer function for the output.
trainable	Whether the layer weights will be updated during training.

kernel_posterior_fn	Function which creates tfd\$Distribution instance representing the surrogate posterior of the kernel parameter. Default value: default_mean_field_normal_fn().
kernel_posterior_tensor_fn	Function which takes a tfd\$Distribution instance and returns a representative value. Default value: function(d) d %>% tfd_sample().
kernel_prior_fn	Function which creates tfd\$Distribution instance. See default_mean_field_normal_fn docstring for required parameter signature. Default value: tfd_normal(loc = 0, scale = 1).
kernel_divergence_fn	Function which takes the surrogate posterior distribution, prior distribution and random variate sample(s) from the surrogate posterior and computes or approximates the KL divergence. The distributions are tfd\$Distribution-like instances and the sample is a Tensor.
bias_posterior_fn	Function which creates a tfd\$Distribution instance representing the surrogate posterior of the bias parameter. Default value: default_mean_field_normal_fn(is_singular = TRUE) (which creates an instance of tfd_deterministic).
bias_posterior_tensor_fn	Function which takes a tfd\$Distribution instance and returns a representative value. Default value: function(d) d %>% tfd_sample().
bias_prior_fn	Function which creates tfd instance. See default_mean_field_normal_fn docstring for required parameter signature. Default value: NULL (no prior, no variational inference)
bias_divergence_fn	Function which takes the surrogate posterior distribution, prior distribution and random variate sample(s) from the surrogate posterior and computes or approximates the KL divergence. The distributions are tfd\$Distribution-like instances and the sample is a Tensor.
...	Additional keyword arguments passed to the keras::layer_dense constructed by this layer.

Details

This layer implements the Bayesian variational inference analogue to a dense layer by assuming the kernel and/or the bias are drawn from distributions.

By default, the layer implements a stochastic forward pass via sampling from the kernel and bias posteriors,

$$\text{outputs} = f(\text{inputs}; \text{kernel}, \text{bias}), \text{kernel}, \text{bias} \sim \text{posterior}$$

where f denotes the layer's calculation. It uses the reparameterization estimator (Kingma and Welling, 2014), which performs a Monte Carlo approximation of the distribution integrating over the kernel and bias.

The arguments permit separate specification of the surrogate posterior ($q(W|x)$), prior ($p(W)$), and divergence for both the kernel and bias distributions.

Upon being built, this layer adds losses (accessible via the losses property) representing the divergences of kernel and/or bias surrogate posteriors and their respective priors. When doing minibatch stochastic optimization, make sure to scale this loss such that it is applied just once per epoch (e.g. if kl is the sum of losses for each element of the batch, you should pass kl / num_examples_per_epoch to your optimizer). You can access the kernel and/or bias posterior and prior distributions after the layer is built via the kernel_posterior, kernel_prior, bias_posterior and bias_prior properties.

Value

a Keras layer

References

- [Diederik Kingma and Max Welling. Auto-Encoding Variational Bayes. In *International Conference on Learning Representations*, 2014.](#)

See Also

Other layers: [layer_autoregressive\(\)](#), [layer_conv_1d_flipout\(\)](#), [layer_conv_1d_reparameterization\(\)](#), [layer_conv_2d_flipout\(\)](#), [layer_conv_3d_flipout\(\)](#), [layer_conv_3d_reparameterization\(\)](#), [layer_dense_flipout\(\)](#), [layer_dense_local_reparameterization\(\)](#), [layer_dense_reparameterization\(\)](#), [layer_dense_variational\(\)](#), [layer_variable\(\)](#)

layer_conv_3d_flipout *3D convolution layer (e.g. spatial convolution over volumes) with Flipout*

Description

This layer creates a convolution kernel that is convolved (actually cross-correlated) with the layer input to produce a tensor of outputs. It may also include a bias addition and activation function on the outputs. It assumes the kernel and/or bias are drawn from distributions.

Usage

```
layer_conv_3d_flipout(
    object,
    filters,
    kernel_size,
    strides = 1,
    padding = "valid",
    data_format = "channels_last",
    dilation_rate = 1,
    activation = NULL,
    activity_regularizer = NULL,
    trainable = TRUE,
    kernel_posterior_fn = tfp$layers$util$default_mean_field_normal_fn(),
```

```

kernel_posterior_tensor_fn = function(d) d %>% tfd_sample(),
kernel_prior_fn = tfp$layers$util$default_multivariate_normal_fn,
kernel_divergence_fn = function(q, p, ignore) tfd_kl_divergence(q, p),
bias_posterior_fn = tfp$layers$util$default_mean_field_normal_fn(is_singular = TRUE),
bias_posterior_tensor_fn = function(d) d %>% tfd_sample(),
bias_prior_fn = NULL,
bias_divergence_fn = function(q, p, ignore) tfd_kl_divergence(q, p),
...
)

```

Arguments

object	What to compose the new Layer instance with. Typically a Sequential model or a Tensor (e.g., as returned by layer_input()). The return value depends on object. If object is: <ul style="list-style-type: none"> • missing or NULL, the Layer instance is returned. • a Sequential model, the model with an additional layer is returned. • a Tensor, the output tensor from layer_instance(object) is returned.
filters	Integer, the dimensionality of the output space (i.e. the number of filters in the convolution).
kernel_size	An integer or list of a single integer, specifying the length of the 1D convolution window.
strides	An integer or list of a single integer, specifying the stride length of the convolution. Specifying any stride value != 1 is incompatible with specifying any dilation_rate value != 1.
padding	One of "valid" or "same" (case-insensitive).
data_format	A string, one of channels_last (default) or channels_first. The ordering of the dimensions in the inputs. channels_last corresponds to inputs with shape (batch, length, channels) while channels_first corresponds to inputs with shape (batch, channels, length).
dilation_rate	An integer or tuple/list of a single integer, specifying the dilation rate to use for dilated convolution. Currently, specifying any dilation_rate value != 1 is incompatible with specifying any strides value != 1.
activation	Activation function. Set it to None to maintain a linear activation.
activity_regularizer	Regularizer function for the output.
trainable	Whether the layer weights will be updated during training.
kernel_posterior_fn	Function which creates tfd\$Distribution instance representing the surrogate posterior of the kernel parameter. Default value: default_mean_field_normal_fn().
kernel_posterior_tensor_fn	Function which takes a tfd\$Distribution instance and returns a representative value. Default value: function(d) d %>% tfd_sample().

kernel_prior_fn	Function which creates <code>tfd\$Distribution</code> instance. See <code>default_mean_field_normal_fn</code> docstring for required parameter signature. Default value: <code>tfd_normal(loc = 0, scale = 1)</code> .
kernel_divergence_fn	Function which takes the surrogate posterior distribution, prior distribution and random variate sample(s) from the surrogate posterior and computes or approximates the KL divergence. The distributions are <code>tfd\$Distribution</code> -like instances and the sample is a <code>Tensor</code> .
bias_posterior_fn	Function which creates a <code>tfd\$Distribution</code> instance representing the surrogate posterior of the bias parameter. Default value: <code>default_mean_field_normal_fn(is_singular = TRUE)</code> (which creates an instance of <code>tfd_deterministic</code>).
bias_posterior_tensor_fn	Function which takes a <code>tfd\$Distribution</code> instance and returns a representative value. Default value: <code>function(d) d %>% tfd_sample()</code> .
bias_prior_fn	Function which creates <code>tfd</code> instance. See <code>default_mean_field_normal_fn</code> docstring for required parameter signature. Default value: <code>NULL</code> (no prior, no variational inference)
bias_divergence_fn	Function which takes the surrogate posterior distribution, prior distribution and random variate sample(s) from the surrogate posterior and computes or approximates the KL divergence. The distributions are <code>tfd\$Distribution</code> -like instances and the sample is a <code>Tensor</code> .
...	Additional keyword arguments passed to the <code>keras::layer_dense</code> constructed by this layer.

Details

This layer implements the Bayesian variational inference analogue to a dense layer by assuming the kernel and/or the bias are drawn from distributions.

By default, the layer implements a stochastic forward pass via sampling from the kernel and bias posteriors,

$$\text{outputs} = f(\text{inputs}; \text{kernel}, \text{bias}), \text{kernel}, \text{bias} \sim \text{posterior}$$

where f denotes the layer's calculation. It uses the Flipout estimator (Wen et al., 2018), which performs a Monte Carlo approximation of the distribution integrating over the kernel and bias. Flipout uses roughly twice as many floating point operations as the reparameterization estimator but has the advantage of significantly lower variance.

The arguments permit separate specification of the surrogate posterior ($q(W|x)$), prior ($p(W)$), and divergence for both the kernel and bias distributions.

Upon being built, this layer adds losses (accessible via the `losses` property) representing the divergences of kernel and/or bias surrogate posteriors and their respective priors. When doing minibatch stochastic optimization, make sure to scale this loss such that it is applied just once per epoch (e.g. if kl is the sum of losses for each element of the batch, you should pass kl

/ num_examples_per_epoch to your optimizer). You can access the kernel and/or bias posterior and prior distributions after the layer is built via the kernel_posterior, kernel_prior, bias_posterior and bias_prior properties.

Value

a Keras layer

References

- Yeming Wen, Paul Vicol, Jimmy Ba, Dustin Tran, and Roger Grosse. Flipout: Efficient Pseudo-Independent Weight Perturbations on Mini-Batches. In *International Conference on Learning Representations*, 2018.

See Also

Other layers: [layer_autoregressive\(\)](#), [layer_conv_1d_flipout\(\)](#), [layer_conv_1d_reparameterization\(\)](#), [layer_conv_2d_flipout\(\)](#), [layer_conv_2d_reparameterization\(\)](#), [layer_conv_3d_reparameterization\(\)](#), [layer_dense_flipout\(\)](#), [layer_dense_local_reparameterization\(\)](#), [layer_dense_reparameterization\(\)](#), [layer_dense_variational\(\)](#), [layer_variable\(\)](#)

layer_conv_3d_reparameterization

3D convolution layer (e.g. spatial convolution over volumes)

Description

This layer creates a convolution kernel that is convolved (actually cross-correlated) with the layer input to produce a tensor of outputs. It may also include a bias addition and activation function on the outputs. It assumes the kernel and/or bias are drawn from distributions.

Usage

```
layer_conv_3d_reparameterization(
    object,
    filters,
    kernel_size,
    strides = 1,
    padding = "valid",
    data_format = "channels_last",
    dilation_rate = 1,
    activation = NULL,
    activity_regularizer = NULL,
    trainable = TRUE,
    kernel_posterior_fn = tfp$layers$util$default_mean_field_normal_fn(),
    kernel_posterior_tensor_fn = function(d) d %>% tfd_sample(),
    kernel_prior_fn = tfp$layers$util$default_multivariate_normal_fn,
```

```

kernel_divergence_fn = function(q, p, ignore) tfd_kl_divergence(q, p),
bias_posterior_fn = tfp$layers$util$default_mean_field_normal_fn(is_singular = TRUE),
bias_posterior_tensor_fn = function(d) d %>% tfd_sample(),
bias_prior_fn = NULL,
bias_divergence_fn = function(q, p, ignore) tfd_kl_divergence(q, p),
...
)

```

Arguments

object	What to compose the new Layer instance with. Typically a Sequential model or a Tensor (e.g., as returned by <code>layer_input()</code>). The return value depends on object. If object is: <ul style="list-style-type: none"> • missing or NULL, the Layer instance is returned. • a Sequential model, the model with an additional layer is returned. • a Tensor, the output tensor from <code>layer_instance(object)</code> is returned.
filters	Integer, the dimensionality of the output space (i.e. the number of filters in the convolution).
kernel_size	An integer or list of a single integer, specifying the length of the 1D convolution window.
strides	An integer or list of a single integer, specifying the stride length of the convolution. Specifying any stride value $\neq 1$ is incompatible with specifying any <code>dilation_rate</code> value $\neq 1$.
padding	One of "valid" or "same" (case-insensitive).
data_format	A string, one of <code>channels_last</code> (default) or <code>channels_first</code> . The ordering of the dimensions in the inputs. <code>channels_last</code> corresponds to inputs with shape (batch, length, channels) while <code>channels_first</code> corresponds to inputs with shape (batch, channels, length).
dilation_rate	An integer or tuple/list of a single integer, specifying the dilation rate to use for dilated convolution. Currently, specifying any <code>dilation_rate</code> value $\neq 1$ is incompatible with specifying any <code>strides</code> value $\neq 1$.
activation	Activation function. Set it to None to maintain a linear activation.
activity_regularizer	Regularizer function for the output.
trainable	Whether the layer weights will be updated during training.
kernel_posterior_fn	Function which creates <code>tfd\$Distribution</code> instance representing the surrogate posterior of the kernel parameter. Default value: <code>default_mean_field_normal_fn()</code> .
kernel_posterior_tensor_fn	Function which takes a <code>tfd\$Distribution</code> instance and returns a representative value. Default value: <code>function(d) d %>% tfd_sample()</code> .
kernel_prior_fn	Function which creates <code>tfd\$Distribution</code> instance. See <code>default_mean_field_normal_fn</code> docstring for required parameter signature. Default value: <code>tfd_normal(loc = 0, scale = 1)</code> .

kernel_divergence_fn	Function which takes the surrogate posterior distribution, prior distribution and random variate sample(s) from the surrogate posterior and computes or approximates the KL divergence. The distributions are tfd\$Distribution-like instances and the sample is a Tensor.
bias_posterior_fn	Function which creates a tfd\$Distribution instance representing the surrogate posterior of the bias parameter. Default value: default_mean_field_normal_fn(is_singular = TRUE) (which creates an instance of tfd_deterministic).
bias_posterior_tensor_fn	Function which takes a tfd\$Distribution instance and returns a representative value. Default value: function(d) d %>% tfd_sample().
bias_prior_fn	Function which creates tfd instance. See default_mean_field_normal_fn docstring for required parameter signature. Default value: NULL (no prior, no variational inference)
bias_divergence_fn	Function which takes the surrogate posterior distribution, prior distribution and random variate sample(s) from the surrogate posterior and computes or approximates the KL divergence. The distributions are tfd\$Distribution-like instances and the sample is a Tensor.
...	Additional keyword arguments passed to the keras::layer_dense constructed by this layer.

Details

This layer implements the Bayesian variational inference analogue to a dense layer by assuming the kernel and/or the bias are drawn from distributions.

By default, the layer implements a stochastic forward pass via sampling from the kernel and bias posteriors,

$$\text{outputs} = f(\text{inputs}; \text{kernel}, \text{bias}), \text{kernel}, \text{bias} \sim \text{posterior}$$

where f denotes the layer's calculation. It uses the reparameterization estimator (Kingma and Welling, 2014), which performs a Monte Carlo approximation of the distribution integrating over the kernel and bias.

The arguments permit separate specification of the surrogate posterior ($q(W|x)$), prior ($p(W)$), and divergence for both the kernel and bias distributions.

Upon being built, this layer adds losses (accessible via the losses property) representing the divergences of kernel and/or bias surrogate posteriors and their respective priors. When doing minibatch stochastic optimization, make sure to scale this loss such that it is applied just once per epoch (e.g. if kl is the sum of losses for each element of the batch, you should pass $kl / \text{num_examples_per_epoch}$ to your optimizer). You can access the kernel and/or bias posterior and prior distributions after the layer is built via the kernel_posterior, kernel_prior, bias_posterior and bias_prior properties.

Value

a Keras layer

References

- Diederik Kingma and Max Welling. Auto-Encoding Variational Bayes. In *International Conference on Learning Representations*, 2014.

See Also

Other layers: `layer_autoregressive()`, `layer_conv_1d_flipout()`, `layer_conv_1d_reparameterization()`, `layer_conv_2d_flipout()`, `layer_conv_2d_reparameterization()`, `layer_conv_3d_flipout()`, `layer_dense_flipout()`, `layer_dense_local_reparameterization()`, `layer_dense_reparameterization()`, `layer_dense_variational()`, `layer_variable()`

layer_dense_flipout *Densely-connected layer class with Flipout estimator.*

Description

This layer implements the Bayesian variational inference analogue to a dense layer by assuming the kernel and/or the bias are drawn from distributions.

Usage

```
layer_dense_flipout(
  object,
  units,
  activation = NULL,
  activity_regularizer = NULL,
  trainable = TRUE,
  kernel_posterior_fn = tfp$layers$util$default_mean_field_normal_fn(),
  kernel_posterior_tensor_fn = function(d) d %>% tfd_sample(),
  kernel_prior_fn = tfp$layers$util$default_multivariate_normal_fn(),
  kernel_divergence_fn = function(q, p, ignore) tfd_kl_divergence(q, p),
  bias_posterior_fn = tfp$layers$util$default_mean_field_normal_fn(is_singular = TRUE),
  bias_posterior_tensor_fn = function(d) d %>% tfd_sample(),
  bias_prior_fn = NULL,
  bias_divergence_fn = function(q, p, ignore) tfd_kl_divergence(q, p),
  seed = NULL,
  ...
)
```

Arguments

object What to compose the new Layer instance with. Typically a Sequential model or a Tensor (e.g., as returned by `layer_input()`). The return value depends on object. If object is:

- missing or NULL, the Layer instance is returned.
- a Sequential model, the model with an additional layer is returned.

- a Tensor, the output tensor from `layer_instance(object)` is returned.

<code>units</code>	integer dimensionality of the output space
<code>activation</code>	Activation function. Set it to <code>None</code> to maintain a linear activation.
<code>activity_regularizer</code>	Regularizer function for the output.
<code>trainable</code>	Whether the layer weights will be updated during training.
<code>kernel_posterior_fn</code>	Function which creates <code>tfd\$Distribution</code> instance representing the surrogate posterior of the kernel parameter. Default value: <code>default_mean_field_normal_fn()</code> .
<code>kernel_posterior_tensor_fn</code>	Function which takes a <code>tfd\$Distribution</code> instance and returns a representative value. Default value: <code>function(d) d %>% tfd_sample()</code> .
<code>kernel_prior_fn</code>	Function which creates <code>tfd\$Distribution</code> instance. See <code>default_mean_field_normal_fn</code> docstring for required parameter signature. Default value: <code>tfd_normal(loc = 0, scale = 1)</code> .
<code>kernel_divergence_fn</code>	Function which takes the surrogate posterior distribution, prior distribution and random variate sample(s) from the surrogate posterior and computes or approximates the KL divergence. The distributions are <code>tfd\$Distribution</code> -like instances and the sample is a Tensor.
<code>bias_posterior_fn</code>	Function which creates a <code>tfd\$Distribution</code> instance representing the surrogate posterior of the bias parameter. Default value: <code>default_mean_field_normal_fn(is_singular = TRUE)</code> (which creates an instance of <code>tfd_deterministic</code>).
<code>bias_posterior_tensor_fn</code>	Function which takes a <code>tfd\$Distribution</code> instance and returns a representative value. Default value: <code>function(d) d %>% tfd_sample()</code> .
<code>bias_prior_fn</code>	Function which creates <code>tfd</code> instance. See <code>default_mean_field_normal_fn</code> docstring for required parameter signature. Default value: <code>NULL</code> (no prior, no variational inference)
<code>bias_divergence_fn</code>	Function which takes the surrogate posterior distribution, prior distribution and random variate sample(s) from the surrogate posterior and computes or approximates the KL divergence. The distributions are <code>tfd\$Distribution</code> -like instances and the sample is a Tensor.
<code>seed</code>	scalar integer which initializes the random number generator. Default value: <code>NULL</code> (i.e., use global seed).
<code>...</code>	Additional keyword arguments passed to the <code>keras::layer_dense</code> constructed by this layer.

Details

By default, the layer implements a stochastic forward pass via sampling from the kernel and bias posteriors,

```
kernel, bias ~ posterior
outputs = activation(matmul(inputs, kernel) + bias)
```

It uses the Flipout estimator (Wen et al., 2018), which performs a Monte Carlo approximation of the distribution integrating over the kernel and bias. Flipout uses roughly twice as many floating point operations as the reparameterization estimator but has the advantage of significantly lower variance.

The arguments permit separate specification of the surrogate posterior ($q(W|x)$), prior ($p(W)$), and divergence for both the kernel and bias distributions.

Upon being built, this layer adds losses (accessible via the `losses` property) representing the divergences of kernel and/or bias surrogate posteriors and their respective priors. When doing minibatch stochastic optimization, make sure to scale this loss such that it is applied just once per epoch (e.g. if `kl` is the sum of losses for each element of the batch, you should pass `kl / num_examples_per_epoch` to your optimizer).

Value

a Keras layer

References

- Yeming Wen, Paul Vicol, Jimmy Ba, Dustin Tran, and Roger Grosse. Flipout: Efficient Pseudo-Independent Weight Perturbations on Mini-Batches. In *International Conference on Learning Representations*, 2018.

See Also

Other layers: [layer_autoregressive\(\)](#), [layer_conv_1d_flipout\(\)](#), [layer_conv_1d_reparameterization\(\)](#), [layer_conv_2d_flipout\(\)](#), [layer_conv_2d_reparameterization\(\)](#), [layer_conv_3d_flipout\(\)](#), [layer_conv_3d_reparameterization\(\)](#), [layer_dense_local_reparameterization\(\)](#), [layer_dense_reparameterization\(\)](#), [layer_dense_variational\(\)](#), [layer_variable\(\)](#)

layer_dense_local_reparameterization

Densely-connected layer class with local reparameterization estimator.

Description

This layer implements the Bayesian variational inference analogue to a dense layer by assuming the kernel and/or the bias are drawn from distributions.

Usage

```

layer_dense_local_reparameterization(
  object,
  units,
  activation = NULL,
  activity_regularizer = NULL,
  trainable = TRUE,
  kernel_posterior_fn = tfp$layers$util$default_mean_field_normal_fn(),
  kernel_posterior_tensor_fn = function(d) d %>% tfd_sample(),
  kernel_prior_fn = tfp$layers$util$default_multivariate_normal_fn,
  kernel_divergence_fn = function(q, p, ignore) tfd_kl_divergence(q, p),
  bias_posterior_fn = tfp$layers$util$default_mean_field_normal_fn(is_singular = TRUE),
  bias_posterior_tensor_fn = function(d) d %>% tfd_sample(),
  bias_prior_fn = NULL,
  bias_divergence_fn = function(q, p, ignore) tfd_kl_divergence(q, p),
  ...
)

```

Arguments

object	What to compose the new Layer instance with. Typically a Sequential model or a Tensor (e.g., as returned by layer_input()). The return value depends on object. If object is: <ul style="list-style-type: none"> • missing or NULL, the Layer instance is returned. • a Sequential model, the model with an additional layer is returned. • a Tensor, the output tensor from layer_instance(object) is returned.
units	integer dimensionality of the output space
activation	Activation function. Set it to None to maintain a linear activation.
activity_regularizer	Regularizer function for the output.
trainable	Whether the layer weights will be updated during training.
kernel_posterior_fn	Function which creates tfd\$Distribution instance representing the surrogate posterior of the kernel parameter. Default value: default_mean_field_normal_fn().
kernel_posterior_tensor_fn	Function which takes a tfd\$Distribution instance and returns a representative value. Default value: function(d) d %>% tfd_sample().
kernel_prior_fn	Function which creates tfd\$Distribution instance. See default_mean_field_normal_fn docstring for required parameter signature. Default value: tfd_normal(loc = 0, scale = 1).
kernel_divergence_fn	Function which takes the surrogate posterior distribution, prior distribution and random variate sample(s) from the surrogate posterior and computes or approximates the KL divergence. The distributions are tfd\$Distribution-like instances and the sample is a Tensor.

bias_posterior_fn	Function which creates a tfd\$Distribution instance representing the surrogate posterior of the bias parameter. Default value: default_mean_field_normal_fn(is_singular = TRUE) (which creates an instance of tfd_deterministic).
bias_posterior_tensor_fn	Function which takes a tfd\$Distribution instance and returns a representative value. Default value: function(d) d %>% tfd_sample().
bias_prior_fn	Function which creates tfd instance. See default_mean_field_normal_fn docstring for required parameter signature. Default value: NULL (no prior, no variational inference)
bias_divergence_fn	Function which takes the surrogate posterior distribution, prior distribution and random variate sample(s) from the surrogate posterior and computes or approximates the KL divergence. The distributions are tfd\$Distribution-like instances and the sample is a Tensor.
...	Additional keyword arguments passed to the keras::layer_dense constructed by this layer.

Details

By default, the layer implements a stochastic forward pass via sampling from the kernel and bias posteriors,

```
kernel, bias ~ posterior
outputs = activation(matmul(inputs, kernel) + bias)
```

It uses the local reparameterization estimator (Kingma et al., 2015), which performs a Monte Carlo approximation of the distribution on the hidden units induced by the kernel and bias. The default kernel_posterior_fn is a normal distribution which factorizes across all elements of the weight matrix and bias vector. Unlike that paper's multiplicative parameterization, this distribution has trainable location and scale parameters which is known as an additive noise parameterization (Molchanov et al., 2017).

The arguments permit separate specification of the surrogate posterior ($q(W|x)$), prior ($p(W)$), and divergence for both the kernel and bias distributions.

Upon being built, this layer adds losses (accessible via the losses property) representing the divergences of kernel and/or bias surrogate posteriors and their respective priors. When doing minibatch stochastic optimization, make sure to scale this loss such that it is applied just once per epoch (e.g. if kl is the sum of losses for each element of the batch, you should pass $kl / \text{num_examples_per_epoch}$ to your optimizer). You can access the kernel and/or bias posterior and prior distributions after the layer is built via the kernel_posterior, kernel_prior, bias_posterior and bias_prior properties.

Value

a Keras layer

References

- Diederik Kingma, Tim Salimans, and Max Welling. Variational Dropout and the Local Reparameterization Trick. In *Neural Information Processing Systems*, 2015.
- Dmitry Molchanov, Arsenii Ashukha, Dmitry Vetrov. Variational Dropout Sparsifies Deep Neural Networks. In *International Conference on Machine Learning*, 2017.

See Also

Other layers: `layer_autoregressive()`, `layer_conv_1d_flipout()`, `layer_conv_1d_reparameterization()`, `layer_conv_2d_flipout()`, `layer_conv_2d_reparameterization()`, `layer_conv_3d_flipout()`, `layer_conv_3d_reparameterization()`, `layer_dense_flipout()`, `layer_dense_reparameterization()`, `layer_dense_variational()`, `layer_variable()`

layer_dense_reparameterization

Densely-connected layer class with reparameterization estimator.

Description

This layer implements the Bayesian variational inference analogue to a dense layer by assuming the kernel and/or the bias are drawn from distributions.

Usage

```
layer_dense_reparameterization(
    object,
    units,
    activation = NULL,
    activity_regularizer = NULL,
    trainable = TRUE,
    kernel_posterior_fn = tfp$layers$util$default_mean_field_normal_fn(),
    kernel_posterior_tensor_fn = function(d) d %>% tfd_sample(),
    kernel_prior_fn = tfp$layers$util$default_multivariate_normal_fn,
    kernel_divergence_fn = function(q, p, ignore) tfd_kl_divergence(q, p),
    bias_posterior_fn = tfp$layers$util$default_mean_field_normal_fn(is_singular = TRUE),
    bias_posterior_tensor_fn = function(d) d %>% tfd_sample(),
    bias_prior_fn = NULL,
    bias_divergence_fn = function(q, p, ignore) tfd_kl_divergence(q, p),
    ...
)
```

Arguments

object What to compose the new Layer instance with. Typically a Sequential model or a Tensor (e.g., as returned by `layer_input()`). The return value depends on object. If object is:

- missing or NULL, the Layer instance is returned.
- a Sequential model, the model with an additional layer is returned.
- a Tensor, the output tensor from layer_instance(object) is returned.

units integer dimensionality of the output space

activation Activation function. Set it to None to maintain a linear activation.

activity_regularizer Regularizer function for the output.

trainable Whether the layer weights will be updated during training.

kernel_posterior_fn Function which creates tfd\$Distribution instance representing the surrogate posterior of the kernel parameter. Default value: default_mean_field_normal_fn().

kernel_posterior_tensor_fn Function which takes a tfd\$Distribution instance and returns a representative value. Default value: function(d) d %>% tfd_sample().

kernel_prior_fn Function which creates tfd\$Distribution instance. See default_mean_field_normal_fn docstring for required parameter signature. Default value: tfd_normal(loc = 0, scale = 1).

kernel_divergence_fn Function which takes the surrogate posterior distribution, prior distribution and random variate sample(s) from the surrogate posterior and computes or approximates the KL divergence. The distributions are tfd\$Distribution-like instances and the sample is a Tensor.

bias_posterior_fn Function which creates a tfd\$Distribution instance representing the surrogate posterior of the bias parameter. Default value: default_mean_field_normal_fn(is_singular = TRUE) (which creates an instance of tfd_deterministic).

bias_posterior_tensor_fn Function which takes a tfd\$Distribution instance and returns a representative value. Default value: function(d) d %>% tfd_sample().

bias_prior_fn Function which creates tfd instance. See default_mean_field_normal_fn docstring for required parameter signature. Default value: NULL (no prior, no variational inference)

bias_divergence_fn Function which takes the surrogate posterior distribution, prior distribution and random variate sample(s) from the surrogate posterior and computes or approximates the KL divergence. The distributions are tfd\$Distribution-like instances and the sample is a Tensor.

... Additional keyword arguments passed to the keras::layer_dense constructed by this layer.

Details

By default, the layer implements a stochastic forward pass via sampling from the kernel and bias posteriors,

```
kernel, bias ~ posterior
outputs = activation(matmul(inputs, kernel) + bias)
```

It uses the reparameterization estimator (Kingma and Welling, 2014) which performs a Monte Carlo approximation of the distribution integrating over the kernel and bias.

The arguments permit separate specification of the surrogate posterior ($q(W|x)$), prior ($p(W)$), and divergence for both the kernel and bias distributions.

Upon being built, this layer adds losses (accessible via the losses property) representing the divergences of kernel and/or bias surrogate posteriors and their respective priors. When doing minibatch stochastic optimization, make sure to scale this loss such that it is applied just once per epoch (e.g. if kl is the sum of losses for each element of the batch, you should pass $kl / \text{num_examples_per_epoch}$ to your optimizer). You can access the kernel and/or bias posterior and prior distributions after the layer is built via the kernel_posterior, kernel_prior, bias_posterior and bias_prior properties.

Value

a Keras layer

References

- Diederik Kingma and Max Welling. Auto-Encoding Variational Bayes. In *International Conference on Learning Representations*, 2014.

See Also

Other layers: [layer_autoregressive\(\)](#), [layer_conv_1d_flipout\(\)](#), [layer_conv_1d_reparameterization\(\)](#), [layer_conv_2d_flipout\(\)](#), [layer_conv_2d_reparameterization\(\)](#), [layer_conv_3d_flipout\(\)](#), [layer_conv_3d_reparameterization\(\)](#), [layer_dense_flipout\(\)](#), [layer_dense_local_reparameterization\(\)](#), [layer_dense_variational\(\)](#), [layer_variable\(\)](#)

layer_dense_variational

Dense Variational Layer

Description

This layer uses variational inference to fit a "surrogate" posterior to the distribution over both the kernel matrix and the bias terms which are otherwise used in a manner similar to `layer_dense()`. This layer fits the "weights posterior" according to the following generative process:

```
[K, b] ~ Prior()
M = matmul(X, K) + b
Y ~ Likelihood(M)
```

Usage

```
layer_dense_variational(
    object,
    units,
    make_posterior_fn,
    make_prior_fn,
    kl_weight = NULL,
    kl_use_exact = FALSE,
    activation = NULL,
    use_bias = TRUE,
    ...
)
```

Arguments

object	What to compose the new Layer instance with. Typically a Sequential model or a Tensor (e.g., as returned by <code>layer_input()</code>). The return value depends on object. If object is: <ul style="list-style-type: none"> • missing or NULL, the Layer instance is returned. • a Sequential model, the model with an additional layer is returned. • a Tensor, the output tensor from <code>layer_instance(object)</code> is returned.
units	Positive integer, dimensionality of the output space.
make_posterior_fn	function taking <code>tf\$size(kernel)</code> , <code>tf\$size(bias)</code> , <code>dtype</code> and returns another callable which takes an input and produces a <code>tfd\$Distribution</code> instance.
make_prior_fn	function taking <code>tf\$size(kernel)</code> , <code>tf\$size(bias)</code> , <code>dtype</code> and returns another callable which takes an input and produces a <code>tfd\$Distribution</code> instance.
kl_weight	Amount by which to scale the KL divergence loss between prior and posterior.
kl_use_exact	Logical indicating that the analytical KL divergence should be used rather than a Monte Carlo approximation.
activation	An activation function. See <code>keras::layer_dense</code> . Default: NULL.
use_bias	Whether or not the dense layers constructed in this layer should have a bias term. See <code>keras::layer_dense</code> . Default: TRUE.
...	Additional keyword arguments passed to the <code>keras::layer_dense</code> constructed by this layer.

Value

a Keras layer

See Also

Other layers: [layer_autoregressive\(\)](#), [layer_conv_1d_flipout\(\)](#), [layer_conv_1d_reparameterization\(\)](#), [layer_conv_2d_flipout\(\)](#), [layer_conv_2d_reparameterization\(\)](#), [layer_conv_3d_flipout\(\)](#), [layer_conv_3d_reparameterization\(\)](#), [layer_dense_flipout\(\)](#), [layer_dense_local_reparameterization\(\)](#), [layer_dense_reparameterization\(\)](#), [layer_variable\(\)](#)

 layer_distribution_lambda

Keras layer enabling plumbing TFP distributions through Keras models

Description

Keras layer enabling plumbing TFP distributions through Keras models

Usage

```
layer_distribution_lambda(
    object,
    make_distribution_fn,
    convert_to_tensor_fn = tfp$distributions$Distribution$sample,
    ...
)
```

Arguments

object	What to compose the new Layer instance with. Typically a Sequential model or a Tensor (e.g., as returned by <code>layer_input()</code>). The return value depends on object. If object is: <ul style="list-style-type: none"> • missing or NULL, the Layer instance is returned. • a Sequential model, the model with an additional layer is returned. • a Tensor, the output tensor from <code>layer_instance(object)</code> is returned.
make_distribution_fn	A callable that takes previous layer outputs and returns a <code>tfd\$Distributions\$Distribution</code> instance.
convert_to_tensor_fn	A callable that takes a <code>tfd\$Distribution</code> instance and returns a <code>tf\$Tensor</code> -like object. Default value: <code>tfd\$distributions\$Distribution\$sample</code> .
...	Additional arguments passed to <code>args of keras::create_layer</code> .

Value

a Keras layer

See Also

For an example how to use in a Keras model, see [layer_independent_normal\(\)](#).

Other `distribution_layers`: [layer_categorical_mixture_of_one_hot_categorical\(\)](#), [layer_independent_bernoulli](#), [layer_independent_logistic\(\)](#), [layer_independent_normal\(\)](#), [layer_independent_poisson\(\)](#), [layer_kl_divergence_add_loss\(\)](#), [layer_kl_divergence_regularizer\(\)](#), [layer_mixture_logistic\(\)](#), [layer_mixture_normal\(\)](#), [layer_mixture_same_family\(\)](#), [layer_multivariate_normal_tri_l\(\)](#), [layer_one_hot_categorical\(\)](#)

layer_independent_bernoulli

An Independent-Bernoulli Keras layer from prod(event_shape) params

Description

An Independent-Bernoulli Keras layer from prod(event_shape) params

Usage

```
layer_independent_bernoulli(
    object,
    event_shape,
    convert_to_tensor_fn = tfp$distributions$Distribution$sample,
    sample_dtype = NULL,
    validate_args = FALSE,
    ...
)
```

Arguments

object	What to compose the new Layer instance with. Typically a Sequential model or a Tensor (e.g., as returned by layer_input()). The return value depends on object. If object is: <ul style="list-style-type: none"> • missing or NULL, the Layer instance is returned. • a Sequential model, the model with an additional layer is returned. • a Tensor, the output tensor from layer_instance(object) is returned.
event_shape	Scalar integer representing the size of single draw from this distribution.
convert_to_tensor_fn	A callable that takes a tfd\$Distribution instance and returns a tf\$Tensor-like object. Default value: tfd\$distributions\$Distribution\$sample.
sample_dtype	dtype of samples produced by this distribution. Default value: NULL (i.e., previous layer's dtype).
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE. @param ...
...	Additional arguments passed to args of keras::create_layer.

Value

a Keras layer

See Also

For an example how to use in a Keras model, see [layer_independent_normal\(\)](#).

Other distribution_layers: [layer_categorical_mixture_of_one_hot_categorical\(\)](#), [layer_distribution_lambda\(\)](#), [layer_independent_logistic\(\)](#), [layer_independent_normal\(\)](#), [layer_independent_poisson\(\)](#), [layer_kl_divergence_add_loss\(\)](#), [layer_kl_divergence_regularizer\(\)](#), [layer_mixture_logistic\(\)](#), [layer_mixture_normal\(\)](#), [layer_mixture_same_family\(\)](#), [layer_multivariate_normal_tri_l\(\)](#), [layer_one_hot_categorical\(\)](#)

layer_independent_logistic

An independent Logistic Keras layer.

Description

An independent Logistic Keras layer.

Usage

```
layer_independent_logistic(
    object,
    event_shape,
    convert_to_tensor_fn = tfp$distributions$Distribution$sample,
    validate_args = FALSE,
    ...
)
```

Arguments

object	What to compose the new Layer instance with. Typically a Sequential model or a Tensor (e.g., as returned by <code>layer_input()</code>). The return value depends on object. If object is: <ul style="list-style-type: none"> • missing or NULL, the Layer instance is returned. • a Sequential model, the model with an additional layer is returned. • a Tensor, the output tensor from <code>layer_instance(object)</code> is returned.
event_shape	Scalar integer representing the size of single draw from this distribution.
convert_to_tensor_fn	A callable that takes a <code>tfd\$Distribution</code> instance and returns a <code>tf\$Tensor</code> -like object. Default value: <code>tfd\$distributions\$Distribution\$sample</code> .
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE. @param ... Additional arguments passed to <code>args</code> of <code>keras::create_layer</code> .
...	Additional arguments passed to <code>args</code> of <code>keras::create_layer</code> .

Value

a Keras layer

See Also

For an example how to use in a Keras model, see [layer_independent_normal\(\)](#).

Other distribution_layers: [layer_categorical_mixture_of_one_hot_categorical\(\)](#), [layer_distribution_lambda\(\)](#), [layer_independent_bernoulli\(\)](#), [layer_independent_normal\(\)](#), [layer_independent_poisson\(\)](#), [layer_kl_divergence_add_loss\(\)](#), [layer_kl_divergence_regularizer\(\)](#), [layer_mixture_logistic\(\)](#), [layer_mixture_normal\(\)](#), [layer_mixture_same_family\(\)](#), [layer_multivariate_normal_tri_l\(\)](#), [layer_one_hot_categorical\(\)](#)

layer_independent_normal

An independent Normal Keras layer.

Description

An independent Normal Keras layer.

Usage

```
layer_independent_normal(
    object,
    event_shape,
    convert_to_tensor_fn = tfp$distributions$Distribution$sample,
    validate_args = FALSE,
    ...
)
```

Arguments

object	What to compose the new Layer instance with. Typically a Sequential model or a Tensor (e.g., as returned by <code>layer_input()</code>). The return value depends on object. If object is: <ul style="list-style-type: none"> • missing or NULL, the Layer instance is returned. • a Sequential model, the model with an additional layer is returned. • a Tensor, the output tensor from <code>layer_instance(object)</code> is returned.
event_shape	Scalar integer representing the size of single draw from this distribution.
convert_to_tensor_fn	A callable that takes a <code>tfd\$Distribution</code> instance and returns a <code>tf\$Tensor</code> -like object. Default value: <code>tfd\$distributions\$Distribution\$sample</code> .
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE. @param ...
...	Additional arguments passed to <code>args</code> of <code>keras::create_layer</code> .

Value

a Keras layer

See Also

Other distribution_layers: [layer_categorical_mixture_of_one_hot_categorical\(\)](#), [layer_distribution_lambda\(\)](#), [layer_independent_bernoulli\(\)](#), [layer_independent_logistic\(\)](#), [layer_independent_poisson\(\)](#), [layer_kl_divergence_add_loss\(\)](#), [layer_kl_divergence_regularizer\(\)](#), [layer_mixture_logistic\(\)](#), [layer_mixture_normal\(\)](#), [layer_mixture_same_family\(\)](#), [layer_multivariate_normal_tri_l\(\)](#), [layer_one_hot_categorical\(\)](#)

Examples

```
library(keras)
input_shape <- c(28, 28, 1)
encoded_shape <- 2
n <- 2
model <- keras_model_sequential(
  list(
    layer_input(shape = input_shape),
    layer_flatten(),
    layer_dense(units = n),
    layer_dense(units = params_size_independent_normal(encoded_shape)),
    layer_independent_normal(event_shape = encoded_shape)
  )
)
```

layer_independent_poisson

An independent Poisson Keras layer.

Description

An independent Poisson Keras layer.

Usage

```
layer_independent_poisson(
  object,
  event_shape,
  convert_to_tensor_fn = tfp$distributions$Distribution$sample,
  validate_args = FALSE,
  ...
)
```

Arguments

object	What to compose the new Layer instance with. Typically a Sequential model or a Tensor (e.g., as returned by <code>layer_input()</code>). The return value depends on object. If object is: <ul style="list-style-type: none"> • missing or NULL, the Layer instance is returned. • a Sequential model, the model with an additional layer is returned. • a Tensor, the output tensor from <code>layer_instance(object)</code> is returned.
event_shape	Scalar integer representing the size of single draw from this distribution.
convert_to_tensor_fn	A callable that takes a <code>tfd\$Distribution</code> instance and returns a <code>tf\$Tensor</code> -like object. Default value: <code>tfd\$distributions\$Distribution\$sample</code> .
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE. @param ... Additional arguments passed to <code>args</code> of <code>keras::create_layer</code> .
...	Additional arguments passed to <code>args</code> of <code>keras::create_layer</code> .

Value

a Keras layer

See Also

For an example how to use in a Keras model, see [layer_independent_normal\(\)](#).

Other `distribution_layers`: [layer_categorical_mixture_of_one_hot_categorical\(\)](#), [layer_distribution_lambda\(\)](#), [layer_independent_bernoulli\(\)](#), [layer_independent_logistic\(\)](#), [layer_independent_normal\(\)](#), [layer_kl_divergence_add_loss\(\)](#), [layer_kl_divergence_regularizer\(\)](#), [layer_mixture_logistic\(\)](#), [layer_mixture_normal\(\)](#), [layer_mixture_same_family\(\)](#), [layer_multivariate_normal_tri_l\(\)](#), [layer_one_hot_categorical\(\)](#)

layer_kl_divergence_add_loss

Pass-through layer that adds a KL divergence penalty to the model loss

Description

Pass-through layer that adds a KL divergence penalty to the model loss

Usage

```
layer_kl_divergence_add_loss(
    object,
    distribution_b,
    use_exact_kl = FALSE,
    test_points_reduce_axis = NULL,
    test_points_fn = tf$convert_to_tensor,
    weight = NULL,
    ...
)
```

Arguments

<code>object</code>	What to compose the new Layer instance with. Typically a Sequential model or a Tensor (e.g., as returned by <code>layer_input()</code>). The return value depends on object. If object is: <ul style="list-style-type: none"> • missing or NULL, the Layer instance is returned. • a Sequential model, the model with an additional layer is returned. • a Tensor, the output tensor from <code>layer_instance(object)</code> is returned.
<code>distribution_b</code>	Distribution instance corresponding to b as in $KL[a,b]$. The previous layer's output is presumed to be a Distribution instance and is a.
<code>use_exact_kl</code>	Logical indicating if KL divergence should be calculated exactly via <code>tfp\$distributions\$kl_divergence</code> or via Monte Carlo approximation. Default value: FALSE.
<code>test_points_reduce_axis</code>	Integer vector or scalar representing dimensions over which to reduce_mean while calculating the Monte Carlo approximation of the KL divergence. As is with all <code>tf\$reduce_*</code> ops, NULL means reduce over all dimensions; () means reduce over none of them. Default value: () (i.e., no reduction).
<code>test_points_fn</code>	A callable taking a <code>tfp\$distributions\$Distribution</code> instance and returning a tensor used for random test points to approximate the KL divergence. Default value: <code>tf\$convert_to_tensor</code> .
<code>weight</code>	Multiplier applied to the calculated KL divergence for each Keras batch member. Default value: NULL (i.e., do not weight each batch member).
<code>...</code>	Additional arguments passed to args of <code>keras::create_layer</code> .

Value

a Keras layer

See Also

For an example how to use in a Keras model, see [layer_independent_normal\(\)](#).

Other `distribution_layers`: [layer_categorical_mixture_of_one_hot_categorical\(\)](#), [layer_distribution_lambda\(\)](#), [layer_independent_bernoulli\(\)](#), [layer_independent_logistic\(\)](#), [layer_independent_normal\(\)](#), [layer_independent_poisson\(\)](#), [layer_kl_divergence_regularizer\(\)](#), [layer_mixture_logistic\(\)](#), [layer_mixture_normal\(\)](#), [layer_mixture_same_family\(\)](#), [layer_multivariate_normal_tri_l\(\)](#), [layer_one_hot_categorical\(\)](#)

 layer_kl_divergence_regularizer

Regularizer that adds a KL divergence penalty to the model loss

Description

When using Monte Carlo approximation (e.g., `use_exact = FALSE`), it is presumed that the input distribution's concretization (i.e., `tf$convert_to_tensor(distribution)`) corresponds to a random sample. To override this behavior, set `test_points_fn`.

Usage

```
layer_kl_divergence_regularizer(
  object,
  distribution_b,
  use_exact_kl = FALSE,
  test_points_reduce_axis = NULL,
  test_points_fn = tf$convert_to_tensor,
  weight = NULL,
  ...
)
```

Arguments

<code>object</code>	What to compose the new <code>Layer</code> instance with. Typically a <code>Sequential</code> model or a <code>Tensor</code> (e.g., as returned by <code>layer_input()</code>). The return value depends on <code>object</code> . If <code>object</code> is: <ul style="list-style-type: none"> • missing or <code>NULL</code>, the <code>Layer</code> instance is returned. • a <code>Sequential</code> model, the model with an additional layer is returned. • a <code>Tensor</code>, the output tensor from <code>layer_instance(object)</code> is returned.
<code>distribution_b</code>	Distribution instance corresponding to <code>b</code> as in $KL[a,b]$. The previous layer's output is presumed to be a <code>Distribution</code> instance and is <code>a</code> .
<code>use_exact_kl</code>	Logical indicating if KL divergence should be calculated exactly via <code>tfp\$distributions\$kl_divergence</code> or via Monte Carlo approximation. Default value: <code>FALSE</code> .
<code>test_points_reduce_axis</code>	Integer vector or scalar representing dimensions over which to <code>reduce_mean</code> while calculating the Monte Carlo approximation of the KL divergence. As is with all <code>tf\$reduce_*</code> ops, <code>NULL</code> means reduce over all dimensions; <code>()</code> means reduce over none of them. Default value: <code>()</code> (i.e., no reduction).
<code>test_points_fn</code>	A callable taking a <code>tfp\$distributions\$Distribution</code> instance and returning a tensor used for random test points to approximate the KL divergence. Default value: <code>tf\$convert_to_tensor</code> .
<code>weight</code>	Multiplier applied to the calculated KL divergence for each Keras batch member. Default value: <code>NULL</code> (i.e., do not weight each batch member).
<code>...</code>	Additional arguments passed to <code>args</code> of <code>keras::create_layer</code> .

Value

a Keras layer

See Also

For an example how to use in a Keras model, see [layer_independent_normal\(\)](#).

Other distribution_layers: [layer_categorical_mixture_of_one_hot_categorical\(\)](#), [layer_distribution_lambda\(\)](#), [layer_independent_bernoulli\(\)](#), [layer_independent_logistic\(\)](#), [layer_independent_normal\(\)](#), [layer_independent_poisson\(\)](#), [layer_kl_divergence_add_loss\(\)](#), [layer_mixture_logistic\(\)](#), [layer_mixture_normal\(\)](#), [layer_mixture_same_family\(\)](#), [layer_multivariate_normal_tri_l\(\)](#), [layer_one_hot_categorical\(\)](#)

layer_mixture_logistic

A mixture distribution Keras layer, with independent logistic components.

Description

A mixture distribution Keras layer, with independent logistic components.

Usage

```
layer_mixture_logistic(
    object,
    num_components,
    event_shape = list(),
    convert_to_tensor_fn = tfp$distributions$Distribution$sample,
    validate_args = FALSE,
    ...
)
```

Arguments

object	What to compose the new Layer instance with. Typically a Sequential model or a Tensor (e.g., as returned by <code>layer_input()</code>). The return value depends on object. If object is: <ul style="list-style-type: none"> • missing or NULL, the Layer instance is returned. • a Sequential model, the model with an additional layer is returned. • a Tensor, the output tensor from <code>layer_instance(object)</code> is returned.
num_components	Number of component distributions in the mixture distribution.
event_shape	integer vector Tensor representing the shape of single draw from this distribution.
convert_to_tensor_fn	A callable that takes a <code>tfd\$Distribution</code> instance and returns a <code>tf\$Tensor</code> -like object. Default value: <code>tfd\$distributions\$Distribution\$sample</code> .

`validate_args` Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE. @param ... Additional arguments passed to args of `keras::create_layer`.

... Additional arguments passed to args of `keras::create_layer`.

Value

a Keras layer

See Also

For an example how to use in a Keras model, see `layer_independent_normal()`.

Other `distribution_layers`: `layer_categorical_mixture_of_one_hot_categorical()`, `layer_distribution_lambda()`, `layer_independent_bernoulli()`, `layer_independent_logistic()`, `layer_independent_normal()`, `layer_independent_poisson()`, `layer_kl_divergence_add_loss()`, `layer_kl_divergence_regularizer()`, `layer_mixture_normal()`, `layer_mixture_same_family()`, `layer_multivariate_normal_tri_l()`, `layer_one_hot_categorical()`

`layer_mixture_normal` *A mixture distribution Keras layer, with independent normal components.*

Description

A mixture distribution Keras layer, with independent normal components.

Usage

```
layer_mixture_normal(
  object,
  num_components,
  event_shape = list(),
  convert_to_tensor_fn = tfp$distributions$Distribution$sample,
  validate_args = FALSE,
  ...
)
```

Arguments

`object` What to compose the new Layer instance with. Typically a Sequential model or a Tensor (e.g., as returned by `layer_input()`). The return value depends on `object`. If `object` is:

- missing or NULL, the Layer instance is returned.
- a Sequential model, the model with an additional layer is returned.
- a Tensor, the output tensor from `layer_instance(object)` is returned.

num_components	Number of component distributions in the mixture distribution.
event_shape	integer vector Tensor representing the shape of single draw from this distribution.
convert_to_tensor_fn	A callable that takes a tfd\$Distribution instance and returns a tf\$Tensor-like object. Default value: tfd\$distributions\$Distribution\$sample.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE. @param ... Additional arguments passed to args of keras::create_layer.
...	Additional arguments passed to args of keras::create_layer.

Value

a Keras layer

See Also

For an example how to use in a Keras model, see [layer_independent_normal\(\)](#).

Other distribution_layers: [layer_categorical_mixture_of_one_hot_categorical\(\)](#), [layer_distribution_lambda\(\)](#), [layer_independent_bernoulli\(\)](#), [layer_independent_logistic\(\)](#), [layer_independent_normal\(\)](#), [layer_independent_poisson\(\)](#), [layer_kl_divergence_add_loss\(\)](#), [layer_kl_divergence_regularizer\(\)](#), [layer_mixture_logistic\(\)](#), [layer_mixture_same_family\(\)](#), [layer_multivariate_normal_tri_l\(\)](#), [layer_one_hot_categorical\(\)](#)

layer_mixture_same_family

A mixture (same-family) Keras layer.

Description

A mixture (same-family) Keras layer.

Usage

```
layer_mixture_same_family(
  object,
  num_components,
  component_layer,
  convert_to_tensor_fn = tfp$distributions$Distribution$sample,
  validate_args = FALSE,
  ...
)
```

Arguments

object	What to compose the new Layer instance with. Typically a Sequential model or a Tensor (e.g., as returned by <code>layer_input()</code>). The return value depends on object. If object is: <ul style="list-style-type: none"> • missing or NULL, the Layer instance is returned. • a Sequential model, the model with an additional layer is returned. • a Tensor, the output tensor from <code>layer_instance(object)</code> is returned.
num_components	Number of component distributions in the mixture distribution.
component_layer	Function that, given a tensor of shape <code>batch_shape + [num_components, component_params_size]</code> , returns a <code>tfd.Distribution</code> -like instance that implements the component distribution (with batch shape <code>batch_shape + [num_components]</code>) – e.g., a TFP distribution layer.
convert_to_tensor_fn	A callable that takes a <code>tfd.Distribution</code> instance and returns a <code>tf.Tensor</code> -like object. Default value: <code>tfd.distributions.Distribution.sample</code> .
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE. @param ... Additional arguments passed to <code>args</code> of <code>keras::create_layer</code> .
...	Additional arguments passed to <code>args</code> of <code>keras::create_layer</code> .

Value

a Keras layer

See Also

For an example how to use in a Keras model, see [layer_independent_normal\(\)](#).

Other distribution_layers: [layer_categorical_mixture_of_one_hot_categorical\(\)](#), [layer_distribution_lambda\(\)](#), [layer_independent_bernoulli\(\)](#), [layer_independent_logistic\(\)](#), [layer_independent_normal\(\)](#), [layer_independent_poisson\(\)](#), [layer_kl_divergence_add_loss\(\)](#), [layer_kl_divergence_regularizer\(\)](#), [layer_mixture_logistic\(\)](#), [layer_mixture_normal\(\)](#), [layer_multivariate_normal_tri_l\(\)](#), [layer_one_hot_categorical\(\)](#)

layer_multivariate_normal_tri_l

A d-variate Multivariate Normal TriL Keras layer from $d+d(d+1)/2$ params*

Description

A d-variate Multivariate Normal TriL Keras layer from $d+d*(d+1)/2$ params

Usage

```
layer_multivariate_normal_tri_l(
    object,
    event_size,
    convert_to_tensor_fn = tfp$distributions$Distribution$sample,
    validate_args = FALSE,
    ...
)
```

Arguments

object	What to compose the new Layer instance with. Typically a Sequential model or a Tensor (e.g., as returned by <code>layer_input()</code>). The return value depends on object. If object is: <ul style="list-style-type: none"> • missing or NULL, the Layer instance is returned. • a Sequential model, the model with an additional layer is returned. • a Tensor, the output tensor from <code>layer_instance(object)</code> is returned.
event_size	Integer vector tensor representing the shape of single draw from this distribution.
convert_to_tensor_fn	A callable that takes a <code>tfd\$Distribution</code> instance and returns a <code>tf\$Tensor</code> -like object. Default value: <code>tfd\$distributions\$Distribution\$sample</code> .
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
...	Additional arguments passed to <code>args</code> of <code>keras::create_layer</code> .

Value

a Keras layer

See Also

For an example how to use in a Keras model, see [layer_independent_normal\(\)](#).

Other `distribution_layers`: [layer_categorical_mixture_of_one_hot_categorical\(\)](#), [layer_distribution_lambda\(\)](#), [layer_independent_bernoulli\(\)](#), [layer_independent_logistic\(\)](#), [layer_independent_normal\(\)](#), [layer_independent_poisson\(\)](#), [layer_kl_divergence_add_loss\(\)](#), [layer_kl_divergence_regularizer\(\)](#), [layer_mixture_logistic\(\)](#), [layer_mixture_normal\(\)](#), [layer_mixture_same_family\(\)](#), [layer_one_hot_categorical\(\)](#)

layer_one_hot_categorical

A d-variate OneHotCategorical Keras layer from d params.

Description

Typical choices for `convert_to_tensor_fn` include:

- `tfp$distributions$Distribution$sample`
- `tfp$distributions$Distribution$mean`
- `tfp$distributions$Distribution$mode`
- `tfp$distributions$OneHotCategorical$logits`

Usage

```
layer_one_hot_categorical(
  object,
  event_size,
  convert_to_tensor_fn = tfp$distributions$Distribution$sample,
  sample_dtype = NULL,
  validate_args = FALSE,
  ...
)
```

Arguments

<code>object</code>	What to compose the new <code>Layer</code> instance with. Typically a <code>Sequential</code> model or a <code>Tensor</code> (e.g., as returned by <code>layer_input()</code>). The return value depends on <code>object</code> . If <code>object</code> is: <ul style="list-style-type: none"> • missing or <code>NULL</code>, the <code>Layer</code> instance is returned. • a <code>Sequential</code> model, the model with an additional layer is returned. • a <code>Tensor</code>, the output tensor from <code>layer_instance(object)</code> is returned.
<code>event_size</code>	Scalar integer representing the size of single draw from this distribution.
<code>convert_to_tensor_fn</code>	A callable that takes a <code>tfd\$Distribution</code> instance and returns a <code>tf\$Tensor</code> -like object. Default value: <code>tfp\$distributions\$Distribution\$sample</code> .
<code>sample_dtype</code>	<code>dtype</code> of samples produced by this distribution. Default value: <code>NULL</code> (i.e., previous layer's <code>dtype</code>).
<code>validate_args</code>	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
<code>...</code>	Additional arguments passed to <code>args</code> of <code>keras::create_layer</code> .

Value

a Keras layer

See Also

For an example how to use in a Keras model, see [layer_independent_normal\(\)](#).

Other distribution_layers: [layer_categorical_mixture_of_one_hot_categorical\(\)](#), [layer_distribution_lambda\(\)](#), [layer_independent_bernoulli\(\)](#), [layer_independent_logistic\(\)](#), [layer_independent_normal\(\)](#), [layer_independent_poisson\(\)](#), [layer_kl_divergence_add_loss\(\)](#), [layer_kl_divergence_regularizer\(\)](#), [layer_mixture_logistic\(\)](#), [layer_mixture_normal\(\)](#), [layer_mixture_same_family\(\)](#), [layer_multivariate_normal\(\)](#)

layer_variable	<i>Variable Layer</i>
----------------	-----------------------

Description

Simply returns a (trainable) variable, regardless of input. This layer implements the mathematical function $f(x) = c$ where c is a constant, i.e., unchanged for all x . Like other Keras layers, the constant is trainable. This layer can also be interpreted as the special case of [layer_dense\(\)](#) when the kernel is forced to be the zero matrix (`tf.zeros`).

Usage

```
layer_variable(
    object,
    shape,
    dtype = NULL,
    activation = NULL,
    initializer = "zeros",
    regularizer = NULL,
    constraint = NULL,
    ...
)
```

Arguments

object	What to compose the new Layer instance with. Typically a <code>Sequential</code> model or a <code>Tensor</code> (e.g., as returned by <code>layer_input()</code>). The return value depends on object. If object is: <ul style="list-style-type: none"> • missing or <code>NULL</code>, the Layer instance is returned. • a <code>Sequential</code> model, the model with an additional layer is returned. • a <code>Tensor</code>, the output tensor from <code>layer_instance(object)</code> is returned.
shape	integer or integer vector specifying the shape of the output of this layer.
dtype	TensorFlow dtype of the variable created by this layer.
activation	An activation function. See <code>keras::layer_dense</code> . Default: <code>NULL</code> .
initializer	Initializer for the constant vector.
regularizer	Regularizer function applied to the constant vector.
constraint	Constraint function applied to the constant vector.
...	Additional keyword arguments passed to the <code>keras::layer_dense</code> constructed by this layer.

Value

a Keras layer

See Also

Other layers: [layer_autoregressive\(\)](#), [layer_conv_1d_flipout\(\)](#), [layer_conv_1d_reparameterization\(\)](#), [layer_conv_2d_flipout\(\)](#), [layer_conv_2d_reparameterization\(\)](#), [layer_conv_3d_flipout\(\)](#), [layer_conv_3d_reparameterization\(\)](#), [layer_dense_flipout\(\)](#), [layer_dense_local_reparameterization\(\)](#), [layer_dense_reparameterization\(\)](#), [layer_dense_variational\(\)](#)

layer_variational_gaussian_process

A Variational Gaussian Process Layer.

Description

Create a Variational Gaussian Process distribution whose `index_points` are the inputs to the layer. Parameterized by number of inducing points and a `kernel_provider`, which should be a `tf.keras.Layer` with an `@property` that late-binds variable parameters to a `tfp.positive_semidefinite_kernel.PositiveSemidefiniteKernel` instance (this requirement has to do with the way that variables must be created in a keras model). The `mean_fn` is an optional argument which, if omitted, will be automatically configured to be a constant function with trainable variable output.

Usage

```
layer_variational_gaussian_process(
    object,
    num_inducing_points,
    kernel_provider,
    event_shape = 1,
    inducing_index_points_initializer = NULL,
    unconstrained_observation_noise_variance_initializer = NULL,
    mean_fn = NULL,
    jitter = 1e-06,
    name = NULL
)
```

Arguments

<code>object</code>	<p>What to compose the new Layer instance with. Typically a Sequential model or a Tensor (e.g., as returned by <code>layer_input()</code>). The return value depends on <code>object</code>. If <code>object</code> is:</p> <ul style="list-style-type: none"> • missing or NULL, the Layer instance is returned. • a Sequential model, the model with an additional layer is returned. • a Tensor, the output tensor from <code>layer_instance(object)</code> is returned.
---------------------	--

num_inducing_points	number of inducing points in the Variational Gaussian Process distribution.
kernel_provider	a Layer instance equipped with an @property, which yields a PositiveSemidefiniteKernel instance. The latter is used to parametrize the constructed Variational Gaussian Process distribution returned by calling the layer.
event_shape	the shape of the output of the layer. This translates to a batch of underlying Variational Gaussian Process distributions. For example, event_shape = 3 means we are modelling a batch of 3 distributions over functions. We can think of this as a distribution over 3-dimensional vector-valued functions.
inducing_index_points_initializer	a tf.keras.initializer.Initializer used to initialize the trainable inducing_index_points variables. Training VGP's is pretty sensitive to choice of initial inducing index point locations. A reasonable heuristic is to scatter them near the data, not too close to each other.
unconstrained_observation_noise_variance_initializer	a tf.keras.initializer.Initializer used to initialize the unconstrained observation noise variable. The observation noise variance is computed from this variable via the tf.nn.softplus function.
mean_fn	a callable that maps layer inputs to mean function values. Passed to the mean_fn parameter of Variational Gaussian Process distribution. If omitted, defaults to a constant function with trainable variable value.
jitter	a small term added to the diagonal of various kernel matrices for numerical stability.
name	name to give to this layer and the scope of ops and variables it contains.

Value

a Keras layer

mcmc_dual_averaging_step_size_adaptation

Adapts the inner kernel's step_size based on log_accept_prob.

Description

The dual averaging policy uses a noisy step size for exploration, while averaging over tuning steps to provide a smoothed estimate of an optimal value. It is based on section 3.2 of Hoffman and Gelman (2013), which modifies the [stochastic convex optimization scheme of Nesterov (2009)]. The modified algorithm applies extra weight to recent iterations while keeping the convergence guarantees of Robbins-Monro, and takes care not to make the step size too small too quickly when maintaining a constant trajectory length, to avoid expensive early iterations. A good target acceptance probability depends on the inner kernel. If this kernel is HamiltonianMonteCarlo, then 0.6-0.9 is a good range to aim for. For RandomWalkMetropolis this should be closer to 0.25. See the individual kernels' docstrings for guidance.

Usage

```
mcmc_dual_averaging_step_size_adaptation(
    inner_kernel,
    num_adaptation_steps,
    target_accept_prob = 0.75,
    exploration_shrinkage = 0.05,
    step_count_smoothing = 10,
    decay_rate = 0.75,
    step_size_setter_fn = NULL,
    step_size_getter_fn = NULL,
    log_accept_prob_getter_fn = NULL,
    validate_args = FALSE,
    name = NULL
)
```

Arguments

`inner_kernel` TransitionKernel-like object.

`num_adaptation_steps` Scalar integer Tensor number of initial steps to during which to adjust the step size. This may be greater, less than, or equal to the number of burnin steps.

`target_accept_prob` A floating point Tensor representing desired acceptance probability. Must be a positive number less than 1. This can either be a scalar, or have shape `[num_chains]`. Default value: 0.75 (the center of asymptotically optimal rate for HMC).

`exploration_shrinkage` Floating point scalar Tensor. How strongly the exploration rate is biased towards the shrinkage target.

`step_count_smoothing` Int32 scalar Tensor. Number of "pseudo-steps" added to the number of steps taken to prevents noisy exploration during the early samples.

`decay_rate` Floating point scalar Tensor. How much to favor recent iterations over earlier ones. A value of 1 gives equal weight to all history.

`step_size_setter_fn` A function with the signature `(kernel_results, new_step_size) -> new_kernel_results` where `kernel_results` are the results of the `inner_kernel`, `new_step_size` is a Tensor or a nested collection of Tensors with the same structure as returned by the `step_size_getter_fn`, and `new_kernel_results` are a copy of `kernel_results` with the step size(s) set.

`step_size_getter_fn` A callable with the signature `(kernel_results) -> step_size` where `kernel_results` are the results of the `inner_kernel`, and `step_size` is a floating point Tensor or a nested collection of such Tensors.

`log_accept_prob_getter_fn` A callable with the signature `(kernel_results) -> log_accept_prob` where `kernel_results` are the results of the `inner_kernel`, and `log_accept_prob` is a floating point Tensor. `log_accept_prob` can either be a scalar, or have shape

	[num_chains]. If it's the latter, step_size should also have the same leading dimension.
validate_args	logical. When TRUE kernel parameters are checked for validity. When FALSE invalid inputs may silently render incorrect outputs.
name	name prefixed to Ops created by this function. Default value: NULL (i.e., 'dual_averaging_step_size_adaptation')

Details

In general, adaptation prevents the chain from reaching a stationary distribution, so obtaining consistent samples requires num_adaptation_steps be set to a value somewhat smaller than the number of burnin steps. However, it may sometimes be helpful to set num_adaptation_steps to a larger value during development in order to inspect the behavior of the chain during adaptation. The step size is assumed to broadcast with the chain state, potentially having leading dimensions corresponding to multiple chains. When there are fewer of those leading dimensions than there are chain dimensions, the corresponding dimensions in the log_accept_prob are averaged (in the direct space, rather than the log space) before being used to adjust the step size. This means that this kernel can do both cross-chain adaptation, or per-chain step size adaptation, depending on the shape of the step size. For example, if your problem has a state with shape [S], your chain state has shape [C0, C1, S] (meaning that there are $C0 * C1$ total chains) and log_accept_prob has shape [C0, C1] (one acceptance probability per chain), then depending on the shape of the step size, the following will happen:

- Step size has shape [], [S] or [1], the log_accept_prob will be averaged across its C0 and C1 dimensions. This means that you will learn a shared step size based on the mean acceptance probability across all chains. This can be useful if you don't have a lot of steps to adapt and want to average away the noise.
- Step size has shape [C1, 1] or [C1, S], the log_accept_prob will be averaged across its C0 dimension. This means that you will learn a shared step size based on the mean acceptance probability across chains that share the coordinate across the C1 dimension. This can be useful when the C1 dimension indexes different distributions, while C0 indexes replicas of a single distribution, all sampled in parallel.
- Step size has shape [C0, C1, 1] or [C0, C1, S], then no averaging will happen. This means that each chain will learn its own step size. This can be useful when all chains are sampling from different distributions. Even when all chains are for the same distribution, this can help during the initial warmup period.
- Step size has shape [C0, 1, 1] or [C0, 1, S], the log_accept_prob will be averaged across its C1 dimension. This means that you will learn a shared step size based on the mean acceptance probability across chains that share the coordinate across the C0 dimension. This can be useful when the C0 dimension indexes different distributions, while C1 indexes replicas of a single distribution, all sampled in parallel.

Value

a Monte Carlo sampling kernel

References

- Matthew D. Hoffman, Andrew Gelman. The No-U-Turn Sampler: Adaptively Setting Path Lengths in Hamiltonian Monte Carlo. In *Journal of Machine Learning Research*, 15(1):1593-1623, 2014.
- Yurii Nesterov. Primal-dual subgradient methods for convex problems. *Mathematical programming* 120.1 (2009): 221-259
- <https://statmodeling.stat.columbia.edu/2017/12/15/burn-vs-warm-iterative-simulation-algorithms>

See Also

For an example how to use see `mcmc_no_u_turn_sampler()`.

Other `mcmc_kernels`: `mcmc_hamiltonian_monte_carlo()`, `mcmc_metropolis_adjusted_langevin_algorithm()`, `mcmc_metropolis_hastings()`, `mcmc_no_u_turn_sampler()`, `mcmc_random_walk_metropolis()`, `mcmc_replica_exchange_mc()`, `mcmc_simple_step_size_adaptation()`, `mcmc_slice_sampler()`, `mcmc_transformed_transition_kernel()`, `mcmc_uncalibrated_hamiltonian_monte_carlo()`, `mcmc_uncalibrated_langevin()`, `mcmc_uncalibrated_random_walk()`

`mcmc_effective_sample_size`

Estimate a lower bound on effective sample size for each independent chain.

Description

Roughly speaking, "effective sample size" (ESS) is the size of an iid sample with the same variance as state.

Usage

```
mcmc_effective_sample_size(
  states,
  filter_threshold = 0,
  filter_beyond_lag = NULL,
  name = NULL
)
```

Arguments

`states` Tensor or list of Tensor objects. Dimension zero should index identically distributed states.

`filter_threshold` Tensor or list of Tensor objects. Must broadcast with `state`. The auto-correlation sequence is truncated after the first appearance of a term less than `filter_threshold`. Setting to `NULL` means we use no threshold filter. Since $|R_k| \leq 1$, setting to any number less than -1 has the same effect.

`filter_beyond_lag` Tensor or list of Tensor objects. Must be int-like and scalar valued. The auto-correlation sequence is truncated to this length. Setting to NULL means we do not filter based on number of lags.

`name` name to prepend to created ops.

Details

More precisely, given a stationary sequence of possibly correlated random variables X_1, X_2, \dots, X_N , each identically distributed ESS is the number such that $\text{Variance}\{N^{-1} * \text{Sum}\{X_i\}\} = \text{ESS}^{-1} * \text{Variance}\{X_1\}$.

If the sequence is uncorrelated, $\text{ESS} = N$. In general, one should expect $\text{ESS} \leq N$, with more highly correlated sequences having smaller ESS.

Value

Tensor or list of Tensor objects. The effective sample size of each component of states. Shape will be `states.shape[1:]`.

See Also

Other `mcmc_functions`: [mcmc_potential_scale_reduction\(\)](#), [mcmc_sample_annealed_importance_chain\(\)](#), [mcmc_sample_chain\(\)](#), [mcmc_sample_halton_sequence\(\)](#)

`mcmc_hamiltonian_monte_carlo`

Runs one step of Hamiltonian Monte Carlo.

Description

Hamiltonian Monte Carlo (HMC) is a Markov chain Monte Carlo (MCMC) algorithm that takes a series of gradient-informed steps to produce a Metropolis proposal. This class implements one random HMC step from a given `current_state`. Mathematical details and derivations can be found in Neal (2011).

Usage

```
mcmc_hamiltonian_monte_carlo(
    target_log_prob_fn,
    step_size,
    num_leapfrog_steps,
    state_gradients_are_stopped = FALSE,
    step_size_update_fn = NULL,
    seed = NULL,
    store_parameters_in_results = FALSE,
    name = NULL
)
```

Arguments

<code>target_log_prob_fn</code>	Function which takes an argument like <code>current_state</code> (if it's a list <code>current_state</code> will be unpacked) and returns its (possibly unnormalized) log-density under the target distribution.
<code>step_size</code>	Tensor or list of Tensors representing the step size for the leapfrog integrator. Must broadcast with the shape of <code>current_state</code> . Larger step sizes lead to faster progress, but too-large step sizes make rejection exponentially more likely. When possible, it's often helpful to match per-variable step sizes to the standard deviations of the target distribution in each variable.
<code>num_leapfrog_steps</code>	Integer number of steps to run the leapfrog integrator for. Total progress per HMC step is roughly proportional to <code>step_size * num_leapfrog_steps</code> .
<code>state_gradients_are_stopped</code>	logical indicating that the proposed new state be run through <code>tf\$stop_gradient</code> . This is particularly useful when combining optimization over samples from the HMC chain. Default value: FALSE (i.e., do not apply <code>stop_gradient</code>).
<code>step_size_update_fn</code>	Function taking <code>current_step_size</code> (typically a <code>tf\$Variable</code>) and <code>kernel_results</code> (typically <code>collections\$namedtuple</code>) and returns updated <code>step_size</code> (Tensors). Default value: NULL (i.e., do not update <code>step_size</code> automatically).
<code>seed</code>	integer to seed the random number generator.
<code>store_parameters_in_results</code>	If TRUE, then <code>step_size</code> and <code>num_leapfrog_steps</code> are written to and read from eponymous fields in the kernel results objects returned from <code>one_step</code> and <code>bootstrap_results</code> . This allows wrapper kernels to adjust those parameters on the fly. This is incompatible with <code>step_size_update_fn</code> , which must be set to NULL.
<code>name</code>	string prefixed to Ops created by this function. Default value: NULL (i.e., 'hmc_kernel').

Details

The `one_step` function can update multiple chains in parallel. It assumes that all leftmost dimensions of `current_state` index independent chain states (and are therefore updated independently). The output of `target_log_prob_fn(current_state)` should sum log-probabilities across all event dimensions. Slices along the rightmost dimensions may have different target distributions; for example, `current_state[0, :]` could have a different target distribution from `current_state[1, :]`. These semantics are governed by `target_log_prob_fn(current_state)`. (The number of independent chains is `tf$size(target_log_prob_fn(current_state))`.)

Value

a Monte Carlo sampling kernel

References

- Radford Neal. *MCMC Using Hamiltonian Dynamics*. *Handbook of Markov Chain Monte Carlo*, 2011.

- Bernard Delyon, Marc Lavielle, Eric Moulines. *Convergence of a stochastic approximation version of the EM algorithm*, Ann. Statist. 27 (1999), no. 1, 94–128.

See Also

Other mcmc_kernels: `mcmc_dual_averaging_step_size_adaptation()`, `mcmc_metropolis_adjusted_langevin_algorithm()`, `mcmc_metropolis_hastings()`, `mcmc_no_u_turn_sampler()`, `mcmc_random_walk_metropolis()`, `mcmc_replica_exchange_mc()`, `mcmc_simple_step_size_adaptation()`, `mcmc_slice_sampler()`, `mcmc_transformed_transition_kernel()`, `mcmc_uncalibrated_hamiltonian_monte_carlo()`, `mcmc_uncalibrated_langevin()`, `mcmc_uncalibrated_random_walk()`

mcmc_metropolis_adjusted_langevin_algorithm

Runs one step of Metropolis-adjusted Langevin algorithm.

Description

Metropolis-adjusted Langevin algorithm (MALA) is a Markov chain Monte Carlo (MCMC) algorithm that takes a step of a discretised Langevin diffusion as a proposal. This class implements one step of MALA using Euler-Maruyama method for a given `current_state` and diagonal preconditioning volatility matrix.

Usage

```
mcmc_metropolis_adjusted_langevin_algorithm(
    target_log_prob_fn,
    step_size,
    volatility_fn = NULL,
    seed = NULL,
    parallel_iterations = 10,
    name = NULL
)
```

Arguments

<code>target_log_prob_fn</code>	Function which takes an argument like <code>current_state</code> (if it's a list <code>current_state</code> will be unpacked) and returns its (possibly unnormalized) log-density under the target distribution.
<code>step_size</code>	Tensor or list of Tensors representing the step size for the leapfrog integrator. Must broadcast with the shape of <code>current_state</code> . Larger step sizes lead to faster progress, but too-large step sizes make rejection exponentially more likely. When possible, it's often helpful to match per-variable step sizes to the standard deviations of the target distribution in each variable.
<code>volatility_fn</code>	function which takes an argument like <code>current_state</code> (or <code>*current_state</code> if it's a list) and returns volatility value at <code>current_state</code> . Should return a Tensor or list of Tensors that must broadcast with the shape of <code>current_state</code> . Defaults to the identity function.

seed integer to seed the random number generator.

parallel_iterations the number of coordinates for which the gradients of the volatility matrix volatility_fn can be computed in parallel.

name String prefixed to Ops created by this function. Default value: NULL (i.e., 'mala_kernel').

Details

Mathematical details and derivations can be found in Roberts and Rosenthal (1998) and Xifara et al. (2013).

The one_step function can update multiple chains in parallel. It assumes that all leftmost dimensions of current_state index independent chain states (and are therefore updated independently). The output of target_log_prob_fn(current_state) should reduce log-probabilities across all event dimensions. Slices along the rightmost dimensions may have different target distributions; for example, current_state[0, :] could have a different target distribution from current_state[1, :]. These semantics are governed by target_log_prob_fn(current_state). (The number of independent chains is tf.size(target_log_prob_fn(current_state)).)

References

- Gareth Roberts and Jeffrey Rosenthal. Optimal Scaling of Discrete Approximations to Langevin Diffusions. *Journal of the Royal Statistical Society: Series B (Statistical Methodology)*, 60: 255-268, 1998.
- T. Xifara et al. Langevin diffusions and the Metropolis-adjusted Langevin algorithm. *arXiv preprint arXiv:1309.2983*, 2013.

See Also

Other mcmc_kernels: [mcmc_dual_averaging_step_size_adaptation\(\)](#), [mcmc_hamiltonian_monte_carlo\(\)](#), [mcmc_metropolis_hastings\(\)](#), [mcmc_no_u_turn_sampler\(\)](#), [mcmc_random_walk_metropolis\(\)](#), [mcmc_replica_exchange_mc\(\)](#), [mcmc_simple_step_size_adaptation\(\)](#), [mcmc_slice_sampler\(\)](#), [mcmc_transformed_transition_kernel\(\)](#), [mcmc_uncalibrated_hamiltonian_monte_carlo\(\)](#), [mcmc_uncalibrated_langevin\(\)](#), [mcmc_uncalibrated_random_walk\(\)](#)

mcmc_metropolis_hastings

Runs one step of the Metropolis-Hastings algorithm.

Description

The Metropolis-Hastings algorithm is a Markov chain Monte Carlo (MCMC) technique which uses a proposal distribution to eventually sample from a target distribution.

Usage

```
mcmc_metropolis_hastings(inner_kernel, seed = NULL, name = NULL)
```

Arguments

inner_kernel	TransitionKernel-like object which has collections\$namedtuple kernel_results and which contains a target_log_prob member and optionally a log_acceptance_correction member.
seed	integer to seed the random number generator.
name	string prefixed to Ops created by this function. Default value: NULL (i.e., "mh_kernel").

Details

Note: inner_kernel\$one_step must return kernel_results as a collections\$namedtuple which must:

- have a target_log_prob field,
- optionally have a log_acceptance_correction field, and,
- have only fields which are Tensor-valued.

The Metropolis-Hastings log acceptance-probability is computed as:

```
log_accept_ratio = (current_kernel_results.target_log_prob
                    - previous_kernel_results.target_log_prob
                    + current_kernel_results.log_acceptance_correction)
```

If current_kernel_results\$log_acceptance_correction does not exist, it is presumed 0 (i.e., that the proposal distribution is symmetric). The most common use-case for log_acceptance_correction is in the Metropolis-Hastings algorithm, i.e.,

$$\text{accept_prob}(x' \mid x) = p(x') / p(x) (g(x \mid x') / g(x' \mid x))$$

where,

p represents the target distribution,

g represents the proposal (conditional) distribution,

x' is the proposed state, and,

x is current state

The log of the parenthetical term is the log_acceptance_correction. The log_acceptance_correction may not necessarily correspond to the ratio of proposal distributions, e.g, log_acceptance_correction has a different interpretation in Hamiltonian Monte Carlo.

Value

a Monte Carlo sampling kernel

See Also

Other mcmc_kernels: [mcmc_dual_averaging_step_size_adaptation\(\)](#), [mcmc_hamiltonian_monte_carlo\(\)](#), [mcmc_metropolis_adjusted_langevin_algorithm\(\)](#), [mcmc_no_u_turn_sampler\(\)](#), [mcmc_random_walk_metropolis\(\)](#), [mcmc_replica_exchange_mc\(\)](#), [mcmc_simple_step_size_adaptation\(\)](#), [mcmc_slice_sampler\(\)](#), [mcmc_transformed_transition_kernel\(\)](#), [mcmc_uncalibrated_hamiltonian_monte_carlo\(\)](#), [mcmc_uncalibrated_langevin\(\)](#), [mcmc_uncalibrated_random_walk\(\)](#)

mcmc_no_u_turn_sampler

Runs one step of the No U-Turn Sampler

Description

The No U-Turn Sampler (NUTS) is an adaptive variant of the Hamiltonian Monte Carlo (HMC) method for MCMC. NUTS adapts the distance traveled in response to the curvature of the target density. Conceptually, one proposal consists of reversibly evolving a trajectory through the sample space, continuing until that trajectory turns back on itself (hence the name, 'No U-Turn'). This class implements one random NUTS step from a given `current_state`. Mathematical details and derivations can be found in Hoffman & Gelman (2011).

Usage

```
mcmc_no_u_turn_sampler(
    target_log_prob_fn,
    step_size,
    max_tree_depth = 10,
    max_energy_diff = 1000,
    unrolled_leapfrog_steps = 1,
    seed = NULL,
    name = NULL
)
```

Arguments

<code>target_log_prob_fn</code>	function which takes an argument like <code>current_state</code> and returns its (possibly unnormalized) log-density under the target distribution.
<code>step_size</code>	Tensor or list of Tensors representing the step size for the leapfrog integrator. Must broadcast with the shape of <code>current_state</code> . Larger step sizes lead to faster progress, but too-large step sizes make rejection exponentially more likely. When possible, it's often helpful to match per-variable step sizes to the standard deviations of the target distribution in each variable.
<code>max_tree_depth</code>	Maximum depth of the tree implicitly built by NUTS. The maximum number of leapfrog steps is bounded by $2 \times \text{max_tree_depth}$ i.e. the number of nodes in a binary tree <code>max_tree_depth</code> nodes deep. The default setting of 10 takes up to 1024 leapfrog steps.
<code>max_energy_diff</code>	Scaler threshold of energy differences at each leapfrog, divergence samples are defined as leapfrog steps that exceed this threshold. Default to 1000.
<code>unrolled_leapfrog_steps</code>	The number of leapfrogs to unroll per tree expansion step. Applies a direct linear multiplier to the maximum trajectory length implied by <code>max_tree_depth</code> . Defaults to 1.

seed integer to seed the random number generator.
 name name prefixed to Ops created by this function. Default value: NULL (i.e., 'nuts_kernel').

Details

The `one_step` function can update multiple chains in parallel. It assumes that a prefix of leftmost dimensions of `current_state` index independent chain states (and are therefore updated independently). The output of `target_log_prob_fn(current_state)` should sum log-probabilities across all event dimensions. Slices along the rightmost dimensions may have different target distributions; for example, `current_state[0][0, ...]` could have a different target distribution from `current_state[0][1, ...]`. These semantics are governed by `target_log_prob_fn(*current_state)`. (The number of independent chains is `tf$size(target_log_prob_fn(current_state))`.)

Value

a Monte Carlo sampling kernel

References

- [Matthew D. Hoffman, Andrew Gelman. The No-U-Turn Sampler: Adaptively Setting Path Lengths in Hamiltonian Monte Carlo. 2011.](#)

See Also

Other `mcmc_kernels`: `mcmc_dual_averaging_step_size_adaptation()`, `mcmc_hamiltonian_monte_carlo()`, `mcmc_metropolis_adjusted_langevin_algorithm()`, `mcmc_metropolis_hastings()`, `mcmc_random_walk_metropolis()`, `mcmc_replica_exchange_mc()`, `mcmc_simple_step_size_adaptation()`, `mcmc_slice_sampler()`, `mcmc_transformed_transition_kernel()`, `mcmc_uncalibrated_hamiltonian_monte_carlo()`, `mcmc_uncalibrated_langevin()`, `mcmc_uncalibrated_random_walk()`

Examples

```
predictors <- tf$cast( c(201,244, 47,287,203,58,210,202,198,158,165,201,157,
  131,166,160,186,125,218,146),tf$float32)
obs <- tf$cast(c(592,401,583,402,495,173,479,504,510,416,393,442,317,311,400,
  337,423,334,533,344),tf$float32)
y_sigma <- tf$cast(c(61,25,38,15,21,15,27,14,30,16,14,25,52,16,34,31,42,26,
  16,22),tf$float32)

# Robust linear regression model
robust_lm <- tfd_joint_distribution_sequential(
  list(
    tfd_normal(loc = 0, scale = 1, name = "b0"),
    tfd_normal(loc = 0, scale = 1, name = "b1"),
    tfd_half_normal(5, name = "df"),
    function(df, b1, b0)
      tfd_independent(
        tfd_student_t(
          # Likelihood
          df = tf$expand_dims(df, axis = -1L),
```

```

    loc = tf$expand_dims(b0, axis = -1L) +
          tf$expand_dims(b1, axis = -1L) * predictors[tf$newaxis, ],
    scale = y_sigma,
    name = "st"
  ), name = "ind")), validate_args = TRUE)

log_prob <-function(b0, b1, df) {robust_lm %>%
  tfd_log_prob(list(b0, b1, df, obs))}
step_size0 <- Map(function(x) tf$cast(x, tf$float32), c(1, .2, .5))

number_of_steps <- 10
burnin <- 5
nchain <- 50

run_chain <- function() {
# random initialization of the starting position of each chain
samples <- robust_lm %>% tfd_sample(nchain)
b0 <- samples[[1]]
b1 <- samples[[2]]
df <- samples[[3]]

# bijector to map constrained parameters to real
unconstraining_bijectors <- list(
  tfb_identity(), tfb_identity(), tfb_exp())

trace_fn <- function(x, pkr) {
  list(pkr$inner_results$inner_results$step_size,
       pkr$inner_results$inner_results$log_accept_ratio)
}

nuts <- mcmc_no_u_turn_sampler(
  target_log_prob_fn = log_prob,
  step_size = step_size0
) %>%
mcmc_transformed_transition_kernel(bijector = unconstraining_bijectors) %>%
mcmc_dual_averaging_step_size_adaptation(
  num_adaptation_steps = burnin,
  step_size_setter_fn = function(pkr, new_step_size)
  pkr`_replace`(
    inner_results = pkr$inner_results`_replace`(step_size = new_step_size)),
  step_size_getter_fn = function(pkr) pkr$inner_results$step_size,
  log_accept_prob_getter_fn = function(pkr) pkr$inner_results$log_accept_ratio
)

nuts %>% mcmc_sample_chain(
  num_results = number_of_steps,
  num_burnin_steps = burnin,
  current_state = list(b0, b1, df),
  trace_fn = trace_fn)
}

run_chain <- tensorflow::tf_function(run_chain)
res <- run_chain()

```

```
mcmc_potential_scale_reduction
```

Gelman and Rubin (1992)'s potential scale reduction for chain convergence.

Description

Given $N > 1$ states from each of $C > 1$ independent chains, the potential scale reduction factor, commonly referred to as \hat{R} , measures convergence of the chains (to the same target) by testing for equality of means.

Usage

```
mcmc_potential_scale_reduction(
  chains_states,
  independent_chain_ndims = 1,
  name = NULL
)
```

Arguments

chains_states Tensor or list of Tensors representing the state(s) of a Markov Chain at each result step. The i th state is assumed to have shape $[N_i, C_{i1}, C_{i2}, \dots, C_{iD}] + A$. Dimension 0 indexes the $N_i > 1$ result steps of the Markov Chain. Dimensions 1 through D index the $C_{i1} \times \dots \times C_{iD}$ independent chains to be tested for convergence to the same target. The remaining dimensions, A , can have any shape (even empty).

independent_chain_ndims Integer type Tensor with value ≥ 1 giving the number of giving the number of dimensions, from $\text{dim} = 1$ to $\text{dim} = D$, holding independent chain results to be tested for convergence.

name name to prepend to created tf. Default: `potential_scale_reduction`.

Details

Specifically, \hat{R} measures the degree to which variance (of the means) between chains exceeds what one would expect if the chains were identically distributed. See Gelman and Rubin (1992), Brooks and Gelman (1998)].

Some guidelines:

- The initial state of the chains should be drawn from a distribution overdispersed with respect to the target.
- If all chains converge to the target, then as $N \rightarrow \infty$, $\hat{R} \rightarrow 1$. Before that, $\hat{R} > 1$ (except in pathological cases, e.g. if the chain paths were identical).

- The above holds for any number of chains $C > 1$. Increasing C improves effectiveness of the diagnostic.
- Sometimes, $R\text{-hat} < 1.2$ is used to indicate approximate convergence, but of course this is problem dependent. See Brooks and Gelman (1998).
- $R\text{-hat}$ only measures non-convergence of the mean. If higher moments, or other statistics are desired, a different diagnostic should be used. See Brooks and Gelman (1998).

To see why $R\text{-hat}$ is reasonable, let X be a random variable drawn uniformly from the combined states (combined over all chains). Then, in the limit $N, C \rightarrow \infty$, with E, Var denoting expectation and variance, $R\text{-hat} = (E[\text{Var}[X | \text{chain}]] + \text{Var}[E[X | \text{chain}]]) / E[\text{Var}[X | \text{chain}]]$. Using the law of total variance, the numerator is the variance of the combined states, and the denominator is the total variance minus the variance of the the individual chain means. If the chains are all drawing from the same distribution, they will have the same mean, and thus the ratio should be one.

Value

Tensor or list of Tensors representing the $R\text{-hat}$ statistic for the state(s). Same dtype as state, and shape equal to `state.shape[1 + independent_chain_ndims:]`.

References

- Stephen P. Brooks and Andrew Gelman. General Methods for Monitoring Convergence of Iterative Simulations. *Journal of Computational and Graphical Statistics*, 7(4), 1998.
- Andrew Gelman and Donald B. Rubin. Inference from Iterative Simulation Using Multiple Sequences. *Statistical Science*, 7(4):457-472, 1992.

See Also

Other mcmc_functions: `mcmc_effective_sample_size()`, `mcmc_sample_annealed_importance_chain()`, `mcmc_sample_chain()`, `mcmc_sample_halton_sequence()`

mcmc_random_walk_metropolis

Runs one step of the RWM algorithm with symmetric proposal.

Description

Random Walk Metropolis is a gradient-free Markov chain Monte Carlo (MCMC) algorithm. The algorithm involves a proposal generating step `proposal_state = current_state + perturb` by a random perturbation, followed by Metropolis-Hastings accept/reject step. For more details see Section 2.1 of Roberts and Rosenthal (2004).

Usage

```
mcmc_random_walk_metropolis(
  target_log_prob_fn,
  new_state_fn = NULL,
  seed = NULL,
  name = NULL
)
```

Arguments

<code>target_log_prob_fn</code>	Function which takes an argument like <code>current_state</code> (if it's a list <code>current_state</code> will be unpacked) and returns its (possibly unnormalized) log-density under the target distribution.
<code>new_state_fn</code>	Function which takes a list of state parts and a seed; returns a same-type list of Tensors, each being a perturbation of the input state parts. The perturbation distribution is assumed to be a symmetric distribution centered at the input state part. Default value: <code>NULL</code> which is mapped to <code>tfp\$mcmc\$random_walk_normal_fn()</code> .
<code>seed</code>	integer to seed the random number generator.
<code>name</code>	String name prefixed to Ops created by this function. Default value: <code>NULL</code> (i.e., <code>'rwm_kernel'</code>).

Details

The current class implements RWM for normal and uniform proposals. Alternatively, the user can supply any custom proposal generating function. The function `one_step` can update multiple chains in parallel. It assumes that all leftmost dimensions of `current_state` index independent chain states (and are therefore updated independently). The output of `target_log_prob_fn(current_state)` should sum log-probabilities across all event dimensions. Slices along the rightmost dimensions may have different target distributions; for example, `current_state[0, :]` could have a different target distribution from `current_state[1, :]`. These semantics are governed by `target_log_prob_fn(current_state)`. (The number of independent chains is `tf$size(target_log_prob_fn(current_state))`.)

Value

a Monte Carlo sampling kernel

See Also

Other `mcmc_kernels`: [mcmc_dual_averaging_step_size_adaptation\(\)](#), [mcmc_hamiltonian_monte_carlo\(\)](#), [mcmc_metropolis_adjusted_langevin_algorithm\(\)](#), [mcmc_metropolis_hastings\(\)](#), [mcmc_no_u_turn_sampler\(\)](#), [mcmc_replica_exchange_mc\(\)](#), [mcmc_simple_step_size_adaptation\(\)](#), [mcmc_slice_sampler\(\)](#), [mcmc_transformed_transition_kernel\(\)](#), [mcmc_uncalibrated_hamiltonian_monte_carlo\(\)](#), [mcmc_uncalibrated_langevin\(\)](#), [mcmc_uncalibrated_random_walk\(\)](#)

mcmc_replica_exchange_mc

Runs one step of the Replica Exchange Monte Carlo

Description

Replica Exchange Monte Carlo is a Markov chain Monte Carlo (MCMC) algorithm that is also known as Parallel Tempering. This algorithm performs multiple sampling with different temperatures in parallel, and exchanges those samplings according to the Metropolis-Hastings criterion. The K replicas are parameterized in terms of inverse_temperature's, ($\beta[0]$, $\beta[1]$, ..., $\beta[K-1]$). If the target distribution has probability density $p(x)$, the k th replica has density $p(x)^{\beta_k}$.

Usage

```
mcmc_replica_exchange_mc(
  target_log_prob_fn,
  inverse_temperatures,
  make_kernel_fn,
  swap_proposal_fn = tfp.mcmc.replica_exchange_mc.default_swap_proposal_fn(1),
  state_includes_replicas = FALSE,
  seed = NULL,
  name = NULL
)
```

Arguments

target_log_prob_fn	Function which takes an argument like current_state (if it's a list current_state will be unpacked) and returns its (possibly unnormalized) log-density under the target distribution.
inverse_temperatures	1D Tensor of inverse temperatures to perform samplings with each replica. Must have statically known shape. inverse_temperatures[0] produces the states returned by samplers, and is typically == 1.
make_kernel_fn	Function which takes target_log_prob_fn and seed args and returns a TransitionKernel instance.
swap_proposal_fn	function which take a number of replicas, and return combinations of replicas for exchange.
state_includes_replicas	Boolean indicating whether the leftmost dimension of each state sample should index replicas. If TRUE, the leftmost dimension of the current_state kwarg to tfp.mcmc.sample_chain will be interpreted as indexing replicas.
seed	integer to seed the random number generator.
name	string prefixed to Ops created by this function. Default value: NULL (i.e., "remc_kernel").

Details

Typically $\beta[0] = 1.0$, and $1.0 > \beta[1] > \beta[2] > \dots > 0.0$.

- $\beta[0] == 1$ ==> First replicas samples from the target density, p .
- $\beta[k] < 1$, for $k = 1, \dots, K-1$ ==> Other replicas sample from "flattened" versions of p (peak is less high, valley less low). These distributions are somewhat closer to a uniform on the support of p . Samples from adjacent replicas $i, i + 1$ are used as proposals for each other in a Metropolis step. This allows the lower beta samples, which explore less dense areas of p , to occasionally be used to help the $\beta == 1$ chain explore new regions of the support. Samples from replica 0 are returned, and the others are discarded.

Value

list of `next_state` (Tensor or Python list of Tensors representing the state(s) of the Markov chain(s) at each result step. Has same shape as `current_state`.) and `kernel_results` (collections.namedtuple of internal calculations used to 'advance the chain').

See Also

Other `mcmc_kernels`: `mcmc_dual_averaging_step_size_adaptation()`, `mcmc_hamiltonian_monte_carlo()`, `mcmc_metropolis_adjusted_langevin_algorithm()`, `mcmc_metropolis_hastings()`, `mcmc_no_u_turn_sampler()`, `mcmc_random_walk_metropolis()`, `mcmc_simple_step_size_adaptation()`, `mcmc_slice_sampler()`, `mcmc_transformed_transition_kernel()`, `mcmc_uncalibrated_hamiltonian_monte_carlo()`, `mcmc_uncalibrated_langevin()`, `mcmc_uncalibrated_random_walk()`

`mcmc_sample_annealed_importance_chain`

Runs annealed importance sampling (AIS) to estimate normalizing constants.

Description

This function uses an MCMC transition operator (e.g., Hamiltonian Monte Carlo) to sample from a series of distributions that slowly interpolates between an initial "proposal" distribution: $\exp(\text{proposal_log_prob_fn}(x) - \text{proposal_log_normalizer})$ and the target distribution: $\exp(\text{target_log_prob_fn}(x) - \text{target_log_normalizer})$, accumulating importance weights along the way. The product of these importance weights gives an unbiased estimate of the ratio of the normalizing constants of the initial distribution and the target distribution: $E[\exp(\text{ais_weights})] = \exp(\text{target_log_normalizer} - \text{proposal_log_normalizer})$.

Usage

```
mcmc_sample_annealed_importance_chain(
    num_steps,
    proposal_log_prob_fn,
    target_log_prob_fn,
    current_state,
    make_kernel_fn,
```

```

    parallel_iterations = 10,
    name = NULL
)

```

Arguments

<code>num_steps</code>	Integer number of Markov chain updates to run. More iterations means more expense, but smoother annealing between q and p, which in turn means exponentially lower variance for the normalizing constant estimator.
<code>proposal_log_prob_fn</code>	function that returns the log density of the initial distribution.
<code>target_log_prob_fn</code>	function which takes an argument like <code>current_state</code> and returns its (possibly unnormalized) log-density under the target distribution.
<code>current_state</code>	Tensor or list of Tensors representing the current state(s) of the Markov chain(s). The first r dimensions index independent chains, $r = \text{tf\$rank}(\text{target_log_prob_fn}(\text{current_state}))$.
<code>make_kernel_fn</code>	function which returns a TransitionKernel-like object. Must take one argument representing the TransitionKernel's <code>target_log_prob_fn</code> . The <code>target_log_prob_fn</code> argument represents the TransitionKernel's target log distribution. Note: <code>sample_annealed_importance_chain</code> creates a new <code>target_log_prob_fn</code> which is an interpolation between the supplied <code>target_log_prob_fn</code> and <code>proposal_log_prob_fn</code> ; it is this interpolated function which is used as an argument to <code>make_kernel_fn</code> .
<code>parallel_iterations</code>	The number of iterations allowed to run in parallel. It must be a positive integer. See <code>tf\$while_loop</code> for more details.
<code>name</code>	string prefixed to Ops created by this function. Default value: NULL (i.e., "sample_annealed_importance_chain").

Details

Note: When running in graph mode, `proposal_log_prob_fn` and `target_log_prob_fn` are called exactly three times (although this may be reduced to two times in the future).

Value

list of `next_state` (Tensor or Python list of Tensors representing the state(s) of the Markov chain(s) at the final iteration. Has same shape as input `current_state`), `ais_weights` (Tensor with the estimated weight(s). Has shape matching `target_log_prob_fn(current_state)`), and `kernel_results` (collections.namedtuple of internal calculations used to advance the chain).

See Also

For an example how to use see [mcmc_sample_chain\(\)](#).

Other mcmc_functions: [mcmc_effective_sample_size\(\)](#), [mcmc_potential_scale_reduction\(\)](#), [mcmc_sample_chain\(\)](#), [mcmc_sample_halton_sequence\(\)](#)

mcmc_sample_chain	<i>Implements Markov chain Monte Carlo via repeated TransitionKernel steps.</i>
-------------------	---

Description

This function samples from an Markov chain at `current_state` and whose stationary distribution is governed by the supplied `TransitionKernel` instance (`kernel`).

Usage

```
mcmc_sample_chain(
  kernel = NULL,
  num_results,
  current_state,
  previous_kernel_results = NULL,
  num_burnin_steps = 0,
  num_steps_between_results = 0,
  trace_fn = NULL,
  return_final_kernel_results = FALSE,
  parallel_iterations = 10,
  seed = NULL,
  name = NULL
)
```

Arguments

<code>kernel</code>	An instance of <code>tfp\$mcmc\$TransitionKernel</code> which implements one step of the Markov chain.
<code>num_results</code>	Integer number of Markov chain draws.
<code>current_state</code>	Tensor or list of Tensors representing the current state(s) of the Markov chain(s).
<code>previous_kernel_results</code>	A Tensor or a nested collection of Tensors representing internal calculations made within the previous call to this function (or as returned by <code>bootstrap_results</code>).
<code>num_burnin_steps</code>	Integer number of chain steps to take before starting to collect results. Default value: 0 (i.e., no burn-in).
<code>num_steps_between_results</code>	Integer number of chain steps between collecting a result. Only one out of every <code>num_steps_between_samples + 1</code> steps is included in the returned results. The number of returned chain states is still equal to <code>num_results</code> . Default value: 0 (i.e., no thinning).
<code>trace_fn</code>	A function that takes in the current chain state and the previous kernel results and return a Tensor or a nested collection of Tensors that is then traced along with the chain state.

return_final_kernel_results	If TRUE, then the final kernel results are returned alongside the chain state and the trace specified by the trace_fn.
parallel_iterations	The number of iterations allowed to run in parallel. It must be a positive integer. See <code>tf\$while_loop</code> for more details.
seed	Optional, a seed for reproducible sampling.
name	string prefixed to Ops created by this function. Default value: NULL, (i.e., "mcmc_sample_chain").

Details

This function can sample from multiple chains, in parallel. (Whether or not there are multiple chains is dictated by the kernel.)

The `current_state` can be represented as a single Tensor or a list of Tensors which collectively represent the current state. Since MCMC states are correlated, it is sometimes desirable to produce additional intermediate states, and then discard them, ending up with a set of states with decreased autocorrelation. See Owen (2017). Such "thinning" is made possible by setting `num_steps_between_results > 0`. The chain then takes `num_steps_between_results` extra steps between the steps that make it into the results. The extra steps are never materialized (in calls to `sess$run`), and thus do not increase memory requirements.

Warning: when setting a seed in the kernel, ensure that `sample_chain`'s `parallel_iterations=1`, otherwise results will not be reproducible. In addition to returning the chain state, this function supports tracing of auxiliary variables used by the kernel. The traced values are selected by specifying `trace_fn`. By default, all kernel results are traced but in the future the default will be changed to no results being traced, so plan accordingly. See below for some examples of this feature.

Value

list of:

- `checkpointable_states_and_trace`: if `return_final_kernel_results` is TRUE. The return value is an instance of `CheckpointableStatesAndTrace`.
- `all_states`: if `return_final_kernel_results` is FALSE and `trace_fn` is NULL. The return value is a Tensor or Python list of Tensors representing the state(s) of the Markov chain(s) at each result step. Has same shape as input `current_state` but with a prepended `num_results-size` dimension.
- `states_and_trace`: if `return_final_kernel_results` is FALSE and `trace_fn` is not NULL. The return value is an instance of `StatesAndTrace`.

References

- Art B. Owen. *Statistically efficient thinning of a Markov chain sampler*. *Technical Report*, 2017.

See Also

Other `mcmc_functions`: `mcmc_effective_sample_size()`, `mcmc_potential_scale_reduction()`, `mcmc_sample_annealed_importance_chain()`, `mcmc_sample_halton_sequence()`

Examples

```

dims <- 10
true_stddev <- sqrt(seq(1, 3, length.out = dims))
likelihood <- tfd_multivariate_normal_diag(scale_diag = true_stddev)

kernel <- mcmc_hamiltonian_monte_carlo(
  target_log_prob_fn = likelihood$log_prob,
  step_size = 0.5,
  num_leapfrog_steps = 2
)

states <- kernel %>% mcmc_sample_chain(
  num_results = 1000,
  num_burnin_steps = 500,
  current_state = rep(0, dims),
  trace_fn = NULL
)

sample_mean <- tf$reduce_mean(states, axis = 0L)
sample_stddev <- tf$sqrt(
  tf$reduce_mean(tf$math$ssquared_difference(states, sample_mean), axis = 0L))

```

mcmc_sample_halton_sequence

Returns a sample from the dim dimensional Halton sequence.

Description

Warning: The sequence elements take values only between 0 and 1. Care must be taken to appropriately transform the domain of a function if it differs from the unit cube before evaluating integrals using Halton samples. It is also important to remember that quasi-random numbers without randomization are not a replacement for pseudo-random numbers in every context. Quasi random numbers are completely deterministic and typically have significant negative autocorrelation unless randomization is used.

Usage

```

mcmc_sample_halton_sequence(
  dim,
  num_results = NULL,
  sequence_indices = NULL,
  dtype = tf$float32,
  randomized = TRUE,
  seed = NULL,
  name = NULL
)

```

Arguments

<code>dim</code>	Positive integer representing each sample's <code>event_size</code> . Must not be greater than 1000.
<code>num_results</code>	(Optional) Positive scalar Tensor of dtype <code>int32</code> . The number of samples to generate. Either this parameter or <code>sequence_indices</code> must be specified but not both. If this parameter is <code>None</code> , then the behaviour is determined by the <code>sequence_indices</code> . Default value: <code>NULL</code> .
<code>sequence_indices</code>	(Optional) Tensor of dtype <code>int32</code> and rank 1. The elements of the sequence to compute specified by their position in the sequence. The entries index into the Halton sequence starting with 0 and hence, must be whole numbers. For example, <code>sequence_indices=[0, 5, 6]</code> will produce the first, sixth and seventh elements of the sequence. If this parameter is <code>None</code> , then the <code>num_results</code> parameter must be specified which gives the number of desired samples starting from the first sample. Default value: <code>NULL</code> .
<code>dtype</code>	(Optional) The dtype of the sample. One of: <code>float16</code> , <code>float32</code> or <code>float64</code> . Default value: <code>tf\$float32</code> .
<code>randomized</code>	(Optional) bool indicating whether to produce a randomized Halton sequence. If <code>TRUE</code> , applies the randomization described in Owen (2017). Default value: <code>TRUE</code> .
<code>seed</code>	(Optional) integer to seed the random number generator. Only used if <code>randomized</code> is <code>TRUE</code> . If not supplied and <code>randomized</code> is <code>TRUE</code> , no seed is set. Default value: <code>NULL</code> .
<code>name</code>	(Optional) string describing ops managed by this function. If not supplied the name of this function is used. Default value: <code>"sample_halton_sequence"</code> .

Details

Computes the members of the low discrepancy Halton sequence in dimension `dim`. The `dim`-dimensional sequence takes values in the unit hypercube in `dim` dimensions. Currently, only dimensions up to 1000 are supported. The prime base for the `k`-th axes is the `k`-th prime starting from 2. For example, if `dim = 3`, then the bases will be `[2, 3, 5]` respectively and the first element of the non-randomized sequence will be: `[0.5, 0.333, 0.2]`. For a more complete description of the Halton sequences see [here](#). For low discrepancy sequences and their applications see [here](#).

If `randomized` is true, this function produces a scrambled version of the Halton sequence introduced by Owen (2017). For the advantages of randomization of low discrepancy sequences see [here](#).

The number of samples produced is controlled by the `num_results` and `sequence_indices` parameters. The user must supply either `num_results` or `sequence_indices` but not both. The former is the number of samples to produce starting from the first element. If `sequence_indices` is given instead, the specified elements of the sequence are generated. For example, `sequence_indices=tf$range(10)` is equivalent to specifying `n=10`.

Value

`halton_elements` Elements of the Halton sequence. Tensor of supplied dtype and shape `[num_results, dim]` if `num_results` was specified or shape `[s, dim]` where `s` is the size of `sequence_indices` if `sequence_indices` were specified.

References

- [Art B. Owen. A randomized Halton algorithm in R. *arXiv preprint arXiv:1706.02808*, 2017.](#)

See Also

For an example how to use see `mcmc_sample_chain()`.

Other mcmc_functions: `mcmc_effective_sample_size()`, `mcmc_potential_scale_reduction()`, `mcmc_sample_annealed_importance_chain()`, `mcmc_sample_chain()`

mcmc_simple_step_size_adaptation

Adapts the inner kernel's step_size based on log_accept_prob.

Description

The simple policy multiplicatively increases or decreases the `step_size` of the inner kernel based on the value of `log_accept_prob`. It is based on equation 19 of Andrieu and Thoms (2008). Given enough steps and small enough `adaptation_rate` the median of the distribution of the acceptance probability will converge to the `target_accept_prob`. A good target acceptance probability depends on the inner kernel. If this kernel is HamiltonianMonteCarlo, then 0.6-0.9 is a good range to aim for. For RandomWalkMetropolis this should be closer to 0.25. See the individual kernels' docstrings for guidance.

Usage

```
mcmc_simple_step_size_adaptation(
    inner_kernel,
    num_adaptation_steps,
    target_accept_prob = 0.75,
    adaptation_rate = 0.01,
    step_size_setter_fn = NULL,
    step_size_getter_fn = NULL,
    log_accept_prob_getter_fn = NULL,
    validate_args = FALSE,
    name = NULL
)
```

Arguments

`inner_kernel` TransitionKernel-like object.

`num_adaptation_steps`

Scalar integer Tensor number of initial steps to during which to adjust the step size. This may be greater, less than, or equal to the number of burnin steps.

`target_accept_prob`

A floating point Tensor representing desired acceptance probability. Must be a positive number less than 1. This can either be a scalar, or have shape `list(num_chains)`. Default value: 0.75 (the center of asymptotically optimal rate for HMC).

<code>adaptation_rate</code>	Tensor representing amount to scale the current <code>step_size</code> .
<code>step_size_setter_fn</code>	A function with the signature <code>(kernel_results, new_step_size) -> new_kernel_results</code> where <code>kernel_results</code> are the results of the <code>inner_kernel</code> , <code>new_step_size</code> is a Tensor or a nested collection of Tensors with the same structure as returned by the <code>step_size_getter_fn</code> , and <code>new_kernel_results</code> are a copy of <code>kernel_results</code> with the step size(s) set.
<code>step_size_getter_fn</code>	A function with the signature <code>(kernel_results) -> step_size</code> where <code>kernel_results</code> are the results of the <code>inner_kernel</code> , and <code>step_size</code> is a floating point Tensor or a nested collection of such Tensors.
<code>log_accept_prob_getter_fn</code>	A function with the signature <code>(kernel_results) -> log_accept_prob</code> where <code>kernel_results</code> are the results of the <code>inner_kernel</code> , and <code>log_accept_prob</code> is a floating point Tensor. <code>log_accept_prob</code> can either be a scalar, or have shape <code>list(num_chains)</code> . If it's the latter, <code>step_size</code> should also have the same leading dimension.
<code>validate_args</code>	Logical. When True kernel parameters are checked for validity. When False invalid inputs may silently render incorrect outputs.
<code>name</code>	string prefixed to Ops created by this class. Default: "simple_step_size_adaptation".

Details

In general, adaptation prevents the chain from reaching a stationary distribution, so obtaining consistent samples requires `num_adaptation_steps` be set to a value somewhat smaller than the number of burnin steps. However, it may sometimes be helpful to set `num_adaptation_steps` to a larger value during development in order to inspect the behavior of the chain during adaptation.

The step size is assumed to broadcast with the chain state, potentially having leading dimensions corresponding to multiple chains. When there are fewer of those leading dimensions than there are chain dimensions, the corresponding dimensions in the `log_accept_prob` are averaged (in the direct space, rather than the log space) before being used to adjust the step size. This means that this kernel can do both cross-chain adaptation, or per-chain step size adaptation, depending on the shape of the step size.

For example, if your problem has a state with shape `[S]`, your chain state has shape `[C0, C1, Y]` (meaning that there are $C0 * C1$ total chains) and `log_accept_prob` has shape `[C0, C1]` (one acceptance probability per chain), then depending on the shape of the step size, the following will happen:

- Step size has shape `[], [S]` or `[1]`, the `log_accept_prob` will be averaged across its `C0` and `C1` dimensions. This means that you will learn a shared step size based on the mean acceptance probability across all chains. This can be useful if you don't have a lot of steps to adapt and want to average away the noise.
- Step size has shape `[C1, 1]` or `[C1, S]`, the `log_accept_prob` will be averaged across its `C0` dimension. This means that you will learn a shared step size based on the mean acceptance probability across chains that share the coordinate across the `C1` dimension. This can be useful when the `C1` dimension indexes different distributions, while `C0` indexes replicas of a single distribution, all sampled in parallel.

- Step size has shape [C0, C1, 1] or [C0, C1, S], then no averaging will happen. This means that each chain will learn its own step size. This can be useful when all chains are sampling from different distributions. Even when all chains are for the same distribution, this can help during the initial warmup period.
- Step size has shape [C0, 1, 1] or [C0, 1, S], the `log_accept_prob` will be averaged across its C1 dimension. This means that you will learn a shared step size based on the mean acceptance probability across chains that share the coordinate across the C0 dimension. This can be useful when the C0 dimension indexes different distributions, while C1 indexes replicas of a single distribution, all sampled in parallel.

Value

a Monte Carlo sampling kernel

References

- [Andrieu, Christophe, Thoms, Johannes. A tutorial on adaptive MCMC. *Statistics and Computing*, 2008.](#)
- [Betancourt, M. J., Byrne, S., & Girolami, M. \(2014\). *Optimizing The Integrator Step Size for Hamiltonian Monte Carlo*.](#)

See Also

Other `mcmc_kernels`: [mcmc_dual_averaging_step_size_adaptation\(\)](#), [mcmc_hamiltonian_monte_carlo\(\)](#), [mcmc_metropolis_adjusted_langevin_algorithm\(\)](#), [mcmc_metropolis_hastings\(\)](#), [mcmc_no_u_turn_sampler\(\)](#), [mcmc_random_walk_metropolis\(\)](#), [mcmc_replica_exchange_mc\(\)](#), [mcmc_slice_sampler\(\)](#), [mcmc_transformed_trans](#), [mcmc_uncalibrated_hamiltonian_monte_carlo\(\)](#), [mcmc_uncalibrated_langevin\(\)](#), [mcmc_uncalibrated_random_w](#)

Examples

```
target_log_prob_fn <- tfd_normal(loc = 0, scale = 1)$log_prob
num_burnin_steps <- 500
num_results <- 500
num_chains <- 64L
step_size <- tf$fill(list(num_chains), 0.1)

kernel <- mcmc_hamiltonian_monte_carlo(
  target_log_prob_fn = target_log_prob_fn,
  num_leapfrog_steps = 2,
  step_size = step_size
) %>%
  mcmc_simple_step_size_adaptation(num_adaptation_steps = round(num_burnin_steps * 0.8))

res <- kernel %>% mcmc_sample_chain(
  num_results = num_results,
  num_burnin_steps = num_burnin_steps,
  current_state = rep(0, num_chains),
  trace_fn = function(x, pkr) {
    list (
```

```

        pkr$inner_results$accepted_results$step_size,
        pkr$inner_results$log_accept_ratio
    )
}
)

samples <- res$all_states
step_size <- res$trace[[1]]
log_accept_ratio <- res$trace[[2]]

```

mcmc_slice_sampler *Runs one step of the slice sampler using a hit and run approach*

Description

Slice Sampling is a Markov Chain Monte Carlo (MCMC) algorithm based, as stated by Neal (2003), on the observation that "...one can sample from a distribution by sampling uniformly from the region under the plot of its density function. A Markov chain that converges to this uniform distribution can be constructed by alternately uniform sampling in the vertical direction with uniform sampling from the horizontal slice defined by the current vertical position, or more generally, with some update that leaves the uniform distribution over this slice invariant". Mathematical details and derivations can be found in Neal (2003). The one dimensional slice sampler is extended to n-dimensions through use of a hit-and-run approach: choose a random direction in n-dimensional space and take a step, as determined by the one-dimensional slice sampling algorithm, along that direction (Belisle et al. 1993).

Usage

```

mcmc_slice_sampler(
  target_log_prob_fn,
  step_size,
  max_doublings,
  seed = NULL,
  name = NULL
)

```

Arguments

target_log_prob_fn	Function which takes an argument like <code>current_state</code> (if it's a list <code>current_state</code> will be unpacked) and returns its (possibly unnormalized) log-density under the target distribution.
step_size	Tensor or list of Tensors representing the step size for the leapfrog integrator. Must broadcast with the shape of <code>current_state</code> . Larger step sizes lead to faster progress, but too-large step sizes make rejection exponentially more likely. When possible, it's often helpful to match per-variable step sizes to the standard deviations of the target distribution in each variable.

max_doublings	Scalar positive int32 tf\$Tensor. The maximum number of doublings to consider.
seed	integer to seed the random number generator.
name	string prefixed to Ops created by this function. Default value: NULL (i.e., 'slice_sampler_kernel').

Details

The `one_step` function can update multiple chains in parallel. It assumes that all leftmost dimensions of `current_state` index independent chain states (and are therefore updated independently). The output of `target_log_prob_fn(*current_state)` should sum log-probabilities across all event dimensions. Slices along the rightmost dimensions may have different target distributions; for example, `current_state[0, :]` could have a different target distribution from `current_state[1, :]`. These semantics are governed by `target_log_prob_fn(*current_state)`. (The number of independent chains is `tf$size(target_log_prob_fn(*current_state))`.)

Note that the sampler only supports states where all components have a common dtype.

Value

list of `next_state` (Tensor or Python list of Tensors representing the state(s) of the Markov chain(s) at each result step. Has same shape as and `current_state`.) and `kernel_results` (collections\$namedtuple of internal calculations used to 'advance the chain').

References

- Radford M. Neal. *Slice Sampling*. *The Annals of Statistics*. 2003, Vol 31, No. 3 , 705-767.
- C.J.P. Belisle, H.E. Romeijn, R.L. Smith. *Hit-and-run algorithms for generating multivariate distributions*. *Math. Oper. Res.*, 18(1993), 225-266.

See Also

Other `mcmc_kernels`: `mcmc_dual_averaging_step_size_adaptation()`, `mcmc_hamiltonian_monte_carlo()`, `mcmc_metropolis_adjusted_langevin_algorithm()`, `mcmc_metropolis_hastings()`, `mcmc_no_u_turn_sampler()`, `mcmc_random_walk_metropolis()`, `mcmc_replica_exchange_mc()`, `mcmc_simple_step_size_adaptation()`, `mcmc_transformed_transition_kernel()`, `mcmc_uncalibrated_hamiltonian_monte_carlo()`, `mcmc_uncalibrated_langevin()`, `mcmc_uncalibrated_random_walk()`

`mcmc_transformed_transition_kernel`

Applies a bijector to the MCMC's state space

Description

The transformed transition kernel enables fitting a bijector which serves to decorrelate the Markov chain Monte Carlo (MCMC) event dimensions thus making the chain mix faster. This is particularly useful when the geometry of the target distribution is unfavorable. In such cases it may take many evaluations of the `target_log_prob_fn` for the chain to mix between faraway states.

Usage

```
mcmc_transformed_transition_kernel(inner_kernel, bijector, name = NULL)
```

Arguments

inner_kernel	TransitionKernel-like object which has a target_log_prob_fn argument.
bijector	bijector or list of bijectors. These bijectors use forward to map the inner_kernel state space to the state expected by inner_kernel\$target_log_prob_fn.
name	string prefixed to Ops created by this function. Default value: NULL (i.e., "transformed_kernel").

Details

The idea of training an affine function to decorrelate chain event dims was presented in Parno and Marzouk (2014). Used in conjunction with the Hamiltonian Monte Carlo transition kernel, the Parno and Marzouk (2014) idea is an instance of Riemannian manifold HMC (Girolami and Calderhead, 2011).

The transformed transition kernel enables arbitrary bijective transformations of arbitrary transition kernels, e.g., one could use bijectors `tfb_affine`, `tfb_real_nvp`, etc. with transition kernels `mcmc_hamiltonian_monte_carlo`, `mcmc_random_walk_metropolis`, etc.

Value

a Monte Carlo sampling kernel

References

- [Matthew Parno and Youssef Marzouk. Transport map accelerated Markov chain Monte Carlo. *arXiv preprint arXiv:1412.5492*, 2014.](#)
- [Mark Girolami and Ben Calderhead. Riemann manifold langevin and hamiltonian monte carlo methods. In *Journal of the Royal Statistical Society*, 2011.](#)

See Also

Other mcmc_kernels: `mcmc_dual_averaging_step_size_adaptation()`, `mcmc_hamiltonian_monte_carlo()`, `mcmc_metropolis_adjusted_langevin_algorithm()`, `mcmc_metropolis_hastings()`, `mcmc_no_u_turn_sampler()`, `mcmc_random_walk_metropolis()`, `mcmc_replica_exchange_mc()`, `mcmc_simple_step_size_adaptation()`, `mcmc_slice_sampler()`, `mcmc_uncalibrated_hamiltonian_monte_carlo()`, `mcmc_uncalibrated_langevin()`, `mcmc_uncalibrated_random_walk()`

```
mcmc_uncalibrated_hamiltonian_monte_carlo
```

Runs one step of Uncalibrated Hamiltonian Monte Carlo

Description

Warning: this kernel will not result in a chain which converges to the `target_log_prob`. To get a convergent MCMC, use `mcmc_hamiltonian_monte_carlo(...)` or `mcmc_metroplis_hastings(mcmc_uncalibrated_hamiltonian_monte_carlo(...))`. For more details on `UncalibratedHamiltonianMonteCarlo`, see `HamiltonianMonteCarlo`.

Usage

```
mcmc_uncalibrated_hamiltonian_monte_carlo(
  target_log_prob_fn,
  step_size,
  num_leapfrog_steps,
  state_gradients_are_stopped = FALSE,
  seed = NULL,
  store_parameters_in_results = FALSE,
  name = NULL
)
```

Arguments

<code>target_log_prob_fn</code>	Function which takes an argument like <code>current_state</code> (if it's a list <code>current_state</code> will be unpacked) and returns its (possibly unnormalized) log-density under the target distribution.
<code>step_size</code>	Tensor or list of Tensors representing the step size for the leapfrog integrator. Must broadcast with the shape of <code>current_state</code> . Larger step sizes lead to faster progress, but too-large step sizes make rejection exponentially more likely. When possible, it's often helpful to match per-variable step sizes to the standard deviations of the target distribution in each variable.
<code>num_leapfrog_steps</code>	Integer number of steps to run the leapfrog integrator for. Total progress per HMC step is roughly proportional to <code>step_size * num_leapfrog_steps</code> .
<code>state_gradients_are_stopped</code>	logical indicating that the proposed new state be run through <code>tf\$stop_gradient</code> . This is particularly useful when combining optimization over samples from the HMC chain. Default value: <code>FALSE</code> (i.e., do not apply <code>stop_gradient</code>).
<code>seed</code>	integer to seed the random number generator.
<code>store_parameters_in_results</code>	If <code>TRUE</code> , then <code>step_size</code> and <code>num_leapfrog_steps</code> are written to and read from eponymous fields in the kernel results objects returned from <code>one_step</code> and <code>bootstrap_results</code> . This allows wrapper kernels to adjust those parameters on the fly. This is incompatible with <code>step_size_update_fn</code> , which must be set to <code>NULL</code> .

name string prefixed to Ops created by this function. Default value: NULL (i.e., 'hmc_kernel').

Value

a Monte Carlo sampling kernel

See Also

Other mcmc_kernels: `mcmc_dual_averaging_step_size_adaptation()`, `mcmc_hamiltonian_monte_carlo()`, `mcmc_metropolis_adjusted_langevin_algorithm()`, `mcmc_metropolis_hastings()`, `mcmc_no_u_turn_sampler()`, `mcmc_random_walk_metropolis()`, `mcmc_replica_exchange_mc()`, `mcmc_simple_step_size_adaptation()`, `mcmc_slice_sampler()`, `mcmc_transformed_transition_kernel()`, `mcmc_uncalibrated_langevin()`, `mcmc_uncalibrated_random_walk()`

mcmc_uncalibrated_langevin

Runs one step of Uncalibrated Langevin discretized diffusion.

Description

The class generates a Langevin proposal using `_euler_method` function and also computes helper `UncalibratedLangevinKernelResults` for the next iteration. Warning: this kernel will not result in a chain which converges to the `target_log_prob`. To get a convergent MCMC, use `MetropolisAdjustedLangevinAlgorithm` or `MetropolisHastings(UncalibratedLangevin(...))`.

Usage

```
mcmc_uncalibrated_langevin(
  target_log_prob_fn,
  step_size,
  volatility_fn = NULL,
  parallel_iterations = 10,
  compute_acceptance = TRUE,
  seed = NULL,
  name = NULL
)
```

Arguments

`target_log_prob_fn` Function which takes an argument like `current_state` (if it's a list `current_state` will be unpacked) and returns its (possibly unnormalized) log-density under the target distribution.

`step_size` Tensor or list of Tensors representing the step size for the leapfrog integrator. Must broadcast with the shape of `current_state`. Larger step sizes lead to faster progress, but too-large step sizes make rejection exponentially more likely. When possible, it's often helpful to match per-variable step sizes to the standard deviations of the target distribution in each variable.

volatility_fn	function which takes an argument like <code>current_state</code> (or <code>*current_state</code> if it's a list) and returns volatility value at <code>current_state</code> . Should return a Tensor or list of Tensors that must broadcast with the shape of <code>current_state</code> . Defaults to the identity function.
parallel_iterations	the number of coordinates for which the gradients of the volatility matrix <code>volatility_fn</code> can be computed in parallel.
compute_acceptance	logical indicating whether to compute the Metropolis log-acceptance ratio used to construct <code>MetropolisAdjustedLangevinAlgorithm</code> kernel.
seed	integer to seed the random number generator.
name	String prefixed to Ops created by this function. Default value: NULL (i.e., 'mala_kernel').

Value

list of `next_state` (Tensor or Python list of Tensors representing the state(s) of the Markov chain(s) at each result step. Has same shape as and `current_state`.) and `kernel_results` (collections.namedtuple of internal calculations used to 'advance the chain').

See Also

Other `mcmc_kernels`: `mcmc_dual_averaging_step_size_adaptation()`, `mcmc_hamiltonian_monte_carlo()`, `mcmc_metropolis_adjusted_langevin_algorithm()`, `mcmc_metropolis_hastings()`, `mcmc_no_u_turn_sampler()`, `mcmc_random_walk_metropolis()`, `mcmc_replica_exchange_mc()`, `mcmc_simple_step_size_adaptation()`, `mcmc_slice_sampler()`, `mcmc_transformed_transition_kernel()`, `mcmc_uncalibrated_hamiltonian_monte_carlo`, `mcmc_uncalibrated_random_walk()`

`mcmc_uncalibrated_random_walk`

Generate proposal for the Random Walk Metropolis algorithm.

Description

Warning: this kernel will not result in a chain which converges to the `target_log_prob`. To get a convergent MCMC, use `mcmc_random_walk_metropolis(...)` or `mcmc_metropolis_hastings(mcmc_uncalibrated_ran`

Usage

```
mcmc_uncalibrated_random_walk(
    target_log_prob_fn,
    new_state_fn = NULL,
    seed = NULL,
    name = NULL
)
```

Arguments

target_log_prob_fn	Function which takes an argument like current_state ((if it's a list current_state will be unpacked) and returns its (possibly unnormalized) log-density under the target distribution.
new_state_fn	Function which takes a list of state parts and a seed; returns a same-type list of Tensors, each being a perturbation of the input state parts. The perturbation distribution is assumed to be a symmetric distribution centered at the input state part. Default value: NULL which is mapped to tfp\$mcmc\$random_walk_normal_fn().
seed	integer to seed the random number generator.
name	String name prefixed to Ops created by this function. Default value: NULL (i.e., 'rwm_kernel').

Value

a Monte Carlo sampling kernel

See Also

Other mcmc_kernels: [mcmc_dual_averaging_step_size_adaptation\(\)](#), [mcmc_hamiltonian_monte_carlo\(\)](#), [mcmc_metropolis_adjusted_langevin_algorithm\(\)](#), [mcmc_metropolis_hastings\(\)](#), [mcmc_no_u_turn_sampler\(\)](#), [mcmc_random_walk_metropolis\(\)](#), [mcmc_replica_exchange_mc\(\)](#), [mcmc_simple_step_size_adaptation\(\)](#), [mcmc_slice_sampler\(\)](#), [mcmc_transformed_transition_kernel\(\)](#), [mcmc_uncalibrated_hamiltonian_monte_carlo\(\)](#), [mcmc_uncalibrated_langevin\(\)](#)

params_size_categorical_mixture_of_one_hot_categorical

number of params needed to create a CategoricalMixtureOfOneHot-Categorical distribution

Description

number of params needed to create a CategoricalMixtureOfOneHotCategorical distribution

Usage

```
params_size_categorical_mixture_of_one_hot_categorical(
  event_size,
  num_components
)
```

Arguments

event_size	event size of this distribution
num_components	number of components in the mixture

Value

a scalar

params_size_independent_bernoulli
number of params needed to create an IndependentBernoulli distribution

Description

number of params needed to create an IndependentBernoulli distribution

Usage

params_size_independent_bernoulli(event_size)

Arguments

event_size event size of this distribution

Value

a scalar

params_size_independent_logistic
number of params needed to create an IndependentLogistic distribution

Description

number of params needed to create an IndependentLogistic distribution

Usage

params_size_independent_logistic(event_size)

Arguments

event_size event size of this distribution

Value

a scalar

params_size_independent_normal

number of params needed to create an IndependentNormal distribution

Description

number of params needed to create an IndependentNormal distribution

Usage

params_size_independent_normal(event_size)

Arguments

event_size event size of this distribution

Value

a scalar

params_size_independent_poisson

number of params needed to create an IndependentPoisson distribution

Description

number of params needed to create an IndependentPoisson distribution

Usage

params_size_independent_poisson(event_size)

Arguments

event_size event size of this distribution

Value

a scalar

params_size_mixture_logistic
number of params needed to create a MixtureLogistic distribution

Description

number of params needed to create a MixtureLogistic distribution

Usage

params_size_mixture_logistic(num_components, event_shape)

Arguments

num_components Number of component distributions in the mixture distribution.
event_shape Number of parameters needed to create a single component distribution.

Value

a scalar

params_size_mixture_normal
number of params needed to create a MixtureNormal distribution

Description

number of params needed to create a MixtureNormal distribution

Usage

params_size_mixture_normal(num_components, event_shape)

Arguments

num_components Number of component distributions in the mixture distribution.
event_shape Number of parameters needed to create a single component distribution.

Value

a scalar

params_size_mixture_same_family
number of params needed to create a MixtureSameFamily distribution

Description

number of params needed to create a MixtureSameFamily distribution

Usage

```
params_size_mixture_same_family(num_components, component_params_size)
```

Arguments

num_components Number of component distributions in the mixture distribution.
 component_params_size
 Number of parameters needed to create a single component distribution.

Value

a scalar

params_size_multivariate_normal_tri_l
number of params needed to create a MultivariateNormalTriL distribution

Description

number of params needed to create a MultivariateNormalTriL distribution

Usage

```
params_size_multivariate_normal_tri_l(event_size)
```

Arguments

event_size event size of this distribution

Value

a scalar

params_size_one_hot_categorical
number of params needed to create a OneHotCategorical distribution

Description

number of params needed to create a OneHotCategorical distribution

Usage

```
params_size_one_hot_categorical(event_size)
```

Arguments

event_size event size of this distribution

Value

a scalar

sts_additive_state_space_model
A state space model representing a sum of component state space models.

Description

A state space model (SSM) posits a set of latent (unobserved) variables that evolve over time with dynamics specified by a probabilistic transition model $p(z[t+1] | z[t])$. At each timestep, we observe a value sampled from an observation model conditioned on the current state, $p(x[t] | z[t])$. The special case where both the transition and observation models are Gaussians with mean specified as a linear function of the inputs, is known as a linear Gaussian state space model and supports tractable exact probabilistic calculations; see `tfd_linear_gaussian_state_space_model` for details.

Usage

```
sts_additive_state_space_model(  
  component_ssms,  
  constant_offset = 0,  
  observation_noise_scale = NULL,  
  initial_state_prior = NULL,  
  initial_step = 0,  
  validate_args = FALSE,  
  allow_nan_stats = TRUE,  
  name = NULL  
)
```

Arguments

<code>component_ssms</code>	list containing one or more <code>tfd_linear_gaussian_state_space_model</code> instances. The components will in general implement different time-series models, with possibly different <code>latent_size</code> , but they must have the same <code>dtype</code> , event shape (<code>num_timesteps</code> and <code>observation_size</code>), and their batch shapes must broadcast to a compatible batch shape.#'
<code>constant_offset</code>	scalar float tensor, or batch of scalars, specifying a constant value added to the sum of outputs from the component models. This allows the components to model the shifted series <code>observed_time_series - constant_offset</code> . Default value: <code>0.#'</code>
<code>observation_noise_scale</code>	Optional scalar float tensor indicating the standard deviation of the observation noise. May contain additional batch dimensions, which must broadcast with the batch shape of elements in <code>component_ssms</code> . If <code>observation_noise_scale</code> is specified for the <code>sts_additive_state_space_model</code> , the observation noise scales of component models are ignored. If <code>NULL</code> , the observation noise scale is derived by summing the noise variances of the component models, i.e., <code>observation_noise_scale = sqrt(sum([ssm.observation_noise_scale**2 for ssm in component_ssms]))</code> .
<code>initial_state_prior</code>	instance of <code>tfd_multivariate_normal</code> representing the prior distribution on latent states. Must have event shape <code>[1]</code> (as <code>tfd_linear_gaussian_state_space_model</code> requires a rank-1 event shape).
<code>initial_step</code>	Optional scalar integer tensor specifying the starting timestep. Default value: <code>0</code> .
<code>validate_args</code>	logical. Whether to validate input with asserts. If <code>validate_args</code> is <code>FALSE</code> , and the inputs are invalid, correct behavior is not guaranteed. Default value: <code>FALSE</code> .
<code>allow_nan_stats</code>	logical. If <code>FALSE</code> , raise an exception if a statistic (e.g. mean/mode/etc...) is undefined for any batch member. If <code>TRUE</code> , batch members with valid parameters leading to undefined statistics will return <code>NaN</code> for this statistic. Default value: <code>TRUE</code> .
<code>name</code>	string prefixed to ops created by this class. Default value: <code>"AdditiveStateSpace-Model"</code> .

Details

The `sts_additive_state_space_model` represents a sum of component state space models. Each of the N components describes a random process generating a distribution on observed time series $x_1[t], x_2[t], \dots, x_N[t]$. The additive model represents the sum of these processes, $y[t] = x_1[t] + x_2[t] + \dots + x_N[t] + \text{eps}[t]$, where $\text{eps}[t] \sim N(0, \text{observation_noise_scale})$ is an observation noise term.

Mathematical Details

The additive model concatenates the latent states of its component models. The generative process runs each component's dynamics in its own subspace of latent space, and then observes the sum of the observation models from the components.

Formally, the transition model is linear Gaussian:

$$p(z[t+1] \mid z[t]) \sim \text{Normal}(\text{loc} = \text{transition_matrix}.\text{matmul}(z[t]), \text{cov} = \text{transition_cov})$$

where each $z[t]$ is a latent state vector concatenating the component state vectors, $z[t] = [z1[t], z2[t], \dots, zN[t]]$, so it has size `latent_size = sum([c.latent_size for c in components])`.

The transition matrix is the block-diagonal composition of transition matrices from the component processes:

```
transition_matrix =
[[ c0.transition_matrix, 0., ..., 0. ],
 [ 0., c1.transition_matrix, ..., 0. ],
 [ ... .. ... ],
 [ 0., 0., ..., cN.transition_matrix ]]
```

and the noise covariance is similarly the block-diagonal composition of component noise covariances:

```
transition_cov =
[[ c0.transition_cov, 0., ..., 0. ],
 [ 0., c1.transition_cov, ..., 0. ],
 [ ... .. ... ],
 [ 0., 0., ..., cN.transition_cov ]]
```

The observation model is also linear Gaussian,

$$p(y[t] \mid z[t]) \sim \text{Normal}(\text{loc} = \text{observation_matrix}.\text{matmul}(z[t]), \text{stddev} = \text{observation_noise_scale})$$

This implementation assumes scalar observations, so `observation_matrix` has shape `[1, latent_size]`. The additive observation matrix simply concatenates the observation matrices from each component:

```
observation_matrix = concat([c0.obs_matrix, c1.obs_matrix, ..., cN.obs_matrix], axis=-1)
```

The effect is that each component observation matrix acts on the dimensions of latent state corresponding to that component, and the overall expected observation is the sum of the expected observations from each component.

If `observation_noise_scale` is not explicitly specified, it is also computed by summing the noise variances of the component processes:

```
observation_noise_scale = sqrt(sum([c.observation_noise_scale**2 for c in components]))
```

Value

an instance of `LinearGaussianStateSpaceModel`.

See Also

Other sts: `sts_autoregressive_state_space_model()`, `sts_autoregressive()`, `sts_constrained_seasonal_state_space_model()`, `sts_dynamic_linear_regression_state_space_model()`, `sts_dynamic_linear_regression()`, `sts_linear_regression()`, `sts_local_level_state_space_model()`, `sts_local_level()`, `sts_local_linear_trend_state_space_model()`, `sts_local_linear_trend()`, `sts_seasonal_state_space_model()`, `sts_seasonal()`, `sts_semi_local_linear_trend_state_space_model()`, `sts_semi_local_linear_trend()`, `sts_smooth_seasonal_state_space_model()`, `sts_smooth_seasonal()`, `sts_sparse_linear_regression()`, `sts_sum()`

sts_autoregressive *Formal representation of an autoregressive model.*

Description

An autoregressive (AR) model posits a latent level whose value at each step is a noisy linear combination of previous steps:

$$\text{level}[t+1] = (\text{sum}(\text{coefficients} * \text{levels}[t:t\text{-order}:-1]) + \text{Normal}(0., \text{level_scale}))$$
Usage

```
sts_autoregressive(
  observed_time_series = NULL,
  order,
  coefficients_prior = NULL,
  level_scale_prior = NULL,
  initial_state_prior = NULL,
  coefficient_constraining_bijector = NULL,
  name = NULL
)
```

Arguments

`observed_time_series` optional float tensor of shape `batch_shape + [T, 1]` (omitting the trailing unit dimension is also supported when `T > 1`), specifying an observed time series. Any priors not explicitly set will be given default values according to the scale of the observed time series (or batch of time series). May optionally be an instance of `sts_masked_time_series`, which includes a mask tensor to specify timesteps with missing observations. Default value: `NULL`.

`order` scalar positive integer specifying the number of past timesteps to regress on.

`coefficients_prior` optional Distribution instance specifying a prior on the coefficients parameter. If `NULL`, a default standard normal (`tfd_multivariate_normal_diag(scale_diag = tf.ones(list(order)))`) prior is used. Default value: `NULL`.

<code>level_scale_prior</code>	optional <code>Distribution</code> instance specifying a prior on the <code>level_scale</code> parameter. If <code>NULL</code> , a heuristic default prior is constructed based on the provided <code>observed_time_series</code> . Default value: <code>NULL</code> .
<code>initial_state_prior</code>	optional <code>Distribution</code> instance specifying a prior on the initial state, corresponding to the values of the process at a set of size order of imagined timesteps before the initial step. If <code>NULL</code> , a heuristic default prior is constructed based on the provided <code>observed_time_series</code> . Default value: <code>NULL</code> .
<code>coefficient_constraining_bijector</code>	optional <code>Bijector</code> instance representing a constraining mapping for the autoregressive coefficients. For example, <code>tfb_tanh()</code> constrains the coefficients to lie in $(-1, 1)$, while <code>tfb_softplus()</code> constrains them to be positive, and <code>tfb_identity()</code> implies no constraint. If <code>NULL</code> , the default behavior constrains the coefficients to lie in $(-1, 1)$ using a <code>tanh</code> bijector. Default value: <code>NULL</code> .
<code>name</code>	the name of this model component. Default value: <code>'Autoregressive'</code> .

Details

The latent state is `levels[t:t-order:-1]`. We observe a noisy realization of the current level: `f[t] = level[t] + Normal(0., observation_noise_scale)` at each timestep.

If `coefficients=[1.]`, the AR process is a simple random walk, equivalent to a `LocalLevel` model. However, a random walk's variance increases with time, while many AR processes (in particular, any first-order process with `abs(coefficient) < 1`) are *stationary*, i.e., they maintain a constant variance over time. This makes AR processes useful models of uncertainty.

Value

an instance of `StructuralTimeSeries`.

See Also

For usage examples see `sts_fit_with_hmc()`, `sts_forecast()`, `sts_decompose_by_component()`.

Other sts: `sts_additive_state_space_model()`, `sts_autoregressive_state_space_model()`, `sts_constrained_seasonal_state_space_model()`, `sts_dynamic_linear_regression_state_space_model()`, `sts_dynamic_linear_regression()`, `sts_linear_regression()`, `sts_local_level_state_space_model()`, `sts_local_level()`, `sts_local_linear_trend_state_space_model()`, `sts_local_linear_trend()`, `sts_seasonal_state_space_model()`, `sts_seasonal()`, `sts_semi_local_linear_trend_state_space_model()`, `sts_semi_local_linear_trend()`, `sts_smooth_seasonal_state_space_model()`, `sts_smooth_seasonal()`, `sts_sparse_linear_regression()`, `sts_sum()`

 sts_autoregressive_state_space_model

State space model for an autoregressive process.

Description

A state space model (SSM) posits a set of latent (unobserved) variables that evolve over time with dynamics specified by a probabilistic transition model $p(z[t+1] | z[t])$. At each timestep, we observe a value sampled from an observation model conditioned on the current state, $p(x[t] | z[t])$. The special case where both the transition and observation models are Gaussians with mean specified as a linear function of the inputs, is known as a linear Gaussian state space model and supports tractable exact probabilistic calculations; see `tfd_linear_gaussian_state_space_model` for details.

Usage

```
sts_autoregressive_state_space_model(
    num_timesteps,
    coefficients,
    level_scale,
    initial_state_prior,
    observation_noise_scale = 0,
    initial_step = 0,
    validate_args = FALSE,
    name = NULL
)
```

Arguments

<code>num_timesteps</code>	Scalar integer tensor number of timesteps to model with this distribution.
<code>coefficients</code>	float tensor of shape <code>tf\$concat(batch_shape, list(order))</code> defining the autoregressive coefficients. The coefficients are defined backwards in time: <code>coefficients[0] * level[t] + coefficients[1] * level[t-1] + ... + coefficients[order-1] * level[t-order+1]</code> .
<code>level_scale</code>	Scalar (any additional dimensions are treated as batch dimensions) float tensor indicating the standard deviation of the transition noise at each step.
<code>initial_state_prior</code>	instance of <code>tfd_multivariate_normal</code> representing the prior distribution on latent states. Must have event shape <code>list(order)</code> .
<code>observation_noise_scale</code>	Scalar (any additional dimensions are treated as batch dimensions) float tensor indicating the standard deviation of the observation noise. Default value: 0.
<code>initial_step</code>	Optional scalar int tensor specifying the starting timestep. Default value: 0.
<code>validate_args</code>	logical. Whether to validate input with asserts. If <code>validate_args</code> is <code>FALSE</code> , and the inputs are invalid, correct behavior is not guaranteed. Default value: <code>FALSE</code> .

name name prefixed to ops created by this class. Default value: "AutoregressiveStateSpaceModel".

Details

In an autoregressive process, the expected level at each timestep is a linear function of previous levels, with added Gaussian noise:

$$\text{level}[t+1] = (\text{sum}(\text{coefficients} * \text{levels}[t:t-\text{order}:-1]) + \text{Normal}(0., \text{level_scale}))$$

The process is characterized by a vector `coefficients` whose size determines the order of the process (how many previous values it looks at), and by `level_scale`, the standard deviation of the noise added at each step. This is formulated as a state space model by letting the latent state encode the most recent values; see 'Mathematical Details' below.

The parameters `level_scale` and `observation_noise_scale` are each (a batch of) scalars, and `coefficients` is a (batch) vector of size `list(order)`. The batch shape of this Distribution is the broadcast batch shape of these parameters and of the `initial_state_prior`.

Mathematical Details

The autoregressive model implements a `tfd_linear_gaussian_state_space_model` with `latent_size = order` and `observation_size = 1`. The latent state vector encodes the recent history of the process, with the current value in the topmost dimension. At each timestep, the transition sums the previous values to produce the new expected value, shifts all other values down by a dimension, and adds noise to the current value. This is formally encoded by the transition model:

$$\text{transition_matrix} = \begin{bmatrix} \text{coefs}[0], & \text{coefs}[1], & \dots, & \text{coefs}[\text{order}] \\ 1., & 0., & \dots, & 0. \\ 0., & 1., & \dots, & 0. \\ \dots & & & \\ 0., & 0., & \dots, & 1., & 0. \end{bmatrix}$$

$$\text{transition_noise} \sim \text{N}(\text{loc}=0., \text{scale}=\text{diag}([\text{level_scale}, 0., 0., \dots, 0.]))$$

The observation model simply extracts the current (topmost) value, and optionally adds independent noise at each step:

$$\begin{aligned} \text{observation_matrix} &= [[1., 0., \dots, 0.]] \\ \text{observation_noise} &\sim \text{N}(\text{loc}=0, \text{scale}=\text{observation_noise_scale}) \end{aligned}$$

Models with `observation_noise_scale = 0` are AR processes in the formal sense. Setting `observation_noise_scale` to a nonzero value corresponds to a latent AR process observed under an iid noise model.

Value

an instance of `LinearGaussianStateSpaceModel`.

See Also

Other sts: `sts_additive_state_space_model()`, `sts_autoregressive()`, `sts_constrained_seasonal_state_space_model()`, `sts_dynamic_linear_regression_state_space_model()`, `sts_dynamic_linear_regression()`, `sts_linear_regression()`, `sts_local_level_state_space_model()`, `sts_local_level()`, `sts_local_linear_trend()`, `sts_local_linear_trend()`, `sts_seasonal_state_space_model()`, `sts_seasonal()`, `sts_semi_local_linear_trend()`, `sts_semi_local_linear_trend()`, `sts_smooth_seasonal_state_space_model()`, `sts_smooth_seasonal()`, `sts_sparse_linear_regression()`, `sts_sum()`

`sts_build_factored_surrogate_posterior`

Build a variational posterior that factors over model parameters.

Description

The surrogate posterior consists of independent Normal distributions for each parameter with trainable loc and scale, transformed using the parameter's bijector to the appropriate support space for that parameter.

Usage

```
sts_build_factored_surrogate_posterior(
  model,
  batch_shape = list(),
  seed = NULL,
  name = NULL
)
```

Arguments

<code>model</code>	An instance of <code>StructuralTimeSeries</code> representing a time-series model. This represents a joint distribution over time-series and their parameters with batch shape <code>[b1, ..., bN].#</code>
<code>batch_shape</code>	Batch shape (<code>list</code> , or <code>integer</code>) of initial states to optimize in parallel. Default value: <code>list()</code> . (i.e., just run a single optimization).
<code>seed</code>	<code>integer</code> to seed the random number generator.
<code>name</code>	string prefixed to ops created by this function. Default value: <code>NULL</code> (i.e., <code>'build_factored_surrogate_posterior'</code>).

Value

variational_posterior `tfd_joint_distribution_named` defining a trainable surrogate posterior over model parameters. Samples from this distribution are named lists with character parameter names as keys.

See Also

Other sts-functions: [sts_build_factored_variational_loss\(\)](#), [sts_decompose_by_component\(\)](#), [sts_decompose_forecast_by_component\(\)](#), [sts_fit_with_hmc\(\)](#), [sts_forecast\(\)](#), [sts_one_step_predictive\(\)](#), [sts_sample_uniform_initial_state\(\)](#)

sts_build_factored_variational_loss

Build a loss function for variational inference in STS models.

Description

Variational inference searches for the distribution within some family of approximate posteriors that minimizes a divergence between the approximate posterior $q(z)$ and true posterior $p(z|\text{observed_time_series})$. By converting inference to optimization, it's generally much faster than sampling-based inference algorithms such as HMC. The tradeoff is that the approximating family rarely contains the true posterior, so it may miss important aspects of posterior structure (in particular, dependence between variables) and should not be blindly trusted. Results may vary; it's generally wise to compare to HMC to evaluate whether inference quality is sufficient for your task at hand.

Usage

```
sts_build_factored_variational_loss(
    observed_time_series,
    model,
    init_batch_shape = list(),
    seed = NULL,
    name = NULL
)
```

Arguments

observed_time_series	float tensor of shape <code>concat([sample_shape, model.batch_shape, [num_timesteps, 1]])</code> where <code>sample_shape</code> corresponds to i.i.d. observations, and the trailing <code>[1]</code> dimension may (optionally) be omitted if <code>num_timesteps > 1</code> . May optionally be an instance of <code>sts_masked_time_series</code> , which includes a mask tensor to specify timesteps with missing observations.
model	An instance of <code>StructuralTimeSeries</code> representing a time-series model. This represents a joint distribution over time-series and their parameters with batch shape <code>[b1, ..., bN]</code> .
init_batch_shape	Batch shape (<code>list</code>) of initial states to optimize in parallel. Default value: <code>list()</code> . (i.e., just run a single optimization).
seed	integer to seed the random number generator.
name	name prefixed to ops created by this function. Default value: <code>NULL</code> (i.e., <code>'build_factored_variational_loss'</code>)

Details

This method constructs a loss function for variational inference using the Kullback-Liebler divergence $KL[q(z) || p(z|observed_time_series)]$, with an approximating family given by independent Normal distributions transformed to the appropriate parameter space for each parameter. Minimizing this loss (the negative ELBO) maximizes a lower bound on the log model evidence $-\log p(observed_time_series)$. This is equivalent to the 'mean-field' method implemented in Kucukelbir et al. (2017) and is a standard approach. The resulting posterior approximations are unimodal; they will tend to underestimate posterior uncertainty when the true posterior contains multiple modes (the $KL[q || p]$ divergence encourages choosing a single mode) or dependence between variables.

Value

list of:

- `variational_loss`: float Tensor of shape `tf$concat([init_batch_shape, model$batch_shape])`, encoding a stochastic estimate of an upper bound on the negative model evidence $-\log p(y)$. Minimizing this loss performs variational inference; the gap between the variational bound and the true (generally unknown) model evidence corresponds to the divergence $KL[q || p]$ between the approximate and true posterior.
- `variational_distributions`: a named list giving the approximate posterior for each model parameter. The keys are character parameter names in order, corresponding to `[param.name for param in model.parameters]`. The values are `tfd$Distribution` instances with batch shape `tf$concat([init_batch_shape, model$batch_shape])`; these will typically be of the form `tfd$TransformedDistribution(tfd.Normal(...), bijector=param.bijector)`

References

- Alp Kucukelbir, Dustin Tran, Rajesh Ranganath, Andrew Gelman, and David M. Blei. Automatic Differentiation Variational Inference. In *Journal of Machine Learning Research*, 2017.

See Also

Other sts-functions: `sts_build_factored_surrogate_posterior()`, `sts_decompose_by_component()`, `sts_decompose_forecast_by_component()`, `sts_fit_with_hmc()`, `sts_forecast()`, `sts_one_step_predictive()`, `sts_sample_uniform_initial_state()`

sts_constrained_seasonal_state_space_model

Seasonal state space model with effects constrained to sum to zero.

Description

Seasonal state space model with effects constrained to sum to zero.

Usage

```

sts_constrained_seasonal_state_space_model(
    num_timesteps,
    num_seasons,
    drift_scale,
    initial_state_prior,
    observation_noise_scale = 1e-04,
    num_steps_per_season = 1,
    initial_step = 0,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = NULL
)

```

Arguments

<code>num_timesteps</code>	Scalar integer tensor number of timesteps to model with this distribution.
<code>num_seasons</code>	Scalar integer number of seasons.
<code>drift_scale</code>	Scalar (any additional dimensions are treated as batch dimensions) float tensor indicating the standard deviation of the change in effect between consecutive occurrences of a given season. This is assumed to be the same for all seasons.
<code>initial_state_prior</code>	instance of <code>tfd_multivariate_normal</code> representing the prior distribution on latent states; must have event shape <code>[num_seasons]</code> .
<code>observation_noise_scale</code>	Scalar (any additional dimensions are treated as batch dimensions) float tensor indicating the standard deviation of the observation noise.
<code>num_steps_per_season</code>	integer number of steps in each season. This may be either a scalar (<code>shape []</code>), in which case all seasons have the same length, or an array of shape <code>[num_seasons]</code> , in which seasons have different length, but remain constant around different cycles, or an array of shape <code>[num_cycles, num_seasons]</code> , in which <code>num_steps_per_season</code> for each season also varies in different cycle (e.g., a 4 years cycle with leap day). Default value: 1.
<code>initial_step</code>	Optional scalar integer tensor specifying the starting timestep. Default value: 0.
<code>validate_args</code>	logical. Whether to validate input with asserts. If <code>validate_args</code> is <code>FALSE</code> , and the inputs are invalid, correct behavior is not guaranteed. Default value: <code>FALSE</code> .
<code>allow_nan_stats</code>	logical. If <code>FALSE</code> , raise an exception if a statistic (e.g. mean/mode/etc...) is undefined for any batch member. If <code>TRUE</code> , batch members with valid parameters leading to undefined statistics will return NaN for this statistic. Default value: <code>TRUE</code> .
<code>name</code>	string prefixed to ops created by this class. Default value: "SeasonalStateSpace-Model".

Value

an instance of `LinearGaussianStateSpaceModel`.

See Also

[sts_seasonal_state_space_model\(\)](#).

Mathematical details

The constrained model implements a reparameterization of the naive `SeasonalStateSpaceModel`. Instead of directly representing the seasonal effects in the latent space, the latent space of the constrained model represents the difference between each effect and the mean effect. The following discussion assumes familiarity with the mathematical details of `SeasonalStateSpaceModel`.

Reparameterization and constraints: let the seasonal effects at a given timestep be $E = [e_1, \dots, e_N]$. The difference between each effect e_i and the mean effect is $z_i = e_i - \sum_i(e_i)/N$. By itself, this transformation is not invertible because recovering the absolute effects requires that we know the mean as well. To fix this, we'll define $z_N = \sum_i(e_i)/N$ as the mean effect. It's easy to see that this is invertible: given the mean effect and the differences of the first $N-1$ effects from the mean, it's easy to solve for all N effects. Formally, we've defined the invertible linear reparameterization $Z = R E$, where

$$R = \begin{bmatrix} 1 - 1/N, & -1/N, & \dots, & -1/N \\ -1/N, & 1 - 1/N, & \dots, & -1/N, \\ \dots & & & \\ 1/N, & 1/N, & \dots, & 1/N \end{bmatrix}$$

represents the change of basis from 'effect coordinates' E to 'residual coordinates' Z . The Z s form the latent space of the `ConstrainedSeasonalStateSpaceModel`. To constrain the mean effect z_N to zero, we fix the prior to zero, $p(z_N) \sim N(0, \theta)$, and after the transition at each timestep we project z_N back to zero. Note that this projection is linear: to set the N th dimension to zero, we simply multiply by the identity matrix with a missing element in the bottom right, i.e., $Z_{\text{constrained}} = P Z$, where $P = \text{eye}(N) - \text{scatter}(N-1, N-1, 1)$.

Model: concretely, suppose a naive seasonal effect model has initial state prior $N(m, S)$, transition matrix F and noise covariance Q , and observation matrix H . Then the corresponding constrained seasonal effect model has initial state prior $N(P R m, P R S R' P')$, transition matrix $P R F R^{-1}$ and noise covariance $F R Q R' F'$, and observation matrix $H R^{-1}$, where the change-of-basis matrix R and constraint projection matrix P are as defined above. This follows directly from applying the reparameterization $Z = R E$, and then enforcing the zero-sum constraint on the prior and transition noise covariances. In practice, because the sum of effects z_N is constrained to be zero, it will never contribute a term to any linear operation on the latent space, so we can drop that dimension from the model entirely. `ConstrainedSeasonalStateSpaceModel` does this, so that it implements the $N-1$ dimension latent space $z_1, \dots, z_{[N-1]}$. Note that since we constrained the mean effect to be zero, the latent z_i 's now recover their interpretation as the *actual* effects, $z_i = e_i$ for $i = 1, \dots, N-1$, even though they were originally defined as residuals. The N th effect is represented only implicitly, as the nonzero mean of the first $N-1$ effects. Although the computational representation is not symmetric across all N effects, we derived the `ConstrainedSeasonalStateSpaceModel` by starting with a symmetric representation and imposing only a symmetric constraint (the zero-sum constraint), so the probability model remains symmetric over all N seasonal effects.

Other sts: `sts_additive_state_space_model()`, `sts_autoregressive_state_space_model()`, `sts_autoregressive()`, `sts_dynamic_linear_regression_state_space_model()`, `sts_dynamic_linear_regression()`, `sts_linear_regression()`, `sts_local_level_state_space_model()`, `sts_local_level()`, `sts_local_linear_trend()`, `sts_local_linear_trend()`, `sts_seasonal_state_space_model()`, `sts_seasonal()`, `sts_semi_local_linear_trend()`, `sts_semi_local_linear_trend()`, `sts_smooth_seasonal_state_space_model()`, `sts_smooth_seasonal()`, `sts_sparse_linear_regression()`, `sts_sum()`

sts_decompose_by_component

Decompose an observed time series into contributions from each component.

Description

This method decomposes a time series according to the posterior representation of a structural time series model. In particular, it:

- Computes the posterior marginal mean and covariances over the additive model's latent space.
- Decomposes the latent posterior into the marginal blocks for each model component.
- Maps the per-component latent posteriors back through each component's observation model, to generate the time series modeled by that component.

Usage

```
sts_decompose_by_component(observed_time_series, model, parameter_samples)
```

Arguments

observed_time_series

float tensor of shape `concat([sample_shape, model.batch_shape, [num_timesteps, 1]])` where `sample_shape` corresponds to i.i.d. observations, and the trailing `[1]` dimension may (optionally) be omitted if `num_timesteps > 1`. May optionally be an instance of `sts_masked_time_series`, which includes a mask tensor to specify timesteps with missing observations.

model

An instance of `sts_sum` representing a structural time series model.

parameter_samples

list of tensors representing posterior samples of model parameters, with shapes `list(tf.concat(list(list(num_posterior_draws), param<1>$prior$batch_shape, param<1>$prior$event_shape), axis=0))` for all model parameters. This may optionally also be a named list mapping parameter names to tensor values.

Value

component_dists A named list mapping component `StructuralTimeSeries` instances (elements of `model$components`) to `Distribution` instances representing the posterior marginal distributions on the process modeled by each component. Each distribution has batch shape matching that of `posterior_means/posterior_covs`, and event shape of `list(num_timesteps)`.

See Also

Other sts-functions: `sts_build_factored_surrogate_posterior()`, `sts_build_factored_variational_loss()`, `sts_decompose_forecast_by_component()`, `sts_fit_with_hmc()`, `sts_forecast()`, `sts_one_step_predictive()`, `sts_sample_uniform_initial_state()`

Examples

```
observed_time_series <- array(rnorm(2 * 1 * 12), dim = c(2, 1, 12))
day_of_week <- observed_time_series %>% sts_seasonal(num_seasons = 7, name = "seasonal")
local_linear_trend <- observed_time_series %>% sts_local_linear_trend(name = "local_linear")
model <- observed_time_series %>%
  sts_sum(components = list(day_of_week, local_linear_trend))
states_and_results <- observed_time_series %>%
  sts_fit_with_hmc(
    model,
    num_results = 10,
    num_warmup_steps = 5,
    num_variational_steps = 15
  )
samples <- states_and_results[[1]]

component_dists <- observed_time_series %>%
  sts_decompose_by_component(model = model, parameter_samples = samples)
```

sts_decompose_forecast_by_component

Decompose a forecast distribution into contributions from each component.

Description

Decompose a forecast distribution into contributions from each component.

Usage

```
sts_decompose_forecast_by_component(model, forecast_dist, parameter_samples)
```

Arguments

model	An instance of <code>sts_sum</code> representing a structural time series model.
forecast_dist	A Distribution instance returned by <code>sts_forecast()</code> . (specifically, must be a <code>tfd.MixtureSameFamily</code> over a <code>tfd_linear_gaussian_state_space_model</code> parameterized by posterior samples).

parameter_samples

list of tensors representing posterior samples of model parameters, with shapes `list(tf$concat(list(list(num_posterior_draws), param<1>$prior$batch_shape, param<1>$prior$event_shape)))` for all model parameters. This may optionally also be a named list mapping parameter names to tensor values.

Value

component_dists A named list mapping component `StructuralTimeSeries` instances (elements of `model$components`) to `Distribution` instances representing the marginal forecast for each component. Each distribution has batch shape matching `forecast_dist` (specifically, the event shape is `[num_steps_forecast]`).

See Also

Other sts-functions: `sts_build_factored_surrogate_posterior()`, `sts_build_factored_variational_loss()`, `sts_decompose_by_component()`, `sts_fit_with_hmc()`, `sts_forecast()`, `sts_one_step_predictive()`, `sts_sample_uniform_initial_state()`

sts_dynamic_linear_regression

Formal representation of a dynamic linear regression model.

Description

The dynamic linear regression model is a special case of a linear Gaussian SSM and a generalization of typical (static) linear regression. The model represents regression weights with a latent state which evolves via a Gaussian random walk:

Usage

```
sts_dynamic_linear_regression(
  observed_time_series = NULL,
  design_matrix,
  drift_scale_prior = NULL,
  initial_weights_prior = NULL,
  name = NULL
)
```

Arguments

observed_time_series

optional float tensor of shape `batch_shape + [T, 1]` (omitting the trailing unit dimension is also supported when `T > 1`), specifying an observed time series. Any priors not explicitly set will be given default values according to the scale of the observed time series (or batch of time series). May optionally be an instance of `sts_masked_time_series`, which includes a mask tensor to specify timesteps with missing observations. Default value: `NULL`.

design_matrix	float tensor of shape <code>tf\$concat(list(batch_shape, list(num_timesteps, num_features)))</code> . This may also optionally be an instance of <code>tf\$linalg\$LinearOperator</code> .
drift_scale_prior	instance of <code>Distribution</code> specifying a prior on the <code>drift_scale</code> parameter. If <code>NULL</code> , a heuristic default prior is constructed based on the provided <code>observed_time_series</code> . Default value: <code>NULL</code> .
initial_weights_prior	instance of <code>tfd_multivariate_normal</code> representing the prior distribution on the latent states (the regression weights). Must have event shape <code>list(num_features)</code> . If <code>NULL</code> , a weakly-informative <code>Normal(0, 10)</code> prior is used. Default value: <code>NULL</code> .
name	the name of this component. Default value: <code>'DynamicLinearRegression'</code> .

Details

`weights[t] ~ Normal(weights[t-1], drift_scale)`

The latent state has dimension `num_features`, while the parameters `drift_scale` and `observation_noise_scale` are each (a batch of) scalars. The batch shape of this distribution is the broadcast batch shape of these parameters, the `initial_state_prior`, and the `design_matrix`. `num_features` is determined from the last dimension of `design_matrix` (equivalent to the number of columns in the design matrix in linear regression).

Value

an instance of `StructuralTimeSeries`.

See Also

For usage examples see `sts_fit_with_hmc()`, `sts_forecast()`, `sts_decompose_by_component()`.

Other sts: `sts_additive_state_space_model()`, `sts_autoregressive_state_space_model()`, `sts_autoregressive()`, `sts_constrained_seasonal_state_space_model()`, `sts_dynamic_linear_regression_state_space_model()`, `sts_linear_regression()`, `sts_local_level_state_space_model()`, `sts_local_level()`, `sts_local_linear_trend()`, `sts_local_linear_trend()`, `sts_seasonal_state_space_model()`, `sts_seasonal()`, `sts_semi_local_linear_trend()`, `sts_semi_local_linear_trend()`, `sts_smooth_seasonal_state_space_model()`, `sts_smooth_seasonal()`, `sts_sparse_linear_regression()`, `sts_sum()`

`sts_dynamic_linear_regression_state_space_model`

State space model for a dynamic linear regression from provided covariates.

Description

A state space model (SSM) posits a set of latent (unobserved) variables that evolve over time with dynamics specified by a probabilistic transition model $p(z[t+1] | z[t])$. At each timestep, we observe a value sampled from an observation model conditioned on the current state, $p(x[t] | z[t])$. The special case where both the transition and observation models are Gaussians with mean

specified as a linear function of the inputs, is known as a linear Gaussian state space model and supports tractable exact probabilistic calculations; see `tfd_linear_gaussian_state_space_model` for details.

Usage

```
sts_dynamic_linear_regression_state_space_model(
    num_timesteps,
    design_matrix,
    drift_scale,
    initial_state_prior,
    observation_noise_scale = 0,
    initial_step = 0,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = NULL
)
```

Arguments

<code>num_timesteps</code>	Scalar integer tensor, number of timesteps to model with this distribution.
<code>design_matrix</code>	float tensor of shape <code>tf\$concat(list(batch_shape, list(num_timesteps, num_features)))</code> . This may also optionally be an instance of <code>tf\$linalg\$LinearOperator</code> .
<code>drift_scale</code>	Scalar (any additional dimensions are treated as batch dimensions) float tensor indicating the standard deviation of the latent state transitions.
<code>initial_state_prior</code>	instance of <code>tfd_multivariate_normal</code> representing the prior distribution on latent states. Must have event shape <code>list(num_features)</code> .
<code>observation_noise_scale</code>	Scalar (any additional dimensions are treated as batch dimensions) float tensor indicating the standard deviation of the observation noise. Default value: 0.
<code>initial_step</code>	scalar integer tensor specifying the starting timestep. Default value: 0.
<code>validate_args</code>	logical. Whether to validate input with asserts. If <code>validate_args</code> is <code>FALSE</code> , and the inputs are invalid, correct behavior is not guaranteed. Default value: <code>FALSE</code> .
<code>allow_nan_stats</code>	logical. If <code>FALSE</code> , raise an exception if a statistic (e.g. mean/mode/etc...) is undefined for any batch member. If <code>TRUE</code> , batch members with valid parameters leading to undefined statistics will return NaN for this statistic. Default value: <code>TRUE</code> .
<code>name</code>	name prefixed to ops created by this class. Default value: <code>'DynamicLinearRegressionStateSpaceModel'</code> .

Details

The dynamic linear regression model is a special case of a linear Gaussian SSM and a generalization of typical (static) linear regression. The model represents regression weights with a latent state which evolves via a Gaussian random walk: $\text{weights}[t] \sim \text{Normal}(\text{weights}[t-1], \text{drift_scale})$

The latent state (the weights) has dimension `num_features`, while the parameters `drift_scale` and `observation_noise_scale` are each (a batch of) scalars. The batch shape of this Distribution is the broadcast batch shape of these parameters, the `initial_state_prior`, and the `design_matrix`. `num_features` is determined from the last dimension of `design_matrix` (equivalent to the number of columns in the design matrix in linear regression).

Mathematical Details

The dynamic linear regression model implements a `tfd_linear_gaussian_state_space_model` with `latent_size = num_features` and `observation_size = 1` following the transition model:

```
transition_matrix = eye(num_features)
transition_noise ~ Normal(0, diag([drift_scale]))
```

which implements the evolution of weights described above. The observation model is:

```
observation_matrix[t] = design_matrix[t]
observation_noise ~ Normal(0, observation_noise_scale)
```

Value

an instance of `LinearGaussianStateSpaceModel`.

See Also

Other sts: [sts_additive_state_space_model\(\)](#), [sts_autoregressive_state_space_model\(\)](#), [sts_autoregressive\(\)](#), [sts_constrained_seasonal_state_space_model\(\)](#), [sts_dynamic_linear_regression\(\)](#), [sts_linear_regression\(\)](#), [sts_local_level_state_space_model\(\)](#), [sts_local_level\(\)](#), [sts_local_linear_trend\(\)](#), [sts_local_linear_trend\(\)](#), [sts_seasonal_state_space_model\(\)](#), [sts_seasonal\(\)](#), [sts_semi_local_linear_trend\(\)](#), [sts_semi_local_linear_trend\(\)](#), [sts_smooth_seasonal_state_space_model\(\)](#), [sts_smooth_seasonal\(\)](#), [sts_sparse_linear_regression\(\)](#), [sts_sum\(\)](#)

sts_fit_with_hmc

Draw posterior samples using Hamiltonian Monte Carlo (HMC)

Description

Markov chain Monte Carlo (MCMC) methods are considered the gold standard of Bayesian inference; under suitable conditions and in the limit of infinitely many draws they generate samples from the true posterior distribution. HMC (Neal, 2011) uses gradients of the model's log-density function to propose samples, allowing it to exploit posterior geometry. However, it is computationally more expensive than variational inference and relatively sensitive to tuning.

Usage

```

sts_fit_with_hmc(
  observed_time_series,
  model,
  num_results = 100,
  num_warmup_steps = 50,
  num_leapfrog_steps = 15,
  initial_state = NULL,
  initial_step_size = NULL,
  chain_batch_shape = list(),
  num_variational_steps = 150,
  variational_optimizer = NULL,
  variational_sample_size = 5,
  seed = NULL,
  name = NULL
)

```

Arguments

observed_time_series float tensor of shape `concat([sample_shape, model.batch_shape, [num_timesteps, 1]])` where `sample_shape` corresponds to i.i.d. observations, and the trailing `[1]` dimension may (optionally) be omitted if `num_timesteps > 1`. May optionally be an instance of `sts_masked_time_series`, which includes a mask tensor to specify timesteps with missing observations.

model An instance of `StructuralTimeSeries` representing a time-series model. This represents a joint distribution over time-series and their parameters with batch shape `[b1, ..., bN]`.

num_results Integer number of Markov chain draws. Default value: 100.

num_warmup_steps Integer number of steps to take before starting to collect results. The warmup steps are also used to adapt the step size towards a target acceptance rate of 0.75. Default value: 50.

num_leapfrog_steps Integer number of steps to run the leapfrog integrator for. Total progress per HMC step is roughly proportional to `step_size * num_leapfrog_steps`. Default value: 15.

initial_state Optional Python list of Tensors, one for each model parameter, representing the initial state(s) of the Markov chain(s). These should have shape `tf.concat(list(chain_batch_shape, sample_shape), 0)`. If NULL, the initial state is set automatically using a sample from a variational posterior. Default value: NULL.

initial_step_size list of tensors, one for each model parameter, representing the step size for the leapfrog integrator. Must broadcast with the shape of `initial_state`. Larger step sizes lead to faster progress, but too-large step sizes make rejection exponentially more likely. If NULL, the step size is set automatically using the standard deviation of a variational posterior. Default value: NULL.

chain_batch_shape	Batch shape (list or int) of chains to run in parallel. Default value: list() (i.e., a single chain).
num_variational_steps	int number of steps to run the variational optimization to determine the initial state and step sizes. Default value: 150.
variational_optimizer	Optional tf\$train\$Optimizer instance to use in the variational optimization. If NULL, defaults to tf\$train\$AdamOptimizer(0.1). Default value: NULL.
variational_sample_size	integer number of Monte Carlo samples to use in estimating the variational divergence. Larger values may stabilize the optimization, but at higher cost per step in time and memory. Default value: 1.
seed	integer to seed the random number generator.
name	name prefixed to ops created by this function. Default value: NULL (i.e., 'fit_with_hmc').

Details

This method attempts to provide a sensible default approach for fitting StructuralTimeSeries models using HMC. It first runs variational inference as a fast posterior approximation, and initializes the HMC sampler from the variational posterior, using the posterior standard deviations to set per-variable step sizes (equivalently, a diagonal mass matrix). During the warmup phase, it adapts the step size to target an acceptance rate of 0.75, which is thought to be in the desirable range for optimal mixing (Betancourt et al., 2014).

Value

list of:

- samples: list of Tensors representing posterior samples of model parameters, with shapes [concat([num_results], chain_batch_shape, param.prior.batch_shape, param.prior.event_shape)] for param in model.parameters
- kernel_results: A (possibly nested) list of Tensors representing internal calculations made within the HMC sampler.

References

- Radford Neal. *MCMC Using Hamiltonian Dynamics*. *Handbook of Markov Chain Monte Carlo*, 2011.
- M.J. Betancourt, Simon Byrne, and Mark Girolami. *Optimizing The Integrator Step Size for Hamiltonian Monte Carlo*.

See Also

Other sts-functions: [sts_build_factored_surrogate_posterior\(\)](#), [sts_build_factored_variational_loss\(\)](#), [sts_decompose_by_component\(\)](#), [sts_decompose_forecast_by_component\(\)](#), [sts_forecast\(\)](#), [sts_one_step_predictive\(\)](#), [sts_sample_uniform_initial_state\(\)](#)

Examples

```

observed_time_series <-
  rep(c(3.5, 4.1, 4.5, 3.9, 2.4, 2.1, 1.2), 5) +
  rep(c(1.1, 1.5, 2.4, 3.1, 4.0), each = 7) %>%
  tensorflow::tf$convert_to_tensor(dtype = tensorflow::tf$float64)
day_of_week <- observed_time_series %>% sts_seasonal(num_seasons = 7)
local_linear_trend <- observed_time_series %>% sts_local_linear_trend()
model <- observed_time_series %>%
  sts_sum(components = list(day_of_week, local_linear_trend))
states_and_results <- observed_time_series %>%
  sts_fit_with_hmc(
    model,
    num_results = 10,
    num_warmup_steps = 5,
    num_variational_steps = 15)

```

sts_forecast

Construct predictive distribution over future observations

Description

Given samples from the posterior over parameters, return the predictive distribution over future observations for `num_steps_forecast` timesteps.

Usage

```

sts_forecast(
  observed_time_series,
  model,
  parameter_samples,
  num_steps_forecast
)

```

Arguments

`observed_time_series` float tensor of shape `concat([sample_shape, model.batch_shape, [num_timesteps, 1]])` where `sample_shape` corresponds to i.i.d. observations, and the trailing `[1]` dimension may (optionally) be omitted if `num_timesteps > 1`. May optionally be an instance of `sts_masked_time_series`, which includes a mask tensor to specify timesteps with missing observations.

`model` An instance of `StructuralTimeSeries` representing a time-series model. This represents a joint distribution over time-series and their parameters with batch shape `[b1, ..., bN]`.

parameter_samples

list of tensors representing posterior samples of model parameters, with shapes `list(tf$concat(list(list(num_posterior_draws), param<1>$prior$batch_shape, param<1>$prior$event_shape))` for all model parameters. This may optionally also be a named list mapping parameter names to tensor values.

num_steps_forecast

scalar integer tensor number of steps to forecast

Value

forecast_dist a `tfd_mixture_same_family` instance with event shape `list(num_steps_forecast, 1)` and batch shape `tf$concat(list(sample_shape, model$batch_shape))`, with `num_posterior_draws` mixture components.

See Also

Other sts-functions: [sts_build_factored_surrogate_posterior\(\)](#), [sts_build_factored_variational_loss\(\)](#), [sts_decompose_by_component\(\)](#), [sts_decompose_forecast_by_component\(\)](#), [sts_fit_with_hmc\(\)](#), [sts_one_step_predictive\(\)](#), [sts_sample_uniform_initial_state\(\)](#)

Examples

```
observed_time_series <-
  rep(c(3.5, 4.1, 4.5, 3.9, 2.4, 2.1, 1.2), 5) +
  rep(c(1.1, 1.5, 2.4, 3.1, 4.0), each = 7) %>%
  tensorflow::tf$convert_to_tensor(dtype = tensorflow::tf$float64)
day_of_week <- observed_time_series %>% sts_seasonal(num_seasons = 7)
local_linear_trend <- observed_time_series %>% sts_local_linear_trend()
model <- observed_time_series %>%
  sts_sum(components = list(day_of_week, local_linear_trend))
states_and_results <- observed_time_series %>%
  sts_fit_with_hmc(
    model,
    num_results = 10,
    num_warmup_steps = 5,
    num_variational_steps = 15)
samples <- states_and_results[[1]]
preds <- observed_time_series %>%
  sts_forecast(model,
    parameter_samples = samples,
    num_steps_forecast = 50)
predictions <- preds %>% tfd_sample(10)
```

sts_linear_regression *Formal representation of a linear regression from provided covariates.*

Description

This model defines a time series given by a linear combination of covariate time series provided in a design matrix:

```
observed_time_series <- tf$matmul(design_matrix, weights)
```

Usage

```
sts_linear_regression(design_matrix, weights_prior = NULL, name = NULL)
```

Arguments

design_matrix	float tensor of shape <code>tf\$concat(list(batch_shape, list(num_timesteps, num_features)))</code> . This may also optionally be an instance of <code>tf\$linalg\$LinearOperator</code> .
weights_prior	Distribution representing a prior over the regression weights. Must have event shape <code>list(num_features)</code> and batch shape broadcastable to the design matrix's <code>batch_shape</code> . Alternately, <code>event_shape</code> may be scalar (<code>list()</code>), in which case the prior is internally broadcast as <code>tfd_transformed_distribution(weights_prior, tfb_i = list(num_features), batch_shape = design_matrix\$batch_shape)</code> . If <code>NULL</code> , defaults to <code>tfd_student_t(df = 5, loc = 0, scale = 10)</code> , a weakly-informative prior loosely inspired by the Stan prior choice recommendations . Default value: <code>NULL</code> .
name	the name of this model component. Default value: <code>'LinearRegression'</code> .

Details

The design matrix has shape `list(num_timesteps, num_features)`. The weights are treated as an unknown random variable of size `list(num_features)` (both components also support batch shape), and are integrated over using the same approximate inference tools as other model parameters, i.e., generally HMC or variational inference.

This component does not itself include observation noise; it defines a deterministic distribution with mass at the point `tf$matmul(design_matrix, weights)`. In practice, it should be combined with observation noise from another component such as `sts_sum`, as demonstrated below.

Value

an instance of `StructuralTimeSeries`.

See Also

For usage examples see `sts_fit_with_hmc()`, `sts_forecast()`, `sts_decompose_by_component()`.

Other sts: `sts_additive_state_space_model()`, `sts_autoregressive_state_space_model()`, `sts_autoregressive()`, `sts_constrained_seasonal_state_space_model()`, `sts_dynamic_linear_regression_state_space_model()`, `sts_dynamic_linear_regression()`, `sts_local_level_state_space_model()`, `sts_local_level()`, `sts_local_linear_trend_state_space_model()`, `sts_local_linear_trend()`, `sts_seasonal_state_space_model()`, `sts_seasonal()`, `sts_semi_local_linear_trend_state_space_model()`, `sts_semi_local_linear_trend()`, `sts_smooth_seasonal_state_space_model()`, `sts_smooth_seasonal()`, `sts_sparse_linear_regression()`, `sts_sum()`

<code>sts_local_level</code>	<i>Formal representation of a local level model</i>
------------------------------	---

Description

The local level model posits a level evolving via a Gaussian random walk:

$$\text{level}[t] = \text{level}[t-1] + \text{Normal}(0., \text{level_scale})$$
Usage

```
sts_local_level(
  observed_time_series = NULL,
  level_scale_prior = NULL,
  initial_level_prior = NULL,
  name = NULL
)
```

Arguments

`observed_time_series` optional float tensor of shape `batch_shape + [T, 1]` (omitting the trailing unit dimension is also supported when `T > 1`), specifying an observed time series. Any priors not explicitly set will be given default values according to the scale of the observed time series (or batch of time series). May optionally be an instance of `sts_masked_time_series`, which includes a mask tensor to specify timesteps with missing observations. Default value: `NULL`.

`level_scale_prior` optional `tfp$distribution` instance specifying a prior on the `level_scale` parameter. If `NULL`, a heuristic default prior is constructed based on the provided `observed_time_series`. Default value: `NULL`.

`initial_level_prior` optional `tfp$distribution` instance specifying a prior on the initial level. If `NULL`, a heuristic default prior is constructed based on the provided `observed_time_series`. Default value: `NULL`.

`name` the name of this model component. Default value: `'LocalLevel'`.

Details

The latent state is [level]. We observe a noisy realization of the current level: $f[t] = \text{level}[t] + \text{Normal}(0., \text{observation_noise_scale})$ at each timestep.

Value

an instance of `StructuralTimeSeries`.

See Also

For usage examples see `sts_fit_with_hmc()`, `sts_forecast()`, `sts_decompose_by_component()`.

Other sts: `sts_additive_state_space_model()`, `sts_autoregressive_state_space_model()`, `sts_autoregressive()`, `sts_constrained_seasonal_state_space_model()`, `sts_dynamic_linear_regression_sts()`, `sts_dynamic_linear_regression()`, `sts_linear_regression()`, `sts_local_level_state_space_model()`, `sts_local_linear_trend_state_space_model()`, `sts_local_linear_trend()`, `sts_seasonal_state_space_model()`, `sts_seasonal()`, `sts_semi_local_linear_trend_state_space_model()`, `sts_semi_local_linear_trend()`, `sts_smooth_seasonal_state_space_model()`, `sts_smooth_seasonal()`, `sts_sparse_linear_regression()`, `sts_sum()`

`sts_local_level_state_space_model`

State space model for a local level

Description

A state space model (SSM) posits a set of latent (unobserved) variables that evolve over time with dynamics specified by a probabilistic transition model $p(z[t+1] | z[t])$. At each timestep, we observe a value sampled from an observation model conditioned on the current state, $p(x[t] | z[t])$. The special case where both the transition and observation models are Gaussians with mean specified as a linear function of the inputs, is known as a linear Gaussian state space model and supports tractable exact probabilistic calculations; see `tfd_linear_gaussian_state_space_model` for details. The local level model is a special case of a linear Gaussian SSM, in which the latent state posits a level evolving via a Gaussian random walk:

$$\text{level}[t] = \text{level}[t-1] + \text{Normal}(0., \text{level_scale})$$
Usage

```
sts_local_level_state_space_model(
    num_timesteps,
    level_scale,
    initial_state_prior,
    observation_noise_scale = 0,
    initial_step = 0,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = NULL
)
```

Arguments

<code>num_timesteps</code>	Scalar integer tensor number of timesteps to model with this distribution.
<code>level_scale</code>	Scalar (any additional dimensions are treated as batch dimensions) float tensor indicating the standard deviation of the level transitions.
<code>initial_state_prior</code>	instance of <code>tfd_multivariate_normal</code> representing the prior distribution on latent states. Must have event shape <code>[1]</code> (as <code>tfd_linear_gaussian_state_space_model</code> requires a rank-1 event shape).
<code>observation_noise_scale</code>	Scalar (any additional dimensions are treated as batch dimensions) float tensor indicating the standard deviation of the observation noise.
<code>initial_step</code>	Optional scalar integer tensor specifying the starting timestep. Default value: 0.
<code>validate_args</code>	logical. Whether to validate input with asserts. If <code>validate_args</code> is <code>FALSE</code> , and the inputs are invalid, correct behavior is not guaranteed. Default value: <code>FALSE</code> .
<code>allow_nan_stats</code>	logical. If <code>FALSE</code> , raise an exception if a statistic (e.g. mean/mode/etc...) is undefined for any batch member. If <code>TRUE</code> , batch members with valid parameters leading to undefined statistics will return <code>NaN</code> for this statistic. Default value: <code>TRUE</code> .
<code>name</code>	string name prefixed to ops created by this class. Default value: "LocalLevel-StateSpaceModel".

Details

The latent state is `[level]` and `[level]` is observed (with noise) at each timestep.

The parameters `level_scale` and `observation_noise_scale` are each (a batch of) scalars. The batch shape of this Distribution is the broadcast batch shape of these parameters and of the `initial_state_prior`.

Mathematical Details

The local level model implements a `tfp$distributions$LinearGaussianStateSpaceModel` with `latent_size = 1` and `observation_size = 1`, following the transition model:

```
transition_matrix = [[1]]
transition_noise ~ N(loc = 0, scale = diag([level_scale]))
```

which implements the evolution of `level` described above, and the observation model:

```
observation_matrix = [[1]]
observation_noise ~ N(loc = 0, scale = observation_noise_scale)
```

Value

an instance of `LinearGaussianStateSpaceModel`.

See Also

Other sts: `sts_additive_state_space_model()`, `sts_autoregressive_state_space_model()`, `sts_autoregressive()`, `sts_constrained_seasonal_state_space_model()`, `sts_dynamic_linear_regression_state_space_model()`, `sts_dynamic_linear_regression()`, `sts_linear_regression()`, `sts_local_level()`, `sts_local_linear_trend_state_space_model()`, `sts_local_linear_trend()`, `sts_seasonal_state_space_model()`, `sts_seasonal()`, `sts_semi_local_linear_trend_state_space_model()`, `sts_semi_local_linear_trend()`, `sts_smooth_seasonal_state_space_model()`, `sts_smooth_seasonal()`, `sts_sparse_linear_regression()`, `sts_sum()`

sts_local_linear_trend

Formal representation of a local linear trend model

Description

The local linear trend model posits a level and slope, each evolving via a Gaussian random walk:

$$\begin{aligned} \text{level}[t] &= \text{level}[t-1] + \text{slope}[t-1] + \text{Normal}(0., \text{level_scale}) \\ \text{slope}[t] &= \text{slope}[t-1] + \text{Normal}(0., \text{slope_scale}) \end{aligned}$$
Usage

```
sts_local_linear_trend(
  observed_time_series = NULL,
  level_scale_prior = NULL,
  slope_scale_prior = NULL,
  initial_level_prior = NULL,
  initial_slope_prior = NULL,
  name = NULL
)
```

Arguments

`observed_time_series`

optional float tensor of shape `batch_shape + [T, 1]` (omitting the trailing unit dimension is also supported when $T > 1$), specifying an observed time series. Any priors not explicitly set will be given default values according to the scale of the observed time series (or batch of time series). May optionally be an instance of `sts_masked_time_series`, which includes a mask tensor to specify timesteps with missing observations. Default value: `NULL`.

`level_scale_prior`

optional `tfp$Distribution` instance specifying a prior on the `level_scale` parameter. If `NULL`, a heuristic default prior is constructed based on the provided `observed_time_series`. Default value: `NULL`.

`slope_scale_prior`

optional `tfd$Distribution` instance specifying a prior on the `slope_scale` parameter. If `NULL`, a heuristic default prior is constructed based on the provided `observed_time_series`. Default value: `NULL`.

<code>initial_level_prior</code>	optional <code>tfp\$Distribution</code> instance specifying a prior on the initial level. If NULL, a heuristic default prior is constructed based on the provided <code>observed_time_series</code> . Default value: NULL.
<code>initial_slope_prior</code>	optional <code>tfd\$Distribution</code> instance specifying a prior on the initial slope. If NULL, a heuristic default prior is constructed based on the provided <code>observed_time_series</code> . Default value: NULL.
<code>name</code>	the name of this model component. Default value: 'LocalLinearTrend'.

Details

The latent state is the two-dimensional tuple [level, slope]. At each timestep we observe a noisy realization of the current level: $f[t] = \text{level}[t] + \text{Normal}(0., \text{observation_noise_scale})$. This model is appropriate for data where the trend direction and magnitude (latent slope) is consistent within short periods but may evolve over time.

Note that this model can produce very high uncertainty forecasts, as uncertainty over the slope compounds quickly. If you expect your data to have nonzero long-term trend, i.e. that slopes tend to revert to some mean, then the `SemiLocalLinearTrend` model may produce sharper forecasts.

Value

an instance of `StructuralTimeSeries`.

See Also

For usage examples see `sts_fit_with_hmc()`, `sts_forecast()`, `sts_decompose_by_component()`.

Other sts: `sts_additive_state_space_model()`, `sts_autoregressive_state_space_model()`, `sts_autoregressive()`, `sts_constrained_seasonal_state_space_model()`, `sts_dynamic_linear_regression_state_space_model()`, `sts_dynamic_linear_regression()`, `sts_linear_regression()`, `sts_local_level_state_space_model()`, `sts_local_level()`, `sts_local_linear_trend_state_space_model()`, `sts_seasonal_state_space_model()`, `sts_seasonal()`, `sts_semi_local_linear_trend_state_space_model()`, `sts_semi_local_linear_trend()`, `sts_smooth_seasonal_state_space_model()`, `sts_smooth_seasonal()`, `sts_sparse_linear_regression()`, `sts_sum()`

`sts_local_linear_trend_state_space_model`

State space model for a local linear trend

Description

A state space model (SSM) posits a set of latent (unobserved) variables that evolve over time with dynamics specified by a probabilistic transition model $p(z[t+1] | z[t])$. At each timestep, we observe a value sampled from an observation model conditioned on the current state, $p(x[t] | z[t])$. The special case where both the transition and observation models are Gaussians with mean specified as a linear function of the inputs, is known as a linear Gaussian state space model and supports tractable exact probabilistic calculations; see `tfd_linear_gaussian_state_space_model` for details.

Usage

```
sts_local_linear_trend_state_space_model(
    num_timesteps,
    level_scale,
    slope_scale,
    initial_state_prior,
    observation_noise_scale = 0,
    initial_step = 0,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = NULL
)
```

Arguments

<code>num_timesteps</code>	Scalar integer tensor number of timesteps to model with this distribution.
<code>level_scale</code>	Scalar (any additional dimensions are treated as batch dimensions) float tensor indicating the standard deviation of the level transitions.
<code>slope_scale</code>	Scalar (any additional dimensions are treated as batch dimensions) float tensor indicating the standard deviation of the slope transitions.
<code>initial_state_prior</code>	instance of <code>tfd_multivariate_normal</code> representing the prior distribution on latent states. Must have event shape [1] (as <code>tfd_linear_gaussian_state_space_model</code> requires a rank-1 event shape).
<code>observation_noise_scale</code>	Scalar (any additional dimensions are treated as batch dimensions) float tensor indicating the standard deviation of the observation noise.
<code>initial_step</code>	Optional scalar integer tensor specifying the starting timestep. Default value: 0.
<code>validate_args</code>	logical. Whether to validate input with asserts. If <code>validate_args</code> is <code>FALSE</code> , and the inputs are invalid, correct behavior is not guaranteed. Default value: <code>FALSE</code> .
<code>allow_nan_stats</code>	logical. If <code>FALSE</code> , raise an exception if a statistic (e.g. mean/mode/etc...) is undefined for any batch member. If <code>TRUE</code> , batch members with valid parameters leading to undefined statistics will return NaN for this statistic. Default value: <code>TRUE</code> .
<code>name</code>	string prefixed to ops created by this class. Default value: "LocalLinearTrend-StateSpaceModel".

Details

The local linear trend model is a special case of a linear Gaussian SSM, in which the latent state posits a level and slope, each evolving via a Gaussian random walk:

```
level[t] = level[t-1] + slope[t-1] + Normal(0., level_scale)
slope[t] = slope[t-1] + Normal(0., slope_scale)
```

The latent state is the two-dimensional tuple [level, slope]. The level is observed at each timestep. The parameters `level_scale`, `slope_scale`, and `observation_noise_scale` are each (a batch of) scalars. The batch shape of this Distribution is the broadcast batch shape of these parameters and of the `initial_state_prior`.

Mathematical Details

The linear trend model implements a `tfd_linear_gaussian_state_space_model` with `latent_size = 2` and `observation_size = 1`, following the transition model:

```
transition_matrix = [[1., 1.]
                    [0., 1.]]
transition_noise ~ N(loc = 0, scale = diag([level_scale, slope_scale]))
```

which implements the evolution of [level, slope] described above, and the observation model:

```
observation_matrix = [[1., 0.]]
observation_noise ~ N(loc= 0 , scale = observation_noise_scale)
```

which picks out the first latent component, i.e., the level, as the observation at each timestep.

Value

an instance of `LinearGaussianStateSpaceModel`.

See Also

Other sts: `sts_additive_state_space_model()`, `sts_autoregressive_state_space_model()`, `sts_autoregressive()`, `sts_constrained_seasonal_state_space_model()`, `sts_dynamic_linear_regression_state_space_model()`, `sts_dynamic_linear_regression()`, `sts_linear_regression()`, `sts_local_level_state_space_model()`, `sts_local_level()`, `sts_local_linear_trend()`, `sts_seasonal_state_space_model()`, `sts_seasonal()`, `sts_semi_local_linear_trend_state_space_model()`, `sts_semi_local_linear_trend()`, `sts_smooth_seasonal_state_space_model()`, `sts_smooth_seasonal()`, `sts_sparse_linear_regression()`, `sts_sum()`

sts_one_step_predictive

Compute one-step-ahead predictive distributions for all timesteps

Description

Given samples from the posterior over parameters, return the predictive distribution over observations at each time T , given observations up through time $T-1$.

Usage

```
sts_one_step_predictive(
    observed_time_series,
    model,
    parameter_samples,
    timesteps_are_event_shape = TRUE
)
```

Arguments

observed_time_series	float tensor of shape <code>concat([sample_shape, model.batch_shape, [num_timesteps, 1]])</code> where <code>sample_shape</code> corresponds to i.i.d. observations, and the trailing <code>[1]</code> dimension may (optionally) be omitted if <code>num_timesteps > 1</code> . May optionally be an instance of <code>sts_masked_time_series</code> , which includes a mask tensor to specify timesteps with missing observations.
model	An instance of <code>StructuralTimeSeries</code> representing a time-series model. This represents a joint distribution over time-series and their parameters with batch shape <code>[b1, ..., bN]</code> .
parameter_samples	list of tensors representing posterior samples of model parameters, with shapes <code>list(tf.concat(list(list(num_posterior_draws), param<1>\$prior\$batch_shape, param<1>\$prior\$event_shape), ...))</code> for all model parameters. This may optionally also be a named list mapping parameter names to tensor values.
timesteps_are_event_shape	Deprecated, for backwards compatibility only. If <code>False</code> , the predictive distribution will return per-timestep probabilities. Default value: <code>TRUE</code> .

Value

`forecast_dist` a `tfd_mixture_same_family` instance with event shape `list(num_timesteps)` and batch shape `tf.concat(list(sample_shape, model$batch_shape))`, with `num_posterior_draws` mixture components. The `t`th step represents the forecast distribution $p(\text{observed_time_series}[t] \mid \text{observed_time_series}[0:t-1], \text{parameter_samples})$.

See Also

Other `sts`-functions: [sts_build_factored_surrogate_posterior\(\)](#), [sts_build_factored_variational_loss\(\)](#), [sts_decompose_by_component\(\)](#), [sts_decompose_forecast_by_component\(\)](#), [sts_fit_with_hmc\(\)](#), [sts_forecast\(\)](#), [sts_sample_uniform_initial_state\(\)](#)

`sts_sample_uniform_initial_state`

Initialize from a uniform $[-2, 2]$ distribution in unconstrained space.

Description

Initialize from a uniform $[-2, 2]$ distribution in unconstrained space.

Usage

```
sts_sample_uniform_initial_state(
  parameter,
  return_constrained = TRUE,
  init_sample_shape = list(),
  seed = NULL
)
```

Arguments

parameter sts\$Parameter named tuple instance.

return_constrained if TRUE, re-applies the constraining bijector to return initializations in the original domain. Otherwise, returns initializations in the unconstrained space. Default value: TRUE.

init_sample_shape sample_shape of the sampled initializations. Default value: list().

seed integer to seed the random number generator.

Value

uniform_initializer Tensor of shape `concat([init_sample_shape, parameter.prior.batch_shape, transformed_event_shape])`, where `transformed_event_shape` is `parameter.prior.event_shape`, if `return_constrained=TRUE`, and otherwise it is `parameter$bijector$inverse_event_shape(parameter$prior$event_shape)`.

See Also

Other sts-functions: [sts_build_factored_surrogate_posterior\(\)](#), [sts_build_factored_variational_loss\(\)](#), [sts_decompose_by_component\(\)](#), [sts_decompose_forecast_by_component\(\)](#), [sts_fit_with_hmc\(\)](#), [sts_forecast\(\)](#), [sts_one_step_predictive\(\)](#)

<code>sts_seasonal</code>	<i>Formal representation of a seasonal effect model.</i>
---------------------------	--

Description

A seasonal effect model posits a fixed set of recurring, discrete 'seasons', each of which is active for a fixed number of timesteps and, while active, contributes a different effect to the time series. These are generally not meteorological seasons, but represent regular recurring patterns such as hour-of-day or day-of-week effects. Each season lasts for a fixed number of timesteps. The effect of each season drifts from one occurrence to the next following a Gaussian random walk:

Usage

```
sts_seasonal(
  observed_time_series = NULL,
  num_seasons,
  num_steps_per_season = 1,
  drift_scale_prior = NULL,
  initial_effect_prior = NULL,
  constrain_mean_effect_to_zero = TRUE,
  name = NULL
)
```


Arguments

- `observed_time_series`
optional float tensor of shape `batch_shape + [T, 1]` (omitting the trailing unit dimension is also supported when $T > 1$), specifying an observed time series. Any priors not explicitly set will be given default values according to the scale of the observed time series (or batch of time series). May optionally be an instance of `sts_masked_time_series`, which includes a mask tensor to specify timesteps with missing observations. Default value: `NULL`.
- `num_seasons` Scalar integer number of seasons.
- `num_steps_per_season`
integer number of steps in each season. This may be either a scalar (shape `[]`), in which case all seasons have the same length, or an array of shape `[num_seasons]`, in which seasons have different length, but remain constant around different cycles, or an array of shape `[num_cycles, num_seasons]`, in which `num_steps_per_season` for each season also varies in different cycle (e.g., a 4 years cycle with leap day). Default value: 1.
- `drift_scale_prior`
optional `tfd$Distribution` instance specifying a prior on the `drift_scale` parameter. If `NULL`, a heuristic default prior is constructed based on the provided `observed_time_series`. Default value: `NULL`.
- `initial_effect_prior`
optional `tfd$Distribution` instance specifying a normal prior on the initial effect of each season. This may be either a scalar `tfd_normal` prior, in which case it applies independently to every season, or it may be multivariate normal (e.g., `tfd_multivariate_normal_diag`) with event shape `[num_seasons]`, in which case it specifies a joint prior across all seasons. If `NULL`, a heuristic default prior is constructed based on the provided `observed_time_series`. Default value: `NULL`.
- `constrain_mean_effect_to_zero`
if `TRUE`, use a model parameterization that constrains the mean effect across all seasons to be zero. This constraint is generally helpful in identifying the contributions of different model components and can lead to more interpretable posterior decompositions. It may be undesirable if you plan to directly examine the latent space of the underlying state space model. Default value: `TRUE`.
- `name` the name of this model component. Default value: `'Seasonal'`.

Details

```
effects[season, occurrence[i]] = (
  effects[season, occurrence[i-1]] + Normal(loc=0., scale=drift_scale))
```

The `drift_scale` parameter governs the standard deviation of the random walk; for example, in a day-of-week model it governs the change in effect from this Monday to next Monday.

Value

an instance of `StructuralTimeSeries`.

See Also

For usage examples see [sts_fit_with_hmc\(\)](#), [sts_forecast\(\)](#), [sts_decompose_by_component\(\)](#).

Other sts: [sts_additive_state_space_model\(\)](#), [sts_autoregressive_state_space_model\(\)](#), [sts_autoregressive\(\)](#), [sts_constrained_seasonal_state_space_model\(\)](#), [sts_dynamic_linear_regression_state_space_model\(\)](#), [sts_dynamic_linear_regression\(\)](#), [sts_linear_regression\(\)](#), [sts_local_level_state_space_model\(\)](#), [sts_local_level\(\)](#), [sts_local_linear_trend_state_space_model\(\)](#), [sts_local_linear_trend\(\)](#), [sts_seasonal_state_space_model\(\)](#), [sts_semi_local_linear_trend_state_space_model\(\)](#), [sts_semi_local_linear_trend\(\)](#), [sts_smooth_seasonal_state_space_model\(\)](#), [sts_smooth_seasonal\(\)](#), [sts_sparse_linear_regression\(\)](#), [sts_sum\(\)](#)

sts_seasonal_state_space_model

State space model for a seasonal effect.

Description

A state space model (SSM) posits a set of latent (unobserved) variables that evolve over time with dynamics specified by a probabilistic transition model $p(z[t+1] | z[t])$. At each timestep, we observe a value sampled from an observation model conditioned on the current state, $p(x[t] | z[t])$. The special case where both the transition and observation models are Gaussians with mean specified as a linear function of the inputs, is known as a linear Gaussian state space model and supports tractable exact probabilistic calculations; see [tfd_linear_gaussian_state_space_model](#) for details.

Usage

```
sts_seasonal_state_space_model(
    num_timesteps,
    num_seasons,
    drift_scale,
    initial_state_prior,
    observation_noise_scale = 0,
    num_steps_per_season = 1,
    initial_step = 0,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = NULL
)
```

Arguments

num_timesteps	Scalar integer tensor number of timesteps to model with this distribution.
num_seasons	Scalar integer number of seasons.
drift_scale	Scalar (any additional dimensions are treated as batch dimensions) float tensor indicating the standard deviation of the change in effect between consecutive occurrences of a given season. This is assumed to be the same for all seasons.

<code>initial_state_prior</code>	instance of <code>tfd_multivariate_normal</code> representing the prior distribution on latent states; must have event shape <code>[num_seasons]</code> .
<code>observation_noise_scale</code>	Scalar (any additional dimensions are treated as batch dimensions) float tensor indicating the standard deviation of the observation noise.
<code>num_steps_per_season</code>	integer number of steps in each season. This may be either a scalar (shape <code>[]</code>), in which case all seasons have the same length, or an array of shape <code>[num_seasons]</code> , in which seasons have different length, but remain constant around different cycles, or an array of shape <code>[num_cycles, num_seasons]</code> , in which <code>num_steps_per_season</code> for each season also varies in different cycle (e.g., a 4 years cycle with leap day). Default value: 1.
<code>initial_step</code>	Optional scalar integer tensor specifying the starting timestep. Default value: 0.
<code>validate_args</code>	logical. Whether to validate input with asserts. If <code>validate_args</code> is <code>FALSE</code> , and the inputs are invalid, correct behavior is not guaranteed. Default value: <code>FALSE</code> .
<code>allow_nan_stats</code>	logical. If <code>FALSE</code> , raise an exception if a statistic (e.g. mean/mode/etc...) is undefined for any batch member. If <code>TRUE</code> , batch members with valid parameters leading to undefined statistics will return <code>NaN</code> for this statistic. Default value: <code>TRUE</code> .
<code>name</code>	string prefixed to ops created by this class. Default value: "SeasonalStateSpace-Model".

Details

A seasonal effect model is a special case of a linear Gaussian SSM. The latent states represent an unknown effect from each of several 'seasons'; these are generally not meteorological seasons, but represent regular recurring patterns such as hour-of-day or day-of-week effects. The effect of each season drifts from one occurrence to the next, following a Gaussian random walk:

$$\text{effects}[\text{season}, \text{occurrence}[i]] = (\text{effects}[\text{season}, \text{occurrence}[i-1]] + \text{Normal}(\text{loc}=0., \text{scale}=\text{drift_scale}))$$

The latent state has dimension `num_seasons`, containing one effect for each seasonal component. The parameters `drift_scale` and `observation_noise_scale` are each (a batch of) scalars. The batch shape of this Distribution is the broadcast batch shape of these parameters and of the `initial_state_prior`. Note: there is no requirement that the effects sum to zero.

Mathematical Details

The seasonal effect model implements a `tfd_linear_gaussian_state_space_model` with `latent_size = num_seasons` and `observation_size = 1`. The latent state is organized so that the *current* seasonal effect is always in the first (zeroth) dimension. The transition model rotates the latent state to shift to a new effect at the end of each season:

$$\text{transition_matrix}[t] = (\text{permutation_matrix}([1, 2, \dots, \text{num_seasons}-1, 0]) \text{ if } \text{season_is_changing}(t))$$

```

        else eye(num_seasons)
transition_noise[t] ~ Normal(loc=0., scale_diag=(
    [drift_scale, 0, ..., 0]
    if season_is_changing(t)
    else [0, 0, ..., 0]))

```

where `season_is_changing(t)` is True if `t % sum(num_steps_per_season)` is in the set of final days for each season, given by `cumsum(num_steps_per_season) - 1`. The observation model always picks out the effect for the current season, i.e., the first element of the latent state:

```

observation_matrix = [[1., 0., ..., 0.]]
observation_noise ~ Normal(loc=0, scale=observation_noise_scale)

```

Value

an instance of `LinearGaussianStateSpaceModel`.

See Also

Other sts: `sts_additive_state_space_model()`, `sts_autoregressive_state_space_model()`, `sts_autoregressive()`, `sts_constrained_seasonal_state_space_model()`, `sts_dynamic_linear_regression_state_space_model()`, `sts_dynamic_linear_regression()`, `sts_linear_regression()`, `sts_local_level_state_space_model()`, `sts_local_level()`, `sts_local_linear_trend_state_space_model()`, `sts_local_linear_trend()`, `sts_seasonal()`, `sts_semi_local_linear_trend_state_space_model()`, `sts_semi_local_linear_trend()`, `sts_smooth_seasonal_state_space_model()`, `sts_smooth_seasonal()`, `sts_sparse_linear_regression()`, `sts_sum()`

sts_semi_local_linear_trend

Formal representation of a semi-local linear trend model.

Description

Like the `sts_local_linear_trend` model, a semi-local linear trend posits a latent level and slope, with the level component updated according to the current slope plus a random walk:

Usage

```

sts_semi_local_linear_trend(
    observed_time_series = NULL,
    level_scale_prior = NULL,
    slope_mean_prior = NULL,
    slope_scale_prior = NULL,
    autoregressive_coef_prior = NULL,
    initial_level_prior = NULL,
    initial_slope_prior = NULL,
    constrain_ar_coef_stationary = TRUE,
    constrain_ar_coef_positive = FALSE,
    name = NULL
)

```

Arguments

<code>observed_time_series</code>	optional float tensor of shape <code>batch_shape + [T, 1]</code> (omitting the trailing unit dimension is also supported when $T > 1$), specifying an observed time series. Any priors not explicitly set will be given default values according to the scale of the observed time series (or batch of time series). May optionally be an instance of <code>sts_masked_time_series</code> , which includes a mask tensor to specify timesteps with missing observations. Default value: <code>NULL</code> .
<code>level_scale_prior</code>	optional <code>tfp\$Distribution</code> instance specifying a prior on the <code>level_scale</code> parameter. If <code>NULL</code> , a heuristic default prior is constructed based on the provided <code>observed_time_series</code> . Default value: <code>NULL</code> .
<code>slope_mean_prior</code>	optional <code>tfd\$Distribution</code> instance specifying a prior on the <code>slope_mean</code> parameter. If <code>NULL</code> , a heuristic default prior is constructed based on the provided <code>observed_time_series</code> . Default value: <code>NULL</code> .
<code>slope_scale_prior</code>	optional <code>tfd\$Distribution</code> instance specifying a prior on the <code>slope_scale</code> parameter. If <code>NULL</code> , a heuristic default prior is constructed based on the provided <code>observed_time_series</code> . Default value: <code>NULL</code> .
<code>autoregressive_coef_prior</code>	optional <code>tfd\$Distribution</code> instance specifying a prior on the <code>autoregressive_coef</code> parameter. If <code>NULL</code> , the default prior is a standard <code>Normal(0, 1)</code> . Note that the prior may be implicitly truncated by <code>constrain_ar_coef_stationary</code> and/or <code>constrain_ar_coef_positive</code> . Default value: <code>NULL</code> .
<code>initial_level_prior</code>	optional <code>tfp\$Distribution</code> instance specifying a prior on the initial level. If <code>NULL</code> , a heuristic default prior is constructed based on the provided <code>observed_time_series</code> . Default value: <code>NULL</code> .
<code>initial_slope_prior</code>	optional <code>tfd\$Distribution</code> instance specifying a prior on the initial slope. If <code>NULL</code> , a heuristic default prior is constructed based on the provided <code>observed_time_series</code> . Default value: <code>NULL</code> .
<code>constrain_ar_coef_stationary</code>	if <code>TRUE</code> , perform inference using a parameterization that restricts <code>autoregressive_coef</code> to the interval <code>(-1, 1)</code> , or <code>(0, 1)</code> if <code>force_positive_ar_coef</code> is also <code>TRUE</code> , corresponding to stationary processes. This will implicitly truncate the support of <code>autoregressive_coef_prior</code> . Default value: <code>TRUE</code> .
<code>constrain_ar_coef_positive</code>	if <code>TRUE</code> , perform inference using a parameterization that restricts <code>autoregressive_coef</code> to be positive, or in <code>(0, 1)</code> if <code>constrain_ar_coef_stationary</code> is also <code>TRUE</code> . This will implicitly truncate the support of <code>autoregressive_coef_prior</code> . Default value: <code>FALSE</code> .
<code>name</code>	the name of this model component. Default value: <code>'SemiLocalLinearTrend'</code> .

Details

`level[t] = level[t-1] + slope[t-1] + Normal(0., level_scale)`

The slope component in a `sts_semi_local_linear_trend` model evolves according to a first-order autoregressive (AR1) process with potentially nonzero mean:

$$\text{slope}[t] = (\text{slope_mean} + \text{autoregressive_coef} * (\text{slope}[t-1] - \text{slope_mean}) + \text{Normal}(0., \text{slope_scale}))$$

Unlike the random walk used in `LocalLinearTrend`, a stationary AR1 process (coefficient in (-1, 1)) maintains bounded variance over time, so a `SemiLocalLinearTrend` model will often produce more reasonable uncertainties when forecasting over long timescales.

Value

an instance of `StructuralTimeSeries`.

See Also

For usage examples see `sts_fit_with_hmc()`, `sts_forecast()`, `sts_decompose_by_component()`.

Other sts: `sts_additive_state_space_model()`, `sts_autoregressive_state_space_model()`, `sts_autoregressive()`, `sts_constrained_seasonal_state_space_model()`, `sts_dynamic_linear_regression_state_space_model()`, `sts_dynamic_linear_regression()`, `sts_linear_regression()`, `sts_local_level_state_space_model()`, `sts_local_level()`, `sts_local_linear_trend_state_space_model()`, `sts_local_linear_trend()`, `sts_seasonal_state_space_model()`, `sts_seasonal()`, `sts_semi_local_linear_trend_state_space_model()`, `sts_smooth_seasonal_state_space_model()`, `sts_smooth_seasonal()`, `sts_sparse_linear_regression()`, `sts_sum()`

`sts_semi_local_linear_trend_state_space_model`

State space model for a semi-local linear trend.

Description

A state space model (SSM) posits a set of latent (unobserved) variables that evolve over time with dynamics specified by a probabilistic transition model $p(z[t+1] | z[t])$. At each timestep, we observe a value sampled from an observation model conditioned on the current state, $p(x[t] | z[t])$. The special case where both the transition and observation models are Gaussians with mean specified as a linear function of the inputs, is known as a linear Gaussian state space model and supports tractable exact probabilistic calculations; see `tfd_linear_gaussian_state_space_model` for details.

Usage

```
sts_semi_local_linear_trend_state_space_model(
    num_timesteps,
    level_scale,
    slope_mean,
    slope_scale,
    autoregressive_coef,
    initial_state_prior,
```

```

    observation_noise_scale = 0,
    initial_step = 0,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = NULL
)

```

Arguments

num_timesteps	Scalar integer tensor number of timesteps to model with this distribution.
level_scale	Scalar (any additional dimensions are treated as batch dimensions) float tensor indicating the standard deviation of the level transitions.
slope_mean	Scalar (any additional dimensions are treated as batch dimensions) float tensor indicating the expected long-term mean of the latent slope.
slope_scale	Scalar (any additional dimensions are treated as batch dimensions) float tensor indicating the standard deviation of the slope transitions.
autoregressive_coef	Scalar (any additional dimensions are treated as batch dimensions) float tensor defining the AR1 process on the latent slope.
initial_state_prior	instance of <code>tfd_multivariate_normal</code> representing the prior distribution on latent states. Must have event shape [1] (as <code>tfd_linear_gaussian_state_space_model</code> requires a rank-1 event shape).
observation_noise_scale	Scalar (any additional dimensions are treated as batch dimensions) float tensor indicating the standard deviation of the observation noise.
initial_step	Optional scalar integer tensor specifying the starting timestep. Default value: 0.
validate_args	logical. Whether to validate input with asserts. If <code>validate_args</code> is <code>FALSE</code> , and the inputs are invalid, correct behavior is not guaranteed. Default value: <code>FALSE</code> .
allow_nan_stats	logical. If <code>FALSE</code> , raise an exception if a statistic (e.g. mean/mode/etc...) is undefined for any batch member. If <code>TRUE</code> , batch members with valid parameters leading to undefined statistics will return NaN for this statistic. Default value: <code>TRUE</code> .
name	string' prefixed to ops created by this class. Default value: "SemiLocalLinearTrend-StateSpaceModel".

Details

The semi-local linear trend model is a special case of a linear Gaussian SSM, in which the latent state posits a level and slope. The level evolves via a Gaussian random walk centered at the current slope, while the slope follows a first-order autoregressive (AR1) process with mean `slope_mean`:

```

level[t] = level[t-1] + slope[t-1] + Normal(0, level_scale)
slope[t] = (slope_mean + autoregressive_coef * (slope[t-1] - slope_mean) +
           Normal(0., slope_scale))

```

The latent state is the two-dimensional tuple [level, slope]. The level is observed at each timestep. The parameters level_scale, slope_mean, slope_scale, autoregressive_coef, and observation_noise_scale are each (a batch of) scalars. The batch shape of this Distribution is the broadcast batch shape of these parameters and of the initial_state_prior.

Mathematical Details

The semi-local linear trend model implements a `tfp.distributions.LinearGaussianStateSpaceModel` with `latent_size = 2` and `observation_size = 1`, following the transition model:

```

transition_matrix = [[1., 1.]
                    [0., autoregressive_coef]]
transition_noise ~ N(loc=slope_mean - autoregressive_coef * slope_mean,
                    scale=diag([level_scale, slope_scale]))

```

which implements the evolution of [level, slope] described above, and the observation model:

```

observation_matrix = [[1., 0.]]
observation_noise ~ N(loc=0, scale=observation_noise_scale)

```

which picks out the first latent component, i.e., the level, as the observation at each timestep.

Value

an instance of `LinearGaussianStateSpaceModel`.

See Also

Other sts: [sts_additive_state_space_model\(\)](#), [sts_autoregressive_state_space_model\(\)](#), [sts_autoregressive\(\)](#), [sts_constrained_seasonal_state_space_model\(\)](#), [sts_dynamic_linear_regression_state_space_model\(\)](#), [sts_dynamic_linear_regression\(\)](#), [sts_linear_regression\(\)](#), [sts_local_level_state_space_model\(\)](#), [sts_local_level\(\)](#), [sts_local_linear_trend_state_space_model\(\)](#), [sts_local_linear_trend\(\)](#), [sts_seasonal_state_space_model\(\)](#), [sts_seasonal\(\)](#), [sts_semi_local_linear_trend\(\)](#), [sts_smooth_seasonal_state_space_model\(\)](#), [sts_smooth_seasonal\(\)](#), [sts_sparse_linear_regression\(\)](#), [sts_sum\(\)](#)

sts_smooth_seasonal *Formal representation of a smooth seasonal effect model*

Description

The smooth seasonal model uses a set of trigonometric terms in order to capture a recurring pattern whereby adjacent (in time) effects are similar. The model uses frequencies calculated via:

Usage

```
sts_smooth_seasonal(
  period,
  frequency_multipliers,
  allow_drift = TRUE,
  drift_scale_prior = NULL,
  initial_state_prior = NULL,
  observed_time_series = NULL,
  name = NULL
)
```

Arguments

period positive scalar float Tensor giving the number of timesteps required for the longest cyclic effect to repeat.

frequency_multipliers One-dimensional float Tensor listing the frequencies (cyclic components) included in the model, as multipliers of the base/fundamental frequency $2 \cdot \pi / \text{period}$. Each component is specified by the number of times it repeats per period, and adds two latent dimensions to the model. A smooth seasonal model that can represent any periodic function is given by `frequency_multipliers = [1, 2, ..., floor(period / 2)]`. However, it is often desirable to enforce a smoothness assumption (and reduce the computational burden) by dropping some of the higher frequencies.

allow_drift optional logical specifying whether the seasonal effects can drift over time. Setting this to `FALSE` removes the `drift_scale` parameter from the model. This is mathematically equivalent to `drift_scale_prior = tfd.Deterministic(0.)`, but removing drift directly is preferred because it avoids the use of a degenerate prior. Default value: `TRUE`.

drift_scale_prior optional `tfd$Distribution` instance specifying a prior on the `drift_scale` parameter. If `NULL`, a heuristic default prior is constructed based on the provided `observed_time_series`. Default value: `NULL`.

initial_state_prior instance of `tfd$MultivariateNormal` representing the prior distribution on the latent states. Must have event shape `[2 * len(frequency_multipliers)]`. If `NULL`, a heuristic default prior is constructed based on the provided `observed_time_series`.

observed_time_series optional float Tensor of shape `batch_shape + [T, 1]` (omitting the trailing unit dimension is also supported when `T > 1`), specifying an observed time series. Any priors not explicitly set will be given default values according to the scale of the observed time series (or batch of time series). May optionally be an instance of `tfpstsMaskedTimeSeries`, which includes a mask Tensor to specify timesteps with missing observations. Default value: `NULL`.

name the name of this model component. Default value: `'LocalLinearTrend'`.

Details

$$\text{frequencies}[j] = 2 \cdot \pi \cdot \text{frequency_multipliers}[j] / \text{period}$$

and then posits two latent states for each frequency. The two latent states associated with frequency j drift over time via:

```
effect[t] = (effect[t-1] * cos(frequencies[j]) +
             auxiliary[t-] * sin(frequencies[j]) +
             Normal(0., drift_scale))
auxiliary[t] = (-effect[t-1] * sin(frequencies[j]) +
               auxiliary[t-] * cos(frequencies[j]) +
               Normal(0., drift_scale))
```

where `effect` is the smooth seasonal effect and `auxiliary` only appears as a matter of construction. The interpretation of `auxiliary` is thus not particularly important.

Value

an instance of `StructuralTimeSeries`.

See Also

For usage examples see `sts_fit_with_hmc()`, `sts_forecast()`, `sts_decompose_by_component()`.

Other sts: `sts_additive_state_space_model()`, `sts_autoregressive_state_space_model()`, `sts_autoregressive()`, `sts_constrained_seasonal_state_space_model()`, `sts_dynamic_linear_regression_state_space_model()`, `sts_dynamic_linear_regression()`, `sts_linear_regression()`, `sts_local_level_state_space_model()`, `sts_local_level()`, `sts_local_linear_trend_state_space_model()`, `sts_local_linear_trend()`, `sts_seasonal_state_space_model()`, `sts_seasonal()`, `sts_semi_local_linear_trend_state_space_model()`, `sts_semi_local_linear_trend()`, `sts_smooth_seasonal_state_space_model()`, `sts_sparse_linear_regression()`, `sts_sum()`

sts_smooth_seasonal_state_space_model

State space model for a smooth seasonal effect

Description

A state space model (SSM) posits a set of latent (unobserved) variables that evolve over time with dynamics specified by a probabilistic transition model $p(z[t+1] | z[t])$. At each timestep, we observe a value sampled from an observation model conditioned on the current state, $p(x[t] | z[t])$. The special case where both the transition and observation models are Gaussians with mean specified as a linear function of the inputs, is known as a linear Gaussian state space model and supports tractable exact probabilistic calculations; see `tfp$distributions$LinearGaussianStateSpaceModel` for details. A smooth seasonal effect model is a special case of a linear Gaussian SSM. It is the sum of a set of "cyclic" components, with one component for each frequency:

$$\text{frequencies}[j] = 2. * \pi * \text{frequency_multipliers}[j] / \text{period}$$

Each cyclic component contains two latent states which we denote `effect` and `auxiliary`. The two latent states for component j drift over time via:

```

effect[t] = (effect[t-1] * cos(frequencies[j]) +
             auxiliary[t-] * sin(frequencies[j]) +
             Normal(0., drift_scale))
auxiliary[t] = (-effect[t-1] * sin(frequencies[j]) +
               auxiliary[t-] * cos(frequencies[j]) +
               Normal(0., drift_scale))

```

Usage

```

sts_smooth_seasonal_state_space_model(
  num_timesteps,
  period,
  frequency_multipliers,
  drift_scale,
  initial_state_prior,
  observation_noise_scale = 0,
  initial_step = 0,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = NULL
)

```

Arguments

num_timesteps Scalar integer Tensor number of timesteps to model with this distribution.

period positive scalar float Tensor giving the number of timesteps required for the longest cyclic effect to repeat.

frequency_multipliers One-dimensional float Tensor listing the frequencies (cyclic components) included in the model, as multipliers of the base/fundamental frequency $2 \cdot \pi / \text{period}$. Each component is specified by the number of times it repeats per period, and adds two latent dimensions to the model. A smooth seasonal model that can represent any periodic function is given by `frequency_multipliers = [1, 2, ..., floor(period / 2)]`. However, it is often desirable to enforce a smoothness assumption (and reduce the computational burden) by dropping some of the higher frequencies.

drift_scale Scalar (any additional dimensions are treated as batch dimensions) float Tensor indicating the standard deviation of the latent state transitions.

initial_state_prior instance of `tfd$MultivariateNormal` representing the prior distribution on latent states. Must have event shape `[num_features]`.

observation_noise_scale Scalar (any additional dimensions are treated as batch dimensions) float Tensor indicating the standard deviation of the observation noise. Default value: `0.`

initial_step scalar integer Tensor specifying the starting timestep. Default value: `0`.

validate_args logical. Whether to validate input with asserts. If `validate_args` is `FALSE`, and the inputs are invalid, correct behavior is not guaranteed. Default value: `FALSE`.

allow_nan_stats	logical. If FALSE, raise an exception if a statistic (e.g. mean/mode/etc...) is undefined for any batch member. If TRUE, batch members with valid parameters leading to undefined statistics will return NaN for this statistic. Default value: TRUE.
name	string prefixed to ops created by this class. Default value: "LocalLinearTrend-StateSpaceModel".

Details

The auxiliary latent state only appears as a matter of construction and thus its interpretation is not particularly important. The total smooth seasonal effect is the sum of the effect values from each of the cyclic components. The parameters `drift_scale` and `observation_noise_scale` are each (a batch of) scalars. The batch shape of this Distribution is the broadcast batch shape of these parameters and of the `initial_state_prior`.

Mathematical Details

The smooth seasonal effect model implements a `tfp$distributions$LinearGaussianStateSpaceModel` with `latent_size = 2 * len(frequency_multipliers)` and `observation_size = 1`. The latent state is the concatenation of the cyclic latent states which themselves comprise an effect and an auxiliary state. The transition matrix is a block diagonal matrix where block `j` is:

$$\text{transition_matrix}[j] = \begin{bmatrix} \cos(\text{frequencies}[j]), & \sin(\text{frequencies}[j]) \\ -\sin(\text{frequencies}[j]), & \cos(\text{frequencies}[j]) \end{bmatrix}$$

The observation model picks out the cyclic effect values from the latent state:

```
observation_matrix = [[1., 0., 1., 0., ..., 1., 0.]]
observation_noise ~ Normal(loc=0, scale=observation_noise_scale)
```

For further mathematical details please see Harvey (1990).

Value

an instance of `LinearGaussianStateSpaceModel`.

references

- Harvey, A. Forecasting, Structural Time Series Models and the Kalman Filter. Cambridge: Cambridge University Press, 1990.

See Also

Other sts: [sts_additive_state_space_model\(\)](#), [sts_autoregressive_state_space_model\(\)](#), [sts_autoregressive\(\)](#), [sts_constrained_seasonal_state_space_model\(\)](#), [sts_dynamic_linear_regression_state_space_model\(\)](#), [sts_dynamic_linear_regression\(\)](#), [sts_linear_regression\(\)](#), [sts_local_level_state_space_model\(\)](#), [sts_local_level\(\)](#), [sts_local_linear_trend_state_space_model\(\)](#), [sts_local_linear_trend\(\)](#), [sts_seasonal_state_space_model\(\)](#), [sts_seasonal\(\)](#), [sts_semi_local_linear_trend_state_space_model\(\)](#), [sts_semi_local_linear_trend\(\)](#), [sts_smooth_seasonal\(\)](#), [sts_sparse_linear_regression\(\)](#), [sts_sum\(\)](#)

 sts_sparse_linear_regression

Formal representation of a sparse linear regression.

Description

This model defines a time series given by a sparse linear combination of covariate time series provided in a design matrix:

Usage

```
sts_sparse_linear_regression(
  design_matrix,
  weights_prior_scale = 0.1,
  weights_batch_shape = NULL,
  name = NULL
)
```

Arguments

design_matrix float tensor of shape `tf$concat(list(batch_shape, list(num_timesteps, num_features)))`. This may also optionally be an instance of `tf$linalg$LinearOperator`.

weights_prior_scale float Tensor defining the scale of the Horseshoe prior on regression weights. Small values encourage the weights to be sparse. The shape must broadcast with `weights_batch_shape`. Default value: 0.1.

weights_batch_shape if NULL, defaults to `design_matrix.batch_shape_tensor()`. Must broadcast with the batch shape of `design_matrix`. Default value: NULL.

name the name of this model component. Default value: 'LinearRegression'.

Details

```
observed_time_series <- tf$matmul(design_matrix, weights)
```

This is identical to `sts_linear_regression`, except that `sts_sparse_linear_regression` uses a parameterization of a Horseshoe prior to encode the assumption that many of the weights are zero, i.e., many of the covariate time series are irrelevant. See the mathematical details section below for further discussion. The prior parameterization used by `sts_sparse_linear_regression` is more suitable for inference than that obtained by simply passing the equivalent `tfd_horseshoe` prior to `sts_linear_regression`; when sparsity is desired, `sts_sparse_linear_regression` will likely yield better results.

This component does not itself include observation noise; it defines a deterministic distribution with mass at the point `tf$matmul(design_matrix, weights)`. In practice, it should be combined with observation noise from another component such as `sts_sum`.

Mathematical Details

The basic horseshoe prior Carvalho et al. (2009) is defined as a Cauchy-normal scale mixture:

```
scales[i] ~ HalfCauchy(loc=0, scale=1)
weights[i] ~ Normal(loc=0., scale=scales[i] * global_scale)`
```

The Cauchy scale parameters puts substantial mass near zero, encouraging weights to be sparse, but their heavy tails allow weights far from zero to be estimated without excessive shrinkage. The horseshoe can be thought of as a continuous relaxation of a traditional 'spike-and-slab' discrete sparsity prior, in which the latent Cauchy scale mixes between 'spike' ($scales[i] \sim 0$) and 'slab' ($scales[i] \gg 0$) regimes.

Following the recommendations in Piironen et al. (2017), `SparseLinearRegression` implements a horseshoe with the following adaptations:

- The Cauchy prior on `scales[i]` is represented as an InverseGamma-Normal compound.
- The `global_scale` parameter is integrated out following a $Cauchy(0., scale=weights_prior_scale)$ hyperprior, which is also represented as an InverseGamma-Normal compound.
- All compound distributions are implemented using a non-centered parameterization. The compound, non-centered representation defines the same marginal prior as the original horseshoe (up to integrating out the global scale), but allows samplers to mix more efficiently through the heavy tails; for variational inference, the compound representation implicitly expands the representational power of the variational model.

Note that we do not yet implement the regularized ('Finnish') horseshoe, proposed in Piironen et al. (2017) for models with weak likelihoods, because the likelihood in STS models is typically Gaussian, where it's not clear that additional regularization is appropriate. If you need this functionality, please email tfprobability@tensorflow.org.

The full prior parameterization implemented in `SparseLinearRegression` is as follows:

```
Sample global_scale from Cauchy(0, scale=weights_prior_scale).
global_scale_variance ~ InverseGamma(alpha=0.5, beta=0.5)
global_scale_noncentered ~ HalfNormal(loc=0, scale=1)
global_scale = (global_scale_noncentered *
sqrt(global_scale_variance) *
weights_prior_scale)
Sample local_scales from Cauchy(0, 1).
local_scale_variances[i] ~ InverseGamma(alpha=0.5, beta=0.5)
local_scales_noncentered[i] ~ HalfNormal(loc=0, scale=1)
local_scales[i] = local_scales_noncentered[i] * sqrt(local_scale_variances[i])
weights[i] ~ Normal(loc=0., scale=local_scales[i] * global_scale)
```

Value

an instance of `StructuralTimeSeries`.

References

- [Carvalho, C., Polson, N. and Scott, J. Handling Sparsity via the Horseshoe. AISTATS \(2009\).](#)
- [Juho Piironen, Aki Vehtari. Sparsity information and regularization in the horseshoe and other shrinkage priors \(2017\).](#)

See Also

For usage examples see [sts_fit_with_hmc\(\)](#), [sts_forecast\(\)](#), [sts_decompose_by_component\(\)](#).

Other sts: [sts_additive_state_space_model\(\)](#), [sts_autoregressive_state_space_model\(\)](#), [sts_autoregressive\(\)](#), [sts_constrained_seasonal_state_space_model\(\)](#), [sts_dynamic_linear_regression_state_space_model\(\)](#), [sts_dynamic_linear_regression\(\)](#), [sts_linear_regression\(\)](#), [sts_local_level_state_space_model\(\)](#), [sts_local_level\(\)](#), [sts_local_linear_trend_state_space_model\(\)](#), [sts_local_linear_trend\(\)](#), [sts_seasonal_state_space_model\(\)](#), [sts_seasonal\(\)](#), [sts_semi_local_linear_trend_state_space_model\(\)](#), [sts_semi_local_linear_trend\(\)](#), [sts_smooth_seasonal_state_space_model\(\)](#), [sts_smooth_seasonal\(\)](#), [sts_sum\(\)](#)

sts_sum

*Sum of structural time series components.***Description**

This class enables compositional specification of a structural time series model from basic components. Given a list of component models, it represents an additive model, i.e., a model of time series that may be decomposed into a sum of terms corresponding to the component models.

Usage

```
sts_sum(
    observed_time_series = NULL,
    components,
    constant_offset = NULL,
    observation_noise_scale_prior = NULL,
    name = NULL
)
```

Arguments

observed_time_series

optional float tensor of shape batch_shape + [T, 1] (omitting the trailing unit dimension is also supported when T > 1), specifying an observed time series. Any priors not explicitly set will be given default values according to the scale of the observed time series (or batch of time series). May optionally be an instance of [sts_masked_time_series](#), which includes a mask tensor to specify timesteps with missing observations. Default value: NULL.

components

list of one or more [StructuralTimeSeries](#) instances. These must have unique names.

constant_offset

optional scalar float tensor, or batch of scalars, specifying a constant value added to the sum of outputs from the component models. This allows the components to model the shifted series `observed_time_series - constant_offset`. If NULL, this is set to the mean of the provided `observed_time_series`. Default value: NULL.

observation_noise_scale_prior
 optional tfd\$Distribution instance specifying a prior on observation_noise_scale.
 If NULL, a heuristic default prior is constructed based on the provided observed_time_series.
 Default value: NULL.

name
 string name of this model component; used as name_scope for ops created by
 this class. Default value: 'Sum'.

Details

Formally, the additive model represents a random process $g[t] = f_1[t] + f_2[t] + \dots + f_N[t] + \text{eps}[t]$, where the f 's are the random processes represented by the components, and $\text{eps}[t] \sim \text{Normal}(\text{loc}=0, \text{scale}=\text{observation_noise_scale})$ is an observation noise term. See the AdditiveStateSpaceModel documentation for mathematical details.

This model inherits the parameters (with priors) of its components, and adds an observation_noise_scale parameter governing the level of noise in the observed time series.

Value

an instance of StructuralTimeSeries.

See Also

For usage examples see [sts_fit_with_hmc\(\)](#), [sts_forecast\(\)](#), [sts_decompose_by_component\(\)](#).

Other sts: [sts_additive_state_space_model\(\)](#), [sts_autoregressive_state_space_model\(\)](#), [sts_autoregressive\(\)](#), [sts_constrained_seasonal_state_space_model\(\)](#), [sts_dynamic_linear_regression_state_space_model\(\)](#), [sts_dynamic_linear_regression\(\)](#), [sts_linear_regression\(\)](#), [sts_local_level_state_space_model\(\)](#), [sts_local_level\(\)](#), [sts_local_linear_trend_state_space_model\(\)](#), [sts_local_linear_trend\(\)](#), [sts_seasonal_state_space_model\(\)](#), [sts_seasonal\(\)](#), [sts_semi_local_linear_trend_state_space_model\(\)](#), [sts_semi_local_linear_trend\(\)](#), [sts_smooth_seasonal_state_space_model\(\)](#), [sts_smooth_seasonal\(\)](#), [sts_sparse_linear_regression\(\)](#)

tfb_absolute_value	<i>Computes</i> $Y = g(X) = \text{Abs}(X)$, <i>element-wise</i>
--------------------	--

Description

This non-injective bijector allows for transformations of scalar distributions with the absolute value function, which maps $(-\infty, \infty)$ to $[0, \infty)$.

- For y in $(0, \infty)$, `tfb_absolute_value$inverse(y)` returns the set inverse $\{x \text{ in } (-\infty, \infty) : |x| = y\}$ as a tuple, $-y, y$. `tfb_absolute_value$inverse(0)` returns $0, 0$, which is not the set inverse (the set inverse is the singleton $\{0\}$), but "works" in conjunction with `TransformedDistribution` to produce a left semi-continuous pdf. For $y < 0$, `tfb_absolute_value$inverse(y)` happily returns the wrong thing, $-y, y$. This is done for efficiency. If `validate_args == TRUE`, $y < 0$ will raise an exception.

Usage

```
tfb_absolute_value(validate_args = FALSE, name = "absolute_value")
```

Arguments

`validate_args` Logical, default FALSE. Whether to validate input with asserts. If `validate_args` is FALSE, and the inputs are invalid, correct behavior is not guaranteed.

`name` name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_template()`, `tfb_masked_autoregressive_flow()`, `tfb_masked_dense()`, `tfb_matrix_inverse_tri_l()`, `tfb_matvec_lu()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_pad()`, `tfb_permute()`, `tfb_power_transform()`, `tfb_rational_quadratic_spline()`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`, `tfb_real_nvp()`, `tfb_reciprocal()`, `tfb_reshape()`, `tfb_scale_matvec_diag()`, `tfb_scale_matvec_linear_operator()`, `tfb_scale_matvec_lu()`, `tfb_scale_matvec_tri_l()`, `tfb_scale_tri_l()`, `tfb_scale()`, `tfb_shifted_gompertz_cdf()`, `tfb_shift()`, `tfb_sigmoid()`, `tfb_sinh_arcsinh()`, `tfb_sinh()`, `tfb_softmax_centered()`, `tfb_softplus()`, `tfb_softsign()`, `tfb_split()`, `tfb_square()`, `tfb_tanh()`, `tfb_transform_diagonal()`, `tfb_transpose()`, `tfb_weibull_cdf()`, `tfb_weibull()`

tfb_affine

Affine bijector

Description

This Bijector is initialized with `shift` Tensor and `scale` arguments, giving the forward operation: $Y = g(X) = \text{scale} @ X + \text{shift}$ where the `scale` term is logically equivalent to: `scale = scale_identity_multiplier * tf.diag(tf.ones(d))`

Usage

```
tfb_affine(
  shift = NULL,
  scale_identity_multiplier = NULL,
  scale_diag = NULL,
```

```

    scale_tril = NULL,
    scale_perturb_factor = NULL,
    scale_perturb_diag = NULL,
    adjoint = FALSE,
    validate_args = FALSE,
    name = "affine",
    dtype = NULL
)

```

Arguments

<code>shift</code>	Floating-point Tensor. If this is set to NULL, no shift is applied.
<code>scale_identity_multiplier</code>	floating point rank 0 Tensor representing a scaling done to the identity matrix. When <code>scale_identity_multiplier = scale_diag = scale_tril = NULL</code> then <code>scale += IdentityMatrix</code> . Otherwise no scaled-identity-matrix is added to <code>scale</code> .
<code>scale_diag</code>	Floating-point Tensor representing the diagonal matrix. <code>scale_diag</code> has shape <code>[N1, N2, ... k]</code> , which represents a <code>k x k</code> diagonal matrix. When NULL no diagonal term is added to <code>scale</code> .
<code>scale_tril</code>	Floating-point Tensor representing the lower triangular matrix. <code>scale_tril</code> has shape <code>[N1, N2, ... k, k]</code> , which represents a <code>k x k</code> lower triangular matrix. When NULL no <code>scale_tril</code> term is added to <code>scale</code> . The upper triangular elements above the diagonal are ignored.
<code>scale_perturb_factor</code>	Floating-point Tensor representing factor matrix with last two dimensions of shape <code>(k, r)</code> When NULL, no rank- <code>r</code> update is added to <code>scale</code> .
<code>scale_perturb_diag</code>	Floating-point Tensor representing the diagonal matrix. <code>scale_perturb_diag</code> has shape <code>[N1, N2, ... r]</code> , which represents an <code>r x r</code> diagonal matrix. When NULL low rank updates will take the form <code>scale_perturb_factor * scale_perturb_factor . T</code> .
<code>adjoint</code>	Logical indicating whether to use the <code>scale</code> matrix as specified or its adjoint. Default value: FALSE.
<code>validate_args</code>	Logical, default FALSE. Whether to validate input with asserts. If <code>validate_args</code> is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
<code>name</code>	name prefixed to Ops created by this class.
<code>dtype</code>	<code>tf\$DType</code> to prefer when converting args to Tensors. Else, we fall back to a common dtype inferred from the args, finally falling back to <code>float32</code> .

Details

If NULL of `scale_identity_multiplier`, `scale_diag`, or `scale_tril` are specified then `scale += IdentityMatrix` Otherwise specifying a `scale` argument has the semantics of `scale += Expand(arg)`, i.e., `scale_diag != NULL` means `scale += tf$diag(scale_diag)`.

Value

a bijector instance.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_template()`, `tfb_masked_autoregressive_flow()`, `tfb_masked_dense()`, `tfb_matrix_inverse_tri_l()`, `tfb_matvec_lu()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_pad()`, `tfb_permute()`, `tfb_power_transform()`, `tfb_rational_quadratic_spline()`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`, `tfb_real_nvp()`, `tfb_reciprocal()`, `tfb_reshape()`, `tfb_scale_matvec_diag()`, `tfb_scale_matvec_linear_operator()`, `tfb_scale_matvec_lu()`, `tfb_scale_matvec_tri_l()`, `tfb_scale_tri_l()`, `tfb_scale()`, `tfb_shifted_gompertz_cdf()`, `tfb_shift()`, `tfb_sigmoid()`, `tfb_sinh_arcsinh()`, `tfb_sinh()`, `tfb_softmax_centered()`, `tfb_softplus()`, `tfb_softsign()`, `tfb_split()`, `tfb_square()`, `tfb_tanh()`, `tfb_transform_diagonal()`, `tfb_transpose()`, `tfb_weibull_cdf()`, `tfb_weibull()`

`tfb_affine_linear_operator`

Computes $Y = g(X; \text{shift}, \text{scale}) = \text{scale} @ X + \text{shift}$

Description

`shift` is a numeric Tensor and `scale` is a LinearOperator. If `X` is a scalar then the forward transformation is: `scale * X + shift` where `*` denotes broadcasted elementwise product.

Usage

```
tfb_affine_linear_operator(
    shift = NULL,
    scale = NULL,
    adjoint = FALSE,
    validate_args = FALSE,
    name = "affine_linear_operator"
)
```

Arguments

<code>shift</code>	Floating-point Tensor.
<code>scale</code>	Subclass of LinearOperator. Represents the (batch) positive definite matrix M in $\mathbb{R}^{[k \times k]}$.
<code>adjoint</code>	Logical indicating whether to use the scale matrix as specified or its adjoint. Default value: FALSE.
<code>validate_args</code>	Logical, default FALSE. Whether to validate input with asserts. If <code>validate_args</code> is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
<code>name</code>	name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_template()`, `tfb_masked_autoregressive_flow()`, `tfb_masked_dense()`, `tfb_matrix_inverse_tri_l()`, `tfb_matvec_lu()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_pad()`, `tfb_permute()`, `tfb_power_transform()`, `tfb_rational_quadratic_spline()`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`, `tfb_real_nvp()`, `tfb_reciprocal()`, `tfb_reshape()`, `tfb_scale_matvec_diag()`, `tfb_scale_matvec_linear_operator()`, `tfb_scale_matvec_lu()`, `tfb_scale_matvec_tri_l()`, `tfb_scale_tri_l()`, `tfb_scale()`, `tfb_shifted_gompertz_cdf()`, `tfb_shift()`, `tfb_sigmoid()`, `tfb_sinh_arcsinh()`, `tfb_sinh()`, `tfb_softmax_centered()`, `tfb_softplus()`, `tfb_softsign()`, `tfb_split()`, `tfb_square()`, `tfb_tanh()`, `tfb_transform_diagonal()`, `tfb_transpose()`, `tfb_weibull_cdf()`, `tfb_weibull()`

tfb_ascending

Maps unconstrained R^n to R^n in ascending order.

Description

Both the domain and the codomain of the mapping is $[-\text{inf}, \text{inf}]^n$, however, the input of the inverse mapping must be strictly increasing. On the last dimension of the tensor, the Ascending bijector performs: $y = \text{tf}\$cumsum([x[0], \text{tf}\$\exp(x[1]), \text{tf}\$\exp(x[2]), \dots, \text{tf}\$\exp(x[-1])])$

Usage

```
tfb_ascending(validate_args = FALSE, name = "ascending")
```

Arguments

`validate_args` Logical, default FALSE. Whether to validate input with asserts. If `validate_args` is FALSE, and the inputs are invalid, correct behavior is not guaranteed.

`name` name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see [tfb_forward\(\)](#), [tfb_inverse\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#).

Other bijectors: [tfb_absolute_value\(\)](#), [tfb_affine_linear_operator\(\)](#), [tfb_affine_scalar\(\)](#), [tfb_affine\(\)](#), [tfb_batch_normalization\(\)](#), [tfb_blockwise\(\)](#), [tfb_chain\(\)](#), [tfb_cholesky_outer_product\(\)](#), [tfb_cholesky_to_inv_cholesky\(\)](#), [tfb_correlation_cholesky\(\)](#), [tfb_cumsum\(\)](#), [tfb_discrete_cosine_transform\(\)](#), [tfb_expm1\(\)](#), [tfb_exp\(\)](#), [tfb_ffjord\(\)](#), [tfb_fill_scale_tri_l\(\)](#), [tfb_fill_triangular\(\)](#), [tfb_glow\(\)](#), [tfb_gompertz_cdf\(\)](#), [tfb_gumbel_cdf\(\)](#), [tfb_gumbel\(\)](#), [tfb_identity\(\)](#), [tfb_inline\(\)](#), [tfb_invert\(\)](#), [tfb_iterated_sigmoid_centered\(\)](#), [tfb_kumaraswamy_cdf\(\)](#), [tfb_kumaraswamy\(\)](#), [tfb_lambert_w_tail\(\)](#), [tfb_masked_autoregressive_default_template\(\)](#), [tfb_masked_autoregressive_flow\(\)](#), [tfb_masked_dense\(\)](#), [tfb_matrix_inverse_tri_l\(\)](#), [tfb_matvec_lu\(\)](#), [tfb_normal_cdf\(\)](#), [tfb_ordered\(\)](#), [tfb_pad\(\)](#), [tfb_permute\(\)](#), [tfb_power_transform\(\)](#), [tfb_rational_quadratic_spline\(\)](#), [tfb_rayleigh_cdf\(\)](#), [tfb_real_nvp_default_template\(\)](#), [tfb_real_nvp\(\)](#), [tfb_reciprocal\(\)](#), [tfb_reshape\(\)](#), [tfb_scale_matvec_diag\(\)](#), [tfb_scale_matvec_linear_operator\(\)](#), [tfb_scale_matvec_lu\(\)](#), [tfb_scale_matvec_tri_l\(\)](#), [tfb_scale_tri_l\(\)](#), [tfb_scale\(\)](#), [tfb_shifted_gompertz_cdf\(\)](#), [tfb_shift\(\)](#), [tfb_sigmoid\(\)](#), [tfb_sinh_arcsinh\(\)](#), [tfb_sinh\(\)](#), [tfb_softmax_centered\(\)](#), [tfb_softplus\(\)](#), [tfb_softsign\(\)](#), [tfb_split\(\)](#), [tfb_square\(\)](#), [tfb_tanh\(\)](#), [tfb_transform_diagonal\(\)](#), [tfb_transpose\(\)](#), [tfb_weibull_cdf\(\)](#), [tfb_weibull\(\)](#)

tfb_batch_normalization

Computes $Y = g(X)$ s.t. $X = g^{-1}(Y) = (Y - \text{mean}(Y)) / \text{std}(Y)$

Description

Applies Batch Normalization (Ioffe and Szegedy, 2015) to samples from a data distribution. This can be used to stabilize training of normalizing flows (Papamakarios et al., 2016; Dinh et al., 2017)

Usage

```
tfb_batch_normalization(
  batchnorm_layer = NULL,
  training = TRUE,
  validate_args = FALSE,
  name = "batch_normalization"
)
```

Arguments

batchnorm_layer	tf\$layers\$BatchNormalization layer object. If NULL, defaults to tf\$layers\$BatchNormalization + 1e-6). This ensures positivity of the scale variable.
training	If TRUE, updates running-average statistics during call to inverse().
validate_args	Logical, default FALSE. Whether to validate input with asserts. If validate_args is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
name	name prefixed to Ops created by this class.

Details

When training Deep Neural Networks (DNNs), it is common practice to normalize or whiten features by shifting them to have zero mean and scaling them to have unit variance.

The `inverse()` method of the `BatchNormalization` bijector, which is used in the log-likelihood computation of data samples, implements the normalization procedure (shift-and-scale) using the mean and standard deviation of the current minibatch.

Conversely, the `forward()` method of the bijector de-normalizes samples (e.g. $X * \text{std}(Y) + \text{mean}(Y)$) with the running-average mean and standard deviation computed at training-time. De-normalization is useful for sampling.

During training time, `BatchNormalization.inverse` and `BatchNormalization.forward` are not guaranteed to be inverses of each other because `inverse(y)` uses statistics of the current minibatch, while `forward(x)` uses running-average statistics accumulated from training. In other words, `tfb_batch_normalization().inverse(tfb_batch_normalization().forward(...))` and `tfb_batch_normalization().forward(tfb_batch_normalization().inverse(...))` will be identical when `training=FALSE` but may be different when `training=TRUE`.

Value

a bijector instance.

References

- [Sergey Ioffe and Christian Szegedy. Batch Normalization: Accelerating Deep Network Training by Reducing Internal Covariate Shift. In *International Conference on Machine Learning*, 2015.](#)
- [Laurent Dinh, Jascha Sohl-Dickstein, and Samy Bengio. Density Estimation using Real NVP. In *International Conference on Learning Representations*, 2017.](#)
- [George Papamakarios, Theo Pavlakou, and Iain Murray. Masked Autoregressive Flow for Density Estimation. In *Neural Information Processing Systems*, 2017.](#)

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_template()`, `tfb_masked_autoregressive_flow()`, `tfb_masked_dense()`, `tfb_matrix_inverse_tri_l()`, `tfb_matvec_lu()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_pad()`, `tfb_permute()`, `tfb_power_transform()`, `tfb_rational_quadratic_spline()`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`, `tfb_real_nvp()`, `tfb_reciprocal()`, `tfb_reshape()`, `tfb_scale_matvec_diag()`, `tfb_scale_matvec_linear_operator()`, `tfb_scale_matvec_lu()`, `tfb_scale_matvec_tri_l()`, `tfb_scale_tri_l()`, `tfb_scale()`, `tfb_shifted_gompertz_cdf()`, `tfb_shift()`, `tfb_sigmoid()`, `tfb_sinh_arcsinh()`, `tfb_sinh()`, `tfb_softmax_centered()`, `tfb_softplus()`, `tfb_softsign()`, `tfb_split()`, `tfb_square()`, `tfb_tanh()`, `tfb_transform_diagonal()`, `tfb_transpose()`, `tfb_weibull_cdf()`, `tfb_weibull()`

tfb_blockwise	<i>Bijector which applies a list of bijectors to blocks of a Tensor</i>
---------------	---

Description

More specifically, given $[F_0, F_1, \dots, F_n]$ which are scalar or vector bijectors this bijector creates a transformation which operates on the vector $[x_0, \dots, x_n]$ with the transformation $[F_0(x_0), F_1(x_1), \dots, F_n(x_n)]$ where x_0, \dots, x_n are blocks (partitions) of the vector.

Usage

```
tfb_blockwise(
    bijectors,
    block_sizes = NULL,
    validate_args = FALSE,
    name = NULL
)
```

Arguments

bijectors	A non-empty list of bijectors.
block_sizes	A 1-D integer Tensor with each element signifying the length of the block of the input vector to pass to the corresponding bijector. The length of block_sizes must be equal to the length of bijectors. If left as NULL, a vector of 1's is used.
validate_args	Logical indicating whether arguments should be checked for correctness.
name	String, name given to ops managed by this object. Default: E.g., <code>tfb_blockwise(list(tfb_exp()), tfb_exp()) == 'blockwise_of_exp_and_softplus'</code> .

Value

a bijector instance.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_template()`, `tfb_masked_autoregressive_flow()`, `tfb_masked_dense()`, `tfb_matrix_inverse_tri_l()`, `tfb_matvec_lu()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_pad()`, `tfb_permute()`, `tfb_power_transform()`, `tfb_rational_quadratic_spline()`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`, `tfb_real_nvp()`, `tfb_reciprocal()`,

```
tfb_reshape(), tfb_scale_matvec_diag(), tfb_scale_matvec_linear_operator(), tfb_scale_matvec_lu(),
tfb_scale_matvec_tri_l(), tfb_scale_tri_l(), tfb_scale(), tfb_shifted_gompertz_cdf(),
tfb_shift(), tfb_sigmoid(), tfb_sinh_arcsinh(), tfb_sinh(), tfb_softmax_centered(),
tfb_softplus(), tfb_softsign(), tfb_split(), tfb_square(), tfb_tanh(), tfb_transform_diagonal(),
tfb_transpose(), tfb_weibull_cdf(), tfb_weibull()
```

 tfb_chain

Bijector which applies a sequence of bijectors

Description

Bijector which applies a sequence of bijectors

Usage

```
tfb_chain(
  bijectors = NULL,
  validate_args = FALSE,
  validate_event_size = TRUE,
  parameters = NULL,
  name = NULL
)
```

Arguments

bijectors list of bijector instances. An empty list makes this bijector equivalent to the Identity bijector.

validate_args Logical indicating whether arguments should be checked for correctness.

validate_event_size Checks that bijectors are not applied to inputs with incomplete support (that is, inputs where one or more elements are a deterministic transformation of the others). For example, the following LDJ would be incorrect: `tfb_chain(list(tfb_scale(), tfb_softmax_centered(= 2))` The jacobian contribution from `tfb_scale()` applies to a 2-dimensional input, but the output from `tfb_softmax_centered()` is a 1-dimensional input embedded in a 2-dimensional space. Setting `validate_event_size=TRUE` (default) prints warnings in these cases. When `validate_args` is also `TRUE`, the warning is promoted to an exception.

parameters Locals dict captured by subclass constructor, to be used for copy/slice re-instantiation operators.

name String, name given to ops managed by this object. Default: E.g., `tfb_chain(list(tfb_exp(), tfb_softmax_centered()))` == "chain_of_exp_of_softmax_centered".

Value

a bijector instance.

See Also

For usage examples see [tfb_forward\(\)](#), [tfb_inverse\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#).

Other bijectors: [tfb_absolute_value\(\)](#), [tfb_affine_linear_operator\(\)](#), [tfb_affine_scalar\(\)](#), [tfb_affine\(\)](#), [tfb_ascending\(\)](#), [tfb_batch_normalization\(\)](#), [tfb_blockwise\(\)](#), [tfb_cholesky_outer_product\(\)](#), [tfb_cholesky_to_inv_cholesky\(\)](#), [tfb_correlation_cholesky\(\)](#), [tfb_cumsum\(\)](#), [tfb_discrete_cosine_transform\(\)](#), [tfb_expm1\(\)](#), [tfb_exp\(\)](#), [tfb_ffjord\(\)](#), [tfb_fill_scale_tri_l\(\)](#), [tfb_fill_triangular\(\)](#), [tfb_glow\(\)](#), [tfb_gompertz_cdf\(\)](#), [tfb_gumbel_cdf\(\)](#), [tfb_gumbel\(\)](#), [tfb_identity\(\)](#), [tfb_inline\(\)](#), [tfb_invert\(\)](#), [tfb_iterated_sigmoid_centered\(\)](#), [tfb_kumaraswamy_cdf\(\)](#), [tfb_kumaraswamy\(\)](#), [tfb_lambert_w_tail\(\)](#), [tfb_masked_autoregressive_default_template\(\)](#), [tfb_masked_autoregressive_flow\(\)](#), [tfb_masked_dense\(\)](#), [tfb_matrix_inverse_tri_l\(\)](#), [tfb_matvec_lu\(\)](#), [tfb_normal_cdf\(\)](#), [tfb_ordered\(\)](#), [tfb_pad\(\)](#), [tfb_permute\(\)](#), [tfb_power_transform\(\)](#), [tfb_rational_quadratic_spline\(\)](#), [tfb_rayleigh_cdf\(\)](#), [tfb_real_nvp_default_template\(\)](#), [tfb_real_nvp\(\)](#), [tfb_reciprocal\(\)](#), [tfb_reshape\(\)](#), [tfb_scale_matvec_diag\(\)](#), [tfb_scale_matvec_linear_operator\(\)](#), [tfb_scale_matvec_lu\(\)](#), [tfb_scale_matvec_tri_l\(\)](#), [tfb_scale_tri_l\(\)](#), [tfb_scale\(\)](#), [tfb_shifted_gompertz_cdf\(\)](#), [tfb_shift\(\)](#), [tfb_sigmoid\(\)](#), [tfb_sinh_arcsinh\(\)](#), [tfb_sinh\(\)](#), [tfb_softmax_centered\(\)](#), [tfb_softplus\(\)](#), [tfb_softsign\(\)](#), [tfb_split\(\)](#), [tfb_square\(\)](#), [tfb_tanh\(\)](#), [tfb_transform_diagonal\(\)](#), [tfb_transpose\(\)](#), [tfb_weibull_cdf\(\)](#), [tfb_weibull\(\)](#)

tfb_cholesky_outer_product

Computes $g(X) = X @ X.T$ where X is lower-triangular; positive-diagonal matrix

Description

Note: the upper-triangular part of X is ignored (whether or not its zero).

Usage

```
tfb_cholesky_outer_product(
    validate_args = FALSE,
    name = "cholesky_outer_product"
)
```

Arguments

`validate_args` Logical, default FALSE. Whether to validate input with asserts. If `validate_args` is FALSE, and the inputs are invalid, correct behavior is not guaranteed.

`name` name prefixed to Ops created by this class.

Details

The surjectivity of g as a map from the set of $n \times n$ positive-diagonal lower-triangular matrices to the set of SPD matrices follows immediately from executing the Cholesky factorization algorithm on an SPD matrix A to produce a positive-diagonal lower-triangular matrix L such that $A = L @ L.T$.

To prove the injectivity of g , suppose that L_1 and L_2 are lower-triangular with positive diagonals and satisfy $A = L_1 @ L_1.T = L_2 @ L_2.T$. Then $inv(L_1) @ A @ inv(L_1).T = [inv(L_1) @ L_2] @ [inv(L_1) @ L_2].T = I$. Setting $L_3 := inv(L_1) @ L_2$, that L_3 is a positive-diagonal lower-triangular matrix follows from $inv(L_1)$ being positive-diagonal lower-triangular (which follows from the diagonal of a triangular matrix being its spectrum), and that the product of two positive-diagonal lower-triangular matrices is another positive-diagonal lower-triangular matrix. A simple inductive argument (proceeding one column of L_3 at a time) shows that, if $I = L_3 @ L_3.T$, with L_3 being lower-triangular with positive-diagonal, then $L_3 = I$. Thus, $L_1 = L_2$, proving injectivity of g .

Value

a bijector instance.

See Also

For usage examples see [tfb_forward\(\)](#), [tfb_inverse\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#).

Other bijectors: [tfb_absolute_value\(\)](#), [tfb_affine_linear_operator\(\)](#), [tfb_affine_scalar\(\)](#), [tfb_affine\(\)](#), [tfb_ascending\(\)](#), [tfb_batch_normalization\(\)](#), [tfb_blockwise\(\)](#), [tfb_chain\(\)](#), [tfb_cholesky_to_inv_cholesky\(\)](#), [tfb_correlation_cholesky\(\)](#), [tfb_cumsum\(\)](#), [tfb_discrete_cosine_transform\(\)](#), [tfb_expm1\(\)](#), [tfb_exp\(\)](#), [tfb_ffjord\(\)](#), [tfb_fill_scale_tri_l\(\)](#), [tfb_fill_triangular\(\)](#), [tfb_glow\(\)](#), [tfb_gompertz_cdf\(\)](#), [tfb_gumbel_cdf\(\)](#), [tfb_gumbel\(\)](#), [tfb_identity\(\)](#), [tfb_inline\(\)](#), [tfb_invert\(\)](#), [tfb_iterated_sigmoid_centered\(\)](#), [tfb_kumaraswamy_cdf\(\)](#), [tfb_kumaraswamy\(\)](#), [tfb_lambert_w_tail\(\)](#), [tfb_masked_autoregressive_default_template\(\)](#), [tfb_masked_autoregressive_flow\(\)](#), [tfb_masked_dense\(\)](#), [tfb_matrix_inverse_tri_l\(\)](#), [tfb_matvec_lu\(\)](#), [tfb_normal_cdf\(\)](#), [tfb_ordered\(\)](#), [tfb_pad\(\)](#), [tfb_permute\(\)](#), [tfb_power_transform\(\)](#), [tfb_rational_quadratic_spline\(\)](#), [tfb_rayleigh_cdf\(\)](#), [tfb_real_nvp_default_template\(\)](#), [tfb_real_nvp\(\)](#), [tfb_reciprocal\(\)](#), [tfb_reshape\(\)](#), [tfb_scale_matvec_diag\(\)](#), [tfb_scale_matvec_linear_operator\(\)](#), [tfb_scale_matvec_lu\(\)](#), [tfb_scale_matvec_tri_l\(\)](#), [tfb_scale_tri_l\(\)](#), [tfb_scale\(\)](#), [tfb_shifted_gompertz_cdf\(\)](#), [tfb_shift\(\)](#), [tfb_sigmoid\(\)](#), [tfb_sinh_arcsinh\(\)](#), [tfb_sinh\(\)](#), [tfb_softmax_centered\(\)](#), [tfb_softplus\(\)](#), [tfb_softsign\(\)](#), [tfb_split\(\)](#), [tfb_square\(\)](#), [tfb_tanh\(\)](#), [tfb_transform_diagonal\(\)](#), [tfb_transpose\(\)](#), [tfb_weibull_cdf\(\)](#), [tfb_weibull\(\)](#)

tfb_cholesky_to_inv_cholesky

Maps the Cholesky factor of M to the Cholesky factor of M^{-1}

Description

The forward and inverse calculations are conceptually identical to: `forward <-function(x) tf$cholesky(tf$linalg$inv(x))`
`inverse = forward` However, the actual calculations exploit the triangular structure of the matrices.

Usage

```
tfb_cholesky_to_inv_cholesky(
  validate_args = FALSE,
  name = "cholesky_to_inv_cholesky"
)
```

Arguments

`validate_args` Logical, default FALSE. Whether to validate input with asserts. If `validate_args` is FALSE, and the inputs are invalid, correct behavior is not guaranteed.

`name` name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_template()`, `tfb_masked_autoregressive_flow()`, `tfb_masked_dense()`, `tfb_matrix_inverse_tri_l()`, `tfb_matvec_lu()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_pad()`, `tfb_permute()`, `tfb_power_transform()`, `tfb_rational_quadratic_spline()`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`, `tfb_real_nvp()`, `tfb_reciprocal()`, `tfb_reshape()`, `tfb_scale_matvec_diag()`, `tfb_scale_matvec_linear_operator()`, `tfb_scale_matvec_lu()`, `tfb_scale_matvec_tri_l()`, `tfb_scale_tri_l()`, `tfb_scale()`, `tfb_shifted_gompertz_cdf()`, `tfb_shift()`, `tfb_sigmoid()`, `tfb_sinh_arcsinh()`, `tfb_sinh()`, `tfb_softmax_centered()`, `tfb_softplus()`, `tfb_softsign()`, `tfb_split()`, `tfb_square()`, `tfb_tanh()`, `tfb_transform_diagonal()`, `tfb_transpose()`, `tfb_weibull_cdf()`, `tfb_weibull()`

tfb_correlation_cholesky

Maps unconstrained reals to Cholesky-space correlation matrices.

Description

This bijector is a mapping between \mathbb{R}^n and the n -dimensional manifold of Cholesky-space correlation matrices embedded in $\mathbb{R}^{\{m^2\}}$, where n is the $(m-1)$ th triangular number; i.e. $n = 1 + 2 + \dots + (m-1)$.

Usage

```
tfb_correlation_cholesky(validate_args = FALSE, name = "correlation_cholesky")
```

Arguments

`validate_args` Logical, default FALSE. Whether to validate input with asserts. If `validate_args` is FALSE, and the inputs are invalid, correct behavior is not guaranteed.

`name` name prefixed to Ops created by this class.

Details

Mathematical Details

The image of unconstrained reals under the CorrelationCholesky bijector is the set of correlation matrices which are positive definite. A **correlation matrix** can be characterized as a symmetric positive semidefinite matrix with 1s on the main diagonal. However, the correlation matrix is positive definite if no component can be expressed as a linear combination of the other components. For a lower triangular matrix L to be a valid Cholesky-factor of a positive definite correlation matrix, it is necessary and sufficient that each row of L have unit Euclidean norm. To see this, observe that if L_i is the i th row of the Cholesky factor corresponding to the correlation matrix R , then the i th diagonal entry of R satisfies:

$$1 = R_{i,i} = L_i \cdot L_i = \|L_i\|^2$$

where \cdot is the dot product of vectors and $\|\dots\|$ denotes the Euclidean norm. Furthermore, observe that $R_{i,j}$ lies in the interval $[-1, 1]$. By the Cauchy-Schwarz inequality:

$$|R_{i,j}| = |L_i \cdot L_j| \leq \|L_i\| \|L_j\| = 1$$

This is a consequence of the fact that R is symmetric positive definite with 1s on the main diagonal. The LKJ distribution with `input_output_cholesky=TRUE` generates samples from (and computes log-densities on) the set of Cholesky factors of positive definite correlation matrices. The CorrelationCholesky bijector provides a bijective mapping from unconstrained reals to the support of the LKJ distribution.

Value

a bijector instance.

References

- [Stan Manual. Section 24.2. Cholesky LKJ Correlation Distribution.](#)
- Daniel Lewandowski, Dorota Kurowicka, and Harry Joe, "Generating random correlation matrices based on vines and extended onion method," Journal of Multivariate Analysis 100 (2009), pp 1989-2001.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_template()`, `tfb_masked_autoregressive_flow()`, `tfb_masked_dense()`, `tfb_matrix_inverse_tri_l()`, `tfb_matvec_lu()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_pad()`, `tfb_permute()`, `tfb_power_transform()`, `tfb_rational_quadratic_spline()`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`, `tfb_real_nvp()`, `tfb_reciprocal()`,

[tfb_reshape\(\)](#), [tfb_scale_matvec_diag\(\)](#), [tfb_scale_matvec_linear_operator\(\)](#), [tfb_scale_matvec_lu\(\)](#), [tfb_scale_matvec_tri_l\(\)](#), [tfb_scale_tri_l\(\)](#), [tfb_scale\(\)](#), [tfb_shifted_gompertz_cdf\(\)](#), [tfb_shift\(\)](#), [tfb_sigmoid\(\)](#), [tfb_sinh_arcsinh\(\)](#), [tfb_sinh\(\)](#), [tfb_softmax_centered\(\)](#), [tfb_softplus\(\)](#), [tfb_softsign\(\)](#), [tfb_split\(\)](#), [tfb_square\(\)](#), [tfb_tanh\(\)](#), [tfb_transform_diagonal\(\)](#), [tfb_transpose\(\)](#), [tfb_weibull_cdf\(\)](#), [tfb_weibull\(\)](#)

tfb_cumsum

Computes the cumulative sum of a tensor along a specified axis.

Description

Computes the cumulative sum of a tensor along a specified axis.

Usage

```
tfb_cumsum(axis = -1, validate_args = FALSE, name = "cumsum")
```

Arguments

axis	int indicating the axis along which to compute the cumulative sum. Note that positive (and zero) values are not supported
validate_args	Logical, default FALSE. Whether to validate input with asserts. If validate_args is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
name	name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see [tfb_forward\(\)](#), [tfb_inverse\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#).

Other bijectors: [tfb_absolute_value\(\)](#), [tfb_affine_linear_operator\(\)](#), [tfb_affine_scalar\(\)](#), [tfb_affine\(\)](#), [tfb_ascending\(\)](#), [tfb_batch_normalization\(\)](#), [tfb_blockwise\(\)](#), [tfb_chain\(\)](#), [tfb_cholesky_outer_product\(\)](#), [tfb_cholesky_to_inv_cholesky\(\)](#), [tfb_correlation_cholesky\(\)](#), [tfb_discrete_cosine_transform\(\)](#), [tfb_expml\(\)](#), [tfb_exp\(\)](#), [tfb_ffjord\(\)](#), [tfb_fill_scale_tri_l\(\)](#), [tfb_fill_triangular\(\)](#), [tfb_glow\(\)](#), [tfb_gompertz_cdf\(\)](#), [tfb_gumbel_cdf\(\)](#), [tfb_gumbel\(\)](#), [tfb_identity\(\)](#), [tfb_inline\(\)](#), [tfb_invert\(\)](#), [tfb_iterated_sigmoid_centered\(\)](#), [tfb_kumaraswamy_cdf\(\)](#), [tfb_kumaraswamy\(\)](#), [tfb_lambert_w_tail\(\)](#), [tfb_masked_autoregressive_default_template\(\)](#), [tfb_masked_autoregressive_flow\(\)](#), [tfb_masked_dense\(\)](#), [tfb_matrix_inverse_tri_l\(\)](#), [tfb_matvec_lu\(\)](#), [tfb_normal_cdf\(\)](#), [tfb_ordered\(\)](#), [tfb_pad\(\)](#), [tfb_permute\(\)](#), [tfb_power_transform\(\)](#), [tfb_rational_quadratic_spline\(\)](#), [tfb_rayleigh_cdf\(\)](#), [tfb_real_nvp_default_template\(\)](#), [tfb_real_nvp\(\)](#), [tfb_reciprocal\(\)](#), [tfb_reshape\(\)](#), [tfb_scale_matvec_diag\(\)](#), [tfb_scale_matvec_linear_operat](#), [tfb_scale_matvec_lu\(\)](#), [tfb_scale_matvec_tri_l\(\)](#), [tfb_scale_tri_l\(\)](#), [tfb_scale\(\)](#), [tfb_shifted_gompertz_c](#), [tfb_shift\(\)](#), [tfb_sigmoid\(\)](#), [tfb_sinh_arcsinh\(\)](#), [tfb_sinh\(\)](#), [tfb_softmax_centered\(\)](#), [tfb_softplus\(\)](#), [tfb_softsign\(\)](#), [tfb_split\(\)](#), [tfb_square\(\)](#), [tfb_tanh\(\)](#), [tfb_transform_diagonal\(\)](#), [tfb_transpose\(\)](#), [tfb_weibull_cdf\(\)](#), [tfb_weibull\(\)](#)

 tfb_discrete_cosine_transform

Computes $Y = g(X) = \text{DCT}(X)$, where DCT type is indicated by the type arg

Description

The **discrete cosine transform** efficiently applies a unitary DCT operator. This can be useful for mixing and decorrelating across the innermost event dimension. The inverse $X = g^{-1}(Y) = \text{IDCT}(Y)$, where IDCT is DCT-III for `type==2`. This bijector can be interleaved with Affine bijectors to build a cascade of structured efficient linear layers as in Moczulski et al., 2016. Note that the operator applied is orthonormal (i.e. `norm='ortho'`).

Usage

```
tfb_discrete_cosine_transform(
  validate_args = FALSE,
  dct_type = 2,
  name = "dct"
)
```

Arguments

<code>validate_args</code>	Logical, default FALSE. Whether to validate input with asserts. If <code>validate_args</code> is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
<code>dct_type</code>	integer, the DCT type performed by the forward transformation. Currently, only 2 and 3 are supported.
<code>name</code>	name prefixed to Ops created by this class.

Value

a bijector instance.

References

- Moczulski M, Denil M, Appleyard J, de Freitas N. ACDC: A structured efficient linear layer. In *International Conference on Learning Representations*, 2016.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_expml()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`,

[tfb_invert\(\)](#), [tfb_iterated_sigmoid_centered\(\)](#), [tfb_kumaraswamy_cdf\(\)](#), [tfb_kumaraswamy\(\)](#),
[tfb_lambert_w_tail\(\)](#), [tfb_masked_autoregressive_default_template\(\)](#), [tfb_masked_autoregressive_flow\(\)](#),
[tfb_masked_dense\(\)](#), [tfb_matrix_inverse_tri_l\(\)](#), [tfb_matvec_lu\(\)](#), [tfb_normal_cdf\(\)](#),
[tfb_ordered\(\)](#), [tfb_pad\(\)](#), [tfb_permute\(\)](#), [tfb_power_transform\(\)](#), [tfb_rational_quadratic_spline\(\)](#),
[tfb_rayleigh_cdf\(\)](#), [tfb_real_nvp_default_template\(\)](#), [tfb_real_nvp\(\)](#), [tfb_reciprocal\(\)](#),
[tfb_reshape\(\)](#), [tfb_scale_matvec_diag\(\)](#), [tfb_scale_matvec_linear_operator\(\)](#), [tfb_scale_matvec_lu\(\)](#),
[tfb_scale_matvec_tri_l\(\)](#), [tfb_scale_tri_l\(\)](#), [tfb_scale\(\)](#), [tfb_shifted_gompertz_cdf\(\)](#),
[tfb_shift\(\)](#), [tfb_sigmoid\(\)](#), [tfb_sinh_arcsinh\(\)](#), [tfb_sinh\(\)](#), [tfb_softmax_centered\(\)](#),
[tfb_softplus\(\)](#), [tfb_softsign\(\)](#), [tfb_split\(\)](#), [tfb_square\(\)](#), [tfb_tanh\(\)](#), [tfb_transform_diagonal\(\)](#),
[tfb_transpose\(\)](#), [tfb_weibull_cdf\(\)](#), [tfb_weibull\(\)](#)

tfb_exp	<i>Computes</i> $Y=g(X)=\exp(X)$
---------	----------------------------------

Description

Computes $Y=g(X)=\exp(X)$

Usage

```
tfb_exp(validate_args = FALSE, name = "exp")
```

Arguments

validate_args	Logical, default FALSE. Whether to validate input with asserts. If validate_args is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
name	name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see [tfb_forward\(\)](#), [tfb_inverse\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#).

Other bijectors: [tfb_absolute_value\(\)](#), [tfb_affine_linear_operator\(\)](#), [tfb_affine_scalar\(\)](#), [tfb_affine\(\)](#), [tfb_ascending\(\)](#), [tfb_batch_normalization\(\)](#), [tfb_blockwise\(\)](#), [tfb_chain\(\)](#), [tfb_cholesky_outer_product\(\)](#), [tfb_cholesky_to_inv_cholesky\(\)](#), [tfb_correlation_cholesky\(\)](#), [tfb_cumsum\(\)](#), [tfb_discrete_cosine_transform\(\)](#), [tfb_expm1\(\)](#), [tfb_ffjord\(\)](#), [tfb_fill_scale_tri_l\(\)](#), [tfb_fill_triangular\(\)](#), [tfb_glow\(\)](#), [tfb_gompertz_cdf\(\)](#), [tfb_gumbel_cdf\(\)](#), [tfb_gumbel\(\)](#), [tfb_identity\(\)](#), [tfb_inline\(\)](#), [tfb_invert\(\)](#), [tfb_iterated_sigmoid_centered\(\)](#), [tfb_kumaraswamy_cdf\(\)](#), [tfb_kumaraswamy\(\)](#), [tfb_lambert_w_tail\(\)](#), [tfb_masked_autoregressive_default_template\(\)](#), [tfb_masked_autoregressive_flow\(\)](#), [tfb_masked_dense\(\)](#), [tfb_matrix_inverse_tri_l\(\)](#), [tfb_matvec_lu\(\)](#), [tfb_normal_cdf\(\)](#), [tfb_ordered\(\)](#), [tfb_pad\(\)](#), [tfb_permute\(\)](#), [tfb_power_transform\(\)](#), [tfb_rational_quadratic_spline\(\)](#), [tfb_rayleigh_cdf\(\)](#), [tfb_real_nvp_default_template\(\)](#), [tfb_real_nvp\(\)](#), [tfb_reciprocal\(\)](#), [tfb_reshape\(\)](#), [tfb_scale_matvec_diag\(\)](#), [tfb_scale_matvec_linear_operat](#), [tfb_scale_matvec_lu\(\)](#), [tfb_scale_matvec_tri_l\(\)](#), [tfb_scale_tri_l\(\)](#), [tfb_scale\(\)](#), [tfb_shifted_gompertz_cdf\(\)](#)

tfb_shift(), tfb_sigmoid(), tfb_sinh_arcsinh(), tfb_sinh(), tfb_softmax_centered(),
 tfb_softplus(), tfb_softsign(), tfb_split(), tfb_square(), tfb_tanh(), tfb_transform_diagonal(),
 tfb_transpose(), tfb_weibull_cdf(), tfb_weibull()

tfb_expml	<i>Computes</i> $Y = g(X) = \exp(X) - 1$
-----------	--

Description

This Bijector is no different from `tfb_chain(list(tfb_affine_scalar(shift=-1), tfb_exp()))`.
 However, this makes use of the more numerically stable routines `tf$math$expm1` and `tf$log1p`.

Usage

```
tfb_expml(validate_args = FALSE, name = "expml")
```

Arguments

`validate_args` Logical, default FALSE. Whether to validate input with asserts. If `validate_args` is FALSE, and the inputs are invalid, correct behavior is not guaranteed.

`name` name prefixed to Ops created by this class.

Details

Note: the `expml(.)` is applied element-wise but the Jacobian is a reduction over the event space.

Value

a bijector instance.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`,
`tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`,
`tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`,
`tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`,
`tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`,
`tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`,
`tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_template()`,
`tfb_masked_autoregressive_flow()`, `tfb_masked_dense()`, `tfb_matrix_inverse_tri_l()`,
`tfb_matvec_lu()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_pad()`, `tfb_permute()`, `tfb_power_transform()`,
`tfb_rational_quadratic_spline()`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`,
`tfb_real_nvp()`, `tfb_reciprocal()`, `tfb_reshape()`, `tfb_scale_matvec_diag()`, `tfb_scale_matvec_linear_operat`
`tfb_scale_matvec_lu()`, `tfb_scale_matvec_tri_l()`, `tfb_scale_tri_l()`, `tfb_scale()`, `tfb_shifted_gompertz_co`
`tfb_shift()`, `tfb_sigmoid()`, `tfb_sinh_arcsinh()`, `tfb_sinh()`, `tfb_softmax_centered()`,
`tfb_softplus()`, `tfb_softsign()`, `tfb_split()`, `tfb_square()`, `tfb_tanh()`, `tfb_transform_diagonal()`,
`tfb_transpose()`, `tfb_weibull_cdf()`, `tfb_weibull()`

tfb_ffjord	<i>Implements a continuous normalizing flow $X \rightarrow Y$ defined via an ODE.</i>
------------	--

Description

This bijector implements a continuous dynamics transformation parameterized by a differential equation, where initial and terminal conditions correspond to domain (X) and image (Y) i.e.

Usage

```
tfb_ffjord(
  state_time_derivative_fn,
  ode_solve_fn = NULL,
  trace_augmentation_fn = tfp$bijectors$ffjord$trace_jacobian_hutchinson,
  initial_time = 0,
  final_time = 1,
  validate_args = FALSE,
  dtype = tf$float32,
  name = "ffjord"
)
```

Arguments

state_time_derivative_fn	function taking arguments time (a scalar representing time) and state (a Tensor representing the state at given time) returning the time derivative of the state at given time.
ode_solve_fn	function taking arguments ode_fn (same as state_time_derivative_fn above), initial_time (a scalar representing the initial time of integration), initial_state (a Tensor of floating dtype represents the initial state) and solution_times (1D Tensor of floating dtype representing time at which to obtain the solution) returning a Tensor of shape [time_axis, initial_state\$shape]. Will take [final_time] as the solution_times argument and state_time_derivative_fn as ode_fn argument. If NULL a DormandPrince solver from <code>tfp\$math\$ode</code> is used. Default value: NULL
trace_augmentation_fn	function taking arguments ode_fn (function same as state_time_derivative_fn above), state_shape (TensorShape of a the state), dtype (same as dtype of the state) and returning a function taking arguments time (a scalar representing the time at which the function is evaluated), state (a Tensor representing the state at given time) that computes a tuple (ode_fn(time, state), jacobian_trace_estimation). jacobian_trace_estimation should represent trace of the jacobian of ode_fn with respect to state. state_time_derivative_fn will be passed as ode_fn argument. Default value: <code>tfp\$bijectors\$ffjord\$trace_jacobian_hutchinson</code>
initial_time	Scalar float representing time to which the x value of the bijector corresponds to. Passed as initial_time to ode_solve_fn. For default solver can be float or floating scalar Tensor. Default value: 0.

<code>final_time</code>	Scalar float representing time to which the y value of the bijector corresponds to. Passed as <code>solution_times</code> to <code>ode_solve_fn</code> . For default solver can be float or floating scalar Tensor. Default value: 1.
<code>validate_args</code>	Logical, default FALSE. Whether to validate input with asserts. If <code>validate_args</code> is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
<code>dtype</code>	<code>tf\$DType</code> to prefer when converting args to Tensors. Else, we fall back to a common dtype inferred from the args, finally falling back to <code>float32</code> .
<code>name</code>	name prefixed to Ops created by this class.

Details

```
d/dt[state(t)] = state_time_derivative_fn(t, state(t))
state(initial_time) = X
state(final_time) = Y
```

For this transformation the value of `log_det_jacobian` follows another differential equation, reducing it to computation of the trace of the jacobian along the trajectory

```
state_time_derivative = state_time_derivative_fn(t, state(t))
d/dt[log_det_jac(t)] = Tr(jacobian(state_time_derivative, state(t)))
```

FFJORD constructor takes two functions `ode_solve_fn` and `trace_augmentation_fn` arguments that customize integration of the differential equation and trace estimation.

Differential equation integration is performed by a call to `ode_solve_fn`.

Custom `ode_solve_fn` must accept the following arguments:

- `ode_fn(time, state)`: Differential equation to be solved.
- `initial_time`: Scalar float or floating Tensor representing the initial time.
- `initial_state`: Floating Tensor representing the initial state.
- `solution_times`: 1D floating Tensor of solution times.

And return a Tensor of shape `[solution_times$shape, initial_state$shape]` representing state values evaluated at `solution_times`. In addition `ode_solve_fn` must support nested structures. For more details see the interface of `tfp$math$ode$Solver$solve()`.

Trace estimation is computed simultaneously with `state_time_derivative` using `augmented_state_time_derivative_fn` that is generated by `trace_augmentation_fn`. `trace_augmentation_fn` takes `state_time_derivative_fn`, `state.shape` and `state.dtype` arguments and returns a `augmented_state_time_derivative_fn` callable that computes both `state_time_derivative` and unreduced `trace_estimation`.

Custom `ode_solve_fn` and `trace_augmentation_fn` examples:

```
# custom_solver_fn: `function(f, t_initial, t_solutions, y_initial, ...)`
# ... : Additional arguments to pass to custom_solver_fn.
ode_solve_fn <- function(ode_fn, initial_time, initial_state, solution_times) {
  custom_solver_fn(ode_fn, initial_time, solution_times, initial_state, ...)
}
ffjord <- tfb_ffjord(state_time_derivative_fn, ode_solve_fn = ode_solve_fn)
```

```

# state_time_derivative_fn: `function(time, state)`
# trace_jac_fn: `function(time, state)` unreduced jacobian trace function
trace_augmentation_fn <- function(ode_fn, state_shape, state_dtype) {
  augmented_ode_fn <- function(time, state) {
    list(ode_fn(time, state), trace_jac_fn(time, state))
  }
  augmented_ode_fn
}
ffjord <- tfb_ffjord(state_time_derivative_fn, trace_augmentation_fn = trace_augmentation_fn)

```

For more details on FFJORD and continuous normalizing flows see Chen et al. (2018), Grathwol et al. (2018).

Value

a bijector instance.

References

- Chen, T. Q., Rubanova, Y., Bettencourt, J., & Duvenaud, D. K. (2018). Neural ordinary differential equations. In Advances in neural information processing systems (pp. 6571-6583)
- Grathwohl, W., Chen, R. T., Bettencourt, J., Sutskever, I., & Duvenaud, D. (2018). Ffjord: Free-form continuous dynamics for scalable reversible generative models. [arXiv:1810.01367](https://arxiv.org/abs/1810.01367).

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_template()`, `tfb_masked_autoregressive_flow()`, `tfb_masked_dense()`, `tfb_matrix_inverse_tri_l()`, `tfb_matvec_lu()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_pad()`, `tfb_permute()`, `tfb_power_transform()`, `tfb_rational_quadratic_spline()`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`, `tfb_real_nvp()`, `tfb_reciprocal()`, `tfb_reshape()`, `tfb_scale_matvec_diag()`, `tfb_scale_matvec_linear_operator()`, `tfb_scale_matvec_lu()`, `tfb_scale_matvec_tri_l()`, `tfb_scale_tri_l()`, `tfb_scale()`, `tfb_shifted_gompertz_cdf()`, `tfb_shift()`, `tfb_sigmoid()`, `tfb_sinh_arcsinh()`, `tfb_sinh()`, `tfb_softmax_centered()`, `tfb_softplus()`, `tfb_softsign()`, `tfb_split()`, `tfb_square()`, `tfb_tanh()`, `tfb_transform_diagonal()`, `tfb_transpose()`, `tfb_weibull_cdf()`, `tfb_weibull()`

tfb_fill_scale_tri_l *Transforms unconstrained vectors to TriL matrices with positive diagonal*

Description

This is implemented as a simple `tfb_chain` of `tfb_fill_triangular` followed by `tfb_transform_diagonal`, and provided mostly as a convenience. The default setup is somewhat opinionated, using a Softplus transformation followed by a small shift ($1e-5$) which attempts to avoid numerical issues from zeros on the diagonal.

Usage

```
tfb_fill_scale_tri_l(
  diag_bijector = NULL,
  diag_shift = 1e-05,
  validate_args = FALSE,
  name = "fill_scale_tril"
)
```

Arguments

<code>diag_bijector</code>	Bijector instance, used to transform the output diagonal to be positive. Default value: <code>NULL</code> (i.e., <code>tfb_softplus()</code>).
<code>diag_shift</code>	Float value broadcastable and added to all diagonal entries after applying the <code>diag_bijector</code> . Setting a positive value forces the output diagonal entries to be positive, but prevents inverting the transformation for matrices with diagonal entries less than this value. Default value: $1e-5$.
<code>validate_args</code>	Logical, default <code>FALSE</code> . Whether to validate input with asserts. If <code>validate_args</code> is <code>FALSE</code> , and the inputs are invalid, correct behavior is not guaranteed.
<code>name</code>	name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expml()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_template()`,

tfb_masked_autoregressive_flow(), tfb_masked_dense(), tfb_matrix_inverse_tri_l(),
 tfb_matvec_lu(), tfb_normal_cdf(), tfb_ordered(), tfb_pad(), tfb_permute(), tfb_power_transform(),
 tfb_rational_quadratic_spline(), tfb_rayleigh_cdf(), tfb_real_nvp_default_template(),
 tfb_real_nvp(), tfb_reciprocal(), tfb_reshape(), tfb_scale_matvec_diag(), tfb_scale_matvec_linear_operat
 tfb_scale_matvec_lu(), tfb_scale_matvec_tri_l(), tfb_scale_tri_l(), tfb_scale(), tfb_shifted_gompertz_c
 tfb_shift(), tfb_sigmoid(), tfb_sinh_arcsinh(), tfb_sinh(), tfb_softmax_centered(),
 tfb_softplus(), tfb_softsign(), tfb_split(), tfb_square(), tfb_tanh(), tfb_transform_diagonal(),
 tfb_transpose(), tfb_weibull_cdf(), tfb_weibull()

tfb_fill_triangular *Transforms vectors to triangular*

Description

Triangular matrix elements are filled in a clockwise spiral. Given input with shape batch_shape + [d], produces output with shape batch_shape + [n, n], where $n = (-1 + \sqrt{1 + 8 * d})/2$. This follows by solving the quadratic equation $d = 1 + 2 + \dots + n = n * (n + 1)/2$.

Usage

```
tfb_fill_triangular(  
    upper = FALSE,  
    validate_args = FALSE,  
    name = "fill_triangular"  
)
```

Arguments

upper	Logical representing whether output matrix should be upper triangular (TRUE) or lower triangular (FALSE, default).
validate_args	Logical, default FALSE. Whether to validate input with asserts. If validate_args is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
name	name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see [tfb_forward\(\)](#), [tfb_inverse\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#).

Other bijectors: [tfb_absolute_value\(\)](#), [tfb_affine_linear_operator\(\)](#), [tfb_affine_scalar\(\)](#), [tfb_affine\(\)](#), [tfb_ascending\(\)](#), [tfb_batch_normalization\(\)](#), [tfb_blockwise\(\)](#), [tfb_chain\(\)](#), [tfb_cholesky_outer_product\(\)](#), [tfb_cholesky_to_inv_cholesky\(\)](#), [tfb_correlation_cholesky\(\)](#), [tfb_cumsum\(\)](#), [tfb_discrete_cosine_transform\(\)](#), [tfb_expm1\(\)](#), [tfb_exp\(\)](#), [tfb_ffjord\(\)](#), [tfb_fill_scale_tri_l\(\)](#), [tfb_glow\(\)](#), [tfb_gompertz_cdf\(\)](#), [tfb_gumbel_cdf\(\)](#), [tfb_gumbel\(\)](#), [tfb_identity\(\)](#), [tfb_inline\(\)](#), [tfb_invert\(\)](#), [tfb_iterated_sigmoid_centered\(\)](#), [tfb_kumaraswamy_cdf\(\)](#),

[tfb_kumaraswamy\(\)](#), [tfb_lambert_w_tail\(\)](#), [tfb_masked_autoregressive_default_template\(\)](#),
[tfb_masked_autoregressive_flow\(\)](#), [tfb_masked_dense\(\)](#), [tfb_matrix_inverse_tri_l\(\)](#),
[tfb_matvec_lu\(\)](#), [tfb_normal_cdf\(\)](#), [tfb_ordered\(\)](#), [tfb_pad\(\)](#), [tfb_permute\(\)](#), [tfb_power_transform\(\)](#),
[tfb_rational_quadratic_spline\(\)](#), [tfb_rayleigh_cdf\(\)](#), [tfb_real_nvp_default_template\(\)](#),
[tfb_real_nvp\(\)](#), [tfb_reciprocal\(\)](#), [tfb_reshape\(\)](#), [tfb_scale_matvec_diag\(\)](#), [tfb_scale_matvec_linear_operat](#)
[tfb_scale_matvec_lu\(\)](#), [tfb_scale_matvec_tri_l\(\)](#), [tfb_scale_tri_l\(\)](#), [tfb_scale\(\)](#), [tfb_shifted_gompertz_c](#)
[tfb_shift\(\)](#), [tfb_sigmoid\(\)](#), [tfb_sinh_arcsinh\(\)](#), [tfb_sinh\(\)](#), [tfb_softmax_centered\(\)](#),
[tfb_softplus\(\)](#), [tfb_softsign\(\)](#), [tfb_split\(\)](#), [tfb_square\(\)](#), [tfb_tanh\(\)](#), [tfb_transform_diagonal\(\)](#),
[tfb_transpose\(\)](#), [tfb_weibull_cdf\(\)](#), [tfb_weibull\(\)](#)

tfb_forward	Returns the forward Bijector evaluation, i.e., $X = g(Y)$.
-------------	---

Description

Returns the forward Bijector evaluation, i.e., $X = g(Y)$.

Usage

```
tfb_forward(bijector, x, name = "forward")
```

Arguments

bijector	The bijector to apply
x	Tensor. The input to the "forward" evaluation.
name	name of the operation

Value

a tensor

See Also

Other bijector_methods: [tfb_forward_log_det_jacobian\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#),
[tfb_inverse\(\)](#)

Examples

```

b <- tfb_affine_scalar(shift = 1, scale = 2)
x <- 10
b %>% tfb_forward(x)

```

`tfb_forward_log_det_jacobian`

Returns the result of the forward evaluation of the log determinant of the Jacobian

Description

Returns the result of the forward evaluation of the log determinant of the Jacobian

Usage

```
tfb_forward_log_det_jacobian(  
  bijector,  
  x,  
  event_ndims,  
  name = "forward_log_det_jacobian"  
)
```

Arguments

<code>bijector</code>	The bijector to apply
<code>x</code>	Tensor. The input to the "forward" Jacobian determinant evaluation.
<code>event_ndims</code>	Number of dimensions in the probabilistic events being transformed. Must be greater than or equal to <code>bijector\$forward_min_event_ndims</code> . The result is summed over the final dimensions to produce a scalar Jacobian determinant for each event, i.e. it has shape <code>x\$shape\$ndims - event_ndims</code> dimensions.
<code>name</code>	name of the operation

Value

a tensor

See Also

Other `bijector_methods`: [tfb_forward\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#), [tfb_inverse\(\)](#)

Examples

```
b <- tfb_affine_scalar(shift = 1, scale = 2)  
x <- 10  
b %>% tfb_forward_log_det_jacobian(x, event_ndims = 0)
```

tfb_glow

*Implements the Glow Bijector from Kingma & Dhariwal (2018).***Description**

Overview: Glow is a chain of bijectors which transforms a rank-1 tensor (vector) into a rank-3 tensor (e.g. an RGB image). Glow does this by chaining together an alternating series of "Blocks," "Squeezes," and "Exits" which are each themselves special chains of other bijectors. The intended use of Glow is as part of a `tfd_transformed_distribution`, in which the base distribution over the vector space is used to generate samples in the image space. In the paper, an Independent Normal distribution is used as the base distribution.

Usage

```
tfb_glow(
  output_shape = c(32, 32, 3),
  num_glow_blocks = 3,
  num_steps_per_block = 32,
  coupling_bijector_fn = NULL,
  exit_bijector_fn = NULL,
  grab_after_block = NULL,
  use_actnorm = TRUE,
  seed = NULL,
  validate_args = FALSE,
  name = "glow"
)
```

Arguments

`output_shape` A list of integers, specifying the event shape of the output, of the bijectors forward pass (the image). Specified as [H, W, C]. Default Value: (32, 32, 3)

`num_glow_blocks` An integer, specifying how many downsampling levels to include in the model. This must divide equally into both H and W, otherwise the bijector would not be invertible. Default Value: 3

`num_steps_per_block` An integer specifying how many Affine Coupling and 1x1 convolution layers to include at each level of the spatial hierarchy. Default Value: 32 (i.e. the value used in the original glow paper).

`coupling_bijector_fn` A function which takes the argument `input_shape` and returns a callable neural network (e.g. a `keras_model_sequential()`). The network should either return a tensor with the same event shape as `input_shape` (this will employ additive coupling), a tensor with the same height and width as `input_shape` but twice the number of channels (this will employ affine coupling), or a bijector which takes in a tensor with event shape `input_shape`, and returns a tensor with shape `input_shape`.

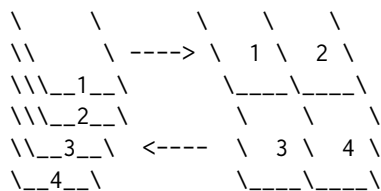
exit_bijector_fn	Similar to coupling_bijector_fn, exit_bijector_fn is a function which takes the argument input_shape and output_chan and returns a callable neural network. The neural network it returns should take a tensor of shape input_shape as the input, and return one of three options: A tensor with output_chan channels, a tensor with 2 * output_chan channels, or a bijector. Additional details can be found in the documentation for ExitBijector.
grab_after_block	A tuple of floats, specifying what fraction of the remaining channels to remove following each glow block. Glow will take the integer floor of this number multiplied by the remaining number of channels. The default is half at each spatial hierarchy. Default value: None (this will take out half of the channels after each block).
use_actnorm	A boolean deciding whether or not to use actnorm. Data-dependent initialization is used to initialize this layer. Default value: FALSE
seed	A seed to control randomness in the 1x1 convolution initialization. Default value: NULL (i.e., non-reproducible sampling).
validate_args	Logical, default FALSE. Whether to validate input with asserts. If validate_args is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
name	name prefixed to Ops created by this class.

Details

A "Block" (implemented as the GlowBlock Bijector) performs much of the transformations which allow glow to produce sophisticated and complex mappings between the image space and the latent space and therefore achieve rich image generation performance. A Block is composed of num_steps_per_block steps, which are each implemented as a Chain containing an ActivationNormalization (ActNorm) bijector, followed by an (invertible) OneByOneConv bijector, and finally a coupling bijector. The coupling bijector is an instance of a RealNVP bijector, and uses the coupling_bijector_fn function to instantiate the coupling bijector function which is given to the RealNVP. This function returns a bijector which defines the coupling (e.g. Shift(Scale) for affine coupling or Shift for additive coupling).

A "Squeeze" converts spatial features into channel features. It is implemented using the Expand bijector. The difference in names is due to the fact that the forward function from glow is meant to ultimately correspond to sampling from a tfp\$util\$TransformedDistribution object, which would use Expand (Squeeze is just Invert(Expand)). The Expand bijector takes a tensor with shape [H, W, C] and returns a tensor with shape [2H, 2W, C / 4], such that each 2x2x1 spatial tile in the output is composed from a single 1x1x4 tile in the input tensor, as depicted in the figure below.

Forward pass (Expand)



Inverse pass (Squeeze) This is implemented using a chain of Reshape -> Transpose -> Reshape bijectors. Note that on an inverse pass through the bijector, each Squeeze will cause the width/height of the image to decrease by a factor of 2. Therefore, the input image must be evenly divisible by 2 at least `num_glow_blocks` times, since it will pass through a Squeeze step that many times.

An "Exit" is simply a junction at which some of the tensor "exits" from the glow bijector and therefore avoids any further alteration. Each exit is implemented as a Blockwise bijector, where some channels are given to the rest of the glow model, and the rest are given to a bypass implemented using the Identity bijector. The fraction of channels to be removed at each exit is determined by the `grab_after_block` arg, indicates the fraction of remaining channels which join the identity bypass. The fraction is converted to an integer number of channels by multiplying by the remaining number of channels and rounding. Additionally, at each exit, glow couples the tensor exiting the highway to the tensor continuing onward. This makes small scale features in the image dependent on larger scale features, since the larger scale features dictate the mean and scale of the distribution over the smaller scale features. This coupling is done similarly to the Coupling bijector in each step of the flow (i.e. using a RealNVP bijector). However for the exit bijector, the coupling is instantiated using `exit_bijector_fn` rather than `coupling_bijector_fn`, allowing for different behaviors between standard coupling and exit coupling. Also note that because the exit utilizes a coupling bijector, there are two special cases (all channels exiting and no channels exiting). The full Glow bijector consists of `num_glow_blocks` Blocks each of which contains `num_steps_per_block` steps. Each step implements a coupling using `bijector_coupling_fn`. Between blocks, glow converts between spatial pixels and channels using the Expand Bijector, and splits channels out of the bijector using the Exit Bijector. The channels which have exited continue onward through Identity bijectors and those which have not exited are given to the next block. After passing through all Blocks, the tensor is reshaped to a rank-1 tensor with the same number of elements. This is where the distribution will be defined. A schematic diagram of Glow is shown below. The forward function of the bijector starts from the bottom and goes upward, while the inverse function starts from the top and proceeds downward.

Value

a bijector instance.

#' ""

```
Glow Schematic Diagram Input Image ##### shape = [H, W, C] \ / <- Expand
Bijector turns spatial \ / dimensions into channels. | XXXXXXXXXXXXXXXXXXXXX |
XXXXXXXXXXXXXXXXXXXXXXXXXX | XXXXXXXXXXXXXXXXXXXXXXXXXXXXX A single step of the flow
consists Glow Block - | XXXXXXXXXXXXXXXXXXXXXXXXXXXXX <- of ActNorm -> 1x1Conv -> Cou-
pling. | XXXXXXXXXXXXXXXXXXXXXXXXXXXXX there are num_steps_per_block | XXXXXXXXXXXXXXXXXXXXXXXXXXXXX
steps of the flow in each block. |_ XXXXXXXXXXXXXXXXXXXXXXXXXXXXX \ / <- Expand bijectors fol-
low each glow \ / block XXXXXXXXXX\ \ \ <- Exit Bijector removes channels _ _ from additional
alteration. | XXXXXXXXXX !!! | XXXXXXXXXX !!! | XXXXXXXXXX !!! After exiting, channels are
passed Glow Block - | XXXXXXXXXX !!! <- downward using the Blockwise and | XXXXXXXXXX
!!! Identify bijectors. | XXXXXXXXXX !!! |_ XXXXXXXXXX !!! \ / <- Expand Bijector \ / XXX\
! <- Exit Bijector _ | XXX !!! | XXX !!! | XXX !!! low Block - | XXX !!! | XXX !!!
! | XXX !!! |_ XXX !!! | XX \ !!! <- (Optional) Exit Bijector ||| v v v Output Distribution
##### shape = [H * W * C]
```

Legend

! **XX** = Step of flow || **X** = Exit bijector || **V** = Expand bijector || **!!!** = Identity bijector || **up** = Forward pass || **dn** = Inverse pass || _____ |

```
[H, W, C]: R:H,%20W,%20C
[2H, 2W, C / 4]: R:2H,%202W,%20C%20/%204
[H, W, C]: R:H,%20W,%20C
[H * W * C]: R:H%20*%20W%20*%20C
```

References

- Diederik P Kingma, Prafulla Dhariwal, Glow: Generative Flow with Invertible 1x1 Convolutions. In *Neural Information Processing Systems*, 2018.
- Laurent Dinh, Jascha Sohl-Dickstein, and Samy Bengio. Density Estimation using Real NVP. In *International Conference on Learning Representations*, 2017.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_template()`, `tfb_masked_autoregressive_flow()`, `tfb_masked_dense()`, `tfb_matrix_inverse_tri_l()`, `tfb_matvec_lu()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_pad()`, `tfb_permute()`, `tfb_power_transform()`, `tfb_rational_quadratic_spline()`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`, `tfb_real_nvp()`, `tfb_reciprocal()`, `tfb_reshape()`, `tfb_scale_matvec_diag()`, `tfb_scale_matvec_linear_operator()`, `tfb_scale_matvec_lu()`, `tfb_scale_matvec_tri_l()`, `tfb_scale_tri_l()`, `tfb_scale()`, `tfb_shifted_gompertz_cdf()`, `tfb_shift()`, `tfb_sigmoid()`, `tfb_sinh_arcsinh()`, `tfb_sinh()`, `tfb_softmax_centered()`, `tfb_softplus()`, `tfb_softsign()`, `tfb_split()`, `tfb_square()`, `tfb_tanh()`, `tfb_transform_diagonal()`, `tfb_transpose()`, `tfb_weibull_cdf()`, `tfb_weibull()`

tfb_gompertz_cdf	Compute $Y = g(X) = 1 - \exp(-c * (\exp(\text{rate} * X) - 1))$, the Gompertz CDF.
------------------	---

Description

This bijector maps inputs from $[-\infty, \infty]$ to $[0, \infty]$. The inverse of the bijector applied to a uniform random variable $X \sim U(0, 1)$ gives back a random variable with the **Gompertz distribution**:

$$Y \sim \text{GompertzCDF}(\text{concentration}, \text{rate})$$

$$\text{pdf}(y; c, r) = r * c * \exp(r * y + c - c * \exp(-c * \exp(r * y)))$$

Note: Because the Gompertz distribution concentrates its mass close to zero, for larger rates or larger concentrations, `bijector.forward` will quickly saturate to 1.

Usage

```
tfb_gompertz_cdf(
    concentration,
    rate,
    validate_args = FALSE,
    name = "gompertz_cdf"
)
```

Arguments

concentration	Positive Float-like Tensor that is the same dtype and is broadcastable with concentration. This is c in $Y = g(X) = 1 - \exp(-c * (\exp(\text{rate} * X) - 1))$.
rate	Positive Float-like Tensor that is the same dtype and is broadcastable with concentration. This is rate in $Y = g(X) = 1 - \exp(-c * (\exp(\text{rate} * X) - 1))$.
validate_args	Logical, default FALSE. Whether to validate input with asserts. If <code>validate_args</code> is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
name	name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see [tfb_forward\(\)](#), [tfb_inverse\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#).

Other bijectors: [tfb_absolute_value\(\)](#), [tfb_affine_linear_operator\(\)](#), [tfb_affine_scalar\(\)](#), [tfb_affine\(\)](#), [tfb_ascending\(\)](#), [tfb_batch_normalization\(\)](#), [tfb_blockwise\(\)](#), [tfb_chain\(\)](#), [tfb_cholesky_outer_product\(\)](#), [tfb_cholesky_to_inv_cholesky\(\)](#), [tfb_correlation_cholesky\(\)](#), [tfb_cumsum\(\)](#), [tfb_discrete_cosine_transform\(\)](#), [tfb_expm1\(\)](#), [tfb_exp\(\)](#), [tfb_ffjord\(\)](#), [tfb_fill_scale_tri_l\(\)](#), [tfb_fill_triangular\(\)](#), [tfb_glow\(\)](#), [tfb_gumbel_cdf\(\)](#), [tfb_gumbel\(\)](#), [tfb_identity\(\)](#), [tfb_inline\(\)](#), [tfb_invert\(\)](#), [tfb_iterated_sigmoid_centered\(\)](#), [tfb_kumaraswamy_cdf\(\)](#), [tfb_kumaraswamy\(\)](#), [tfb_lambert_w_tail\(\)](#), [tfb_masked_autoregressive_default_template\(\)](#), [tfb_masked_autoregressive_flow\(\)](#), [tfb_masked_dense\(\)](#), [tfb_matrix_inverse_tri_l\(\)](#), [tfb_matvec_lu\(\)](#), [tfb_normal_cdf\(\)](#), [tfb_ordered\(\)](#), [tfb_pad\(\)](#), [tfb_permute\(\)](#), [tfb_power_transform\(\)](#), [tfb_rational_quadratic_spline\(\)](#), [tfb_rayleigh_cdf\(\)](#), [tfb_real_nvp_default_template\(\)](#), [tfb_real_nvp\(\)](#), [tfb_reciprocal\(\)](#), [tfb_reshape\(\)](#), [tfb_scale_matvec_diag\(\)](#), [tfb_scale_matvec_linear_operator\(\)](#), [tfb_scale_matvec_lu\(\)](#), [tfb_scale_matvec_tri_l\(\)](#), [tfb_scale_tri_l\(\)](#), [tfb_scale\(\)](#), [tfb_shifted_gompertz_cdf\(\)](#), [tfb_shift\(\)](#), [tfb_sigmoid\(\)](#), [tfb_sinh_arcsinh\(\)](#), [tfb_sinh\(\)](#), [tfb_softmax_centered\(\)](#), [tfb_softplus\(\)](#), [tfb_softsign\(\)](#), [tfb_split\(\)](#), [tfb_square\(\)](#), [tfb_tanh\(\)](#), [tfb_transform_diagonal\(\)](#), [tfb_transpose\(\)](#), [tfb_weibull_cdf\(\)](#), [tfb_weibull\(\)](#)

tfb_gumbel	<i>Computes</i> $Y = g(X) = \exp(-\exp(-(X - \text{loc}) / \text{scale}))$
------------	--

Description

This bijector maps inputs from $[-\text{inf}, \text{inf}]$ to $[0, 1]$. The inverse of the bijector applied to a uniform random variable $X \sim U(0, 1)$ gives back a random variable with the [Gumbel distribution](#):

Usage

```
tfb_gumbel(loc = 0, scale = 1, validate_args = FALSE, name = "gumbel")
```

Arguments

loc	Float-like Tensor that is the same dtype and is broadcastable with scale. This is loc in $Y = g(X) = \exp(-\exp(-(X - \text{loc}) / \text{scale}))$.
scale	Positive Float-like Tensor that is the same dtype and is broadcastable with loc. This is scale in $Y = g(X) = \exp(-\exp(-(X - \text{loc}) / \text{scale}))$.
validate_args	Logical, default FALSE. Whether to validate input with asserts. If validate_args is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
name	name prefixed to Ops created by this class.

Details

$Y \sim \text{Gumbel}(\text{loc}, \text{scale})$ pdf(y ; loc, scale) = $\exp(-((y - \text{loc}) / \text{scale} + \exp(-(y - \text{loc}) / \text{scale}))) / \text{scale}$

Value

a bijector instance.

See Also

For usage examples see [tfb_forward\(\)](#), [tfb_inverse\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#).

Other bijectors: [tfb_absolute_value\(\)](#), [tfb_affine_linear_operator\(\)](#), [tfb_affine_scalar\(\)](#), [tfb_affine\(\)](#), [tfb_ascending\(\)](#), [tfb_batch_normalization\(\)](#), [tfb_blockwise\(\)](#), [tfb_chain\(\)](#), [tfb_cholesky_outer_product\(\)](#), [tfb_cholesky_to_inv_cholesky\(\)](#), [tfb_correlation_cholesky\(\)](#), [tfb_cumsum\(\)](#), [tfb_discrete_cosine_transform\(\)](#), [tfb_expm1\(\)](#), [tfb_exp\(\)](#), [tfb_ffjord\(\)](#), [tfb_fill_scale_tri_l\(\)](#), [tfb_fill_triangular\(\)](#), [tfb_glow\(\)](#), [tfb_gompertz_cdf\(\)](#), [tfb_gumbel_cdf\(\)](#), [tfb_identity\(\)](#), [tfb_inline\(\)](#), [tfb_invert\(\)](#), [tfb_iterated_sigmoid_centered\(\)](#), [tfb_kumaraswamy_cdf\(\)](#), [tfb_kumaraswamy\(\)](#), [tfb_lambert_w_tail\(\)](#), [tfb_masked_autoregressive_default_template\(\)](#), [tfb_masked_autoregressive_flow\(\)](#), [tfb_masked_dense\(\)](#), [tfb_matrix_inverse_tri_l\(\)](#), [tfb_matvec_lu\(\)](#), [tfb_normal_cdf\(\)](#), [tfb_ordered\(\)](#), [tfb_pad\(\)](#), [tfb_permute\(\)](#), [tfb_power_transform\(\)](#), [tfb_rational_quadratic_spline\(\)](#), [tfb_rayleigh_cdf\(\)](#), [tfb_real_nvp_default_template\(\)](#), [tfb_real_nvp\(\)](#), [tfb_reciprocal\(\)](#), [tfb_reshape\(\)](#), [tfb_scale_matvec_diag\(\)](#), [tfb_scale_matvec_linear_operat](#), [tfb_scale_matvec_lu\(\)](#), [tfb_scale_matvec_tri_l\(\)](#), [tfb_scale_tri_l\(\)](#), [tfb_scale\(\)](#), [tfb_shifted_gompertz_co](#), [tfb_shift\(\)](#), [tfb_sigmoid\(\)](#), [tfb_sinh_arcsinh\(\)](#), [tfb_sinh\(\)](#), [tfb_softmax_centered\(\)](#),

[tfb_softplus\(\)](#), [tfb_softsign\(\)](#), [tfb_split\(\)](#), [tfb_square\(\)](#), [tfb_tanh\(\)](#), [tfb_transform_diagonal\(\)](#), [tfb_transpose\(\)](#), [tfb_weibull_cdf\(\)](#), [tfb_weibull\(\)](#)

tfb_gumbel_cdf	<i>Compute $Y = g(X) = \exp(-\exp(-(X - \text{loc}) / \text{scale}))$, the Gumbel CDF.</i>
----------------	---

Description

This bijector maps inputs from $[-\infty, \infty]$ to $[0, 1]$. The inverse of the bijector applied to a uniform random variable $X \sim U(0, 1)$ gives back a random variable with the **Gumbel distribution**:

Usage

```
tfb_gumbel_cdf(loc = 0, scale = 1, validate_args = FALSE, name = "gumbel_cdf")
```

Arguments

loc	Float-like Tensor that is the same dtype and is broadcastable with scale. This is loc in $Y = g(X) = \exp(-\exp(-(X - \text{loc}) / \text{scale}))$.
scale	Positive Float-like Tensor that is the same dtype and is broadcastable with loc. This is scale in $Y = g(X) = \exp(-\exp(-(X - \text{loc}) / \text{scale}))$.
validate_args	Logical, default FALSE. Whether to validate input with asserts. If validate_args is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
name	name prefixed to Ops created by this class.

Details

$Y \sim \text{GumbelCDF}(\text{loc}, \text{scale})$
 $\text{pdf}(y; \text{loc}, \text{scale}) = \exp(-((y - \text{loc}) / \text{scale} + \exp(-(y - \text{loc}) / \text{scale}))) / \text{scale}$

Value

a bijector instance.

See Also

For usage examples see [tfb_forward\(\)](#), [tfb_inverse\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#).

Other bijectors: [tfb_absolute_value\(\)](#), [tfb_affine_linear_operator\(\)](#), [tfb_affine_scalar\(\)](#), [tfb_affine\(\)](#), [tfb_ascending\(\)](#), [tfb_batch_normalization\(\)](#), [tfb_blockwise\(\)](#), [tfb_chain\(\)](#), [tfb_cholesky_outer_product\(\)](#), [tfb_cholesky_to_inv_cholesky\(\)](#), [tfb_correlation_cholesky\(\)](#), [tfb_cumsum\(\)](#), [tfb_discrete_cosine_transform\(\)](#), [tfb_expm1\(\)](#), [tfb_exp\(\)](#), [tfb_ffjord\(\)](#), [tfb_fill_scale_tri_l\(\)](#), [tfb_fill_triangular\(\)](#), [tfb_glow\(\)](#), [tfb_gompertz_cdf\(\)](#), [tfb_gumbel\(\)](#), [tfb_identity\(\)](#), [tfb_inline\(\)](#), [tfb_invert\(\)](#), [tfb_iterated_sigmoid_centered\(\)](#), [tfb_kumaraswamy_cdf\(\)](#), [tfb_kumaraswamy\(\)](#), [tfb_lambert_w_tail\(\)](#), [tfb_masked_autoregressive_default_template\(\)](#), [tfb_masked_autoregressive_flow\(\)](#), [tfb_masked_dense\(\)](#), [tfb_matrix_inverse_tri_l\(\)](#), [tfb_matvec_lu\(\)](#), [tfb_normal_cdf\(\)](#), [tfb_ordered\(\)](#), [tfb_pad\(\)](#), [tfb_permute\(\)](#), [tfb_power_transform\(\)](#),

tfb_rational_quadratic_spline(), tfb_rayleigh_cdf(), tfb_real_nvp_default_template(),
 tfb_real_nvp(), tfb_reciprocal(), tfb_reshape(), tfb_scale_matvec_diag(), tfb_scale_matvec_linear_operat
 tfb_scale_matvec_lu(), tfb_scale_matvec_tri_l(), tfb_scale_tri_l(), tfb_scale(), tfb_shifted_gompertz_c
 tfb_shift(), tfb_sigmoid(), tfb_sinh_arcsinh(), tfb_sinh(), tfb_softmax_centered(),
 tfb_softplus(), tfb_softsign(), tfb_split(), tfb_square(), tfb_tanh(), tfb_transform_diagonal(),
 tfb_transpose(), tfb_weibull_cdf(), tfb_weibull()

tfb_identity	<i>Computes</i> $Y = g(X) = X$
--------------	--------------------------------

Description

Computes $Y = g(X) = X$

Usage

```
tfb_identity(validate_args = FALSE, name = "identity")
```

Arguments

`validate_args` Logical, default FALSE. Whether to validate input with asserts. If `validate_args` is FALSE, and the inputs are invalid, correct behavior is not guaranteed.

`name` name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see [tfb_forward\(\)](#), [tfb_inverse\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#).

Other bijectors: [tfb_absolute_value\(\)](#), [tfb_affine_linear_operator\(\)](#), [tfb_affine_scalar\(\)](#), [tfb_affine\(\)](#), [tfb_ascending\(\)](#), [tfb_batch_normalization\(\)](#), [tfb_blockwise\(\)](#), [tfb_chain\(\)](#), [tfb_cholesky_outer_product\(\)](#), [tfb_cholesky_to_inv_cholesky\(\)](#), [tfb_correlation_cholesky\(\)](#), [tfb_cumsum\(\)](#), [tfb_discrete_cosine_transform\(\)](#), [tfb_expm1\(\)](#), [tfb_exp\(\)](#), [tfb_ffjord\(\)](#), [tfb_fill_scale_tri_l\(\)](#), [tfb_fill_triangular\(\)](#), [tfb_glow\(\)](#), [tfb_gompertz_cdf\(\)](#), [tfb_gumbel_cdf\(\)](#), [tfb_gumbel\(\)](#), [tfb_inline\(\)](#), [tfb_invert\(\)](#), [tfb_iterated_sigmoid_centered\(\)](#), [tfb_kumaraswamy_cdf\(\)](#), [tfb_kumaraswamy\(\)](#), [tfb_lambert_w_tail\(\)](#), [tfb_masked_autoregressive_default_template\(\)](#), [tfb_masked_autoregressive_flow\(\)](#), [tfb_masked_dense\(\)](#), [tfb_matrix_inverse_tri_l\(\)](#), [tfb_matvec_lu\(\)](#), [tfb_normal_cdf\(\)](#), [tfb_ordered\(\)](#), [tfb_pad\(\)](#), [tfb_permute\(\)](#), [tfb_power_transform\(\)](#), [tfb_rational_quadratic_spline\(\)](#), [tfb_rayleigh_cdf\(\)](#), [tfb_real_nvp_default_template\(\)](#), [tfb_real_nvp\(\)](#), [tfb_reciprocal\(\)](#), [tfb_reshape\(\)](#), [tfb_scale_matvec_diag\(\)](#), [tfb_scale_matvec_linear_operat](#), [tfb_scale_matvec_lu\(\)](#), [tfb_scale_matvec_tri_l\(\)](#), [tfb_scale_tri_l\(\)](#), [tfb_scale\(\)](#), [tfb_shifted_gompertz_c](#), [tfb_shift\(\)](#), [tfb_sigmoid\(\)](#), [tfb_sinh_arcsinh\(\)](#), [tfb_sinh\(\)](#), [tfb_softmax_centered\(\)](#), [tfb_softplus\(\)](#), [tfb_softsign\(\)](#), [tfb_split\(\)](#), [tfb_square\(\)](#), [tfb_tanh\(\)](#), [tfb_transform_diagonal\(\)](#), [tfb_transpose\(\)](#), [tfb_weibull_cdf\(\)](#), [tfb_weibull\(\)](#)

 tfb_inline

Bijector constructed from custom functions

Description

Bijector constructed from custom functions

Usage

```
tfb_inline(
  forward_fn = NULL,
  inverse_fn = NULL,
  inverse_log_det_jacobian_fn = NULL,
  forward_log_det_jacobian_fn = NULL,
  forward_event_shape_fn = NULL,
  forward_event_shape_tensor_fn = NULL,
  inverse_event_shape_fn = NULL,
  inverse_event_shape_tensor_fn = NULL,
  is_constant_jacobian = NULL,
  validate_args = FALSE,
  forward_min_event_ndims = NULL,
  inverse_min_event_ndims = NULL,
  name = "inline"
)
```

Arguments

`forward_fn` Function implementing the forward transformation.

`inverse_fn` Function implementing the inverse transformation.

`inverse_log_det_jacobian_fn`
 Function implementing the `log_det_jacobian` of the forward transformation.

`forward_log_det_jacobian_fn`
 Function implementing the `log_det_jacobian` of the inverse transformation.

`forward_event_shape_fn`
 Function implementing non-identical static event shape changes. Default: shape is assumed unchanged.

`forward_event_shape_tensor_fn`
 Function implementing non-identical event shape changes. Default: shape is assumed unchanged.

`inverse_event_shape_fn`
 Function implementing non-identical static event shape changes. Default: shape is assumed unchanged.

`inverse_event_shape_tensor_fn`
 Function implementing non-identical event shape changes. Default: shape is assumed unchanged.

<code>is_constant_jacobian</code>	Logical indicating that the Jacobian is constant for all input arguments.
<code>validate_args</code>	Logical, default FALSE. Whether to validate input with asserts. If <code>validate_args</code> is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
<code>forward_min_event_ndims</code>	Integer indicating the minimal dimensionality this bijector acts on.
<code>inverse_min_event_ndims</code>	Integer indicating the minimal dimensionality this bijector acts on.
<code>name</code>	name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_template()`, `tfb_masked_autoregressive_flow()`, `tfb_masked_dense()`, `tfb_matrix_inverse_tri_l()`, `tfb_matvec_lu()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_pad()`, `tfb_permute()`, `tfb_power_transform()`, `tfb_rational_quadratic_spline()`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`, `tfb_real_nvp()`, `tfb_reciprocal()`, `tfb_reshape()`, `tfb_scale_matvec_diag()`, `tfb_scale_matvec_linear_operator()`, `tfb_scale_matvec_lu()`, `tfb_scale_matvec_tri_l()`, `tfb_scale_tri_l()`, `tfb_scale()`, `tfb_shifted_gompertz_cdf()`, `tfb_shift()`, `tfb_sigmoid()`, `tfb_sinh_arcsinh()`, `tfb_sinh()`, `tfb_softmax_centered()`, `tfb_softplus()`, `tfb_softsign()`, `tfb_split()`, `tfb_square()`, `tfb_tanh()`, `tfb_transform_diagonal()`, `tfb_transpose()`, `tfb_weibull_cdf()`, `tfb_weibull()`

<code>tfb_inverse</code>	<i>Returns the inverse Bijector evaluation, i.e., $X = g^{-1}(Y)$.</i>
--------------------------	---

Description

Returns the inverse Bijector evaluation, i.e., $X = g^{-1}(Y)$.

Usage

```
tfb_inverse(bijector, y, name = "inverse")
```

Arguments

bijector	The bijector to apply
y	Tensor. The input to the "inverse" evaluation.
name	name of the operation

Value

a tensor

See Also

Other bijector_methods: [tfb_forward_log_det_jacobian\(\)](#), [tfb_forward\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#)

Examples

```
b <- tfb_affine_scalar(shift = 1, scale = 2)
x <- 10
y <- b %>% tfb_forward(x)
b %>% tfb_inverse(y)
```

tfb_inverse_log_det_jacobian

Returns the result of the inverse evaluation of the log determinant of the Jacobian

Description

Returns the result of the inverse evaluation of the log determinant of the Jacobian

Usage

```
tfb_inverse_log_det_jacobian(
  bijector,
  y,
  event_ndims,
  name = "inverse_log_det_jacobian"
)
```

Arguments

bijector	The bijector to apply
y	Tensor. The input to the "inverse" Jacobian determinant evaluation.
event_ndims	Number of dimensions in the probabilistic events being transformed. Must be greater than or equal to <code>bijector\$inverse_min_event_ndims</code> . The result is summed over the final dimensions to produce a scalar Jacobian determinant for each event, i.e. it has shape <code>x\$shape\$ndims - event_ndims</code> dimensions.
name	name of the operation

Value

a tensor

See Also

Other bijector_methods: `tfb_forward_log_det_jacobian()`, `tfb_forward()`, `tfb_inverse()`

Examples

```
b <- tfb_affine_scalar(shift = 1, scale = 2)
x <- 10
y <- b %>% tfb_forward(x)
b %>% tfb_inverse_log_det_jacobian(y, event_ndims = 0)
```

tfb_invert

Bijector which inverts another Bijector

Description

Creates a Bijector which swaps the meaning of inverse and forward. Note: An inverted bijector's `inverse_log_det_jacobian` is often more efficient if the base bijector implements `_forward_log_det_jacobian`. If `_forward_log_det_jacobian` is not implemented then the following code is used: `y = b$inverse(x)`
`-b$inverse_log_det_jacobian(y)`

Usage

```
tfb_invert(bijector, validate_args = FALSE, name = NULL)
```

Arguments

<code>bijector</code>	Bijector instance.
<code>validate_args</code>	Logical, default FALSE. Whether to validate input with asserts. If <code>validate_args</code> is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
<code>name</code>	name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_template()`, `tfb_masked_autoregressive_flow()`, `tfb_masked_dense()`, `tfb_matrix_inverse_tri_l()`, `tfb_matvec_lu()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_pad()`, `tfb_permute()`, `tfb_power_transform()`, `tfb_rational_quadratic_spline()`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`, `tfb_real_nvp()`, `tfb_reciprocal()`, `tfb_reshape()`, `tfb_scale_matvec_diag()`, `tfb_scale_matvec_linear_operat`, `tfb_scale_matvec_lu()`, `tfb_scale_matvec_tri_l()`, `tfb_scale_tri_l()`, `tfb_scale()`, `tfb_shifted_gompertz_cdf()`, `tfb_shift()`, `tfb_sigmoid()`, `tfb_sinh_arcsinh()`, `tfb_sinh()`, `tfb_softmax_centered()`, `tfb_softplus()`, `tfb_softsign()`, `tfb_split()`, `tfb_square()`, `tfb_tanh()`, `tfb_transform_diagonal()`, `tfb_transpose()`, `tfb_weibull_cdf()`, `tfb_weibull()`

`tfb_iterated_sigmoid_centered`

Bijector which applies a Stick Breaking procedure.

Description

Bijector which applies a Stick Breaking procedure.

Usage

```
tfb_iterated_sigmoid_centered(validate_args = FALSE, name = "iterated_sigmoid")
```

Arguments

<code>validate_args</code>	Logical, default FALSE. Whether to validate input with asserts. If <code>validate_args</code> is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
<code>name</code>	name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`,

```
tfb_cumsum(), tfb_discrete_cosine_transform(), tfb_expm1(), tfb_exp(), tfb_ffjord(),
tfb_fill_scale_tri_l(), tfb_fill_triangular(), tfb_glow(), tfb_gompertz_cdf(), tfb_gumbel_cdf(),
tfb_gumbel(), tfb_identity(), tfb_inline(), tfb_invert(), tfb_kumaraswamy_cdf(), tfb_kumaraswamy(),
tfb_lambert_w_tail(), tfb_masked_autoregressive_default_template(), tfb_masked_autoregressive_flow(),
tfb_masked_dense(), tfb_matrix_inverse_tri_l(), tfb_matvec_lu(), tfb_normal_cdf(),
tfb_ordered(), tfb_pad(), tfb_permute(), tfb_power_transform(), tfb_rational_quadratic_spline(),
tfb_rayleigh_cdf(), tfb_real_nvp_default_template(), tfb_real_nvp(), tfb_reciprocal(),
tfb_reshape(), tfb_scale_matvec_diag(), tfb_scale_matvec_linear_operator(), tfb_scale_matvec_lu(),
tfb_scale_matvec_tri_l(), tfb_scale_tri_l(), tfb_scale(), tfb_shifted_gompertz_cdf(),
tfb_shift(), tfb_sigmoid(), tfb_sinh_arcsinh(), tfb_sinh(), tfb_softmax_centered(),
tfb_softplus(), tfb_softsign(), tfb_split(), tfb_square(), tfb_tanh(), tfb_transform_diagonal(),
tfb_transpose(), tfb_weibull_cdf(), tfb_weibull()
```

tfb_kumaraswamy	<i>Computes</i> $Y = g(X) = (1 - (1 - X)^{(1/b)})^{1/a}$, with X in $[0, 1]$
-----------------	---

Description

This bijector maps inputs from $[0, 1]$ to $[0, 1]$. The inverse of the bijector applied to a uniform random variable $X \sim U(0, 1)$ gives back a random variable with the **Kumaraswamy distribution**: $Y \sim \text{Kumaraswamy}(a, b)$ $\text{pdf}(y; a, b, 0 \leq y \leq 1) = a * b * y^{a-1} * (1 - y^b)^{b-1}$

Usage

```
tfb_kumaraswamy(
  concentration1 = NULL,
  concentration0 = NULL,
  validate_args = FALSE,
  name = "kumaraswamy"
)
```

Arguments

concentration1	float scalar indicating the transform power, i.e., $Y = g(X) = (1 - (1 - X)^{(1/b)})^{1/a}$ where a is concentration1.
concentration0	float scalar indicating the transform power, i.e., $Y = g(X) = (1 - (1 - X)^{(1/b)})^{1/a}$ where b is concentration0.
validate_args	Logical, default FALSE. Whether to validate input with asserts. If <code>validate_args</code> is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
name	name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_template()`, `tfb_masked_autoregressive_flow()`, `tfb_masked_dense()`, `tfb_matrix_inverse_tri_l()`, `tfb_matvec_lu()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_pad()`, `tfb_permute()`, `tfb_power_transform()`, `tfb_rational_quadratic_spline()`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`, `tfb_real_nvp()`, `tfb_reciprocal()`, `tfb_reshape()`, `tfb_scale_matvec_diag()`, `tfb_scale_matvec_linear_operator()`, `tfb_scale_matvec_lu()`, `tfb_scale_matvec_tri_l()`, `tfb_scale_tri_l()`, `tfb_scale()`, `tfb_shifted_gompertz_cdf()`, `tfb_shift()`, `tfb_sigmoid()`, `tfb_sinh_arcsinh()`, `tfb_sinh()`, `tfb_softmax_centered()`, `tfb_softplus()`, `tfb_softsign()`, `tfb_split()`, `tfb_square()`, `tfb_tanh()`, `tfb_transform_diagonal()`, `tfb_transpose()`, `tfb_weibull_cdf()`, `tfb_weibull()`

<code>tfb_kumaraswamy_cdf</code>	<i>Computes</i> $Y = g(X) = (1 - (1 - X)^{(1/b)})^{(1/a)}$, with X in $[0, 1]$
----------------------------------	---

Description

This bijector maps inputs from $[0, 1]$ to $[0, 1]$. The inverse of the bijector applied to a uniform random variable $X \sim U(0, 1)$ gives back a random variable with the **Kumaraswamy distribution**: $Y \sim \text{Kumaraswamy}(a, b)$ $\text{pdf}(y; a, b, 0 \leq y \leq 1) = a * b * y^{a-1} * (1 - y^b)^{b-1}$

Usage

```
tfb_kumaraswamy_cdf(
  concentration1 = 1,
  concentration0 = 1,
  validate_args = FALSE,
  name = "kumaraswamy_cdf"
)
```

Arguments

<code>concentration1</code>	float scalar indicating the transform power, i.e., $Y = g(X) = (1 - (1 - X)^{(1/b)})^{(1/a)}$ where a is <code>concentration1</code> .
<code>concentration0</code>	float scalar indicating the transform power, i.e., $Y = g(X) = (1 - (1 - X)^{(1/b)})^{(1/a)}$ where b is <code>concentration0</code> .
<code>validate_args</code>	Logical, default <code>FALSE</code> . Whether to validate input with asserts. If <code>validate_args</code> is <code>FALSE</code> , and the inputs are invalid, correct behavior is not guaranteed.
<code>name</code>	name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_template()`, `tfb_masked_autoregressive_flow()`, `tfb_masked_dense()`, `tfb_matrix_inverse_tri_l()`, `tfb_matvec_lu()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_pad()`, `tfb_permute()`, `tfb_power_transform()`, `tfb_rational_quadratic_spline()`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`, `tfb_real_nvp()`, `tfb_reciprocal()`, `tfb_reshape()`, `tfb_scale_matvec_diag()`, `tfb_scale_matvec_linear_operator()`, `tfb_scale_matvec_lu()`, `tfb_scale_matvec_tri_l()`, `tfb_scale_tri_l()`, `tfb_scale()`, `tfb_shifted_gompertz_cdf()`, `tfb_shift()`, `tfb_sigmoid()`, `tfb_sinh_arcsinh()`, `tfb_sinh()`, `tfb_softmax_centered()`, `tfb_softplus()`, `tfb_softsign()`, `tfb_split()`, `tfb_square()`, `tfb_tanh()`, `tfb_transform_diagonal()`, `tfb_transpose()`, `tfb_weibull_cdf()`, `tfb_weibull()`

<code>tfb_lambert_w_tail</code>	<i>LambertWTail transformation for heavy-tail Lambert W x F random variables.</i>
---------------------------------	---

Description

A random variable Y has a Lambert $W \times F$ distribution if $W_{\text{tau}}(Y) = X$ has distribution F , where $\text{tau} = (\text{shift}, \text{scale}, \text{tail})$ parameterizes the inverse transformation.

Usage

```
tfb_lambert_w_tail(
  shift = NULL,
  scale = NULL,
  tailweight = NULL,
  validate_args = FALSE,
  name = "lambertw_tail"
)
```

Arguments

<code>shift</code>	Floating point tensor; the shift for centering (uncentering) the input (output) random variable(s).
<code>scale</code>	Floating point tensor; the scaling (unscaling) of the input (output) random variable(s). Must contain only positive values.

tailweight	Floating point tensor; the tail behaviors of the output random variable(s). Must contain only non-negative values.
validate_args	Logical, default FALSE. Whether to validate input with asserts. If validate_args is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
name	name prefixed to Ops created by this class.

Details

This bijector defines the transformation underlying Lambert W x F distributions that transform an input random variable to an output random variable with heavier tails. It is defined as $Y = (U * \exp(0.5 * \text{tail} * U^2)) * \text{scale} + \text{shift}$, $\text{tail} \geq 0$ where $U = (X - \text{shift}) / \text{scale}$ is a shifted/scaled input random variable, and $\text{tail} \geq 0$ is the tail parameter.

Attributes: shift: shift to center (uncenter) the input data. scale: scale to normalize (de-normalize) the input data. tailweight: Tail parameter δ of heavy-tail transformation; must be ≥ 0 .

Value

a bijector instance.

See Also

For usage examples see [tfb_forward\(\)](#), [tfb_inverse\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#).

Other bijectors: [tfb_absolute_value\(\)](#), [tfb_affine_linear_operator\(\)](#), [tfb_affine_scalar\(\)](#), [tfb_affine\(\)](#), [tfb_ascending\(\)](#), [tfb_batch_normalization\(\)](#), [tfb_blockwise\(\)](#), [tfb_chain\(\)](#), [tfb_cholesky_outer_product\(\)](#), [tfb_cholesky_to_inv_cholesky\(\)](#), [tfb_correlation_cholesky\(\)](#), [tfb_cumsum\(\)](#), [tfb_discrete_cosine_transform\(\)](#), [tfb_expm1\(\)](#), [tfb_exp\(\)](#), [tfb_ffjord\(\)](#), [tfb_fill_scale_tri_l\(\)](#), [tfb_fill_triangular\(\)](#), [tfb_glow\(\)](#), [tfb_gompertz_cdf\(\)](#), [tfb_gumbel_cdf\(\)](#), [tfb_gumbel\(\)](#), [tfb_identity\(\)](#), [tfb_inline\(\)](#), [tfb_invert\(\)](#), [tfb_iterated_sigmoid_centered\(\)](#), [tfb_kumaraswamy_cdf\(\)](#), [tfb_kumaraswamy\(\)](#), [tfb_masked_autoregressive_default_template\(\)](#), [tfb_masked_autoregressive_flow\(\)](#), [tfb_masked_dense\(\)](#), [tfb_matrix_inverse_tri_l\(\)](#), [tfb_matvec_lu\(\)](#), [tfb_normal_cdf\(\)](#), [tfb_ordered\(\)](#), [tfb_pad\(\)](#), [tfb_permute\(\)](#), [tfb_power_transform\(\)](#), [tfb_rational_quadratic_spline\(\)](#), [tfb_rayleigh_cdf\(\)](#), [tfb_real_nvp_default_template\(\)](#), [tfb_real_nvp\(\)](#), [tfb_reciprocal\(\)](#), [tfb_reshape\(\)](#), [tfb_scale_matvec_diag\(\)](#), [tfb_scale_matvec_linear_operator\(\)](#), [tfb_scale_matvec_lu\(\)](#), [tfb_scale_matvec_tri_l\(\)](#), [tfb_scale_tri_l\(\)](#), [tfb_scale\(\)](#), [tfb_shifted_gompertz_cdf\(\)](#), [tfb_shift\(\)](#), [tfb_sigmoid\(\)](#), [tfb_sinh_arcsinh\(\)](#), [tfb_sinh\(\)](#), [tfb_softmax_centered\(\)](#), [tfb_softplus\(\)](#), [tfb_softsign\(\)](#), [tfb_split\(\)](#), [tfb_square\(\)](#), [tfb_tanh\(\)](#), [tfb_transform_diagonal\(\)](#), [tfb_transpose\(\)](#), [tfb_weibull_cdf\(\)](#), [tfb_weibull\(\)](#)

tfb_masked_autoregressive_default_template

Masked Autoregressive Density Estimator

Description

This will be wrapped in a `make_template` to ensure the variables are only created once. It takes the input and returns the loc ("mu" in Germain et al. (2015)) and `log_scale` ("alpha" in Germain et al. (2015)) from the MADE network.

Usage

```
tfb_masked_autoregressive_default_template(
    hidden_layers,
    shift_only = FALSE,
    activation = tf$nn$relu,
    log_scale_min_clip = -5,
    log_scale_max_clip = 3,
    log_scale_clip_gradient = FALSE,
    name = NULL,
    ...
)
```

Arguments

<code>hidden_layers</code>	list-like of non-negative integer, scalars indicating the number of units in each hidden layer. Default: <code>list(512, 512)</code> .
<code>shift_only</code>	logical indicating if only the shift term shall be computed. Default: <code>FALSE</code> .
<code>activation</code>	Activation function (callable). Explicitly setting to <code>NULL</code> implies a linear activation.
<code>log_scale_min_clip</code>	float-like scalar Tensor, or a Tensor with the same shape as <code>log_scale</code> . The minimum value to clip by. Default: <code>-5</code> .
<code>log_scale_max_clip</code>	float-like scalar Tensor, or a Tensor with the same shape as <code>log_scale</code> . The maximum value to clip by. Default: <code>3</code> .
<code>log_scale_clip_gradient</code>	logical indicating that the gradient of <code>tf\$clip_by_value</code> should be preserved. Default: <code>FALSE</code> .
<code>name</code>	A name for ops managed by this function. Default: <code>"tfb_masked_autoregressive_default_template"</code> .
<code>...</code>	<code>tf\$layers\$dense</code> arguments

Details

Warning: This function uses `masked_dense` to create randomly initialized `tf$Variables`. It is presumed that these will be fit, just as you would any other neural architecture which uses `tf$layers$dense`.

About Hidden Layers Each element of `hidden_layers` should be greater than the `input_depth` (i.e., `input_depth = tf$shape(input)[-1]` where `input` is the input to the neural network). This is necessary to ensure the autoregressivity property.

About Clipping This function also optionally clips the `log_scale` (but possibly not its gradient). This is useful because if `log_scale` is too small/large it might underflow/overflow making it impossible for the `MaskedAutoregressiveFlow` bijector to implement a bijection. Additionally, the `log_scale_clip_gradient` bool indicates whether the gradient should also be clipped. The default does not clip the gradient; this is useful because it still provides gradient information (for fitting) yet solves the numerical stability problem. I.e., `log_scale_clip_gradient = FALSE` means $\text{grad}[\exp(\text{clip}(x))] = \text{grad}[x] \exp(\text{clip}(x))$ rather than the usual $\text{grad}[\text{clip}(x)] \exp(\text{clip}(x))$.

Value

list of:

- shift: Float-like Tensor of shift terms
- log_scale: Float-like Tensor of log(scale) terms

References

- Mathieu Germain, Karol Gregor, Iain Murray, and Hugo Larochelle. MADE: Masked Auto-encoder for Distribution Estimation. In *International Conference on Machine Learning, 2015*.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_flow()`, `tfb_masked_dense()`, `tfb_matrix_inverse_tri_l()`, `tfb_matvec_lu()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_pad()`, `tfb_permute()`, `tfb_power_transform()`, `tfb_rational_quadratic_spline()`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`, `tfb_real_nvp()`, `tfb_reciprocal()`, `tfb_reshape()`, `tfb_scale_matvec_diag()`, `tfb_scale_matvec_linear_operator()`, `tfb_scale_matvec_lu()`, `tfb_scale_matvec_tri_l()`, `tfb_scale_tri_l()`, `tfb_scale()`, `tfb_shifted_gompertz_cdf()`, `tfb_shift()`, `tfb_sigmoid()`, `tfb_sinh_arcsinh()`, `tfb_sinh()`, `tfb_softmax_centered()`, `tfb_softplus()`, `tfb_softsign()`, `tfb_split()`, `tfb_square()`, `tfb_tanh()`, `tfb_transform_diagonal()`, `tfb_transpose()`, `tfb_weibull_cdf()`, `tfb_weibull()`

tfb_masked_autoregressive_flow

Affine MaskedAutoregressiveFlow bijector

Description

The affine autoregressive flow (Papamakarios et al., 2016) provides a relatively simple framework for user-specified (deep) architectures to learn a distribution over continuous events. Regarding terminology,

Usage

```
tfb_masked_autoregressive_flow(
    shift_and_log_scale_fn,
    is_constant_jacobian = FALSE,
```

```

    unroll_loop = FALSE,
    event_ndims = 1L,
    validate_args = FALSE,
    name = NULL
)

```

Arguments

`shift_and_log_scale_fn`

Function which computes `shift` and `log_scale` from both the forward domain (`x`) and the inverse domain (`y`). Calculation must respect the "autoregressive property". Suggested default: `tfb_masked_autoregressive_default_template(hidden_layers=...)`. Typically the function contains `tf$Variables` and is wrapped using `tf$make_template`. Returning `NULL` for either (both) `shift`, `log_scale` is equivalent to (but more efficient than) returning zero.

`is_constant_jacobian`

Logical, default: `FALSE`. When `TRUE` the implementation assumes `log_scale` does not depend on the forward domain (`x`) or inverse domain (`y`) values. (No validation is made; `is_constant_jacobian=FALSE` is always safe but possibly computationally inefficient.)

`unroll_loop`

Logical indicating whether the `tf$while_loop` in `_forward` should be replaced with a static for loop. Requires that the final dimension of `x` be known at graph construction time. Defaults to `FALSE`.

`event_ndims`

integer, the intrinsic dimensionality of this bijector. 1 corresponds to a simple vector autoregressive bijector as implemented by the `tfb_masked_autoregressive_default_template`. 2 might be useful for a 2D convolutional `shift_and_log_scale_fn` and so on.

`validate_args`

Logical, default `FALSE`. Whether to validate input with asserts. If `validate_args` is `FALSE`, and the inputs are invalid, correct behavior is not guaranteed.

`name`

name prefixed to Ops created by this class.

Details

"Autoregressive models decompose the joint density as a product of conditionals, and model each conditional in turn. Normalizing flows transform a base density (e.g. a standard Gaussian) into the target density by an invertible transformation with tractable Jacobian." (Papamakarios et al., 2016)

In other words, the "autoregressive property" is equivalent to the decomposition, $p(x) = \prod_{i=0}^{d-1} p(x[\text{perm}[i]] \mid x[\text{perm}[0:i]])$ where `perm` is some permutation of $\{0, \dots, d\}$. In the simple case where the permutation is identity this reduces to:

$p(x) = \prod_{i=0}^{d-1} p(x[i] \mid x[0:i])$. The provided `shift_and_log_scale_fn`, `tfb_masked_autoregressive_default_template`, achieves this property by zeroing out weights in its `masked_dense` layers. In TensorFlow Probability, "normalizing flows" are implemented as `tfp.bijectors.Bijectors`. The forward "autoregression" is implemented using a `tf.while_loop` and a deep neural network (DNN) with masked weights such that the autoregressive property is automatically met in the inverse. A `TransformedDistribution` using `MaskedAutoregressiveFlow(...)` uses the (expensive) forward-mode calculation to draw samples and the (cheap) reverse-mode calculation to compute log-probabilities. Conversely, a `TransformedDistribution` using `Invert(MaskedAutoregressiveFlow(...))` uses the (expensive) forward-mode calculation to compute log-probabilities and the (cheap) reverse-mode calculation to compute samples.

Given a `shift_and_log_scale_fn`, the forward and inverse transformations are (a sequence of) affine transformations. A "valid" `shift_and_log_scale_fn` must compute each shift (aka μ or "mu" in Germain et al. (2015)) and $\log(\text{scale})$ (aka "alpha" in Germain et al. (2015)) such that each are broadcastable with the arguments to forward and inverse, i.e., such that the calculations in forward, inverse below are possible.

For convenience, `tfb_masked_autoregressive_default_template` is offered as a possible `shift_and_log_scale_fn` function. It implements the MADE architecture (Germain et al., 2015). MADE is a feed-forward network that computes a shift and $\log(\text{scale})$ using `masked_dense` layers in a deep neural network. Weights are masked to ensure the autoregressive property. It is possible that this architecture is suboptimal for your task. To build alternative networks, either change the arguments to `tfb_masked_autoregressive_default_template`, use the `masked_dense` function to roll-out your own, or use some other architecture, e.g., using `tf.layers`. Warning: no attempt is made to validate that the `shift_and_log_scale_fn` enforces the "autoregressive property".

Assuming `shift_and_log_scale_fn` has valid shape and autoregressive semantics, the forward transformation is

```
def forward(x):
    y = zeros_like(x)
    event_size = x.shape[-event_dims:].num_elements()
    for _ in range(event_size):
        shift, log_scale = shift_and_log_scale_fn(y)
        y = x * tf.exp(log_scale) + shift
    return y
```

and the inverse transformation is

```
def inverse(y):
    shift, log_scale = shift_and_log_scale_fn(y)
    return (y - shift) / tf.exp(log_scale)
```

Notice that the inverse does not need a for-loop. This is because in the forward pass each calculation of `shift` and `log_scale` is based on the `y` calculated so far (not `x`). In the inverse, the `y` is fully known, thus is equivalent to the scaling used in forward after `event_size` passes, i.e., the "last" `y` used to compute `shift`, `log_scale`. (Roughly speaking, this also proves the transform is bijective.)

Value

a bijector instance.

References

- Mathieu Germain, Karol Gregor, Iain Murray, and Hugo Larochelle. MADE: Masked Autoencoder for Distribution Estimation. In *International Conference on Machine Learning*, 2015.
- Diederik P. Kingma, Tim Salimans, Rafal Jozefowicz, Xi Chen, Ilya Sutskever, and Max Welling. Improving Variational Inference with Inverse Autoregressive Flow. In *Neural Information Processing Systems*, 2016.
- George Papamakarios, Theo Pavlakou, and Iain Murray. Masked Autoregressive Flow for Density Estimation. In *Neural Information Processing Systems*, 2017.

See Also

For usage examples see [tfb_forward\(\)](#), [tfb_inverse\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#).

Other bijectors: [tfb_absolute_value\(\)](#), [tfb_affine_linear_operator\(\)](#), [tfb_affine_scalar\(\)](#), [tfb_affine\(\)](#), [tfb_ascending\(\)](#), [tfb_batch_normalization\(\)](#), [tfb_blockwise\(\)](#), [tfb_chain\(\)](#), [tfb_cholesky_outer_product\(\)](#), [tfb_cholesky_to_inv_cholesky\(\)](#), [tfb_correlation_cholesky\(\)](#), [tfb_cumsum\(\)](#), [tfb_discrete_cosine_transform\(\)](#), [tfb_expm1\(\)](#), [tfb_exp\(\)](#), [tfb_ffjord\(\)](#), [tfb_fill_scale_tri_l\(\)](#), [tfb_fill_triangular\(\)](#), [tfb_glow\(\)](#), [tfb_gompertz_cdf\(\)](#), [tfb_gumbel_cdf\(\)](#), [tfb_gumbel\(\)](#), [tfb_identity\(\)](#), [tfb_inline\(\)](#), [tfb_invert\(\)](#), [tfb_iterated_sigmoid_centered\(\)](#), [tfb_kumaraswamy_cdf\(\)](#), [tfb_kumaraswamy\(\)](#), [tfb_lambert_w_tail\(\)](#), [tfb_masked_autoregressive_default_template\(\)](#), [tfb_masked_dense\(\)](#), [tfb_matrix_inverse_tri_l\(\)](#), [tfb_matvec_lu\(\)](#), [tfb_normal_cdf\(\)](#), [tfb_ordered\(\)](#), [tfb_pad\(\)](#), [tfb_permute\(\)](#), [tfb_power_transform\(\)](#), [tfb_rational_quadratic_spline\(\)](#), [tfb_rayleigh_cdf\(\)](#), [tfb_real_nvp_default_template\(\)](#), [tfb_real_nvp\(\)](#), [tfb_reciprocal\(\)](#), [tfb_reshape\(\)](#), [tfb_scale_matvec_diag\(\)](#), [tfb_scale_matvec_linear_operator\(\)](#), [tfb_scale_matvec_lu\(\)](#), [tfb_scale_matvec_tri_l\(\)](#), [tfb_scale_tri_l\(\)](#), [tfb_scale\(\)](#), [tfb_shifted_gompertz_cdf\(\)](#), [tfb_shift\(\)](#), [tfb_sigmoid\(\)](#), [tfb_sinh_arcsinh\(\)](#), [tfb_sinh\(\)](#), [tfb_softmax_centered\(\)](#), [tfb_softplus\(\)](#), [tfb_softsign\(\)](#), [tfb_split\(\)](#), [tfb_square\(\)](#), [tfb_tanh\(\)](#), [tfb_transform_diagonal\(\)](#), [tfb_transpose\(\)](#), [tfb_weibull_cdf\(\)](#), [tfb_weibull\(\)](#)

tfb_masked_dense	<i>Autoregressively masked dense layer</i>
------------------	--

Description

Analogous to `tf.layers.dense`.

Usage

```
tfb_masked_dense(
    inputs,
    units,
    num_blocks = NULL,
    exclusive = FALSE,
    kernel_initializer = NULL,
    reuse = NULL,
    name = NULL,
    ...
)
```

Arguments

<code>inputs</code>	Tensor input.
<code>units</code>	integer scalar representing the dimensionality of the output space.
<code>num_blocks</code>	integer scalar representing the number of blocks for the MADE masks.
<code>exclusive</code>	logical scalar representing whether to zero the diagonal of the mask, used for the first layer of a MADE.

kernel_initializer	Initializer function for the weight matrix. If NULL (default), weights are initialized using the <code>tf\$glorot_random_initializer</code>
reuse	logical scalar representing whether to reuse the weights of a previous layer by the same name.
name	string used to describe ops managed by this function.
...	tf\$layers\$dense arguments

Details

See Germain et al. (2015) for detailed explanation.

Value

tensor

References

- Mathieu Germain, Karol Gregor, Iain Murray, and Hugo Larochelle. *MADE: Masked Autoencoder for Distribution Estimation*. In *International Conference on Machine Learning*, 2015.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_template()`, `tfb_masked_autoregressive_flow()`, `tfb_matrix_inverse_tri_l()`, `tfb_matvec_lu()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_pad()`, `tfb_permute()`, `tfb_power_transform()`, `tfb_rational_quadratic_spline()`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`, `tfb_real_nvp()`, `tfb_reciprocal()`, `tfb_reshape()`, `tfb_scale_matvec_diag()`, `tfb_scale_matvec_linear_operator()`, `tfb_scale_matvec_lu()`, `tfb_scale_matvec_tri_l()`, `tfb_scale_tri_l()`, `tfb_scale()`, `tfb_shifted_gompertz_cdf()`, `tfb_shift()`, `tfb_sigmoid()`, `tfb_sinh_arcsinh()`, `tfb_sinh()`, `tfb_softmax_centered()`, `tfb_softplus()`, `tfb_softsign()`, `tfb_split()`, `tfb_square()`, `tfb_tanh()`, `tfb_transform_diagonal()`, `tfb_transpose()`, `tfb_weibull_cdf()`, `tfb_weibull()`

tfb_matrix_inverse_tri_l

Computes $g(L) = \text{inv}(L)$, where L is a lower-triangular matrix

Description

L must be nonsingular; equivalently, all diagonal entries of L must be nonzero. The input must have rank ≥ 2 . The input is treated as a batch of matrices with batch shape `input.shape[:-2]`, where each matrix has dimensions `input.shape[-2]` by `input.shape[-1]` (hence `input.shape[-2]` must equal `input.shape[-1]`).

Usage

```
tfb_matrix_inverse_tri_l(validate_args = FALSE, name = "matrix_inverse_tril")
```

Arguments

`validate_args` Logical, default FALSE. Whether to validate input with asserts. If `validate_args` is FALSE, and the inputs are invalid, correct behavior is not guaranteed.

`name` name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_template()`, `tfb_masked_autoregressive_flow()`, `tfb_masked_dense()`, `tfb_matvec_lu()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_pad()`, `tfb_permute()`, `tfb_power_transform()`, `tfb_rational_quadratic_spline()`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`, `tfb_real_nvp()`, `tfb_reciprocal()`, `tfb_reshape()`, `tfb_scale_matvec_diag()`, `tfb_scale_matvec_linear_operator()`, `tfb_scale_matvec_lu()`, `tfb_scale_matvec_tri_l()`, `tfb_scale_tri_l()`, `tfb_scale()`, `tfb_shifted_gompertz_cdf()`, `tfb_shift()`, `tfb_sigmoid()`, `tfb_sinh_arcsinh()`, `tfb_sinh()`, `tfb_softmax_centered()`, `tfb_softplus()`, `tfb_softsign()`, `tfb_split()`, `tfb_square()`, `tfb_tanh()`, `tfb_transform_diagonal()`, `tfb_transpose()`, `tfb_weibull_cdf()`, `tfb_weibull()`

tfb_matvec_lu	<i>Matrix-vector multiply using LU decomposition</i>
---------------	--

Description

This bijector is identical to the "Convolution1x1" used in Glow (Kingma and Dhariwal, 2018).

Usage

```
tfb_matvec_lu(lower_upper, permutation, validate_args = FALSE, name = NULL)
```

Arguments

lower_upper	The LU factorization as returned by <code>tf\$linalg\$lu</code> .
permutation	The LU factorization permutation as returned by <code>tf\$linalg\$lu</code> .
validate_args	Logical, default FALSE. Whether to validate input with asserts. If <code>validate_args</code> is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
name	name prefixed to Ops created by this class.

Details

Warning: this bijector never verifies the scale matrix (as parameterized by LU decomposition) is invertible. Ensuring this is the case is the caller's responsibility.

Value

a bijector instance.

References

- [Diederik P. Kingma, Prafulla Dhariwal. Glow: Generative Flow with Invertible 1x1 Convolutions. *arXiv preprint arXiv:1807.03039*, 2018.](#)

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_template()`, `tfb_masked_autoregressive_flow()`, `tfb_masked_dense()`, `tfb_matrix_inverse_tri_l()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_pad()`, `tfb_permute()`, `tfb_power_transform()`, `tfb_rational_quadratic()`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`, `tfb_real_nvp()`, `tfb_reciprocal()`,

tfb_reshape(), tfb_scale_matvec_diag(), tfb_scale_matvec_linear_operator(), tfb_scale_matvec_lu(),
 tfb_scale_matvec_tri_l(), tfb_scale_tri_l(), tfb_scale(), tfb_shifted_gompertz_cdf(),
 tfb_shift(), tfb_sigmoid(), tfb_sinh_arcsinh(), tfb_sinh(), tfb_softmax_centered(),
 tfb_softplus(), tfb_softsign(), tfb_split(), tfb_square(), tfb_tanh(), tfb_transform_diagonal(),
 tfb_transpose(), tfb_weibull_cdf(), tfb_weibull()

tfb_normal_cdf	<i>Computes</i> $Y = g(X) = \text{NormalCDF}(x)$
----------------	--

Description

This bijector maps inputs from $[-\infty, \infty]$ to $[0, 1]$. The inverse of the bijector applied to a uniform random variable $X \sim U(0, 1)$ gives back a random variable with the **Normal distribution**:

Usage

```
tfb_normal_cdf(validate_args = FALSE, name = "normal")
```

Arguments

validate_args	Logical, default FALSE. Whether to validate input with asserts. If validate_args is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
name	name prefixed to Ops created by this class.

Details

$Y \sim \text{Normal}(0, 1)$ pdf($y; 0., 1.$) = $1 / \sqrt{2 * \pi} * \exp(-y ** 2 / 2)$

Value

a bijector instance.

See Also

For usage examples see [tfb_forward\(\)](#), [tfb_inverse\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#).

Other bijectors: [tfb_absolute_value\(\)](#), [tfb_affine_linear_operator\(\)](#), [tfb_affine_scalar\(\)](#), [tfb_affine\(\)](#), [tfb_ascending\(\)](#), [tfb_batch_normalization\(\)](#), [tfb_blockwise\(\)](#), [tfb_chain\(\)](#), [tfb_cholesky_outer_product\(\)](#), [tfb_cholesky_to_inv_cholesky\(\)](#), [tfb_correlation_cholesky\(\)](#), [tfb_cumsum\(\)](#), [tfb_discrete_cosine_transform\(\)](#), [tfb_expm1\(\)](#), [tfb_exp\(\)](#), [tfb_ffjord\(\)](#), [tfb_fill_scale_tri_l\(\)](#), [tfb_fill_triangular\(\)](#), [tfb_glow\(\)](#), [tfb_gompertz_cdf\(\)](#), [tfb_gumbel_cdf\(\)](#), [tfb_gumbel\(\)](#), [tfb_identity\(\)](#), [tfb_inline\(\)](#), [tfb_invert\(\)](#), [tfb_iterated_sigmoid_centered\(\)](#), [tfb_kumaraswamy_cdf\(\)](#), [tfb_kumaraswamy\(\)](#), [tfb_lambert_w_tail\(\)](#), [tfb_masked_autoregressive_default_template\(\)](#), [tfb_masked_autoregressive_flow\(\)](#), [tfb_masked_dense\(\)](#), [tfb_matrix_inverse_tri_l\(\)](#), [tfb_matvec_lu\(\)](#), [tfb_ordered\(\)](#), [tfb_pad\(\)](#), [tfb_permute\(\)](#), [tfb_power_transform\(\)](#), [tfb_rational_quadratic_spline\(\)](#), [tfb_rayleigh_cdf\(\)](#), [tfb_real_nvp_default_template\(\)](#), [tfb_real_nvp\(\)](#), [tfb_reciprocal\(\)](#), [tfb_reshape\(\)](#), [tfb_scale_matvec_diag\(\)](#), [tfb_scale_matvec_linear_operator\(\)](#), [tfb_scale_matvec_lu\(\)](#), [tfb_scale_matvec_tri_l\(\)](#), [tfb_scale_tri_l\(\)](#), [tfb_scale\(\)](#), [tfb_shifted_gompertz_cdf\(\)](#),

[tfb_shift\(\)](#), [tfb_sigmoid\(\)](#), [tfb_sinh_arcsinh\(\)](#), [tfb_sinh\(\)](#), [tfb_softmax_centered\(\)](#), [tfb_softplus\(\)](#), [tfb_softsign\(\)](#), [tfb_split\(\)](#), [tfb_square\(\)](#), [tfb_tanh\(\)](#), [tfb_transform_diagonal\(\)](#), [tfb_transpose\(\)](#), [tfb_weibull_cdf\(\)](#), [tfb_weibull\(\)](#)

tfb_ordered	<i>Bijector which maps a tensor x_k that has increasing elements in the last dimension to an unconstrained tensor y_k</i>
-------------	---

Description

Both the domain and the codomain of the mapping is $[-\infty, \infty]$, however, the input of the forward mapping must be strictly increasing. The inverse of the bijector applied to a normal random vector $y \sim N(0, 1)$ gives back a sorted random vector with the same distribution $x \sim N(0, 1)$ where $x = \text{sort}(y)$

Usage

```
tfb_ordered(validate_args = FALSE, name = "ordered")
```

Arguments

validate_args	Logical, default FALSE. Whether to validate input with asserts. If validate_args is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
name	name prefixed to Ops created by this class.

Details

On the last dimension of the tensor, Ordered bijector performs: $y[0] = x[0]$ $y[1:] = \text{tf}\$log(x[1:] - x[:-1])$

Value

a bijector instance.

See Also

For usage examples see [tfb_forward\(\)](#), [tfb_inverse\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#).

Other bijectors: [tfb_absolute_value\(\)](#), [tfb_affine_linear_operator\(\)](#), [tfb_affine_scalar\(\)](#), [tfb_affine\(\)](#), [tfb_ascending\(\)](#), [tfb_batch_normalization\(\)](#), [tfb_blockwise\(\)](#), [tfb_chain\(\)](#), [tfb_cholesky_outer_product\(\)](#), [tfb_cholesky_to_inv_cholesky\(\)](#), [tfb_correlation_cholesky\(\)](#), [tfb_cumsum\(\)](#), [tfb_discrete_cosine_transform\(\)](#), [tfb_expm1\(\)](#), [tfb_exp\(\)](#), [tfb_ffjord\(\)](#), [tfb_fill_scale_tri_l\(\)](#), [tfb_fill_triangular\(\)](#), [tfb_glow\(\)](#), [tfb_gompertz_cdf\(\)](#), [tfb_gumbel_cdf\(\)](#), [tfb_gumbel\(\)](#), [tfb_identity\(\)](#), [tfb_inline\(\)](#), [tfb_invert\(\)](#), [tfb_iterated_sigmoid_centered\(\)](#), [tfb_kumaraswamy_cdf\(\)](#), [tfb_kumaraswamy\(\)](#), [tfb_lambert_w_tail\(\)](#), [tfb_masked_autoregressive_default_template\(\)](#), [tfb_masked_autoregressive_flow\(\)](#), [tfb_masked_dense\(\)](#), [tfb_matrix_inverse_tri_l\(\)](#), [tfb_matvec_lu\(\)](#), [tfb_normal_cdf\(\)](#), [tfb_pad\(\)](#), [tfb_permute\(\)](#), [tfb_power_transform\(\)](#), [tfb_rational_quadratic_spline\(\)](#), [tfb_rayleigh_cdf\(\)](#), [tfb_real_nvp_default_template\(\)](#),

tfb_real_nvp(), tfb_reciprocal(), tfb_reshape(), tfb_scale_matvec_diag(), tfb_scale_matvec_linear_operat
 tfb_scale_matvec_lu(), tfb_scale_matvec_tri_l(), tfb_scale_tri_l(), tfb_scale(), tfb_shifted_gompertz_co
 tfb_shift(), tfb_sigmoid(), tfb_sinh_arcsinh(), tfb_sinh(), tfb_softmax_centered(),
 tfb_softplus(), tfb_softsign(), tfb_split(), tfb_square(), tfb_tanh(), tfb_transform_diagonal(),
 tfb_transpose(), tfb_weibull_cdf(), tfb_weibull()

tfb_pad *Pads a value to the event_shape of a Tensor.*

Description

The semantics of `bijector_pad` generally follow that of `tf$pad()` except that `bijector_pad`'s `paddings` argument applies to the rightmost dimensions. Additionally, the new argument `axis` enables overriding the dimensions to which `paddings` is applied. Like `paddings`, the `axis` argument is also relative to the rightmost dimension and must therefore be negative. The argument `paddings` is a vector of integer pairs each representing the number of left and/or right `constant_values` to pad to the corresponding rightmost dimensions. That is, unless `axis` is specified, specifying `k` different `paddings` means the rightmost `k` dimensions will be "grown" by the sum of the respective `paddings` row. When `axis` is specified, it indicates the dimension to which the corresponding `paddings` element is applied. By default `axis` is `NULL` which means it is logically equivalent to `range(start=-len(paddings), limit=0)`, i.e., the rightmost dimensions.

Usage

```
tfb_pad(
  paddings = list(c(0, 1)),
  mode = "CONSTANT",
  constant_values = 0,
  axis = NULL,
  validate_args = FALSE,
  name = NULL
)
```

Arguments

<code>paddings</code>	A vector-shaped Tensor of integer pairs representing the number of elements to pad on the left and right, respectively. Default value: <code>list(reticulate::tuple(0L, 1L))</code> .
<code>mode</code>	One of 'CONSTANT', 'REFLECT', or 'SYMMETRIC' (case-insensitive). For more details, see <code>tf\$pad</code> .
<code>constant_values</code>	In "CONSTANT" mode, the scalar pad value to use. Must be same type as tensor. For more details, see <code>tf\$pad</code> .
<code>axis</code>	The dimensions for which <code>paddings</code> are applied. Must be 1:1 with <code>paddings</code> or <code>NULL</code> . Default value: <code>NULL</code> (i.e., <code>tf\$range(start = -length(paddings), limit = 0)</code>).
<code>validate_args</code>	Logical, default <code>FALSE</code> . Whether to validate input with asserts. If <code>validate_args</code> is <code>FALSE</code> , and the inputs are invalid, correct behavior is not guaranteed.
<code>name</code>	name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_template()`, `tfb_masked_autoregressive_flow()`, `tfb_masked_dense()`, `tfb_matrix_inverse_tri_l()`, `tfb_matvec_lu()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_permute()`, `tfb_power_transform()`, `tfb_rational_quadratic_spline()`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`, `tfb_real_nvp()`, `tfb_reciprocal()`, `tfb_reshape()`, `tfb_scale_matvec_diag()`, `tfb_scale_matvec_linear_operator()`, `tfb_scale_matvec_lu()`, `tfb_scale_matvec_tri_l()`, `tfb_scale_tri_l()`, `tfb_scale()`, `tfb_shifted_gompertz_cdf()`, `tfb_shift()`, `tfb_sigmoid()`, `tfb_sinh_arcsinh()`, `tfb_sinh()`, `tfb_softmax_centered()`, `tfb_softplus()`, `tfb_softsign()`, `tfb_split()`, `tfb_square()`, `tfb_tanh()`, `tfb_transform_diagonal()`, `tfb_transpose()`, `tfb_weibull_cdf()`, `tfb_weibull()`

tfb_permute

Permutes the rightmost dimension of a Tensor

Description

Permutes the rightmost dimension of a Tensor

Usage

```
tfb_permute(permutation, axis = -1L, validate_args = FALSE, name = NULL)
```

Arguments

permutation	An integer-like vector-shaped Tensor representing the permutation to apply to the axis dimension of the transformed Tensor.
axis	Scalar integer Tensor representing the dimension over which to <code>tf\$gather</code> . <code>axis</code> must be relative to the end (reading left to right) thus must be negative. Default value: -1 (i.e., right-most).
validate_args	Logical, default FALSE. Whether to validate input with asserts. If <code>validate_args</code> is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
name	name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expml()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_template()`, `tfb_masked_autoregressive_flow()`, `tfb_masked_dense()`, `tfb_matrix_inverse_tri_l()`, `tfb_matvec_lu()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_pad()`, `tfb_power_transform()`, `tfb_rational_quadratic_spline()`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`, `tfb_real_nvp()`, `tfb_reciprocal()`, `tfb_reshape()`, `tfb_scale_matvec_diag()`, `tfb_scale_matvec_linear_operator()`, `tfb_scale_matvec_lu()`, `tfb_scale_matvec_tri_l()`, `tfb_scale_tri_l()`, `tfb_scale()`, `tfb_shifted_gompertz_cdf()`, `tfb_shift()`, `tfb_sigmoid()`, `tfb_sinh_arcsinh()`, `tfb_sinh()`, `tfb_softmax_centered()`, `tfb_softplus()`, `tfb_softsign()`, `tfb_split()`, `tfb_square()`, `tfb_tanh()`, `tfb_transform_diagonal()`, `tfb_transpose()`, `tfb_weibull_cdf()`, `tfb_weibull()`

`tfb_power_transform` *Computes* $Y = g(X) = (1 + X * c)**(1 / c)$, where $X \geq -1 / c$

Description

The **power transform** maps inputs from $[0, \infty]$ to $[-1/c, \infty]$; this is equivalent to the inverse of this bijector. This bijector is equivalent to the Exp bijector when $c=0$.

Usage

```
tfb_power_transform(power, validate_args = FALSE, name = "power_transform")
```

Arguments

<code>power</code>	float scalar indicating the transform power, i.e., $Y = g(X) = (1 + X * c)**(1 / c)$ where c is the power.
<code>validate_args</code>	Logical, default FALSE. Whether to validate input with asserts. If <code>validate_args</code> is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
<code>name</code>	name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_exp1()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_template()`, `tfb_masked_autoregressive_flow()`, `tfb_masked_dense()`, `tfb_matrix_inverse_tri_l()`, `tfb_matvec_lu()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_pad()`, `tfb_permute()`, `tfb_rational_quadratic_spline`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`, `tfb_real_nvp()`, `tfb_reciprocal()`, `tfb_reshape()`, `tfb_scale_matvec_diag()`, `tfb_scale_matvec_linear_operator()`, `tfb_scale_matvec_lu()`, `tfb_scale_matvec_tri_l()`, `tfb_scale_tri_l()`, `tfb_scale()`, `tfb_shifted_gompertz_cdf()`, `tfb_shift()`, `tfb_sigmoid()`, `tfb_sinh_arcsinh()`, `tfb_sinh()`, `tfb_softmax_centered()`, `tfb_softplus()`, `tfb_softsign()`, `tfb_split()`, `tfb_square()`, `tfb_tanh()`, `tfb_transform_diagonal()`, `tfb_transpose()`, `tfb_weibull_cdf()`, `tfb_weibull()`

tfb_rational_quadratic_spline

A piecewise rational quadratic spline, as developed in Conor et al.(2019).

Description

This transformation represents a monotonically increasing piecewise rational quadratic function. Outside of the bounds of `knot_x/knot_y`, the transform behaves as an identity function.

Usage

```
tfb_rational_quadratic_spline(
    bin_widths,
    bin_heights,
    knot_slopes,
    range_min = -1,
    validate_args = FALSE,
    name = NULL
)
```

Arguments

`bin_widths` The widths of the spans between subsequent knot x positions, a floating point Tensor. Must be positive, and at least 1-D. Innermost axis must sum to the same value as `bin_heights`. The knot x positions will be a first at `range_min`, followed by knots at `range_min + cumsum(bin_widths, axis=-1)`.

bin_heights	The heights of the spans between subsequent knot y positions, a floating point Tensor. Must be positive, and at least 1-D. Innermost axis must sum to the same value as bin_widths. The knot y positions will be a first at range_min, followed by knots at range_min + cumsum(bin_heights,axis=-1).
knot_slopes	The slope of the spline at each knot, a floating point Tensor. Must be positive. 1s are implicitly padded for the first and last implicit knots corresponding to range_min and range_min + sum(bin_widths,axis=-1). Innermost axis size should be 1 less than that of bin_widths/bin_heights, or 1 for broadcasting.
range_min	The x/y position of the first knot, which has implicit slope 1. range_max is implicit, and can be computed as range_min + sum(bin_widths,axis=-1). Scalar floating point Tensor.
validate_args	Logical, default FALSE. Whether to validate input with asserts. If validate_args is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
name	name prefixed to Ops created by this class.

Details

Typically this bijector will be used as part of a chain, with splines for trailing x dimensions conditioned on some of the earlier x dimensions, and with the inverse then solved first for unconditioned dimensions, then using conditioning derived from those inverses, and so forth.

For each argument, the innermost axis indexes bins/knots and batch axes index axes of x/y spaces. A RationalQuadraticSpline with a separate transform for each of three dimensions might have bin_widths shaped [3, 32]. To use the same spline for each of x's three dimensions we may broadcast against x and use a bin_widths parameter shaped [32].

Parameters will be broadcast against each other and against the input x/ys, so if we want fixed slopes, we can use kwarg knot_slopes=1. A typical recipe for acquiring compatible bin widths and heights would be:

```
nbins <- unconstrained_vector$shape[-1]
range_min <- 1
range_max <- 1
min_bin_size = 1e-2
scale <- range_max - range_min - nbins * min_bin_size
bin_widths = tf$math$softmax(unconstrained_vector) * scale + min_bin_size
```

Value

a bijector instance.

References

- [Conor Durkan, Artur Bekasov, Iain Murray, George Papamakarios. Neural Spline Flows. arXiv preprint arXiv:1906.04032, 2019.](#)

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_temp`, `tfb_masked_autoregressive_flow()`, `tfb_masked_dense()`, `tfb_matrix_inverse_tri_l()`, `tfb_matvec_lu()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_pad()`, `tfb_permute()`, `tfb_power_transform()`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`, `tfb_real_nvp()`, `tfb_reciprocal()`, `tfb_reshape()`, `tfb_scale_matvec_diag()`, `tfb_scale_matvec_linear_operator()`, `tfb_scale_matvec_lu()`, `tfb_scale_matvec_tri_l()`, `tfb_scale_tri_l()`, `tfb_scale()`, `tfb_shifted_gompertz_cdf()`, `tfb_shift()`, `tfb_sigmoid()`, `tfb_sinh_arcsinh()`, `tfb_sinh()`, `tfb_softmax_centered()`, `tfb_softplus()`, `tfb_softsign()`, `tfb_split()`, `tfb_square()`, `tfb_tanh()`, `tfb_transform_diagonal()`, `tfb_transpose()`, `tfb_weibull_cdf()`, `tfb_weibull()`

<code>tfb_rayleigh_cdf</code>	<i>Compute $Y = g(X) = 1 - \exp(-(X/\text{scale})^{**2} / 2)$, $X \geq 0$.</i>
-------------------------------	--

Description

This bijector maps inputs from $[0, \infty]$ to $[0, 1]$. The inverse of the bijector applied to a uniform random variable $X \sim U(0, 1)$ gives back a random variable with the **Rayleigh distribution**:

$Y \sim \text{Rayleigh}(\text{scale})$

$\text{pdf}(y; \text{scale}, y \geq 0) = (1 / \text{scale}) * (y / \text{scale}) * \exp(-(y / \text{scale})^{**2} / 2)$

Usage

```
tfb_rayleigh_cdf(scale, validate_args = FALSE, name = "rayleigh_cdf")
```

Arguments

<code>scale</code>	Positive floating-point tensor. This is l in $Y = g(X) = 1 - \exp(-(X/l)^{**2} / 2)$, $X \geq 0$.
<code>validate_args</code>	Logical, default <code>FALSE</code> . Whether to validate input with asserts. If <code>validate_args</code> is <code>FALSE</code> , and the inputs are invalid, correct behavior is not guaranteed.
<code>name</code>	name prefixed to Ops created by this class.

Details

Likewise, the forward of this bijector is the Rayleigh distribution CDF.

Value

a bijector instance.

See Also

For usage examples see [tfb_forward\(\)](#), [tfb_inverse\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#).

Other bijectors: [tfb_absolute_value\(\)](#), [tfb_affine_linear_operator\(\)](#), [tfb_affine_scalar\(\)](#), [tfb_affine\(\)](#), [tfb_ascending\(\)](#), [tfb_batch_normalization\(\)](#), [tfb_blockwise\(\)](#), [tfb_chain\(\)](#), [tfb_cholesky_outer_product\(\)](#), [tfb_cholesky_to_inv_cholesky\(\)](#), [tfb_correlation_cholesky\(\)](#), [tfb_cumsum\(\)](#), [tfb_discrete_cosine_transform\(\)](#), [tfb_expm1\(\)](#), [tfb_exp\(\)](#), [tfb_ffjord\(\)](#), [tfb_fill_scale_tri_l\(\)](#), [tfb_fill_triangular\(\)](#), [tfb_glow\(\)](#), [tfb_gompertz_cdf\(\)](#), [tfb_gumbel_cdf\(\)](#), [tfb_gumbel\(\)](#), [tfb_identity\(\)](#), [tfb_inline\(\)](#), [tfb_invert\(\)](#), [tfb_iterated_sigmoid_centered\(\)](#), [tfb_kumaraswamy_cdf\(\)](#), [tfb_kumaraswamy\(\)](#), [tfb_lambert_w_tail\(\)](#), [tfb_masked_autoregressive_default_template\(\)](#), [tfb_masked_autoregressive_flow\(\)](#), [tfb_masked_dense\(\)](#), [tfb_matrix_inverse_tri_l\(\)](#), [tfb_matvec_lu\(\)](#), [tfb_normal_cdf\(\)](#), [tfb_ordered\(\)](#), [tfb_pad\(\)](#), [tfb_permute\(\)](#), [tfb_power_transform\(\)](#), [tfb_rational_quadratic_spline\(\)](#), [tfb_real_nvp_default_template\(\)](#), [tfb_real_nvp\(\)](#), [tfb_reciprocal\(\)](#), [tfb_reshape\(\)](#), [tfb_scale_matvec_diag\(\)](#), [tfb_scale_matvec_linear_operator\(\)](#), [tfb_scale_matvec_lu\(\)](#), [tfb_scale_matvec_tri_l\(\)](#), [tfb_scale_tri_l\(\)](#), [tfb_scale\(\)](#), [tfb_shifted_gompertz_cdf\(\)](#), [tfb_shift\(\)](#), [tfb_sigmoid\(\)](#), [tfb_sinh_arcsinh\(\)](#), [tfb_sinh\(\)](#), [tfb_softmax_centered\(\)](#), [tfb_softplus\(\)](#), [tfb_softsign\(\)](#), [tfb_split\(\)](#), [tfb_square\(\)](#), [tfb_tanh\(\)](#), [tfb_transform_diagonal\(\)](#), [tfb_transpose\(\)](#), [tfb_weibull_cdf\(\)](#), [tfb_weibull\(\)](#)

tfb_real_nvp

RealNVP affine coupling layer for vector-valued events

Description

Real NVP models a normalizing flow on a D-dimensional distribution via a single D-d-dimensional conditional distribution (Dinh et al., 2017): $y[d:D] = x[d:D] * \text{tf.exp}(\text{log_scale_fn}(x[0:d])) + \text{shift_fn}(x[0:d])$ $y[0:d] = x[0:d]$ The last D-d units are scaled and shifted based on the first d units only, while the first d units are 'masked' and left unchanged. Real NVP's `shift_and_log_scale_fn` computes vector-valued quantities. For scale-and-shift transforms that do not depend on any masked units, i.e. `d=0`, use the `tfb_affine` bijector with learned parameters instead. Masking is currently only supported for base distributions with `event_ndims=1`. For more sophisticated masking schemes like checkerboard or channel-wise masking (Papamakarios et al., 2016), use the `tfb_permute` bijector to re-order desired masked units into the first d units. For base distributions with `event_ndims > 1`, use the `tfb_reshape` bijector to flatten the event shape.

Usage

```
tfb_real_nvp(
    num_masked,
    shift_and_log_scale_fn,
    is_constant_jacobian = FALSE,
    validate_args = FALSE,
    name = NULL
)
```

Arguments

num_masked	integer indicating that the first d units of the event should be masked. Must be in the closed interval $[1, D-1]$, where D is the event size of the base distribution.
shift_and_log_scale_fn	Function which computes <code>shift</code> and <code>log_scale</code> from both the forward domain (x) and the inverse domain (y). Calculation must respect the "autoregressive property". Suggested default: <code>tfb_real_nvp_default_template(hidden_layers=...)</code> . Typically the function contains <code>tf\$Variables</code> and is wrapped using <code>tf\$make_template</code> . Returning NULL for either (both) <code>shift</code> , <code>log_scale</code> is equivalent to (but more efficient than) returning zero.
is_constant_jacobian	Logical, default: FALSE. When TRUE the implementation assumes <code>log_scale</code> does not depend on the forward domain (x) or inverse domain (y) values. (No validation is made; <code>is_constant_jacobian=FALSE</code> is always safe but possibly computationally inefficient.)
validate_args	Logical, default FALSE. Whether to validate input with asserts. If <code>validate_args</code> is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
name	name prefixed to Ops created by this class.

Details

Recall that the MAF bijector (Papamakarios et al., 2016) implements a normalizing flow via an autoregressive transformation. MAF and IAF have opposite computational tradeoffs - MAF can train all units in parallel but must sample units sequentially, while IAF must train units sequentially but can sample in parallel. In contrast, Real NVP can compute both forward and inverse computations in parallel. However, the lack of an autoregressive transformations makes it less expressive on a per-bijector basis.

A "valid" `shift_and_log_scale_fn` must compute each `shift` (aka `loc` or " μ " in Papamakarios et al. (2016) and `log(scale)` (aka " α " in Papamakarios et al. (2016)) such that each are broadcastable with the arguments to forward and inverse, i.e., such that the calculations in forward, inverse below are possible. For convenience, `real_nvp_default_nvp` is offered as a possible `shift_and_log_scale_fn` function.

NICE (Dinh et al., 2014) is a special case of the Real NVP bijector which discards the scale transformation, resulting in a constant-time inverse-log-determinant-Jacobian. To use a NICE bijector instead of Real NVP, `shift_and_log_scale_fn` should return (`shift`, NULL), and `is_constant_jacobian` should be set to TRUE in the RealNVP constructor. Calling `tfb_real_nvp_default_template` with `shift_only=TRUE` returns one such NICE-compatible `shift_and_log_scale_fn`.

Caching: the scalar input depth D of the base distribution is not known at construction time. The first call to any of `forward(x)`, `inverse(x)`, `inverse_log_det_jacobian(x)`, or `forward_log_det_jacobian(x)` memoizes D , which is re-used in subsequent calls. This shape must be known prior to graph execution (which is the case if using `tf$layers`).

Value

a bijector instance.

References

- George Papamakarios, Theo Pavlakou, and Iain Murray. Masked Autoregressive Flow for Density Estimation. In *Neural Information Processing Systems*, 2017.
- Laurent Dinh, Jascha Sohl-Dickstein, and Samy Bengio. Density Estimation using Real NVP. In *International Conference on Learning Representations*, 2017.
- Laurent Dinh, David Krueger, and Yoshua Bengio. NICE: Non-linear Independent Components Estimation. *arXiv preprint arXiv:1410.8516*, 2014.
- Eric Jang. Normalizing Flows Tutorial, Part 2: Modern Normalizing Flows. Technical Report, 2018.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_template()`, `tfb_masked_autoregressive_flow()`, `tfb_masked_dense()`, `tfb_matrix_inverse_tri_l()`, `tfb_matvec_lu()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_pad()`, `tfb_permute()`, `tfb_power_transform()`, `tfb_rational_quadratic_spline()`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`, `tfb_reciprocal()`, `tfb_reshape()`, `tfb_scale_matvec_diag()`, `tfb_scale_matvec_linear_operator()`, `tfb_scale_matvec_lu()`, `tfb_scale_matvec_tri_l()`, `tfb_scale_tri_l()`, `tfb_scale()`, `tfb_shifted_gompertz_cdf()`, `tfb_shift()`, `tfb_sigmoid()`, `tfb_sinh_arcsinh()`, `tfb_sinh()`, `tfb_softmax_centered()`, `tfb_softplus()`, `tfb_softsign()`, `tfb_split()`, `tfb_square()`, `tfb_tanh()`, `tfb_transform_diagonal()`, `tfb_transpose()`, `tfb_weibull_cdf()`, `tfb_weibull()`

tfb_real_nvp_default_template

Build a scale-and-shift function using a multi-layer neural network

Description

This will be wrapped in a `make_template` to ensure the variables are only created once. It takes the d -dimensional input $x[0:d]$ and returns the $D-d$ dimensional outputs `loc` ("mu") and `log_scale` ("alpha").

Usage

```
tfb_real_nvp_default_template(
    hidden_layers,
    shift_only = FALSE,
    activation = tf$nn$relu,
```

```

    name = NULL,
    ...
)

```

Arguments

hidden_layers	list-like of non-negative integer, scalars indicating the number of units in each hidden layer. Default: list(512, 512).
shift_only	logical indicating if only the shift term shall be computed (i.e. NICE bijector). Default: FALSE.
activation	Activation function (callable). Explicitly setting to NULL implies a linear activation.
name	A name for ops managed by this function. Default: "tfb_real_nvp_default_template".
...	tf\$layers\$dense arguments

Details

The default template does not support conditioning and will raise an exception if `condition_kwargs` are passed to it. To use conditioning in real nvp bijector, implement a conditioned shift/scale template that handles the `condition_kwargs`.

Value

list of:

- shift: Float-like Tensor of shift terms
- log_scale: Float-like Tensor of log(scale) terms

References

- [George Papamakarios, Theo Pavlakou, and Iain Murray. Masked Autoregressive Flow for Density Estimation. In *Neural Information Processing Systems*, 2017.](#)

See Also

For usage examples see [tfb_forward\(\)](#), [tfb_inverse\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#).

Other bijectors: [tfb_absolute_value\(\)](#), [tfb_affine_linear_operator\(\)](#), [tfb_affine_scalar\(\)](#), [tfb_affine\(\)](#), [tfb_ascending\(\)](#), [tfb_batch_normalization\(\)](#), [tfb_blockwise\(\)](#), [tfb_chain\(\)](#), [tfb_cholesky_outer_product\(\)](#), [tfb_cholesky_to_inv_cholesky\(\)](#), [tfb_correlation_cholesky\(\)](#), [tfb_cumsum\(\)](#), [tfb_discrete_cosine_transform\(\)](#), [tfb_expm1\(\)](#), [tfb_exp\(\)](#), [tfb_ffjord\(\)](#), [tfb_fill_scale_tri_l\(\)](#), [tfb_fill_triangular\(\)](#), [tfb_glow\(\)](#), [tfb_gompertz_cdf\(\)](#), [tfb_gumbel_cdf\(\)](#), [tfb_gumbel\(\)](#), [tfb_identity\(\)](#), [tfb_inline\(\)](#), [tfb_invert\(\)](#), [tfb_iterated_sigmoid_centered\(\)](#), [tfb_kumaraswamy_cdf\(\)](#), [tfb_kumaraswamy\(\)](#), [tfb_lambert_w_tail\(\)](#), [tfb_masked_autoregressive_default_template\(\)](#), [tfb_masked_autoregressive_flow\(\)](#), [tfb_masked_dense\(\)](#), [tfb_matrix_inverse_tri_l\(\)](#), [tfb_matvec_lu\(\)](#), [tfb_normal_cdf\(\)](#), [tfb_ordered\(\)](#), [tfb_pad\(\)](#), [tfb_permute\(\)](#), [tfb_power_transform\(\)](#), [tfb_rational_quadratic_spline\(\)](#), [tfb_rayleigh_cdf\(\)](#), [tfb_real_nvp\(\)](#), [tfb_reciprocal\(\)](#), [tfb_reshape\(\)](#), [tfb_scale_matvec_diag\(\)](#), [tfb_scale_matvec_linear_operator\(\)](#), [tfb_scale_matvec_lu\(\)](#), [tfb_scale_matvec_tri_l\(\)](#), [tfb_scale_tri_l\(\)](#), [tfb_scale\(\)](#), [tfb_shifted_gompertz_cdf\(\)](#),

[tfb_shift\(\)](#), [tfb_sigmoid\(\)](#), [tfb_sinh_arcsinh\(\)](#), [tfb_sinh\(\)](#), [tfb_softmax_centered\(\)](#), [tfb_softplus\(\)](#), [tfb_softsign\(\)](#), [tfb_split\(\)](#), [tfb_square\(\)](#), [tfb_tanh\(\)](#), [tfb_transform_diagonal\(\)](#), [tfb_transpose\(\)](#), [tfb_weibull_cdf\(\)](#), [tfb_weibull\(\)](#)

tfb_reciprocal	<i>A Bijector that computes $b(x) = 1. / x$</i>
----------------	--

Description

A Bijector that computes $b(x) = 1. / x$

Usage

```
tfb_reciprocal(validate_args = FALSE, name = "reciprocal")
```

Arguments

`validate_args` Logical, default FALSE. Whether to validate input with asserts. If `validate_args` is FALSE, and the inputs are invalid, correct behavior is not guaranteed.

`name` name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see [tfb_forward\(\)](#), [tfb_inverse\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#).

Other bijectors: [tfb_absolute_value\(\)](#), [tfb_affine_linear_operator\(\)](#), [tfb_affine_scalar\(\)](#), [tfb_affine\(\)](#), [tfb_ascending\(\)](#), [tfb_batch_normalization\(\)](#), [tfb_blockwise\(\)](#), [tfb_chain\(\)](#), [tfb_cholesky_outer_product\(\)](#), [tfb_cholesky_to_inv_cholesky\(\)](#), [tfb_correlation_cholesky\(\)](#), [tfb_cumsum\(\)](#), [tfb_discrete_cosine_transform\(\)](#), [tfb_expm1\(\)](#), [tfb_exp\(\)](#), [tfb_ffjord\(\)](#), [tfb_fill_scale_tri_l\(\)](#), [tfb_fill_triangular\(\)](#), [tfb_glow\(\)](#), [tfb_gompertz_cdf\(\)](#), [tfb_gumbel_cdf\(\)](#), [tfb_gumbel\(\)](#), [tfb_identity\(\)](#), [tfb_inline\(\)](#), [tfb_invert\(\)](#), [tfb_iterated_sigmoid_centered\(\)](#), [tfb_kumaraswamy_cdf\(\)](#), [tfb_kumaraswamy\(\)](#), [tfb_lambert_w_tail\(\)](#), [tfb_masked_autoregressive_default_template\(\)](#), [tfb_masked_autoregressive_flow\(\)](#), [tfb_masked_dense\(\)](#), [tfb_matrix_inverse_tri_l\(\)](#), [tfb_matvec_lu\(\)](#), [tfb_normal_cdf\(\)](#), [tfb_ordered\(\)](#), [tfb_pad\(\)](#), [tfb_permute\(\)](#), [tfb_power_transform\(\)](#), [tfb_rational_quadratic_spline\(\)](#), [tfb_rayleigh_cdf\(\)](#), [tfb_real_nvp_default_template\(\)](#), [tfb_real_nvp\(\)](#), [tfb_reshape\(\)](#), [tfb_scale_matvec_diag\(\)](#), [tfb_scale_matvec_linear_operator\(\)](#), [tfb_scale_matvec_lu\(\)](#), [tfb_scale_matvec_tri_l\(\)](#), [tfb_scale_tri_l\(\)](#), [tfb_scale\(\)](#), [tfb_shifted_gompertz_cdf\(\)](#), [tfb_shift\(\)](#), [tfb_sigmoid\(\)](#), [tfb_sinh_arcsinh\(\)](#), [tfb_sinh\(\)](#), [tfb_softmax_centered\(\)](#), [tfb_softplus\(\)](#), [tfb_softsign\(\)](#), [tfb_split\(\)](#), [tfb_square\(\)](#), [tfb_tanh\(\)](#), [tfb_transform_diagonal\(\)](#), [tfb_transpose\(\)](#), [tfb_weibull_cdf\(\)](#), [tfb_weibull\(\)](#)

tfb_reshape

*Reshapes the event_shape of a Tensor***Description**

The semantics generally follow that of `tf$reshape()`, with a few differences:

- The user must provide both the input and output shape, so that the transformation can be inverted. If an input shape is not specified, the default assumes a vector-shaped input, i.e., `event_shape_in = list(-1)`.
- The Reshape bijector automatically broadcasts over the leftmost dimensions of its input (`sample_shape` and `batch_shape`); only the rightmost `event_ndims_in` dimensions are reshaped. The number of dimensions to reshape is inferred from the provided `event_shape_in` (`event_ndims_in = length(event_shape_in)`).

Usage

```
tfb_reshape(
  event_shape_out,
  event_shape_in = c(-1),
  validate_args = FALSE,
  name = NULL
)
```

Arguments

<code>event_shape_out</code>	An integer-like vector-shaped Tensor representing the event shape of the transformed output.
<code>event_shape_in</code>	An optional integer-like vector-shaped Tensor representing the event shape of the input. This is required in order to define inverse operations; the default of <code>list(-1)</code> assumes a vector-shaped input.
<code>validate_args</code>	Logical, default <code>FALSE</code> . Whether to validate input with asserts. If <code>validate_args</code> is <code>FALSE</code> , and the inputs are invalid, correct behavior is not guaranteed.
<code>name</code>	name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`,

tfb_fill_scale_tri_l(), tfb_fill_triangular(), tfb_glow(), tfb_gompertz_cdf(), tfb_gumbel_cdf(), tfb_gumbel(), tfb_identity(), tfb_inline(), tfb_invert(), tfb_iterated_sigmoid_centered(), tfb_kumaraswamy_cdf(), tfb_kumaraswamy(), tfb_lambert_w_tail(), tfb_masked_autoregressive_default_template(), tfb_masked_autoregressive_flow(), tfb_masked_dense(), tfb_matrix_inverse_tri_l(), tfb_matvec_lu(), tfb_normal_cdf(), tfb_ordered(), tfb_pad(), tfb_permute(), tfb_power_transform(), tfb_rational_quadratic_spline(), tfb_rayleigh_cdf(), tfb_real_nvp_default_template(), tfb_real_nvp(), tfb_reciprocal(), tfb_scale_matvec_diag(), tfb_scale_matvec_linear_operator(), tfb_scale_matvec_lu(), tfb_scale_matvec_tri_l(), tfb_scale_tri_l(), tfb_scale(), tfb_shifted_gompertz_cdf(), tfb_shift(), tfb_sigmoid(), tfb_sinh_arcsinh(), tfb_sinh(), tfb_softmax_centered(), tfb_softplus(), tfb_softsign(), tfb_split(), tfb_square(), tfb_tanh(), tfb_transform_diagonal(), tfb_transpose(), tfb_weibull_cdf(), tfb_weibull()

 tfb_scale

 Compute $Y = g(X; \text{scale}) = \text{scale} * X$.

Description

Examples:

```
Y <- 2 * X
b <- tfb_scale(scale = 2)
```

Usage

```
tfb_scale(
  scale = NULL,
  log_scale = NULL,
  validate_args = FALSE,
  name = "scale"
)
```

Arguments

scale	Floating-point Tensor.
log_scale	Floating-point Tensor. Logarithm of the scale. If this is set to NULL, no scale is applied. This should not be set if scale is set.
validate_args	Logical, default FALSE. Whether to validate input with asserts. If validate_args is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
name	name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_template()`, `tfb_masked_autoregressive_flow()`, `tfb_masked_dense()`, `tfb_matrix_inverse_tri_l()`, `tfb_matvec_lu()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_pad()`, `tfb_permute()`, `tfb_power_transform()`, `tfb_rational_quadratic_spline()`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`, `tfb_real_nvp()`, `tfb_reciprocal()`, `tfb_reshape()`, `tfb_scale_matvec_diag()`, `tfb_scale_matvec_linear_operator()`, `tfb_scale_matvec_lu()`, `tfb_scale_matvec_tri_l()`, `tfb_scale_tri_l()`, `tfb_shifted_gompertz_cdf()`, `tfb_shift()`, `tfb_sigmoid()`, `tfb_sinh_arcsinh()`, `tfb_sinh()`, `tfb_softmax_centered()`, `tfb_softplus()`, `tfb_softsign()`, `tfb_split()`, `tfb_square()`, `tfb_tanh()`, `tfb_transform_diagonal()`, `tfb_transpose()`, `tfb_weibull_cdf()`, `tfb_weibull()`

`tfb_scale_matvec_diag` Compute $Y = g(X; \text{scale}) = \text{scale} @ X$

Description

In TF parlance, the scale term is logically equivalent to:

```
scale = tf$diag(scale_diag)
```

The scale term is applied without materializing a full dense matrix.

Usage

```
tfb_scale_matvec_diag(
  scale_diag,
  adjoint = FALSE,
  validate_args = FALSE,
  name = "scale_matvec_diag",
  dtype = NULL
)
```

Arguments

<code>scale_diag</code>	Floating-point Tensor representing the diagonal matrix. <code>scale_diag</code> has shape <code>[N1, N2, ..., k]</code> , which represents a <code>k x k</code> diagonal matrix.
<code>adjoint</code>	logical indicating whether to use the scale matrix as specified or its adjoint. Default value: <code>FALSE</code> .

validate_args	Logical, default FALSE. Whether to validate input with asserts. If validate_args is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
name	name prefixed to Ops created by this class.
dtype	tf\$DType to prefer when converting args to Tensors. Else, we fall back to a common dtype inferred from the args, finally falling back to float32.

Value

a bijector instance.

See Also

For usage examples see [tfb_forward\(\)](#), [tfb_inverse\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#).

Other bijectors: [tfb_absolute_value\(\)](#), [tfb_affine_linear_operator\(\)](#), [tfb_affine_scalar\(\)](#), [tfb_affine\(\)](#), [tfb_ascending\(\)](#), [tfb_batch_normalization\(\)](#), [tfb_blockwise\(\)](#), [tfb_chain\(\)](#), [tfb_cholesky_outer_product\(\)](#), [tfb_cholesky_to_inv_cholesky\(\)](#), [tfb_correlation_cholesky\(\)](#), [tfb_cumsum\(\)](#), [tfb_discrete_cosine_transform\(\)](#), [tfb_expm1\(\)](#), [tfb_exp\(\)](#), [tfb_ffjord\(\)](#), [tfb_fill_scale_tri_l\(\)](#), [tfb_fill_triangular\(\)](#), [tfb_glow\(\)](#), [tfb_gompertz_cdf\(\)](#), [tfb_gumbel_cdf\(\)](#), [tfb_gumbel\(\)](#), [tfb_identity\(\)](#), [tfb_inline\(\)](#), [tfb_invert\(\)](#), [tfb_iterated_sigmoid_centered\(\)](#), [tfb_kumaraswamy_cdf\(\)](#), [tfb_kumaraswamy\(\)](#), [tfb_lambert_w_tail\(\)](#), [tfb_masked_autoregressive_default_template\(\)](#), [tfb_masked_autoregressive_flow\(\)](#), [tfb_masked_dense\(\)](#), [tfb_matrix_inverse_tri_l\(\)](#), [tfb_matvec_lu\(\)](#), [tfb_normal_cdf\(\)](#), [tfb_ordered\(\)](#), [tfb_pad\(\)](#), [tfb_permute\(\)](#), [tfb_power_transform\(\)](#), [tfb_rational_quadratic_spline\(\)](#), [tfb_rayleigh_cdf\(\)](#), [tfb_real_nvp_default_template\(\)](#), [tfb_real_nvp\(\)](#), [tfb_reciprocal\(\)](#), [tfb_reshape\(\)](#), [tfb_scale_matvec_linear_operator\(\)](#), [tfb_scale_matvec_lu\(\)](#), [tfb_scale_matvec_tri_l\(\)](#), [tfb_scale_tri_l\(\)](#), [tfb_scale\(\)](#), [tfb_shifted_gompertz_cdf\(\)](#), [tfb_shift\(\)](#), [tfb_sigmoid\(\)](#), [tfb_sinh_arcsinh\(\)](#), [tfb_sinh\(\)](#), [tfb_softmax_centered\(\)](#), [tfb_softplus\(\)](#), [tfb_softsign\(\)](#), [tfb_split\(\)](#), [tfb_square\(\)](#), [tfb_tanh\(\)](#), [tfb_transform_diagonal\(\)](#), [tfb_transpose\(\)](#), [tfb_weibull_cdf\(\)](#), [tfb_weibull\(\)](#)

tfb_scale_matvec_linear_operator

Compute $Y = g(X; \text{scale}) = \text{scale} @ X$.

Description

scale is a LinearOperator. If X is a scalar then the forward transformation is: scale * X where * denotes broadcasted elementwise product.

Usage

```
tfb_scale_matvec_linear_operator(
  scale,
  adjoint = FALSE,
  validate_args = FALSE,
  name = "scale_matvec_linear_operator"
)
```

Arguments

scale	Subclass of LinearOperator. Represents the (batch, non-singular) linear transformation by which the Bijector transforms inputs.
adjoint	logical indicating whether to use the scale matrix as specified or its adjoint. Default value: FALSE.
validate_args	Logical, default FALSE. Whether to validate input with asserts. If validate_args is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
name	name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see [tfb_forward\(\)](#), [tfb_inverse\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#).

Other bijectors: [tfb_absolute_value\(\)](#), [tfb_affine_linear_operator\(\)](#), [tfb_affine_scalar\(\)](#), [tfb_affine\(\)](#), [tfb_ascending\(\)](#), [tfb_batch_normalization\(\)](#), [tfb_blockwise\(\)](#), [tfb_chain\(\)](#), [tfb_cholesky_outer_product\(\)](#), [tfb_cholesky_to_inv_cholesky\(\)](#), [tfb_correlation_cholesky\(\)](#), [tfb_cumsum\(\)](#), [tfb_discrete_cosine_transform\(\)](#), [tfb_expm1\(\)](#), [tfb_exp\(\)](#), [tfb_ffjord\(\)](#), [tfb_fill_scale_tri_l\(\)](#), [tfb_fill_triangular\(\)](#), [tfb_glow\(\)](#), [tfb_gompertz_cdf\(\)](#), [tfb_gumbel_cdf\(\)](#), [tfb_gumbel\(\)](#), [tfb_identity\(\)](#), [tfb_inline\(\)](#), [tfb_invert\(\)](#), [tfb_iterated_sigmoid_centered\(\)](#), [tfb_kumaraswamy_cdf\(\)](#), [tfb_kumaraswamy\(\)](#), [tfb_lambert_w_tail\(\)](#), [tfb_masked_autoregressive_default_temp\(\)](#), [tfb_masked_autoregressive_flow\(\)](#), [tfb_masked_dense\(\)](#), [tfb_matrix_inverse_tri_l\(\)](#), [tfb_matvec_lu\(\)](#), [tfb_normal_cdf\(\)](#), [tfb_ordered\(\)](#), [tfb_pad\(\)](#), [tfb_permute\(\)](#), [tfb_power_transform\(\)](#), [tfb_rational_quadratic_spline\(\)](#), [tfb_rayleigh_cdf\(\)](#), [tfb_real_nvp_default_template\(\)](#), [tfb_real_nvp\(\)](#), [tfb_reciprocal\(\)](#), [tfb_reshape\(\)](#), [tfb_scale_matvec_diag\(\)](#), [tfb_scale_matvec_lu\(\)](#), [tfb_scale_matvec_tri_l\(\)](#), [tfb_scale_tri_l\(\)](#), [tfb_scale\(\)](#), [tfb_shifted_gompertz_cdf\(\)](#), [tfb_shift\(\)](#), [tfb_sigmoid\(\)](#), [tfb_sinh_arcsinh\(\)](#), [tfb_sinh\(\)](#), [tfb_softmax_centered\(\)](#), [tfb_softplus\(\)](#), [tfb_softsign\(\)](#), [tfb_split\(\)](#), [tfb_square\(\)](#), [tfb_tanh\(\)](#), [tfb_transform_diagonal\(\)](#), [tfb_transpose\(\)](#), [tfb_weibull_cdf\(\)](#), [tfb_weibull\(\)](#)

tfb_scale_matvec_lu *Matrix-vector multiply using LU decomposition.*

Description

This bijector is identical to the "Convolution1x1" used in Glow (Kingma and Dhariwal, 2018).

Usage

```
tfb_scale_matvec_lu(
  lower_upper,
  permutation,
  validate_args = FALSE,
  name = NULL
)
```

Arguments

lower_upper	The LU factorization as returned by <code>tf.linalg.lu</code> .
permutation	The LU factorization permutation as returned by <code>tf.linalg.lu</code> .
validate_args	Logical, default <code>FALSE</code> . Whether to validate input with asserts. If <code>validate_args</code> is <code>FALSE</code> , and the inputs are invalid, correct behavior is not guaranteed.
name	name prefixed to Ops created by this class.

Value

a bijector instance.

References

- [Diederik P. Kingma, Prafulla Dhariwal. Glow: Generative Flow with Invertible 1x1 Convolutions. *arXiv preprint arXiv:1807.03039*, 2018.](#)

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_template()`, `tfb_masked_autoregressive_flow()`, `tfb_masked_dense()`, `tfb_matrix_inverse_tri_l()`, `tfb_matvec_lu()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_pad()`, `tfb_permute()`, `tfb_power_transform()`, `tfb_rational_quadratic_spline()`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`, `tfb_real_nvp()`, `tfb_reciprocal()`, `tfb_reshape()`, `tfb_scale_matvec_diag()`, `tfb_scale_matvec_linear_operator()`, `tfb_scale_matvec_tri_l()`, `tfb_scale_tri_l()`, `tfb_scale()`, `tfb_shifted_gompertz_cdf()`, `tfb_shift()`, `tfb_sigmoid()`, `tfb_sinh_arcsinh()`, `tfb_sinh()`, `tfb_softmax_centered()`, `tfb_softplus()`, `tfb_softsign()`, `tfb_split()`, `tfb_square()`, `tfb_tanh()`, `tfb_transform_diagonal()`, `tfb_transpose()`, `tfb_weibull_cdf()`, `tfb_weibull()`

`tfb_scale_matvec_tri_l`

Compute $Y = g(X; \text{scale}) = \text{scale} @ X$.

Description

The scale term is presumed lower-triangular and non-singular (ie, no zeros on the diagonal), which permits efficient determinant calculation (linear in matrix dimension, instead of cubic).

Usage

```
tfb_scale_matvec_tri_l(
    scale_tril,
    adjoint = FALSE,
    validate_args = FALSE,
    name = "scale_matvec_tril",
    dtype = NULL
)
```

Arguments

scale_tril	Floating-point Tensor representing the lower triangular matrix. scale_tril has shape [N1, N2, ... k, k], which represents a k x k lower triangular matrix. When NULL no scale_tril term is added to scale. The upper triangular elements above the diagonal are ignored.
adjoint	logical indicating whether to use the scale matrix as specified or its adjoint. Note that lower-triangularity is taken into account first: the region above the diagonal of scale_tril is treated as zero (irrespective of the adjoint setting). A lower-triangular input with adjoint=TRUE will behave like an upper triangular transform. Default value: FALSE.
validate_args	Logical, default FALSE. Whether to validate input with asserts. If validate_args is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
name	name prefixed to Ops created by this class.
dtype	tf\$DType to prefer when converting args to Tensors. Else, we fall back to a common dtype inferred from the args, finally falling back to float32.

Value

a bijector instance.

See Also

For usage examples see [tfb_forward\(\)](#), [tfb_inverse\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#).

Other bijectors: [tfb_absolute_value\(\)](#), [tfb_affine_linear_operator\(\)](#), [tfb_affine_scalar\(\)](#), [tfb_affine\(\)](#), [tfb_ascending\(\)](#), [tfb_batch_normalization\(\)](#), [tfb_blockwise\(\)](#), [tfb_chain\(\)](#), [tfb_cholesky_outer_product\(\)](#), [tfb_cholesky_to_inv_cholesky\(\)](#), [tfb_correlation_cholesky\(\)](#), [tfb_cumsum\(\)](#), [tfb_discrete_cosine_transform\(\)](#), [tfb_expm1\(\)](#), [tfb_exp\(\)](#), [tfb_ffjord\(\)](#), [tfb_fill_scale_tri_l\(\)](#), [tfb_fill_triangular\(\)](#), [tfb_glow\(\)](#), [tfb_gompertz_cdf\(\)](#), [tfb_gumbel_cdf\(\)](#), [tfb_gumbel\(\)](#), [tfb_identity\(\)](#), [tfb_inline\(\)](#), [tfb_invert\(\)](#), [tfb_iterated_sigmoid_centered\(\)](#), [tfb_kumaraswamy_cdf\(\)](#), [tfb_kumaraswamy\(\)](#), [tfb_lambert_w_tail\(\)](#), [tfb_masked_autoregressive_default_template\(\)](#), [tfb_masked_autoregressive_flow\(\)](#), [tfb_masked_dense\(\)](#), [tfb_matrix_inverse_tri_l\(\)](#), [tfb_matvec_lu\(\)](#), [tfb_normal_cdf\(\)](#), [tfb_ordered\(\)](#), [tfb_pad\(\)](#), [tfb_permute\(\)](#), [tfb_power_transform\(\)](#), [tfb_rational_quadratic_spline\(\)](#), [tfb_rayleigh_cdf\(\)](#), [tfb_real_nvp_default_template\(\)](#), [tfb_real_nvp\(\)](#), [tfb_reciprocal\(\)](#), [tfb_reshape\(\)](#), [tfb_scale_matvec_diag\(\)](#), [tfb_scale_matvec_linear_operator\(\)](#), [tfb_scale_matvec_lu\(\)](#), [tfb_scale_tri_l\(\)](#), [tfb_scale\(\)](#), [tfb_shifted_gompertz_cdf\(\)](#), [tfb_shift\(\)](#), [tfb_sigmoid\(\)](#), [tfb_sinh_arcsinh\(\)](#), [tfb_sinh\(\)](#), [tfb_softmax_centered\(\)](#), [tfb_softplus\(\)](#), [tfb_softsign\(\)](#), [tfb_split\(\)](#), [tfb_square\(\)](#), [tfb_tanh\(\)](#), [tfb_transform_diagonal\(\)](#), [tfb_transpose\(\)](#), [tfb_weibull_cdf\(\)](#), [tfb_weibull\(\)](#)

tfb_scale_tri_l	<i>Transforms unconstrained vectors to TriL matrices with positive diagonal</i>
-----------------	---

Description

This is implemented as a simple `tfb_chain` of `tfb_fill_triangular` followed by `tfb_transform_diagonal`, and provided mostly as a convenience. The default setup is somewhat opinionated, using a Softplus transformation followed by a small shift (1e-5) which attempts to avoid numerical issues from zeros on the diagonal.

Usage

```
tfb_scale_tri_l(
  diag_bijector = NULL,
  diag_shift = 1e-05,
  validate_args = FALSE,
  name = "scale_tril"
)
```

Arguments

<code>diag_bijector</code>	Bijector instance, used to transform the output diagonal to be positive. Default value: <code>NULL</code> (i.e., <code>tfb_softplus()</code>).
<code>diag_shift</code>	Float value broadcastable and added to all diagonal entries after applying the <code>diag_bijector</code> . Setting a positive value forces the output diagonal entries to be positive, but prevents inverting the transformation for matrices with diagonal entries less than this value. Default value: <code>1e-5</code> .
<code>validate_args</code>	Logical, default <code>FALSE</code> . Whether to validate input with asserts. If <code>validate_args</code> is <code>FALSE</code> , and the inputs are invalid, correct behavior is not guaranteed.
<code>name</code>	name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expml()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_temp.`

tfb_masked_autoregressive_flow(), tfb_masked_dense(), tfb_matrix_inverse_tri_l(),
 tfb_matvec_lu(), tfb_normal_cdf(), tfb_ordered(), tfb_pad(), tfb_permute(), tfb_power_transform(),
 tfb_rational_quadratic_spline(), tfb_rayleigh_cdf(), tfb_real_nvp_default_template(),
 tfb_real_nvp(), tfb_reciprocal(), tfb_reshape(), tfb_scale_matvec_diag(), tfb_scale_matvec_linear_operat
 tfb_scale_matvec_lu(), tfb_scale_matvec_tri_l(), tfb_scale(), tfb_shifted_gompertz_cdf(),
 tfb_shift(), tfb_sigmoid(), tfb_sinh_arcsinh(), tfb_sinh(), tfb_softmax_centered(),
 tfb_softplus(), tfb_softsign(), tfb_split(), tfb_square(), tfb_tanh(), tfb_transform_diagonal(),
 tfb_transpose(), tfb_weibull_cdf(), tfb_weibull()

tfb_shift	<i>Compute</i> $Y = g(X; \text{shift}) = X + \text{shift}$.
-----------	--

Description

where shift is a numeric Tensor.

Usage

```
tfb_shift(shift, validate_args = FALSE, name = "shift")
```

Arguments

shift	floating-point tensor
validate_args	Logical, default FALSE. Whether to validate input with asserts. If validate_args is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
name	name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see [tfb_forward\(\)](#), [tfb_inverse\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#).

Other bijectors: [tfb_absolute_value\(\)](#), [tfb_affine_linear_operator\(\)](#), [tfb_affine_scalar\(\)](#),
[tfb_affine\(\)](#), [tfb_ascending\(\)](#), [tfb_batch_normalization\(\)](#), [tfb_blockwise\(\)](#), [tfb_chain\(\)](#),
[tfb_cholesky_outer_product\(\)](#), [tfb_cholesky_to_inv_cholesky\(\)](#), [tfb_correlation_cholesky\(\)](#),
[tfb_cumsum\(\)](#), [tfb_discrete_cosine_transform\(\)](#), [tfb_expm1\(\)](#), [tfb_exp\(\)](#), [tfb_ffjord\(\)](#),
[tfb_fill_scale_tri_l\(\)](#), [tfb_fill_triangular\(\)](#), [tfb_glow\(\)](#), [tfb_gompertz_cdf\(\)](#), [tfb_gumbel_cdf\(\)](#),
[tfb_gumbel\(\)](#), [tfb_identity\(\)](#), [tfb_inline\(\)](#), [tfb_invert\(\)](#), [tfb_iterated_sigmoid_centered\(\)](#),
[tfb_kumaraswamy_cdf\(\)](#), [tfb_kumaraswamy\(\)](#), [tfb_lambert_w_tail\(\)](#), [tfb_masked_autoregressive_default_temp](#)
[tfb_masked_autoregressive_flow\(\)](#), [tfb_masked_dense\(\)](#), [tfb_matrix_inverse_tri_l\(\)](#),
[tfb_matvec_lu\(\)](#), [tfb_normal_cdf\(\)](#), [tfb_ordered\(\)](#), [tfb_pad\(\)](#), [tfb_permute\(\)](#), [tfb_power_transform\(\)](#),
[tfb_rational_quadratic_spline\(\)](#), [tfb_rayleigh_cdf\(\)](#), [tfb_real_nvp_default_template\(\)](#),
[tfb_real_nvp\(\)](#), [tfb_reciprocal\(\)](#), [tfb_reshape\(\)](#), [tfb_scale_matvec_diag\(\)](#), [tfb_scale_matvec_linear_operat](#)
[tfb_scale_matvec_lu\(\)](#), [tfb_scale_matvec_tri_l\(\)](#), [tfb_scale_tri_l\(\)](#), [tfb_scale\(\)](#), [tfb_shifted_gompertz_co](#)
[tfb_sigmoid\(\)](#), [tfb_sinh_arcsinh\(\)](#), [tfb_sinh\(\)](#), [tfb_softmax_centered\(\)](#), [tfb_softplus\(\)](#),

tfb_softsign(), tfb_split(), tfb_square(), tfb_tanh(), tfb_transform_diagonal(), tfb_transpose(),
tfb_weibull_cdf(), tfb_weibull()

tfb_shifted_gompertz_cdf

Compute $Y = g(X) = (1 - \exp(-\text{rate} * X)) * \exp(-c * \exp(-\text{rate} * X))$

Description

This bijector maps inputs from $[-\text{inf}, \text{inf}]$ to $[0, \text{inf}]$. The inverse of the bijector applied to a uniform random variable $X \sim U(0, 1)$ gives back a random variable with the **Shifted Gompertz distribution**:

$Y \sim \text{ShiftedGompertzCDF}(\text{concentration}, \text{rate})$
 $\text{pdf}(y; c, r) = r * \exp(-r * y - \exp(-r * y)) / c * (1 + (1 - \exp(-r * y)) / c)$

Usage

```
tfb_shifted_gompertz_cdf(  
  concentration,  
  rate,  
  validate_args = FALSE,  
  name = "shifted_gompertz_cdf"  
)
```

Arguments

concentration	Positive Float-like Tensor that is the same dtype and is broadcastable with concentration. This is c in $Y = g(X) = (1 - \exp(-\text{rate} * X)) * \exp(-c * \exp(-\text{rate} * X))$.
rate	Positive Float-like Tensor that is the same dtype and is broadcastable with concentration. This is rate in $Y = g(X) = (1 - \exp(-\text{rate} * X)) * \exp(-c * \exp(-\text{rate} * X))$.
validate_args	Logical, default FALSE. Whether to validate input with asserts. If <code>validate_args</code> is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
name	name prefixed to Ops created by this class.

Details

Note: Even though this is called `ShiftedGompertzCDF`, when applied to the Uniform distribution, this is not the same as applying a `GompertzCDF` with a `Shift` bijector (i.e. the Shifted Gompertz distribution is not the same as a Gompertz distribution with a location parameter).

Note: Because the Shifted Gompertz distribution concentrates its mass close to zero, for larger rates or larger concentrations, `bijector$forward` will quickly saturate to 1.

Value

a bijector instance.

See Also

For usage examples see [tfb_forward\(\)](#), [tfb_inverse\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#).

Other bijectors: [tfb_absolute_value\(\)](#), [tfb_affine_linear_operator\(\)](#), [tfb_affine_scalar\(\)](#), [tfb_affine\(\)](#), [tfb_ascending\(\)](#), [tfb_batch_normalization\(\)](#), [tfb_blockwise\(\)](#), [tfb_chain\(\)](#), [tfb_cholesky_outer_product\(\)](#), [tfb_cholesky_to_inv_cholesky\(\)](#), [tfb_correlation_cholesky\(\)](#), [tfb_cumsum\(\)](#), [tfb_discrete_cosine_transform\(\)](#), [tfb_expm1\(\)](#), [tfb_exp\(\)](#), [tfb_ffjord\(\)](#), [tfb_fill_scale_tri_l\(\)](#), [tfb_fill_triangular\(\)](#), [tfb_glow\(\)](#), [tfb_gompertz_cdf\(\)](#), [tfb_gumbel_cdf\(\)](#), [tfb_gumbel\(\)](#), [tfb_identity\(\)](#), [tfb_inline\(\)](#), [tfb_invert\(\)](#), [tfb_iterated_sigmoid_centered\(\)](#), [tfb_kumaraswamy_cdf\(\)](#), [tfb_kumaraswamy\(\)](#), [tfb_lambert_w_tail\(\)](#), [tfb_masked_autoregressive_default_template\(\)](#), [tfb_masked_autoregressive_flow\(\)](#), [tfb_masked_dense\(\)](#), [tfb_matrix_inverse_tri_l\(\)](#), [tfb_matvec_lu\(\)](#), [tfb_normal_cdf\(\)](#), [tfb_ordered\(\)](#), [tfb_pad\(\)](#), [tfb_permute\(\)](#), [tfb_power_transform\(\)](#), [tfb_rational_quadratic_spline\(\)](#), [tfb_rayleigh_cdf\(\)](#), [tfb_real_nvp_default_template\(\)](#), [tfb_real_nvp\(\)](#), [tfb_reciprocal\(\)](#), [tfb_reshape\(\)](#), [tfb_scale_matvec_diag\(\)](#), [tfb_scale_matvec_linear_operator\(\)](#), [tfb_scale_matvec_lu\(\)](#), [tfb_scale_matvec_tri_l\(\)](#), [tfb_scale_tri_l\(\)](#), [tfb_scale\(\)](#), [tfb_shift\(\)](#), [tfb_sigmoid\(\)](#), [tfb_sinh_arcsinh\(\)](#), [tfb_sinh\(\)](#), [tfb_softmax_centered\(\)](#), [tfb_softplus\(\)](#), [tfb_softsign\(\)](#), [tfb_split\(\)](#), [tfb_square\(\)](#), [tfb_tanh\(\)](#), [tfb_transform_diagonal\(\)](#), [tfb_transpose\(\)](#), [tfb_weibull_cdf\(\)](#), [tfb_weibull\(\)](#)

tfb_sigmoid	<i>Computes</i> $Y = g(X) = 1 / (1 + \exp(-X))$
-------------	---

Description

Computes $Y = g(X) = 1 / (1 + \exp(-X))$

Usage

```
tfb_sigmoid(validate_args = FALSE, name = "sigmoid")
```

Arguments

validate_args	Logical, default FALSE. Whether to validate input with asserts. If validate_args is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
name	name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_template()`, `tfb_masked_autoregressive_flow()`, `tfb_masked_dense()`, `tfb_matrix_inverse_tri_l()`, `tfb_matvec_lu()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_pad()`, `tfb_permute()`, `tfb_power_transform()`, `tfb_rational_quadratic_spline()`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`, `tfb_real_nvp()`, `tfb_reciprocal()`, `tfb_reshape()`, `tfb_scale_matvec_diag()`, `tfb_scale_matvec_linear_operator()`, `tfb_scale_matvec_lu()`, `tfb_scale_matvec_tri_l()`, `tfb_scale_tri_l()`, `tfb_scale()`, `tfb_shifted_gompertz_cdf()`, `tfb_shift()`, `tfb_sinh_arcsinh()`, `tfb_sinh()`, `tfb_softmax_centered()`, `tfb_softplus()`, `tfb_softsign()`, `tfb_split()`, `tfb_square()`, `tfb_tanh()`, `tfb_transform_diagonal()`, `tfb_transpose()`, `tfb_weibull_cdf()`, `tfb_weibull()`

tfb_sinh

Bijector that computes $Y = \sinh(X)$.

Description

Bijector that computes $Y = \sinh(X)$.

Usage

```
tfb_sinh(validate_args = FALSE, name = "sinh")
```

Arguments

`validate_args` Logical, default FALSE. Whether to validate input with asserts. If `validate_args` is FALSE, and the inputs are invalid, correct behavior is not guaranteed.

`name` name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`,

tfb_fill_scale_tri_l(), tfb_fill_triangular(), tfb_glow(), tfb_gompertz_cdf(), tfb_gumbel_cdf(), tfb_gumbel(), tfb_identity(), tfb_inline(), tfb_invert(), tfb_iterated_sigmoid_centered(), tfb_kumaraswamy_cdf(), tfb_kumaraswamy(), tfb_lambert_w_tail(), tfb_masked_autoregressive_default_template(), tfb_masked_autoregressive_flow(), tfb_masked_dense(), tfb_matrix_inverse_tri_l(), tfb_matvec_lu(), tfb_normal_cdf(), tfb_ordered(), tfb_pad(), tfb_permute(), tfb_power_transform(), tfb_rational_quadratic_spline(), tfb_rayleigh_cdf(), tfb_real_nvp_default_template(), tfb_real_nvp(), tfb_reciprocal(), tfb_reshape(), tfb_scale_matvec_diag(), tfb_scale_matvec_linear_operator(), tfb_scale_matvec_lu(), tfb_scale_matvec_tri_l(), tfb_scale_tri_l(), tfb_scale(), tfb_shifted_gompertz_cdf(), tfb_shift(), tfb_sigmoid(), tfb_sinh_arcsinh(), tfb_softmax_centered(), tfb_softplus(), tfb_softsign(), tfb_split(), tfb_square(), tfb_tanh(), tfb_transform_diagonal(), tfb_transpose(), tfb_weibull_cdf(), tfb_weibull()

tfb_sinh_arcsinh	<i>Computes</i> $Y = g(X) = \text{Sinh}(\text{Arcsinh}(X) + \text{skewness}) * \text{tailweight}$
------------------	---

Description

For skewness in $(-\infty, \infty)$ and tailweight in $(0, \infty)$, this transformation is a diffeomorphism of the real line $(-\infty, \infty)$. The inverse transform is $X = g^{-1}(Y) = \text{Sinh}(\text{ArcSinh}(Y) / \text{tailweight} - \text{skewness})$. The SinhArcsinh transformation of the Normal is described in [Sinh-arcsinh distributions](#)

Usage

```
tfb_sinh_arcsinh(
  skewness = NULL,
  tailweight = NULL,
  validate_args = FALSE,
  name = "SinhArcsinh"
)
```

Arguments

skewness	Skewness parameter. Float-type Tensor. Default is 0 of type float32.
tailweight	Tailweight parameter. Positive Tensor of same dtype as skewness and broadcastable shape. Default is 1 of type float32.
validate_args	Logical, default FALSE. Whether to validate input with asserts. If validate_args is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
name	name prefixed to Ops created by this class.

Details

This Bijector allows a similar transformation of any distribution supported on $(-\infty, \infty)$.

Value

a bijector instance.

Meaning of the parameters

- If skewness = 0 and tailweight = 1, this transform is the identity.
- Positive (negative) skewness leads to positive (negative) skew.
- positive skew means, for unimodal X centered at zero, the mode of Y is "tilted" to the right.
- positive skew means positive values of Y become more likely, and negative values become less likely.
- Larger (smaller) tailweight leads to fatter (thinner) tails.
- Fatter tails mean larger values of |Y| become more likely.
- If X is a unit Normal, tailweight < 1 leads to a distribution that is "flat" around Y = 0, and a very steep drop-off in the tails.
- If X is a unit Normal, tailweight > 1 leads to a distribution more peaked at the mode with heavier tails. To see the argument about the tails, note that for $|X| \gg 1$ and $|X| \gg (|\text{skewness}| * \text{tailweight})$, we have $Y \approx 0.5 X \text{tailweight} e^{*(\text{sign}(X) \text{skewness} * \text{tailweight})}$.

See Also

For usage examples see [tfb_forward\(\)](#), [tfb_inverse\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#).

Other bijectors: [tfb_absolute_value\(\)](#), [tfb_affine_linear_operator\(\)](#), [tfb_affine_scalar\(\)](#), [tfb_affine\(\)](#), [tfb_ascending\(\)](#), [tfb_batch_normalization\(\)](#), [tfb_blockwise\(\)](#), [tfb_chain\(\)](#), [tfb_cholesky_outer_product\(\)](#), [tfb_cholesky_to_inv_cholesky\(\)](#), [tfb_correlation_cholesky\(\)](#), [tfb_cumsum\(\)](#), [tfb_discrete_cosine_transform\(\)](#), [tfb_expm1\(\)](#), [tfb_exp\(\)](#), [tfb_ffjord\(\)](#), [tfb_fill_scale_tri_l\(\)](#), [tfb_fill_triangular\(\)](#), [tfb_glow\(\)](#), [tfb_gompertz_cdf\(\)](#), [tfb_gumbel_cdf\(\)](#), [tfb_gumbel\(\)](#), [tfb_identity\(\)](#), [tfb_inline\(\)](#), [tfb_invert\(\)](#), [tfb_iterated_sigmoid_centered\(\)](#), [tfb_kumaraswamy_cdf\(\)](#), [tfb_kumaraswamy\(\)](#), [tfb_lambert_w_tail\(\)](#), [tfb_masked_autoregressive_default_template\(\)](#), [tfb_masked_autoregressive_flow\(\)](#), [tfb_masked_dense\(\)](#), [tfb_matrix_inverse_tri_l\(\)](#), [tfb_matvec_lu\(\)](#), [tfb_normal_cdf\(\)](#), [tfb_ordered\(\)](#), [tfb_pad\(\)](#), [tfb_permute\(\)](#), [tfb_power_transform\(\)](#), [tfb_rational_quadratic_spline\(\)](#), [tfb_rayleigh_cdf\(\)](#), [tfb_real_nvp_default_template\(\)](#), [tfb_real_nvp\(\)](#), [tfb_reciprocal\(\)](#), [tfb_reshape\(\)](#), [tfb_scale_matvec_diag\(\)](#), [tfb_scale_matvec_linear_operator\(\)](#), [tfb_scale_matvec_lu\(\)](#), [tfb_scale_matvec_tri_l\(\)](#), [tfb_scale_tri_l\(\)](#), [tfb_scale\(\)](#), [tfb_shifted_gompertz_cdf\(\)](#), [tfb_shift\(\)](#), [tfb_sigmoid\(\)](#), [tfb_sinh\(\)](#), [tfb_softmax_centered\(\)](#), [tfb_softplus\(\)](#), [tfb_softsign\(\)](#), [tfb_split\(\)](#), [tfb_square\(\)](#), [tfb_tanh\(\)](#), [tfb_transform_diagonal\(\)](#), [tfb_transpose\(\)](#), [tfb_weibull_cdf\(\)](#), [tfb_weibull\(\)](#)

tfb_softmax_centered *Computes* $Y = g(X) = \exp([X \ 0]) / \text{sum}(\exp([X \ 0]))$

Description

To implement **softmax** as a bijection, the forward transformation appends a value to the input and the inverse removes this coordinate. The appended coordinate represents a pivot, e.g., $\text{softmax}(x) = \exp(x-c) / \text{sum}(\exp(x-c))$ where c is the implicit last coordinate.

Usage

```
tfb_softmax_centered(validate_args = FALSE, name = "softmax_centered")
```

Arguments

validate_args	Logical, default FALSE. Whether to validate input with asserts. If validate_args is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
name	name prefixed to Ops created by this class.

Details

At first blush it may seem like the **Invariance of domain** theorem implies this implementation is not a bijection. However, the appended dimension makes the (forward) image non-open and the theorem does not directly apply.

Value

a bijector instance.

See Also

For usage examples see [tfb_forward\(\)](#), [tfb_inverse\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#).

Other bijectors: [tfb_absolute_value\(\)](#), [tfb_affine_linear_operator\(\)](#), [tfb_affine_scalar\(\)](#), [tfb_affine\(\)](#), [tfb_ascending\(\)](#), [tfb_batch_normalization\(\)](#), [tfb_blockwise\(\)](#), [tfb_chain\(\)](#), [tfb_cholesky_outer_product\(\)](#), [tfb_cholesky_to_inv_cholesky\(\)](#), [tfb_correlation_cholesky\(\)](#), [tfb_cumsum\(\)](#), [tfb_discrete_cosine_transform\(\)](#), [tfb_expm1\(\)](#), [tfb_exp\(\)](#), [tfb_ffjord\(\)](#), [tfb_fill_scale_tri_l\(\)](#), [tfb_fill_triangular\(\)](#), [tfb_glow\(\)](#), [tfb_gompertz_cdf\(\)](#), [tfb_gumbel_cdf\(\)](#), [tfb_gumbel\(\)](#), [tfb_identity\(\)](#), [tfb_inline\(\)](#), [tfb_invert\(\)](#), [tfb_iterated_sigmoid_centered\(\)](#), [tfb_kumaraswamy_cdf\(\)](#), [tfb_kumaraswamy\(\)](#), [tfb_lambert_w_tail\(\)](#), [tfb_masked_autoregressive_default_template\(\)](#), [tfb_masked_autoregressive_flow\(\)](#), [tfb_masked_dense\(\)](#), [tfb_matrix_inverse_tri_l\(\)](#), [tfb_matvec_lu\(\)](#), [tfb_normal_cdf\(\)](#), [tfb_ordered\(\)](#), [tfb_pad\(\)](#), [tfb_permute\(\)](#), [tfb_power_transform\(\)](#), [tfb_rational_quadratic_spline\(\)](#), [tfb_rayleigh_cdf\(\)](#), [tfb_real_nvp_default_template\(\)](#), [tfb_real_nvp\(\)](#), [tfb_reciprocal\(\)](#), [tfb_reshape\(\)](#), [tfb_scale_matvec_diag\(\)](#), [tfb_scale_matvec_linear_operator\(\)](#), [tfb_scale_matvec_lu\(\)](#), [tfb_scale_matvec_tri_l\(\)](#), [tfb_scale_tri_l\(\)](#), [tfb_scale\(\)](#), [tfb_shifted_gompertz_cdf\(\)](#), [tfb_shift\(\)](#), [tfb_sigmoid\(\)](#), [tfb_sinh_arcsinh\(\)](#), [tfb_sinh\(\)](#), [tfb_softplus\(\)](#), [tfb_softsign\(\)](#), [tfb_split\(\)](#), [tfb_square\(\)](#), [tfb_tanh\(\)](#), [tfb_transform_diagonal\(\)](#), [tfb_transpose\(\)](#), [tfb_weibull_cdf\(\)](#), [tfb_weibull\(\)](#)

tfb_softplus

Computes $Y = g(X) = \text{Log}[1 + \exp(X)]$

Description

The softplus Bijector has the following two useful properties:

- The domain is the positive real numbers
- $\text{softplus}(x) \approx x$, for large x , so it does not overflow as easily as the Exp Bijector.

Usage

```
tfb_softplus(
  hinge_softness = NULL,
  low = NULL,
  validate_args = FALSE,
  name = "softplus"
)
```

Arguments

`hinge_softness` Nonzero floating point Tensor. Controls the softness of what would otherwise be a kink at the origin. Default is 1.0.

`low` Nonzero floating point tensor, lower bound on output values. Implicitly zero if NULL. Otherwise, the transformation $y = \text{softplus}(x) + \text{low}$ is implemented. This is equivalent to a `tfb_chain(list(tfb_shift(low), tfb_softplus()))` bijector and is provided for convenience.

`validate_args` Logical, default FALSE. Whether to validate input with asserts. If `validate_args` is FALSE, and the inputs are invalid, correct behavior is not guaranteed.

`name` name prefixed to Ops created by this class.

Details

The optional nonzero `hinge_softness` parameter changes the transition at zero. With `hinge_softness = c`, the bijector is:

```
f_c(x) := c * g(x / c) = c * Log[1 + exp(x / c)].
'''
```

For large $x \gg 1$,

```
'''
c * Log[1 + exp(x / c)] approx c * Log[exp(x / c)] = x
'''
```

so the behavior for large x is the same as the standard `softplus`.

As $c > 0$ approaches 0 from the right, $f_c(x)$ becomes less and less soft, approaching $\max(0, x)$.

* $c = 1$ is the default.

* $c > 0$ but small means $f(x)$ approx $\text{ReLU}(x) = \max(0, x)$.

* $c < 0$ flips sign and reflects around the y-axis: $f_{\{-c\}}(x) = -f_c(-x)$.

* $c = 0$ results in a non-bijective transformation and triggers an exception.

Note: `log(.)` and `exp(.)` are applied element-wise but the Jacobian is a reduction over the event space.

```
[1 + exp(x / c)]: R:1%20+%20exp(x%20/%20c)
```

```
[1 + exp(x / c)]: R:1%20+%20exp(x%20/%20c)
```

```
[exp(x / c)]: R:exp(x%20/%20c)
```

Value

a bijector instance.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_template()`, `tfb_masked_autoregressive_flow()`, `tfb_masked_dense()`, `tfb_matrix_inverse_tri_l()`, `tfb_matvec_lu()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_pad()`, `tfb_permute()`, `tfb_power_transform()`, `tfb_rational_quadratic_spline()`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`, `tfb_real_nvp()`, `tfb_reciprocal()`, `tfb_reshape()`, `tfb_scale_matvec_diag()`, `tfb_scale_matvec_linear_operator()`, `tfb_scale_matvec_lu()`, `tfb_scale_matvec_tri_l()`, `tfb_scale_tri_l()`, `tfb_scale()`, `tfb_shifted_gompertz_cdf()`, `tfb_shift()`, `tfb_sigmoid()`, `tfb_sinh_arcsinh()`, `tfb_sinh()`, `tfb_softmax_centered()`, `tfb_softsign()`, `tfb_split()`, `tfb_square()`, `tfb_tanh()`, `tfb_transform_diagonal()`, `tfb_transpose()`, `tfb_weibull_cdf()`, `tfb_weibull()`

`tfb_softsign`

Computes $Y = g(X) = X / (1 + |X|)$

Description

The softsign Bijector has the following two useful properties:

- The domain is all real numbers
- `softsign(x)` approx `sgn(x)`, for large `|x|`.

Usage

```
tfb_softsign(validate_args = FALSE, name = "softsign")
```

Arguments

`validate_args` Logical, default `FALSE`. Whether to validate input with asserts. If `validate_args` is `FALSE`, and the inputs are invalid, correct behavior is not guaranteed.

`name` name prefixed to Ops created by this class.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_template()`, `tfb_masked_autoregressive_flow()`, `tfb_masked_dense()`, `tfb_matrix_inverse_tri_l()`, `tfb_matvec_lu()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_pad()`, `tfb_permute()`, `tfb_power_transform()`, `tfb_rational_quadratic_spline()`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`, `tfb_real_nvp()`, `tfb_reciprocal()`, `tfb_reshape()`, `tfb_scale_matvec_diag()`, `tfb_scale_matvec_linear_operator()`, `tfb_scale_matvec_lu()`, `tfb_scale_matvec_tri_l()`, `tfb_scale_tri_l()`, `tfb_scale()`, `tfb_shifted_gompertz_cdf()`, `tfb_shift()`, `tfb_sigmoid()`, `tfb_sinh_arcsinh()`, `tfb_sinh()`, `tfb_softmax_centered()`, `tfb_softplus()`, `tfb_split()`, `tfb_square()`, `tfb_tanh()`, `tfb_transform_diagonal()`, `tfb_transpose()`, `tfb_weibull_cdf()`, `tfb_weibull()`

tfb_split

Split a Tensor event along an axis into a list of Tensors.

Description

The inverse of `split` concatenates a list of Tensors along `axis`.

Usage

```
tfb_split(num_or_size_splits, axis = -1, validate_args = FALSE, name = "split")
```

Arguments

<code>num_or_size_splits</code>	Either an integer indicating the number of splits along <code>axis</code> or a 1-D integer Tensor or Python list containing the sizes of each output tensor along <code>axis</code> . If a list/Tensor, it may contain at most one value of <code>-1</code> , which indicates a split size that is unknown and determined from input.
<code>axis</code>	A negative integer or scalar <code>int32</code> Tensor. The dimension along which to split. Must be negative to enable the bijector to support arbitrary batch dimensions. Defaults to <code>-1</code> (note that this is different from the <code>tf\$Split</code> default of <code>0</code>). Must be statically known.
<code>validate_args</code>	Logical, default <code>FALSE</code> . Whether to validate input with asserts. If <code>validate_args</code> is <code>FALSE</code> , and the inputs are invalid, correct behavior is not guaranteed.
<code>name</code>	name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_template()`, `tfb_masked_autoregressive_flow()`, `tfb_masked_dense()`, `tfb_matrix_inverse_tri_l()`, `tfb_matvec_lu()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_pad()`, `tfb_permute()`, `tfb_power_transform()`, `tfb_rational_quadratic_spline()`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`, `tfb_real_nvp()`, `tfb_reciprocal()`, `tfb_reshape()`, `tfb_scale_matvec_diag()`, `tfb_scale_matvec_linear_operator()`, `tfb_scale_matvec_lu()`, `tfb_scale_matvec_tri_l()`, `tfb_scale_tri_l()`, `tfb_scale()`, `tfb_shifted_gompertz_cdf()`, `tfb_shift()`, `tfb_sigmoid()`, `tfb_sinh_arcsinh()`, `tfb_sinh()`, `tfb_softmax_centered()`, `tfb_softplus()`, `tfb_softsign()`, `tfb_square()`, `tfb_tanh()`, `tfb_transform_diagonal()`, `tfb_transpose()`, `tfb_weibull_cdf()`, `tfb_weibull()`

tfb_square

Computes $g(X) = X^2$; X is a positive real number.

Description

g is a bijection between the non-negative real numbers (\mathbb{R}_+) and the non-negative real numbers.

Usage

```
tfb_square(validate_args = FALSE, name = "square")
```

Arguments

`validate_args` Logical, default FALSE. Whether to validate input with asserts. If `validate_args` is FALSE, and the inputs are invalid, correct behavior is not guaranteed.

`name` name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`,

tfb_fill_scale_tri_l(), tfb_fill_triangular(), tfb_glow(), tfb_gompertz_cdf(), tfb_gumbel_cdf(),
 tfb_gumbel(), tfb_identity(), tfb_inline(), tfb_invert(), tfb_iterated_sigmoid_centered(),
 tfb_kumaraswamy_cdf(), tfb_kumaraswamy(), tfb_lambert_w_tail(), tfb_masked_autoregressive_default_temp,
 tfb_masked_autoregressive_flow(), tfb_masked_dense(), tfb_matrix_inverse_tri_l(),
 tfb_matvec_lu(), tfb_normal_cdf(), tfb_ordered(), tfb_pad(), tfb_permute(), tfb_power_transform(),
 tfb_rational_quadratic_spline(), tfb_rayleigh_cdf(), tfb_real_nvp_default_template(),
 tfb_real_nvp(), tfb_reciprocal(), tfb_reshape(), tfb_scale_matvec_diag(), tfb_scale_matvec_linear_operat
 tfb_scale_matvec_lu(), tfb_scale_matvec_tri_l(), tfb_scale_tri_l(), tfb_scale(), tfb_shifted_gompertz_co
 tfb_shift(), tfb_sigmoid(), tfb_sinh_arcsinh(), tfb_sinh(), tfb_softmax_centered(),
 tfb_softplus(), tfb_softsign(), tfb_split(), tfb_tanh(), tfb_transform_diagonal(),
 tfb_transpose(), tfb_weibull_cdf(), tfb_weibull()

tfb_tanh

Computes $Y = \tanh(X)$ **Description**

$Y = \tanh(X)$, therefore Y in $(-1, 1)$.

Usage

```
tfb_tanh(validate_args = FALSE, name = "tanh")
```

Arguments

`validate_args` Logical, default FALSE. Whether to validate input with asserts. If `validate_args` is FALSE, and the inputs are invalid, correct behavior is not guaranteed.

`name` name prefixed to Ops created by this class.

Details

This can be achieved by an affine transform of the Sigmoid bijector, i.e., it is equivalent to
`tfb_chain(list(tfb_affine(shift = -1, scale = 2), tfb_sigmoid(), tfb_affine(scale = 2)))`
 However, using the Tanh bijector directly is slightly faster and more numerically stable.

Value

a bijector instance.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`,
`tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`,
`tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`,
`tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`,
`tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`,

[tfb_gumbel\(\)](#), [tfb_identity\(\)](#), [tfb_inline\(\)](#), [tfb_invert\(\)](#), [tfb_iterated_sigmoid_centered\(\)](#),
[tfb_kumaraswamy_cdf\(\)](#), [tfb_kumaraswamy\(\)](#), [tfb_lambert_w_tail\(\)](#), [tfb_masked_autoregressive_default_template\(\)](#),
[tfb_masked_autoregressive_flow\(\)](#), [tfb_masked_dense\(\)](#), [tfb_matrix_inverse_tri_l\(\)](#),
[tfb_matvec_lu\(\)](#), [tfb_normal_cdf\(\)](#), [tfb_ordered\(\)](#), [tfb_pad\(\)](#), [tfb_permute\(\)](#), [tfb_power_transform\(\)](#),
[tfb_rational_quadratic_spline\(\)](#), [tfb_rayleigh_cdf\(\)](#), [tfb_real_nvp_default_template\(\)](#),
[tfb_real_nvp\(\)](#), [tfb_reciprocal\(\)](#), [tfb_reshape\(\)](#), [tfb_scale_matvec_diag\(\)](#), [tfb_scale_matvec_linear_operator\(\)](#),
[tfb_scale_matvec_lu\(\)](#), [tfb_scale_matvec_tri_l\(\)](#), [tfb_scale_tri_l\(\)](#), [tfb_scale\(\)](#), [tfb_shifted_gompertz_cdf\(\)](#),
[tfb_shift\(\)](#), [tfb_sigmoid\(\)](#), [tfb_sinh_arcsinh\(\)](#), [tfb_sinh\(\)](#), [tfb_softmax_centered\(\)](#),
[tfb_softplus\(\)](#), [tfb_softsign\(\)](#), [tfb_split\(\)](#), [tfb_square\(\)](#), [tfb_transform_diagonal\(\)](#),
[tfb_transpose\(\)](#), [tfb_weibull_cdf\(\)](#), [tfb_weibull\(\)](#)

tfb_transform_diagonal

Applies a Bijector to the diagonal of a matrix

Description

Applies a Bijector to the diagonal of a matrix

Usage

```

tfb_transform_diagonal(
  diag_bijector,
  validate_args = FALSE,
  name = "transform_diagonal"
)

```

Arguments

diag_bijector	Bijector instance used to transform the diagonal.
validate_args	Logical, default FALSE. Whether to validate input with asserts. If validate_args is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
name	name prefixed to Ops created by this class.

Value

a bijector instance.

See Also

For usage examples see [tfb_forward\(\)](#), [tfb_inverse\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#).

Other bijectors: [tfb_absolute_value\(\)](#), [tfb_affine_linear_operator\(\)](#), [tfb_affine_scalar\(\)](#), [tfb_affine\(\)](#), [tfb_ascending\(\)](#), [tfb_batch_normalization\(\)](#), [tfb_blockwise\(\)](#), [tfb_chain\(\)](#), [tfb_cholesky_outer_product\(\)](#), [tfb_cholesky_to_inv_cholesky\(\)](#), [tfb_correlation_cholesky\(\)](#), [tfb_cumsum\(\)](#), [tfb_discrete_cosine_transform\(\)](#), [tfb_expm1\(\)](#), [tfb_exp\(\)](#), [tfb_ffjord\(\)](#), [tfb_fill_scale_tri_l\(\)](#), [tfb_fill_triangular\(\)](#), [tfb_glow\(\)](#), [tfb_gompertz_cdf\(\)](#), [tfb_gumbel_cdf\(\)](#), [tfb_gumbel\(\)](#), [tfb_identity\(\)](#), [tfb_inline\(\)](#), [tfb_invert\(\)](#), [tfb_iterated_sigmoid_centered\(\)](#),

tfb_kumaraswamy_cdf(), tfb_kumaraswamy(), tfb_lambert_w_tail(), tfb_masked_autoregressive_default_template(),
 tfb_masked_autoregressive_flow(), tfb_masked_dense(), tfb_matrix_inverse_tri_l(),
 tfb_matvec_lu(), tfb_normal_cdf(), tfb_ordered(), tfb_pad(), tfb_permute(), tfb_power_transform(),
 tfb_rational_quadratic_spline(), tfb_rayleigh_cdf(), tfb_real_nvp_default_template(),
 tfb_real_nvp(), tfb_reciprocal(), tfb_reshape(), tfb_scale_matvec_diag(), tfb_scale_matvec_linear_operation(),
 tfb_scale_matvec_lu(), tfb_scale_matvec_tri_l(), tfb_scale_tri_l(), tfb_scale(), tfb_shifted_gompertz_cdf(),
 tfb_shift(), tfb_sigmoid(), tfb_sinh_arcsinh(), tfb_sinh(), tfb_softmax_centered(),
 tfb_softplus(), tfb_softsign(), tfb_split(), tfb_square(), tfb_tanh(), tfb_transpose(),
 tfb_weibull_cdf(), tfb_weibull()

tfb_transpose	<i>Computes</i> $Y = g(X) = \text{transpose_rightmost_dims}(X, \text{rightmost_perm})$
---------------	---

Description

This bijector is semantically similar to `tf.transpose` except that it transposes only the rightmost "event" dimensions. That is, unlike `tf.transpose` the `perm` argument is itself a permutation of `tf.range(rightmost_transposed_ndims)` rather than `tf.rank(x)`, i.e., users specify the (rightmost) dimensions to permute, not all dimensions.

Usage

```
tfb_transpose(
  perm = NULL,
  rightmost_transposed_ndims = NULL,
  validate_args = FALSE,
  name = "transpose"
)
```

Arguments

<code>perm</code>	Positive integer vector-shaped Tensor representing permutation of rightmost dims (for forward transformation). Note that the 0th index represents the first of the rightmost dims and the largest value must be <code>rightmost_transposed_ndims - 1</code> and corresponds to <code>tf.rank(x) - 1</code> . Only one of <code>perm</code> and <code>rightmost_transposed_ndims</code> can (and must) be specified. Default value: <code>tf.range(start=rightmost_transposed_ndims, limit=-1)</code>
<code>rightmost_transposed_ndims</code>	Positive integer scalar-shaped Tensor representing the number of rightmost dimensions to permute. Only one of <code>perm</code> and <code>rightmost_transposed_ndims</code> can (and must) be specified. Default value: <code>tf.size(perm)</code> .
<code>validate_args</code>	Logical, default <code>FALSE</code> . Whether to validate input with asserts. If <code>validate_args</code> is <code>FALSE</code> , and the inputs are invalid, correct behavior is not guaranteed.
<code>name</code>	name prefixed to Ops created by this class.

Details

The actual (forward) transformation is:

```
sample_batch_ndims <-tf$rank(x) -tf$size(perm) perm = tf$concat(list(tf$range(sample_batch_ndims), samp
+ perm),axis=0) tf$transpose(x,perm)
```

Value

a bijector instance.

See Also

For usage examples see [tfb_forward\(\)](#), [tfb_inverse\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#).

Other bijectors: [tfb_absolute_value\(\)](#), [tfb_affine_linear_operator\(\)](#), [tfb_affine_scalar\(\)](#), [tfb_affine\(\)](#), [tfb_ascending\(\)](#), [tfb_batch_normalization\(\)](#), [tfb_blockwise\(\)](#), [tfb_chain\(\)](#), [tfb_cholesky_outer_product\(\)](#), [tfb_cholesky_to_inv_cholesky\(\)](#), [tfb_correlation_cholesky\(\)](#), [tfb_cumsum\(\)](#), [tfb_discrete_cosine_transform\(\)](#), [tfb_expm1\(\)](#), [tfb_exp\(\)](#), [tfb_ffjord\(\)](#), [tfb_fill_scale_tri_l\(\)](#), [tfb_fill_triangular\(\)](#), [tfb_glow\(\)](#), [tfb_gompertz_cdf\(\)](#), [tfb_gumbel_cdf\(\)](#), [tfb_gumbel\(\)](#), [tfb_identity\(\)](#), [tfb_inline\(\)](#), [tfb_invert\(\)](#), [tfb_iterated_sigmoid_centered\(\)](#), [tfb_kumaraswamy_cdf\(\)](#), [tfb_kumaraswamy\(\)](#), [tfb_lambert_w_tail\(\)](#), [tfb_masked_autoregressive_default_template\(\)](#), [tfb_masked_autoregressive_flow\(\)](#), [tfb_masked_dense\(\)](#), [tfb_matrix_inverse_tri_l\(\)](#), [tfb_matvec_lu\(\)](#), [tfb_normal_cdf\(\)](#), [tfb_ordered\(\)](#), [tfb_pad\(\)](#), [tfb_permute\(\)](#), [tfb_power_transform\(\)](#), [tfb_rational_quadratic_spline\(\)](#), [tfb_rayleigh_cdf\(\)](#), [tfb_real_nvp_default_template\(\)](#), [tfb_real_nvp\(\)](#), [tfb_reciprocal\(\)](#), [tfb_reshape\(\)](#), [tfb_scale_matvec_diag\(\)](#), [tfb_scale_matvec_linear_operator\(\)](#), [tfb_scale_matvec_lu\(\)](#), [tfb_scale_matvec_tri_l\(\)](#), [tfb_scale_tri_l\(\)](#), [tfb_scale\(\)](#), [tfb_shifted_gompertz_cdf\(\)](#), [tfb_shift\(\)](#), [tfb_sigmoid\(\)](#), [tfb_sinh_arcsinh\(\)](#), [tfb_sinh\(\)](#), [tfb_softmax_centered\(\)](#), [tfb_softplus\(\)](#), [tfb_softsign\(\)](#), [tfb_split\(\)](#), [tfb_square\(\)](#), [tfb_tanh\(\)](#), [tfb_transform_diagonal\(\)](#), [tfb_weibull_cdf\(\)](#), [tfb_weibull\(\)](#)

tfb_weibull	<i>Computes</i> $Y = g(X) = 1 - \exp((-X / \text{scale}) ** \text{concentration})$ <i>where</i> $X \geq 0$
-------------	---

Description

This bijector maps inputs from $[0, \infty]$ to $[0, 1]$. The inverse of the bijector applied to a uniform random variable $X \sim U(0, 1)$ gives back a random variable with the **Weibull distribution**:

Usage

```
tfb_weibull(
  scale = 1,
  concentration = 1,
  validate_args = FALSE,
  name = "weibull"
)
```

Arguments

scale	Positive Float-type Tensor that is the same dtype and is broadcastable with concentration. This is l in $Y = g(X) = 1 - \exp((-x / l) ** k)$.
concentration	Positive Float-type Tensor that is the same dtype and is broadcastable with scale. This is k in $Y = g(X) = 1 - \exp((-x / l) ** k)$.
validate_args	Logical, default <code>FALSE</code> . Whether to validate input with asserts. If <code>validate_args</code> is <code>FALSE</code> , and the inputs are invalid, correct behavior is not guaranteed.
name	name prefixed to Ops created by this class.

Details

$Y \sim \text{Weibull}(\text{scale}, \text{concentration})$ pdf(y ; scale, concentration, $y \geq 0$) = (concentration / scale) * (y / scale)**(concentration - 1) * exp(-(y / scale)**concentration)

Value

a bijector instance.

See Also

For usage examples see `tfb_forward()`, `tfb_inverse()`, `tfb_inverse_log_det_jacobian()`.

Other bijectors: `tfb_absolute_value()`, `tfb_affine_linear_operator()`, `tfb_affine_scalar()`, `tfb_affine()`, `tfb_ascending()`, `tfb_batch_normalization()`, `tfb_blockwise()`, `tfb_chain()`, `tfb_cholesky_outer_product()`, `tfb_cholesky_to_inv_cholesky()`, `tfb_correlation_cholesky()`, `tfb_cumsum()`, `tfb_discrete_cosine_transform()`, `tfb_expm1()`, `tfb_exp()`, `tfb_ffjord()`, `tfb_fill_scale_tri_l()`, `tfb_fill_triangular()`, `tfb_glow()`, `tfb_gompertz_cdf()`, `tfb_gumbel_cdf()`, `tfb_gumbel()`, `tfb_identity()`, `tfb_inline()`, `tfb_invert()`, `tfb_iterated_sigmoid_centered()`, `tfb_kumaraswamy_cdf()`, `tfb_kumaraswamy()`, `tfb_lambert_w_tail()`, `tfb_masked_autoregressive_default_template()`, `tfb_masked_autoregressive_flow()`, `tfb_masked_dense()`, `tfb_matrix_inverse_tri_l()`, `tfb_matvec_lu()`, `tfb_normal_cdf()`, `tfb_ordered()`, `tfb_pad()`, `tfb_permute()`, `tfb_power_transform()`, `tfb_rational_quadratic_spline()`, `tfb_rayleigh_cdf()`, `tfb_real_nvp_default_template()`, `tfb_real_nvp()`, `tfb_reciprocal()`, `tfb_reshape()`, `tfb_scale_matvec_diag()`, `tfb_scale_matvec_linear_operator()`, `tfb_scale_matvec_lu()`, `tfb_scale_matvec_tri_l()`, `tfb_scale_tri_l()`, `tfb_scale()`, `tfb_shifted_gompertz_cdf()`, `tfb_shift()`, `tfb_sigmoid()`, `tfb_sinh_arcsinh()`, `tfb_sinh()`, `tfb_softmax_centered()`, `tfb_softplus()`, `tfb_softsign()`, `tfb_split()`, `tfb_square()`, `tfb_tanh()`, `tfb_transform_diagonal()`, `tfb_transpose()`, `tfb_weibull_cdf()`

tfb_weibull_cdf

Compute $Y = g(X) = 1 - \exp((-X / \text{scale}) ** \text{concentration})$, $X \geq 0$.

Description

This bijector maps inputs from $[0, \infty]$ to $[0, 1]$. The inverse of the bijector applied to a uniform random variable $X \sim U(0, 1)$ gives back a random variable with the **Weibull distribution**:

```

Y ~ Weibull(scale, concentration)
pdf(y; scale, concentration, y >= 0) =
  (concentration / scale) * (y / scale)**(concentration - 1) *
  exp(-(y / scale)**concentration)

```

Usage

```

tfb_weibull_cdf(
  scale = 1,
  concentration = 1,
  validate_args = FALSE,
  name = "weibull_cdf"
)

```

Arguments

scale	Positive Float-type Tensor that is the same dtype and is broadcastable with concentration. This is l in $Y = g(X) = 1 - \exp((-x / l) ** k)$.
concentration	Positive Float-type Tensor that is the same dtype and is broadcastable with scale. This is k in $Y = g(X) = 1 - \exp((-x / l) ** k)$.
validate_args	Logical, default FALSE. Whether to validate input with asserts. If validate_args is FALSE, and the inputs are invalid, correct behavior is not guaranteed.
name	name prefixed to Ops created by this class.

Details

Likewise, the forward of this bijector is the Weibull distribution CDF.

Value

a bijector instance.

See Also

For usage examples see [tfb_forward\(\)](#), [tfb_inverse\(\)](#), [tfb_inverse_log_det_jacobian\(\)](#).

Other bijectors: [tfb_absolute_value\(\)](#), [tfb_affine_linear_operator\(\)](#), [tfb_affine_scalar\(\)](#), [tfb_affine\(\)](#), [tfb_ascending\(\)](#), [tfb_batch_normalization\(\)](#), [tfb_blockwise\(\)](#), [tfb_chain\(\)](#), [tfb_cholesky_outer_product\(\)](#), [tfb_cholesky_to_inv_cholesky\(\)](#), [tfb_correlation_cholesky\(\)](#), [tfb_cumsum\(\)](#), [tfb_discrete_cosine_transform\(\)](#), [tfb_expm1\(\)](#), [tfb_exp\(\)](#), [tfb_ffjord\(\)](#), [tfb_fill_scale_tri_l\(\)](#), [tfb_fill_triangular\(\)](#), [tfb_glow\(\)](#), [tfb_gompertz_cdf\(\)](#), [tfb_gumbel_cdf\(\)](#), [tfb_gumbel\(\)](#), [tfb_identity\(\)](#), [tfb_inline\(\)](#), [tfb_invert\(\)](#), [tfb_iterated_sigmoid_centered\(\)](#), [tfb_kumaraswamy_cdf\(\)](#), [tfb_kumaraswamy\(\)](#), [tfb_lambert_w_tail\(\)](#), [tfb_masked_autoregressive_default_template\(\)](#), [tfb_masked_autoregressive_flow\(\)](#), [tfb_masked_dense\(\)](#), [tfb_matrix_inverse_tri_l\(\)](#), [tfb_matvec_lu\(\)](#), [tfb_normal_cdf\(\)](#), [tfb_ordered\(\)](#), [tfb_pad\(\)](#), [tfb_permute\(\)](#), [tfb_power_transform\(\)](#), [tfb_rational_quadratic_spline\(\)](#), [tfb_rayleigh_cdf\(\)](#), [tfb_real_nvp_default_template\(\)](#), [tfb_real_nvp\(\)](#), [tfb_reciprocal\(\)](#), [tfb_reshape\(\)](#), [tfb_scale_matvec_diag\(\)](#), [tfb_scale_matvec_linear_operator\(\)](#), [tfb_scale_matvec_lu\(\)](#), [tfb_scale_matvec_tri_l\(\)](#), [tfb_scale_tri_l\(\)](#), [tfb_scale\(\)](#), [tfb_shifted_gompertz_cdf\(\)](#), [tfb_shift\(\)](#), [tfb_sigmoid\(\)](#), [tfb_sinh_arcsinh\(\)](#), [tfb_sinh\(\)](#), [tfb_softmax_centered\(\)](#),

`tfb_softplus()`, `tfb_softsign()`, `tfb_split()`, `tfb_square()`, `tfb_tanh()`, `tfb_transform_diagonal()`, `tfb_transpose()`, `tfb_weibull()`

tfd_autoregressive *Autoregressive distribution*

Description

The Autoregressive distribution enables learning (often) richer multivariate distributions by repeatedly applying a **diffeomorphic** transformation (such as implemented by Bijectors).

Usage

```
tfd_autoregressive(
  distribution_fn,
  sample0 = NULL,
  num_steps = NULL,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "Autoregressive"
)
```

Arguments

<code>distribution_fn</code>	Function which constructs a <code>tfd\$Distribution</code> -like instance from a Tensor (e.g., <code>sample0</code>). The function must respect the "autoregressive property", i.e., there exists a permutation of event such that each coordinate is a diffeomorphic function of on preceding coordinates.
<code>sample0</code>	Initial input to <code>distribution_fn</code> ; used to build the distribution in <code>__init__</code> which in turn specifies this distribution's properties, e.g., <code>event_shape</code> , <code>batch_shape</code> , <code>dtype</code> . If unspecified, then <code>distribution_fn</code> should be default constructable.
<code>num_steps</code>	Number of times <code>distribution_fn</code> is composed from samples, e.g., <code>num_steps=2</code> implies <code>distribution_fn(distribution_fn(sample0)\$sample(n))\$sample()</code> .
<code>validate_args</code>	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
<code>allow_nan_stats</code>	Logical, default <code>TRUE</code> . When <code>TRUE</code> , statistics (e.g., mean, mode, variance) use the value <code>NaN</code> to indicate the result is undefined. When <code>FALSE</code> , an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Details

Regarding terminology, "Autoregressive models decompose the joint density as a product of conditionals, and model each conditional in turn. Normalizing flows transform a base density (e.g. a standard Gaussian) into the target density by an invertible transformation with tractable Jacobian." (Papamakarios et al., 2016)

In other words, the "autoregressive property" is equivalent to the decomposition, $p(x) = \prod\{ p(x[i] | x[0:i]) : i=0, \dots, d \}$. The provided `shift_and_log_scale_fn`, `tfd_masked_autoregressive_default_template`, achieves this property by zeroing out weights in its `masked_dense` layers. Practically speaking the autoregressive property means that there exists a permutation of the event coordinates such that each coordinate is a diffeomorphic function of only preceding coordinates (van den Oord et al., 2016).

Mathematical Details

The probability function is

$$\text{prob}(x; \text{fn}, n) = \text{fn}(x).\text{prob}(x)$$

And a sample is generated by

$$x = \text{fn}(\dots \text{fn}(\text{fn}(x_0).\text{sample}()).\text{sample}()).\text{sample}()$$

where the ellipses (...) represent $n-2$ composed calls to `fn`, `fn` constructs a `tfd$Distribution`-like instance, and `x0` is a fixed initializing Tensor.

Value

a distribution instance.

References

- George Papamakarios, Theo Pavlakou, and Iain Murray. Masked Autoregressive Flow for Density Estimation. In *Neural Information Processing Systems*, 2017.
- Aaron van den Oord, Nal Kalchbrenner, Oriol Vinyals, Lasse Espeholt, Alex Graves, and Koray Kavukcuoglu. Conditional Image Generation with PixelCNN Decoders. In *Neural Information Processing Systems*, 2016.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space`.


```
tfd_lkj(), tfd_log_logistic(), tfd_log_normal(), tfd_logistic(), tfd_mixture_same_family(),
tfd_mixture(), tfd_multinomial(), tfd_multivariate_normal_diag_plus_low_rank(), tfd_multivariate_normal_diag_plus_low_rank(),
tfd_multivariate_normal_full_covariance(), tfd_multivariate_normal_linear_operator(),
tfd_multivariate_normal_tri_l(), tfd_multivariate_student_t_linear_operator(), tfd_negative_binomial(),
tfd_normal(), tfd_one_hot_categorical(), tfd_pareto(), tfd_pixel_cnn(), tfd_poisson_log_normal_quadrature(),
tfd_poisson(), tfd_power_spherical(), tfd_probit_bernoulli(), tfd_quantized(), tfd_relaxed_bernoulli(),
tfd_relaxed_one_hot_categorical(), tfd_sample_distribution(), tfd_sinh_arcsinh(),
tfd_skellam(), tfd_spherical_uniform(), tfd_student_t_process(), tfd_student_t(), tfd_transformed_distribution(),
tfd_triangular(), tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian(),
tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()
```

tfd_batch_reshape	<i>Batch-Reshaping distribution</i>
-------------------	-------------------------------------

Description

This "meta-distribution" reshapes the batch dimensions of another distribution.

Usage

```
tfd_batch_reshape(
  distribution,
  batch_shape,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = NULL
)
```

Arguments

distribution	The base distribution instance to reshape. Typically an instance of <code>Distribution</code> .
batch_shape	Positive integer-like vector-shaped <code>Tensor</code> representing the new shape of the batch dimensions. Up to one dimension may contain <code>-1</code> , meaning the remainder of the batch size.
validate_args	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
allow_nan_stats	Logical, default <code>TRUE</code> . When <code>TRUE</code> , statistics (e.g., mean, mode, variance) use the value <code>NaN</code> to indicate the result is undefined. When <code>FALSE</code> , an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag_plus_low_rank_batched()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_bates

Bates distribution.

Description

The Bates distribution is the distribution of the average of `total_count` independent samples from `Uniform(low,high)`. It is parameterized by the interval bounds `low` and `high`, and `total_count`, the number of samples. Although some care has been taken to avoid numerical issues, the pdf, cdf, and log versions thereof may still exhibit numerical instability. They are relatively stable near the tails; however near the mode they are unstable if `total_count` is greater than about 75 for `tf$float64`, 25 for `tf$float32`, and 7 for `tf$float16`. Beyond these limits a warning will be shown if `validate_args=FALSE`; otherwise an exception is thrown. For high `total_count`, consider using a Normal approximation.

Usage

```
tfd_bates(
  total_count,
  low = 0,
  high = 1,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "Bates"
)
```

Arguments

<code>total_count</code>	Non-negative integer-valued Tensor with shape broadcastable to the batch shape $[N_1, \dots, N_m]$, $m \geq 0$. This controls the number of samples of <code>Uniform(low, high)</code> to take the mean of.
<code>low</code>	Floating point Tensor representing the lower bounds of the support. Should be broadcastable to $[N_1, \dots, N_m]$ with $m \geq 0$, the same dtype as <code>total_count</code> , and $low < high$ component-wise, after broadcasting. Defaults to 0.
<code>high</code>	Floating point Tensor representing the upper bounds of the support. Should be broadcastable to $[N_1, \dots, N_m]$ with $m \geq 0$, the same dtype as <code>total_count</code> , and $low < high$ component-wise, after broadcasting. Defaults to 1.
<code>validate_args</code>	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
<code>allow_nan_stats</code>	Logical, default <code>TRUE</code> . When <code>TRUE</code> , statistics (e.g., mean, mode, variance) use the value <code>NaN</code> to indicate the result is undefined. When <code>FALSE</code> , an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Details**Mathematical Details**

The probability density function (pdf) is supported in the interval $[low, high]$. If $[low, high]$ is the unit interval $[0, 1]$, the pdf is,

$$\text{pdf}(x; n, 0, 1) = \binom{n}{j} (-1)^j (nx - j)^{n-1-j}$$

where

- `total_count` = n ,
- $j = \text{floor}(nx)$
- $n!$ is the factorial of n ,
- $\binom{n}{k}$ is the binomial coefficient $n! / (k!(n - k)!)$ For arbitrary intervals $[low, high]$, the pdf is,

$$\text{pdf}(x; n, low, high) = \text{pdf}((x - low) / (high - low); n, 0, 1) / (high - low)$$

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_bernoulli

Bernoulli distribution

Description

The Bernoulli distribution with `probs` parameter, i.e., the probability of a 1 outcome (vs a 0 outcome).

Usage

```
tfd_bernoulli(
  logits = NULL,
  probs = NULL,
  dtype = tf$int32,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "Bernoulli"
)
```

Arguments

logits	An N-D Tensor representing the log-odds of a 1 event. Each entry in the Tensor parametrizes an independent Bernoulli distribution where the probability of an event is sigmoid(logits). Only one of logits or probs should be passed in.
probs	An N-D Tensor representing the probability of a 1 event. Each entry in the Tensor parameterizes an independent Bernoulli distribution. Only one of logits or probs should be passed in.
dtype	The type of the event samples. Default: int32.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Value

a distribution instance.

See Also

For usage examples see e.g. [tfd_sample\(\)](#), [tfd_log_prob\(\)](#), [tfd_mean\(\)](#).

Other distributions: [tfd_autoregressive\(\)](#), [tfd_batch_reshape\(\)](#), [tfd_bates\(\)](#), [tfd_beta_binomial\(\)](#), [tfd_beta\(\)](#), [tfd_binomial\(\)](#), [tfd_categorical\(\)](#), [tfd_cauchy\(\)](#), [tfd_chi2\(\)](#), [tfd_chi\(\)](#), [tfd_cholesky_lkj\(\)](#), [tfd_continuous_bernoulli\(\)](#), [tfd_deterministic\(\)](#), [tfd_dirichlet_multinomial\(\)](#), [tfd_dirichlet\(\)](#), [tfd_empirical\(\)](#), [tfd_exp_gamma\(\)](#), [tfd_exp_inverse_gamma\(\)](#), [tfd_exponential\(\)](#), [tfd_gamma_gamma\(\)](#), [tfd_gamma\(\)](#), [tfd_gaussian_process_regression_model\(\)](#), [tfd_gaussian_process\(\)](#), [tfd_generalized_normal\(\)](#), [tfd_geometric\(\)](#), [tfd_gumbel\(\)](#), [tfd_half_cauchy\(\)](#), [tfd_half_normal\(\)](#), [tfd_hidden_markov_model\(\)](#), [tfd_horseshoe\(\)](#), [tfd_independent\(\)](#), [tfd_inverse_gamma\(\)](#), [tfd_inverse_gaussian\(\)](#), [tfd_johnson_s_u\(\)](#), [tfd_joint_distribution_named_auto_batched\(\)](#), [tfd_joint_distribution_named\(\)](#), [tfd_joint_distribution_sequential_auto_batched\(\)](#), [tfd_joint_distribution_sequential\(\)](#), [tfd_kumaraswamy\(\)](#), [tfd_laplace\(\)](#), [tfd_linear_gaussian_state_space\(\)](#), [tfd_lkj\(\)](#), [tfd_log_logistic\(\)](#), [tfd_log_normal\(\)](#), [tfd_logistic\(\)](#), [tfd_mixture_same_family\(\)](#), [tfd_mixture\(\)](#), [tfd_multinomial\(\)](#), [tfd_multivariate_normal_diag_plus_low_rank\(\)](#), [tfd_multivariate_normal\(\)](#), [tfd_multivariate_normal_full_covariance\(\)](#), [tfd_multivariate_normal_linear_operator\(\)](#), [tfd_multivariate_normal_tri_l\(\)](#), [tfd_multivariate_student_t_linear_operator\(\)](#), [tfd_negative_binomial\(\)](#), [tfd_normal\(\)](#), [tfd_one_hot_categorical\(\)](#), [tfd_pareto\(\)](#), [tfd_pixel_cnn\(\)](#), [tfd_poisson_log_normal_quadrature\(\)](#), [tfd_poisson\(\)](#), [tfd_power_spherical\(\)](#), [tfd_probit_bernoulli\(\)](#), [tfd_quantized\(\)](#), [tfd_relaxed_bernoulli\(\)](#), [tfd_relaxed_one_hot_categorical\(\)](#), [tfd_sample_distribution\(\)](#), [tfd_sinh_arcsinh\(\)](#), [tfd_skellam\(\)](#), [tfd_spherical_uniform\(\)](#), [tfd_student_t_process\(\)](#), [tfd_student_t\(\)](#), [tfd_transformed_distribution\(\)](#), [tfd_triangular\(\)](#), [tfd_truncated_cauchy\(\)](#), [tfd_truncated_normal\(\)](#), [tfd_uniform\(\)](#), [tfd_variational_gaussian\(\)](#), [tfd_vector_diffeomixture\(\)](#), [tfd_vector_exponential_diag\(\)](#), [tfd_vector_exponential_linear_operator\(\)](#), [tfd_vector_laplace_diag\(\)](#), [tfd_vector_laplace_linear_operator\(\)](#), [tfd_vector_sinh_arcsinh_diag\(\)](#), [tfd_von_mises_fisher\(\)](#), [tfd_von_mises\(\)](#), [tfd_weibull\(\)](#), [tfd_wishart_linear_operator\(\)](#), [tfd_wishart_tri_l\(\)](#), [tfd_wishart\(\)](#), [tfd_zipf\(\)](#)

tfd_beta

*Beta distribution***Description**

The Beta distribution is defined over the (0, 1) interval using parameters `concentration1` (aka "alpha") and `concentration0` (aka "beta").

Usage

```
tfd_beta(
  concentration1 = NULL,
  concentration0 = NULL,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "Beta"
)
```

Arguments

`concentration1` Positive floating-point Tensor indicating mean number of successes; aka "alpha". Implies `self.dtype` and `self.batch_shape`, i.e., `concentration1.shape = [N1, N2, ..., Nm] = self.batch_shape`.

`concentration0` Positive floating-point Tensor indicating mean number of failures; aka "beta". Otherwise has same semantics as `concentration1`.

`validate_args` Logical, default `FALSE`. When `TRUE` distribution parameters are checked for validity despite possibly degrading runtime performance. When `FALSE` invalid inputs may silently render incorrect outputs. Default value: `FALSE`.

`allow_nan_stats` Logical, default `TRUE`. When `TRUE`, statistics (e.g., mean, mode, variance) use the value `NaN` to indicate the result is undefined. When `FALSE`, an exception is raised if one or more of the statistic's batch members are undefined.

`name` name prefixed to Ops created by this class.

Details**Mathematical Details**

The probability density function (pdf) is,

$$\text{pdf}(x; \alpha, \beta) = \frac{x^{\alpha-1} (1-x)^{\beta-1}}{Z}$$

$$Z = \frac{\Gamma(\alpha)\Gamma(\beta)}{\Gamma(\alpha+\beta)}$$

where:

- `concentration1` = `alpha`,
- `concentration0` = `beta`,
- `Z` is the normalization constant, and,

- Gamma is the [gamma function](#). The concentration parameters represent mean total counts of a 1 or a 0, i.e.,

```
concentration1 = alpha = mean * total_concentration
concentration0 = beta = (1. - mean) * total_concentration
```

where mean in (0, 1) and total_concentration is a positive real number representing a mean total_count = concentration1 + concentration0. Distribution parameters are automatically broadcast in all functions; see examples for details. Warning: The samples can be zero due to finite precision. This happens more often when some of the concentrations are very small. Make sure to round the samples to `np.finfo(dtype).tiny` before computing the density. Samples of this distribution are reparameterized (pathwise differentiable). The derivatives are computed using the approach described in the paper [Michael Figurnov, Shakir Mohamed, Andriy Mnih. Implicit Reparameterization Gradients, 2018](#)

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_beta_binomial	<i>Beta-Binomial compound distribution</i>
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Description

The Beta-Binomial distribution is parameterized by (a batch of) `total_count` parameters, the number of trials per draw from Binomial distributions where the probabilities of success per trial are drawn from underlying Beta distributions; the Beta distributions are parameterized by `concentration1` (aka 'alpha') and `concentration0` (aka 'beta'). Mathematically, it is (equivalent to) a special case of the Dirichlet-Multinomial over two classes, although the computational representation is slightly different: while the Beta-Binomial is a distribution over the number of successes in `total_count` trials, the two-class Dirichlet-Multinomial is a distribution over the number of successes and failures.

Usage

```
tfd_beta_binomial(
    total_count,
    concentration1,
    concentration0,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = "BetaBinomial"
)
```

Arguments

<code>total_count</code>	Non-negative integer-valued tensor, whose dtype is the same as <code>concentration1</code> and <code>concentration0</code> . The shape is broadcastable to $[N_1, \dots, N_m]$ with $m \geq 0$. When <code>total_count</code> is broadcast with <code>concentration1</code> and <code>concentration0</code> , it defines the distribution as a batch of $N_1 \times \dots \times N_m$ different Beta-Binomial distributions. Its components should be equal to integer values.
<code>concentration1</code>	Positive floating-point Tensor indicating mean number of successes. Specifically, the expected number of successes is $\text{total_count} * \text{concentration1} / (\text{concentration1} + \text{concentration0})$.
<code>concentration0</code>	Positive floating-point Tensor indicating mean number of failures; see description of <code>concentration1</code> for details.
<code>validate_args</code>	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
<code>allow_nan_stats</code>	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Details

Mathematical Details

The Beta-Binomial is a distribution over the number of successes in `total_count` independent Binomial trials, with each trial having the same probability of success, the underlying probability being unknown but drawn from a Beta distribution with known parameters. The probability mass function (pmf) is,

$$\text{pmf}(k; n, a, b) = \text{Beta}(k + a, n - k + b) / Z$$

$$Z = (k! (n - k)! / n!) * \text{Beta}(a, b)$$

where:

- `concentration1 = a > 0`,
- `concentration0 = b > 0`,
- `total_count = n`, `n` a positive integer,
- `n!` is `n` factorial,
- $\text{Beta}(x, y) = \text{Gamma}(x) \text{Gamma}(y) / \text{Gamma}(x + y)$ is the **beta function**, and
- `Gamma` is the **gamma function**.

Dirichlet-Multinomial is a **compound distribution**, i.e., its samples are generated as follows.

1. Choose success probabilities: `probs ~ Beta(concentration1, concentration0)`
 2. Draw integers representing the number of successes: `counts ~ Binomial(total_count, probs)`
- Distribution parameters are automatically broadcast in all functions; see examples for details.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature`

```
tfd_poisson(), tfd_power_spherical(), tfd_probit_bernoulli(), tfd_quantized(), tfd_relaxed_bernoulli(),
tfd_relaxed_one_hot_categorical(), tfd_sample_distribution(), tfd_sinh_arcsinh(),
tfd_skellam(), tfd_spherical_uniform(), tfd_student_t_process(), tfd_student_t(), tfd_transformed_distribution(),
tfd_triangular(), tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian(),
tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()
```

tfd_binomial	<i>Binomial distribution</i>
--------------	------------------------------

Description

This distribution is parameterized by `probs`, a (batch of) probabilities for drawing a 1 and `total_count`, the number of trials per draw from the Binomial.

Usage

```
tfd_binomial(
    total_count,
    logits = NULL,
    probs = NULL,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = "Beta"
)
```

Arguments

<code>total_count</code>	Non-negative floating point tensor with shape broadcastable to $[N_1, \dots, N_m]$ with $m \geq 0$ and the same dtype as <code>probs</code> or <code>logits</code> . Defines this as a batch of $N_1 \times \dots \times N_m$ different Binomial distributions. Its components should be equal to integer values.
<code>logits</code>	Floating point tensor representing the log-odds of a positive event with shape broadcastable to $[N_1, \dots, N_m]$ $m \geq 0$, and the same dtype as <code>total_count</code> . Each entry represents logits for the probability of success for independent Binomial distributions. Only one of <code>logits</code> or <code>probs</code> should be passed in.
<code>probs</code>	Positive floating point tensor with shape broadcastable to $[N_1, \dots, N_m]$ $m \geq 0$, <code>probs</code> in $[0, 1]$. Each entry represents the probability of success for independent Binomial distributions. Only one of <code>logits</code> or <code>probs</code> should be passed in.
<code>validate_args</code>	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .

<code>allow_nan_stats</code>	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Details

Mathematical Details

The Binomial is a distribution over the number of 1's in `total_count` independent trials, with each trial having the same probability of 1, i.e., `probs`.

The probability mass function (pmf) is,

$$\text{pmf}(k; n, p) = p^k (1 - p)^{(n - k)} / Z$$

$$Z = \sum_{k=0}^n \binom{n}{k} p^k (1 - p)^{n - k} / n!$$

where:

- `total_count = n`,
- `probs = p`,
- `Z` is the normalizing constant, and,
- `n!` is the factorial of `n`.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`.

[tfd_triangular\(\)](#), [tfd_truncated_cauchy\(\)](#), [tfd_truncated_normal\(\)](#), [tfd_uniform\(\)](#), [tfd_variational_gaussian\(\)](#), [tfd_vector_diffeomixture\(\)](#), [tfd_vector_exponential_diag\(\)](#), [tfd_vector_exponential_linear_operator\(\)](#), [tfd_vector_laplace_diag\(\)](#), [tfd_vector_laplace_linear_operator\(\)](#), [tfd_vector_sinh_arcsinh_diag\(\)](#), [tfd_von_mises_fisher\(\)](#), [tfd_von_mises\(\)](#), [tfd_weibull\(\)](#), [tfd_wishart_linear_operator\(\)](#), [tfd_wishart_tri_l\(\)](#), [tfd_wishart\(\)](#), [tfd_zipf\(\)](#)

tfd_blockwise	<i>Blockwise distribution</i>
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Description

Blockwise distribution

Usage

```
tfd_blockwise(
    distributions,
    dtype_override = NULL,
    validate_args = FALSE,
    allow_nan_stats = FALSE,
    name = "Blockwise"
)
```

Arguments

distributions list of Distribution instances. All distribution instances must have the same `batch_shape` and all must have `'event_ndims==1'`, i.e., be vector-variate distributions.

dtype_override samples of distributions will be cast to this dtype. If unspecified, all distributions must have the same dtype. Default value: NULL (i.e., do not cast).

validate_args Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.

allow_nan_stats Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.

name name prefixed to Ops created by this class.

Value

a distribution instance.

See Also

For usage examples see e.g. [tfd_sample\(\)](#), [tfd_log_prob\(\)](#), [tfd_mean\(\)](#).

tfd_categorical	<i>Categorical distribution over integers</i>
-----------------	---

Description

The Categorical distribution is parameterized by either probabilities or log-probabilities of a set of K classes. It is defined over the integers $\{0, 1, \dots, K-1\}$.

Usage

```
tfd_categorical(
  logits = NULL,
  probs = NULL,
  dtype = tf.int32,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "Categorical"
)
```

Arguments

<code>logits</code>	An N-D Tensor, $N \geq 1$, representing the log probabilities of a set of Categorical distributions. The first $N-1$ dimensions index into a batch of independent distributions and the last dimension represents a vector of logits for each class. Only one of <code>logits</code> or <code>probs</code> should be passed in.
<code>probs</code>	An N-D Tensor, $N \geq 1$, representing the probabilities of a set of Categorical distributions. The first $N-1$ dimensions index into a batch of independent distributions and the last dimension represents a vector of probabilities for each class. Only one of <code>logits</code> or <code>probs</code> should be passed in.
<code>dtype</code>	The type of the event samples (default: <code>int32</code>).
<code>validate_args</code>	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
<code>allow_nan_stats</code>	Logical, default <code>TRUE</code> . When <code>TRUE</code> , statistics (e.g., mean, mode, variance) use the value <code>NaN</code> to indicate the result is undefined. When <code>FALSE</code> , an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Details

The Categorical distribution is closely related to the `OneHotCategorical` and `Multinomial` distributions. The Categorical distribution can be intuited as generating samples according to `argmax{ OneHotCategorical(probs) }` itself being identical to `argmax{ Multinomial(probs, total_count=1) }`.

Mathematical Details

The probability mass function (pmf) is,

```
pmf(k; pi) = prod_j pi_j**[k == j]
```

Pitfalls

The number of classes, K, must not exceed:

- the largest integer representable by self.dtype, i.e., $2^{*(\text{mantissa_bits}+1)}$ (IEEE 754),
- the maximum Tensor index, i.e., $2^{*31}-1$.

Note: This condition is validated only when `validate_args = TRUE`.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_cauchy

Cauchy distribution with location loc and scale scale

Description

Mathematical details

Usage

```
tfd_cauchy(
  loc,
  scale,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "Cauchy"
)
```

Arguments

loc	Floating point tensor; the modes of the distribution(s).
scale	Floating point tensor; the locations of the distribution(s). Must contain only positive values.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

The probability density function (pdf) is,

$$\text{pdf}(x; \text{loc}, \text{scale}) = 1 / (\pi \text{scale} (1 + z^{**2}))$$

$$z = (x - \text{loc}) / \text{scale}$$

where loc is the location, and scale is the scale. The Cauchy distribution is a member of the **location-scale family**, i.e. $Y \sim \text{Cauchy}(\text{loc}, \text{scale})$ is equivalent to,

$$X \sim \text{Cauchy}(\text{loc}=0, \text{scale}=1)$$

$$Y = \text{loc} + \text{scale} * X$$

Value

a distribution instance.

See Also

For usage examples see e.g. [tfd_sample\(\)](#), [tfd_log_prob\(\)](#), [tfd_mean\(\)](#).

Other distributions: [tfd_autoregressive\(\)](#), [tfd_batch_reshape\(\)](#), [tfd_bates\(\)](#), [tfd_bernoulli\(\)](#), [tfd_beta_binomial\(\)](#), [tfd_beta\(\)](#), [tfd_binomial\(\)](#), [tfd_categorical\(\)](#), [tfd_chi2\(\)](#), [tfd_chi\(\)](#), [tfd_cholesky_lkj\(\)](#), [tfd_continuous_bernoulli\(\)](#), [tfd_deterministic\(\)](#), [tfd_dirichlet_multinomial\(\)](#), [tfd_dirichlet\(\)](#), [tfd_empirical\(\)](#), [tfd_exp_gamma\(\)](#), [tfd_exp_inverse_gamma\(\)](#), [tfd_exponential\(\)](#),

[tfd_gamma_gamma\(\)](#), [tfd_gamma\(\)](#), [tfd_gaussian_process_regression_model\(\)](#), [tfd_gaussian_process\(\)](#),
[tfd_generalized_normal\(\)](#), [tfd_geometric\(\)](#), [tfd_gumbel\(\)](#), [tfd_half_cauchy\(\)](#), [tfd_half_normal\(\)](#),
[tfd_hidden_markov_model\(\)](#), [tfd_horseshoe\(\)](#), [tfd_independent\(\)](#), [tfd_inverse_gamma\(\)](#),
[tfd_inverse_gaussian\(\)](#), [tfd_johnson_s_u\(\)](#), [tfd_joint_distribution_named_auto_batched\(\)](#),
[tfd_joint_distribution_named\(\)](#), [tfd_joint_distribution_sequential_auto_batched\(\)](#),
[tfd_joint_distribution_sequential\(\)](#), [tfd_kumaraswamy\(\)](#), [tfd_laplace\(\)](#), [tfd_linear_gaussian_state_space\(\)](#),
[tfd_lkj\(\)](#), [tfd_log_logistic\(\)](#), [tfd_log_normal\(\)](#), [tfd_logistic\(\)](#), [tfd_mixture_same_family\(\)](#),
[tfd_mixture\(\)](#), [tfd_multinomial\(\)](#), [tfd_multivariate_normal_diag_plus_low_rank\(\)](#), [tfd_multivariate_normal\(\)](#),
[tfd_multivariate_normal_full_covariance\(\)](#), [tfd_multivariate_normal_linear_operator\(\)](#),
[tfd_multivariate_normal_tri_l\(\)](#), [tfd_multivariate_student_t_linear_operator\(\)](#), [tfd_negative_binomial\(\)](#),
[tfd_normal\(\)](#), [tfd_one_hot_categorical\(\)](#), [tfd_pareto\(\)](#), [tfd_pixel_cnn\(\)](#), [tfd_poisson_log_normal_quadrature\(\)](#),
[tfd_poisson\(\)](#), [tfd_power_spherical\(\)](#), [tfd_probit_bernoulli\(\)](#), [tfd_quantized\(\)](#), [tfd_relaxed_bernoulli\(\)](#),
[tfd_relaxed_one_hot_categorical\(\)](#), [tfd_sample_distribution\(\)](#), [tfd_sinh_arcsinh\(\)](#),
[tfd_skellam\(\)](#), [tfd_spherical_uniform\(\)](#), [tfd_student_t_process\(\)](#), [tfd_student_t\(\)](#), [tfd_transformed_distribution\(\)](#),
[tfd_triangular\(\)](#), [tfd_truncated_cauchy\(\)](#), [tfd_truncated_normal\(\)](#), [tfd_uniform\(\)](#), [tfd_variational_gaussian\(\)](#),
[tfd_vector_diffeomixture\(\)](#), [tfd_vector_exponential_diag\(\)](#), [tfd_vector_exponential_linear_operator\(\)](#),
[tfd_vector_laplace_diag\(\)](#), [tfd_vector_laplace_linear_operator\(\)](#), [tfd_vector_sinh_arcsinh_diag\(\)](#),
[tfd_von_mises_fisher\(\)](#), [tfd_von_mises\(\)](#), [tfd_weibull\(\)](#), [tfd_wishart_linear_operator\(\)](#),
[tfd_wishart_tri_l\(\)](#), [tfd_wishart\(\)](#), [tfd_zipf\(\)](#)

tfd_cdf

Cumulative distribution function. Given random variable X, the cumulative distribution function cdf is: $\text{cdf}(x) := P[X \leq x]$

Description

Cumulative distribution function. Given random variable X, the cumulative distribution function cdf is: $\text{cdf}(x) := P[X \leq x]$

Usage

```
tfd_cdf(distribution, value, ...)
```

Arguments

distribution	The distribution being used.
value	float or double Tensor.
...	Additional parameters passed to Python.

Value

a Tensor of shape $\text{sample_shape}(x) + \text{self}\batch_shape with values of type $\text{self}\$\text{dtype}$.

See Also

Other `distribution_methods`: [tfd_covariance\(\)](#), [tfd_cross_entropy\(\)](#), [tfd_entropy\(\)](#), [tfd_kl_divergence\(\)](#), [tfd_log_cdf\(\)](#), [tfd_log_prob\(\)](#), [tfd_log_survival_function\(\)](#), [tfd_mean\(\)](#), [tfd_mode\(\)](#), [tfd_prob\(\)](#), [tfd_quantile\(\)](#), [tfd_sample\(\)](#), [tfd_stddev\(\)](#), [tfd_survival_function\(\)](#), [tfd_variance\(\)](#)

Examples

```
d <- tfd_normal(loc = c(1, 2), scale = c(1, 0.5))
x <- d %>% tfd_sample()
d %>% tfd_cdf(x)
```

tfd_chi	<i>Chi distribution</i>
---------	-------------------------

Description

The Chi distribution is defined over nonnegative real numbers and uses a degrees of freedom ("df") parameter.

Usage

```
tfd_chi(df, validate_args = FALSE, allow_nan_stats = TRUE, name = "Chi")
```

Arguments

df	Floating point tensor, the degrees of freedom of the distribution(s). df must contain only positive values.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details**Mathematical Details**

The probability density function (pdf) is,

$$\text{pdf}(x; \text{df}, x \geq 0) = x^{(\text{df} - 1)} \exp(-0.5 x^2) / Z$$

$$Z = 2^{(0.5 \text{df} - 1)} \Gamma(0.5 \text{df})$$

where:

- df denotes the degrees of freedom,
- Z is the normalization constant, and,
- Gamma is the [gamma function](#).

The Chi distribution is a transformation of the Chi2 distribution; it is the distribution of the positive square root of a variable obeying a Chi distribution.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag_plus_low_rank_linear_operator()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_chi2

Chi Square distribution

Description

The Chi2 distribution is defined over positive real numbers using a degrees of freedom ("df") parameter.

Usage

```
tfd_chi2(df, validate_args = FALSE, allow_nan_stats = TRUE, name = "Chi2")
```

Arguments

`df` Floating point tensor, the degrees of freedom of the distribution(s). `df` must contain only positive values.

<code>validate_args</code>	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
<code>allow_nan_stats</code>	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Details

Mathematical Details

The probability density function (pdf) is,

$$\text{pdf}(x; \text{df}, x > 0) = x^{(0.5 \text{df} - 1)} \exp(-0.5 x) / Z$$

$$Z = 2^{(0.5 \text{df})} \text{Gamma}(0.5 \text{df})$$

where

- `df` denotes the degrees of freedom,
- `Z` is the normalization constant, and,
- `Gamma` is the [gamma function](#). The Chi2 distribution is a special case of the Gamma distribution, i.e.,

$$\text{Chi2}(\text{df}) = \text{Gamma}(\text{concentration}=\text{df} * 0.5, \text{rate}=0.5)$$

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`

```
tfd_poisson(), tfd_power_spherical(), tfd_probit_bernoulli(), tfd_quantized(), tfd_relaxed_bernoulli(),
tfd_relaxed_one_hot_categorical(), tfd_sample_distribution(), tfd_sinh_arcsinh(),
tfd_skellam(), tfd_spherical_uniform(), tfd_student_t_process(), tfd_student_t(), tfd_transformed_distribution(),
tfd_triangular(), tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian(),
tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()
```

tfd_cholesky_lkj	<i>The CholeskyLKJ distribution on cholesky factors of correlation matrices</i>
------------------	---

Description

This is a one-parameter family of distributions on cholesky factors of correlation matrices. In other words, if $X \sim \text{CholeskyLKJ}(c)$, then $X @ X^T \sim \text{LKJ}(c)$. For more details on the LKJ distribution, see `tfd_lkj`.

Usage

```
tfd_cholesky_lkj(
    dimension,
    concentration,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = "CholeskyLKJ"
)
```

Arguments

<code>dimension</code>	integer. The dimension of the correlation matrices to sample.
<code>concentration</code>	float or double Tensor. The positive concentration parameter of the CholeskyLKJ distributions.
<code>validate_args</code>	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
<code>allow_nan_stats</code>	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_continuous_bernoulli

Continuous Bernoulli distribution.

Description

This distribution is parameterized by `probs`, a (batch of) parameters taking values in $(0, 1)$. Note that, unlike in the Bernoulli case, `probs` does not correspond to a probability, but the same name is used due to the similarity with the Bernoulli.

Usage

```
tfd_continuous_bernoulli(
  logits = NULL,
  probs = NULL,
  dtype = tf$float32,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "ContinuousBernoulli"
)
```

Arguments

<code>logits</code>	An N-D Tensor. Each entry in the Tensor parameterizes an independent continuous Bernoulli distribution with parameter $\text{sigmoid}(\text{logits})$. Only one of <code>logits</code> or <code>probs</code> should be passed in. Note that this does not correspond to the log-odds as in the Bernoulli case.
<code>probs</code>	An N-D Tensor representing the parameter of a continuous Bernoulli. Each entry in the Tensor parameterizes an independent continuous Bernoulli distribution. Only one of <code>logits</code> or <code>probs</code> should be passed in. Note that this also does not correspond to a probability as in the Bernoulli case.
<code>dtype</code>	The type of the event samples. Default: <code>float32</code> .
<code>validate_args</code>	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
<code>allow_nan_stats</code>	Logical, default <code>TRUE</code> . When <code>TRUE</code> , statistics (e.g., mean, mode, variance) use the value <code>NaN</code> to indicate the result is undefined. When <code>FALSE</code> , an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Details**Mathematical Details**

The continuous Bernoulli is a distribution over the interval $[0, 1]$, parameterized by `probs` in $(0, 1)$. The probability density function (pdf) is,

$$\text{pdf}(x; \text{probs}) = \text{probs}^x * (1 - \text{probs})^{(1 - x)} * C(\text{probs})$$

$$C(\text{probs}) = (2 * \text{atanh}(1 - 2 * \text{probs}) / (1 - 2 * \text{probs})) \text{ if } \text{probs} \neq 0.5 \text{ else } 2.$$

While the normalizing constant $C(\text{probs})$ is a continuous function of `probs` (even at `probs = 0.5`), computing it at values close to 0.5 can result in numerical instabilities due to 0/0 errors. A Taylor approximation of $C(\text{probs})$ is thus used for values of `probs` in a small interval $[\text{lims}[0], \text{lims}[1]]$ around 0.5. For more details, see Loaiza-Ganem and Cunningham (2019). NOTE: Unlike the Bernoulli, numerical instabilities can happen for `probs` very close to 0 or 1. Current implementation allows any value in $(0, 1)$, but this could be changed to $(1e-6, 1-1e-6)$ to avoid these issues.

Value

a distribution instance.

References

- Loaiza-Ganem G and Cunningham JP. The continuous Bernoulli: fixing a pervasive error in variational autoencoders. NeurIPS2019. <https://arxiv.org/abs/1907.06845>

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_covariance

Covariance.

Description

Covariance is (possibly) defined only for non-scalar-event distributions. For example, for a length- k , vector-valued distribution, it is calculated as, $\text{Cov}[i, j] = \text{Covariance}(X_i, X_j) = E[(X_i - E[X_i])(X_j - E[X_j])]$ where Cov is a (batch of) $k \times k$ matrix, $0 \leq (i, j) < k$, and E denotes expectation.

Usage

```
tfd_covariance(distribution, ...)
```

Arguments

`distribution` The distribution being used.
`...` Additional parameters passed to Python.

Details

Alternatively, for non-vector, multivariate distributions (e.g., matrix-valued, Wishart), Covariance shall return a (batch of) matrices under some vectorization of the events, i.e., $\text{Cov}[i, j] = \text{Covariance}(\text{Vec}(X)_i, \text{Vec}(X)_j) = [\text{as above}]$ where Cov is a (batch of) $k \times k$ matrices, $0 \leq (i, j) < k = \text{reduce_prod}(\text{event_shape})$, and Vec is some function mapping indices of this distribution's event dimensions to indices of a length- k vector.

Value

Floating-point Tensor with shape $[B_1, \dots, B_n, k, k]$ where the first n dimensions are batch coordinates and $k = \text{reduce_prod}(\text{self.event_shape})$.

See Also

Other distribution_methods: [tfd_cdf\(\)](#), [tfd_cross_entropy\(\)](#), [tfd_entropy\(\)](#), [tfd_kl_divergence\(\)](#), [tfd_log_cdf\(\)](#), [tfd_log_prob\(\)](#), [tfd_log_survival_function\(\)](#), [tfd_mean\(\)](#), [tfd_mode\(\)](#), [tfd_prob\(\)](#), [tfd_quantile\(\)](#), [tfd_sample\(\)](#), [tfd_stddev\(\)](#), [tfd_survival_function\(\)](#), [tfd_variance\(\)](#)

Examples

```
d <- tfd_normal(loc = c(1, 2), scale = c(1, 0.5))
d %>% tfd_variance()
```

tfd_cross_entropy	<i>Computes the (Shannon) cross entropy.</i>
-------------------	--

Description

Denote this distribution (self) by P and the other distribution by Q . Assuming P, Q are absolutely continuous with respect to one another and permit densities $p(x) \text{ dr}(x)$ and $q(x) \text{ dr}(x)$, (Shannon) cross entropy is defined as: $H[P, Q] = E_p[-\log q(X)] = -\int_F p(x) \log q(x) \text{ dr}(x)$ where F denotes the support of the random variable $X \sim P$.

Usage

```
tfd_cross_entropy(distribution, other, name = "cross_entropy")
```

Arguments

distribution	The distribution being used.
other	tfp\$distributions\$Distribution instance.
name	String prepended to names of ops created by this function.

Value

cross_entropy: self.dtype Tensor with shape [B1, ..., Bn] representing n different calculations of (Shannon) cross entropy.

See Also

Other distribution_methods: [tfd_cdf\(\)](#), [tfd_covariance\(\)](#), [tfd_entropy\(\)](#), [tfd_kl_divergence\(\)](#), [tfd_log_cdf\(\)](#), [tfd_log_prob\(\)](#), [tfd_log_survival_function\(\)](#), [tfd_mean\(\)](#), [tfd_mode\(\)](#), [tfd_prob\(\)](#), [tfd_quantile\(\)](#), [tfd_sample\(\)](#), [tfd_stddev\(\)](#), [tfd_survival_function\(\)](#), [tfd_variance\(\)](#)

Examples

```
d1 <- tfd_normal(loc = 1, scale = 1)
d2 <- tfd_normal(loc = 2, scale = 1)
d1 %>% tfd_cross_entropy(d2)
```

tfd_deterministic *Scalar Deterministic distribution on the real line*

Description

The scalar Deterministic distribution is parameterized by a (batch) point `loc` on the real line. The distribution is supported at this point only, and corresponds to a random variable that is constant, equal to `loc`. See [Degenerate rv](#).

Usage

```
tfd_deterministic(
  loc,
  atol = NULL,
  rtol = NULL,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "Deterministic"
)
```

Arguments

<code>loc</code>	Numeric Tensor of shape [B1, ..., Bb], with $b \geq 0$. The point (or batch of points) on which this distribution is supported.
<code>atol</code>	Non-negative Tensor of same dtype as <code>loc</code> and broadcastable shape. The absolute tolerance for comparing closeness to <code>loc</code> . Default is 0.
<code>rtol</code>	Non-negative Tensor of same dtype as <code>loc</code> and broadcastable shape. The relative tolerance for comparing closeness to <code>loc</code> . Default is 0.

<code>validate_args</code>	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
<code>allow_nan_stats</code>	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Details

Mathematical Details

The probability mass function (pmf) and cumulative distribution function (cdf) are

$$\begin{aligned} \text{pmf}(x; \text{loc}) &= 1, \text{ if } x == \text{loc}, \text{ else } 0 \\ \text{cdf}(x; \text{loc}) &= 1, \text{ if } x \geq \text{loc}, \text{ else } 0 \end{aligned}$$

Value

a distribution instance.

See Also

For usage examples see e.g. [tfd_sample\(\)](#), [tfd_log_prob\(\)](#), [tfd_mean\(\)](#).

Other distributions: [tfd_autoregressive\(\)](#), [tfd_batch_reshape\(\)](#), [tfd_bates\(\)](#), [tfd_bernoulli\(\)](#), [tfd_beta_binomial\(\)](#), [tfd_beta\(\)](#), [tfd_binomial\(\)](#), [tfd_categorical\(\)](#), [tfd_cauchy\(\)](#), [tfd_chi2\(\)](#), [tfd_chi\(\)](#), [tfd_cholesky_lkj\(\)](#), [tfd_continuous_bernoulli\(\)](#), [tfd_dirichlet_multinomial\(\)](#), [tfd_dirichlet\(\)](#), [tfd_empirical\(\)](#), [tfd_exp_gamma\(\)](#), [tfd_exp_inverse_gamma\(\)](#), [tfd_exponential\(\)](#), [tfd_gamma_gamma\(\)](#), [tfd_gamma\(\)](#), [tfd_gaussian_process_regression_model\(\)](#), [tfd_gaussian_process\(\)](#), [tfd_generalized_normal\(\)](#), [tfd_geometric\(\)](#), [tfd_gumbel\(\)](#), [tfd_half_cauchy\(\)](#), [tfd_half_normal\(\)](#), [tfd_hidden_markov_model\(\)](#), [tfd_horseshoe\(\)](#), [tfd_independent\(\)](#), [tfd_inverse_gamma\(\)](#), [tfd_inverse_gaussian\(\)](#), [tfd_johnson_s_u\(\)](#), [tfd_joint_distribution_named_auto_batched\(\)](#), [tfd_joint_distribution_named\(\)](#), [tfd_joint_distribution_sequential_auto_batched\(\)](#), [tfd_joint_distribution_sequential\(\)](#), [tfd_kumaraswamy\(\)](#), [tfd_laplace\(\)](#), [tfd_linear_gaussian_state_space\(\)](#), [tfd_lkj\(\)](#), [tfd_log_logistic\(\)](#), [tfd_log_normal\(\)](#), [tfd_logistic\(\)](#), [tfd_mixture_same_family\(\)](#), [tfd_mixture\(\)](#), [tfd_multinomial\(\)](#), [tfd_multivariate_normal_diag_plus_low_rank\(\)](#), [tfd_multivariate_normal_full_covariance\(\)](#), [tfd_multivariate_normal_linear_operator\(\)](#), [tfd_multivariate_normal_tri_l\(\)](#), [tfd_multivariate_student_t_linear_operator\(\)](#), [tfd_negative_binomial\(\)](#), [tfd_normal\(\)](#), [tfd_one_hot_categorical\(\)](#), [tfd_pareto\(\)](#), [tfd_pixel_cnn\(\)](#), [tfd_poisson_log_normal_quadrature\(\)](#), [tfd_poisson\(\)](#), [tfd_power_spherical\(\)](#), [tfd_probit_bernoulli\(\)](#), [tfd_quantized\(\)](#), [tfd_relaxed_bernoulli\(\)](#), [tfd_relaxed_one_hot_categorical\(\)](#), [tfd_sample_distribution\(\)](#), [tfd_sinh_arcsinh\(\)](#), [tfd_skellam\(\)](#), [tfd_spherical_uniform\(\)](#), [tfd_student_t_process\(\)](#), [tfd_student_t\(\)](#), [tfd_transformed_distribution\(\)](#), [tfd_triangular\(\)](#), [tfd_truncated_cauchy\(\)](#), [tfd_truncated_normal\(\)](#), [tfd_uniform\(\)](#), [tfd_variational_gaussian\(\)](#), [tfd_vector_diffeomixture\(\)](#), [tfd_vector_exponential_diag\(\)](#), [tfd_vector_exponential_linear_operator\(\)](#), [tfd_vector_laplace_diag\(\)](#), [tfd_vector_laplace_linear_operator\(\)](#), [tfd_vector_sinh_arcsinh_diag\(\)](#), [tfd_von_mises_fisher\(\)](#), [tfd_von_mises\(\)](#), [tfd_weibull\(\)](#), [tfd_wishart_linear_operator\(\)](#), [tfd_wishart_tri_l\(\)](#), [tfd_wishart\(\)](#), [tfd_zipf\(\)](#)

tfd_dirichlet	<i>Dirichlet distribution</i>
---------------	-------------------------------

Description

The Dirichlet distribution is defined over the $(k-1)$ -simplex using a positive, length- k vector concentration ($k > 1$). The Dirichlet is identically the Beta distribution when $k = 2$.

Usage

```
tfd_dirichlet(
  concentration,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "Dirichlet"
)
```

Arguments

concentration	Positive floating-point Tensor indicating mean number of class occurrences; aka "alpha". Implies <code>self.dtype</code> , and <code>self.batch_shape</code> , <code>self.event_shape</code> , i.e., if <code>concentration.shape = [N1, N2, ..., Nm, k]</code> then <code>batch_shape = [N1, N2, ..., Nm]</code> and <code>event_shape = [k]</code> .
validate_args	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
allow_nan_stats	Logical, default <code>TRUE</code> . When <code>TRUE</code> , statistics (e.g., mean, mode, variance) use the value <code>NaN</code> to indicate the result is undefined. When <code>FALSE</code> , an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

Mathematical Details

The Dirichlet is a distribution over the open $(k-1)$ -simplex, i.e.,

$$S^{k-1} = \{ (x_0, \dots, x_{k-1}) \text{ in } \mathbb{R}^k : \sum_j x_j = 1 \text{ and } \forall_j x_j > 0 \}.$$

The probability density function (pdf) is,

$$\text{pdf}(x; \alpha) = \prod_j x_j^{(\alpha_j - 1)} / Z$$

$$Z = \prod_j \Gamma(\alpha_j) / \Gamma(\sum_j \alpha_j)$$

where:

- x in S^{k-1} , i.e., the $(k-1)$ -simplex,

- concentration = alpha = [alpha_0, ..., alpha_{k-1}], alpha_j > 0,
- Z is the normalization constant aka the [multivariate beta function](#), and,
- Gamma is the [gamma function](#).

The concentration represents mean total counts of class occurrence, i.e.,

```
concentration = alpha = mean * total_concentration
```

where mean in S^{k-1} and total_concentration is a positive real number representing a mean total count. Distribution parameters are automatically broadcast in all functions; see examples for details. Warning: Some components of the samples can be zero due to finite precision. This happens more often when some of the concentrations are very small. Make sure to round the samples to `np.finfo(dtype).tiny` before computing the density. Samples of this distribution are reparameterized (pathwise differentiable). The derivatives are computed using the approach described in the paper [Michael Figurnov, Shakir Mohamed, Andriy Mnih. Implicit Reparameterization Gradients, 2018](#)

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag_plus_low_rank_linear_operator()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_dirichlet_multinomial

Dirichlet-Multinomial compound distribution

Description

The Dirichlet-Multinomial distribution is parameterized by a (batch of) length-K concentration vectors ($K > 1$) and a `total_count` number of trials, i.e., the number of trials per draw from the DirichletMultinomial. It is defined over a (batch of) length-K vector counts such that `tf.reduce_sum(counts, -1) = total_count`. The Dirichlet-Multinomial is identically the Beta-Binomial distribution when $K = 2$.

Usage

```
tfd_dirichlet_multinomial(
    total_count,
    concentration,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = "DirichletMultinomial"
)
```

Arguments

<code>total_count</code>	Non-negative floating point tensor, whose dtype is the same as <code>concentration</code> . The shape is broadcastable to $[N_1, \dots, N_m]$ with $m \geq 0$. Defines this as a batch of $N_1 \times \dots \times N_m$ different Dirichlet multinomial distributions. Its components should be equal to integer values.
<code>concentration</code>	Positive floating point tensor, whose dtype is the same as <code>n</code> with shape broadcastable to $[N_1, \dots, N_m, K]$ $m \geq 0$. Defines this as a batch of $N_1 \times \dots \times N_m$ different K class Dirichlet multinomial distributions.
<code>validate_args</code>	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
<code>allow_nan_stats</code>	Logical, default <code>TRUE</code> . When <code>TRUE</code> , statistics (e.g., mean, mode, variance) use the value <code>NaN</code> to indicate the result is undefined. When <code>FALSE</code> , an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Details

Mathematical Details

The Dirichlet-Multinomial is a distribution over K-class counts, i.e., a length-K vector of non-negative integer counts $n = [n_0, \dots, n_{K-1}]$.

The probability mass function (pmf) is,

$$\text{pmf}(n; \text{alpha}, N) = \text{Beta}(\text{alpha} + n) / (\text{prod}_j n_j!) / Z$$

$$Z = \text{Beta}(\text{alpha}) / N!$$

where:

- concentration = alpha = [alpha_0, ..., alpha_{K-1}], alpha_j > 0,
- total_count = N, N a positive integer,
- N! is N factorial, and,
- Beta(x) = prod_j Gamma(x_j) / Gamma(sum_j x_j) is the **multivariate beta function**, and,
- Gamma is the **gamma function**.

Dirichlet-Multinomial is a **compound distribution**, i.e., its samples are generated as follows.

1. Choose class probabilities: probs = [p_0, ..., p_{K-1}] ~ Dir(concentration)
2. Draw integers: counts = [n_0, ..., n_{K-1}] ~ Multinomial(total_count, probs)

The last concentration dimension parametrizes a single Dirichlet-Multinomial distribution. When calling distribution functions (e.g., dist\$prob(counts)), concentration, total_count and counts are broadcast to the same shape. The last dimension of counts corresponds single Dirichlet-Multinomial distributions. Distribution parameters are automatically broadcast in all functions; see examples for details.

Pitfalls The number of classes, K, must not exceed:

- the largest integer representable by self.dtype, i.e., 2**(mantissa_bits+1) (IEEE754),
- the maximum Tensor index, i.e., 2**31-1.

Note: This condition is validated only when validate_args = TRUE.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`

```
tfd_poisson(), tfd_power_spherical(), tfd_probit_bernoulli(), tfd_quantized(), tfd_relaxed_bernoulli(),
tfd_relaxed_one_hot_categorical(), tfd_sample_distribution(), tfd_sinh_arcsinh(),
tfd_skellam(), tfd_spherical_uniform(), tfd_student_t_process(), tfd_student_t(), tfd_transformed_distribution(),
tfd_triangular(), tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian(),
tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()
```

```
tfd_doublesided_maxwell
```

Double-sided Maxwell distribution.

Description

This distribution is useful to compute measure valued derivatives for Gaussian distributions. See Mohamed et al. (2019) for more details.

Usage

```
tfd_doublesided_maxwell(
  loc,
  scale,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "doublesided_maxwell"
)
```

Arguments

<code>loc</code>	Floating point tensor; location of the distribution
<code>scale</code>	Floating point tensor; the scales of the distribution. Must contain only positive values.
<code>validate_args</code>	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
<code>allow_nan_stats</code>	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	string prefixed to Ops created by this class. Default value: 'doublesided_maxwell'.

Details**Mathematical details**

The double-sided Maxwell distribution generalizes the Maxwell distribution to the entire real line.

$$\text{pdf}(x; \mu, \sigma) = 1/(\sigma \sqrt{2\pi}) * ((x-\mu)/\sigma)^2 * \exp(-0.5 ((x-\mu)/\sigma)^2)$$

where $\text{loc} = \mu$ and $\text{scale} = \sigma$. The DoublesidedMaxwell distribution is a member of the **location-scale family**, i.e., it can be constructed as,

$$X \sim \text{DoublesidedMaxwell}(\text{loc}=0, \text{scale}=1)$$

$$Y = \text{loc} + \text{scale} * X$$

The double-sided Maxwell is a symmetric distribution that extends the one-sided maxwell from \mathbb{R}^+ to the entire real line. Their densities are therefore the same up to a factor of 0.5.

It has several methods for generating random variates from it. The version here uses 3 Gaussian variates and a uniform variate to generate the samples The sampling path is:

$$\mu + \sigma * \text{sgn}(U-0.5) * \sqrt{X^2 + Y^2 + Z^2} \quad U \sim \text{Unif}; X, Y, Z \sim N(0,1)$$

In the sampling process above, the random variates generated by $\sqrt{X^2 + Y^2 + Z^2}$ are samples from the one-sided Maxwell (or Maxwell-Boltzmann) distribution.

Value

a distribution instance.

References

- Mohamed, et all, "Monte Carlo Gradient Estimation in Machine Learning.",2019
- B. Heidergott, et al "Sensitivity estimation for Gaussian systems", 2008. European Journal of Operational Research, vol. 187, pp193-207.
- G. Pflug. "Optimization of Stochastic Models: The Interface Between Simulation and Optimization", 2002. Chp. 4.2, pg 247.

See Also

For usage examples see e.g. [tfd_sample\(\)](#), [tfd_log_prob\(\)](#), [tfd_mean\(\)](#).

tfd_empirical

Empirical distribution

Description

The Empirical distribution is parameterized by a (batch) multiset of samples. It describes the empirical measure (observations) of a variable. Note: some methods (`log_prob`, `prob`, `cdf`, `mode`, `entropy`) are not differentiable with regard to samples.

Usage

```
tfd_empirical(
  samples,
  event_ndims = 0,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "Empirical"
)
```

Arguments

<code>samples</code>	Numeric Tensor of shape $[B_1, \dots, B_k, S, E_1, \dots, E_n]$, $k, n \geq 0$. Samples or batches of samples on which the distribution is based. The first k dimensions index into a batch of independent distributions. Length of S dimension determines number of samples in each multiset. The last n dimension represents samples for each distribution. n is specified by argument <code>event_ndims</code> .
<code>event_ndims</code>	<code>int32</code> , default 0 . number of dimensions for each event. When 0 this distribution has scalar samples. When 1 this distribution has vector-like samples.
<code>validate_args</code>	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
<code>allow_nan_stats</code>	Logical, default <code>TRUE</code> . When <code>TRUE</code> , statistics (e.g., mean, mode, variance) use the value <code>NaN</code> to indicate the result is undefined. When <code>FALSE</code> , an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Details**Mathematical Details**

The probability mass function (pmf) and cumulative distribution function (cdf) are

$$\begin{aligned} \text{pmf}(k; s_1, \dots, s_n) &= \sum_i I(k)^{\{k == s_i\}} / n \\ I(k)^{\{k == s_i\}} &= 1, \text{ if } k == s_i, \text{ else } 0. \\ \text{cdf}(k; s_1, \dots, s_n) &= \sum_i I(k)^{\{k \geq s_i\}} / n \\ I(k)^{\{k \geq s_i\}} &= 1, \text{ if } k \geq s_i, \text{ else } 0. \end{aligned}$$
Value

a distribution instance.

See Also

For usage examples see e.g. [tfd_sample\(\)](#), [tfd_log_prob\(\)](#), [tfd_mean\(\)](#).

Other distributions: [tfd_autoregressive\(\)](#), [tfd_batch_reshape\(\)](#), [tfd_bates\(\)](#), [tfd_bernoulli\(\)](#), [tfd_beta_binomial\(\)](#), [tfd_beta\(\)](#), [tfd_binomial\(\)](#), [tfd_categorical\(\)](#), [tfd_cauchy\(\)](#), [tfd_chi2\(\)](#), [tfd_chi\(\)](#), [tfd_cholesky_lkj\(\)](#), [tfd_continuous_bernoulli\(\)](#), [tfd_deterministic\(\)](#), [tfd_dirichlet_multinomial\(\)](#)

[tfd_dirichlet\(\)](#), [tfd_exp_gamma\(\)](#), [tfd_exp_inverse_gamma\(\)](#), [tfd_exponential\(\)](#), [tfd_gamma_gamma\(\)](#),
[tfd_gamma\(\)](#), [tfd_gaussian_process_regression_model\(\)](#), [tfd_gaussian_process\(\)](#), [tfd_generalized_normal\(\)](#),
[tfd_geometric\(\)](#), [tfd_gumbel\(\)](#), [tfd_half_cauchy\(\)](#), [tfd_half_normal\(\)](#), [tfd_hidden_markov_model\(\)](#),
[tfd_horseshoe\(\)](#), [tfd_independent\(\)](#), [tfd_inverse_gamma\(\)](#), [tfd_inverse_gaussian\(\)](#), [tfd_johnson_s_u\(\)](#),
[tfd_joint_distribution_named_auto_batched\(\)](#), [tfd_joint_distribution_named\(\)](#), [tfd_joint_distribution_sequential\(\)](#),
[tfd_kumaraswamy\(\)](#), [tfd_laplace\(\)](#), [tfd_linear_gaussian_state_space\(\)](#), [tfd_lkj\(\)](#), [tfd_log_logistic\(\)](#),
[tfd_log_normal\(\)](#), [tfd_logistic\(\)](#), [tfd_mixture_same_family\(\)](#), [tfd_mixture\(\)](#), [tfd_multinomial\(\)](#),
[tfd_multivariate_normal_diag_plus_low_rank\(\)](#), [tfd_multivariate_normal_full_covariance\(\)](#),
[tfd_multivariate_normal_linear_operator\(\)](#), [tfd_multivariate_normal_tri_l\(\)](#),
[tfd_multivariate_student_t_linear_operator\(\)](#), [tfd_negative_binomial\(\)](#), [tfd_normal\(\)](#),
[tfd_one_hot_categorical\(\)](#), [tfd_pareto\(\)](#), [tfd_pixel_cnn\(\)](#), [tfd_poisson_log_normal_quadrature\(\)](#),
[tfd_poisson\(\)](#), [tfd_power_spherical\(\)](#), [tfd_probit_bernoulli\(\)](#), [tfd_quantized\(\)](#), [tfd_relaxed_bernoulli\(\)](#),
[tfd_relaxed_one_hot_categorical\(\)](#), [tfd_sample_distribution\(\)](#), [tfd_sinh_arcsinh\(\)](#),
[tfd_skellam\(\)](#), [tfd_spherical_uniform\(\)](#), [tfd_student_t_process\(\)](#), [tfd_student_t\(\)](#), [tfd_transformed_distribution\(\)](#),
[tfd_triangular\(\)](#), [tfd_truncated_cauchy\(\)](#), [tfd_truncated_normal\(\)](#), [tfd_uniform\(\)](#), [tfd_variational_gaussian\(\)](#),
[tfd_vector_diffeomixture\(\)](#), [tfd_vector_exponential_diag\(\)](#), [tfd_vector_exponential_linear_operator\(\)](#),
[tfd_vector_laplace_diag\(\)](#), [tfd_vector_laplace_linear_operator\(\)](#), [tfd_vector_sinh_arcsinh_diag\(\)](#),
[tfd_von_mises_fisher\(\)](#), [tfd_von_mises\(\)](#), [tfd_weibull\(\)](#), [tfd_wishart_linear_operator\(\)](#),
[tfd_wishart_tri_l\(\)](#), [tfd_wishart\(\)](#), [tfd_zipf\(\)](#)

tfd_entropy

Shannon entropy in nats.

Description

Shannon entropy in nats.

Usage

```
tfd_entropy(distribution, ...)
```

Arguments

distribution The distribution being used.
... Additional parameters passed to Python.

Value

a Tensor of shape `sample_shape(x) + self$batch_shape` with values of type `self$dtype`.

See Also

Other `distribution_methods`: [tfd_cdf\(\)](#), [tfd_covariance\(\)](#), [tfd_cross_entropy\(\)](#), [tfd_kl_divergence\(\)](#),
[tfd_log_cdf\(\)](#), [tfd_log_prob\(\)](#), [tfd_log_survival_function\(\)](#), [tfd_mean\(\)](#), [tfd_mode\(\)](#),
[tfd_prob\(\)](#), [tfd_quantile\(\)](#), [tfd_sample\(\)](#), [tfd_stddev\(\)](#), [tfd_survival_function\(\)](#), [tfd_variance\(\)](#)

Examples

```
d <- tfd_normal(loc = c(1, 2), scale = c(1, 0.5))
d %>% tfd_entropy()
```

tfd_exponential	<i>Exponential distribution</i>
-----------------	---------------------------------

Description

The Exponential distribution is parameterized by an event rate parameter.

Usage

```
tfd_exponential(
  rate,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "Exponential"
)
```

Arguments

rate	Floating point tensor, equivalent to 1 / mean. Must contain only positive values.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details**Mathematical Details**

The probability density function (pdf) is,

$$\text{pdf}(x; \text{lambda}, x > 0) = \exp(-\text{lambda } x) / Z$$

$$Z = 1 / \text{lambda}$$

where rate = lambda and Z is the normalizing constant.

The Exponential distribution is a special case of the Gamma distribution, i.e.,

$$\text{Exponential}(\text{rate}) = \text{Gamma}(\text{concentration}=1., \text{rate})$$

The Exponential distribution uses a rate parameter, or "inverse scale", which can be intuited as,

```
X ~ Exponential(rate=1)
Y = X / rate
```

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_exp_gamma

ExpGamma distribution.

Description

The ExpGamma distribution is defined over the real line using parameters concentration (aka "alpha") and rate (aka "beta"). This distribution is a transformation of the Gamma distribution such that $X \sim \text{ExpGamma}(\cdot) \Rightarrow \exp(X) \sim \text{Gamma}(\cdot)$.

Usage

```
tfd_exp_gamma(
  concentration,
  rate = NULL,
  log_rate = NULL,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "ExpGamma"
)
```

Arguments

concentration	Floating point tensor, the concentration params of the distribution(s). Must contain only positive values.
rate	Floating point tensor, the inverse scale params of the distribution(s). Must contain only positive values. Mutually exclusive with <code>log_rate</code> .
log_rate	Floating point tensor, natural logarithm of the inverse scale params of the distribution(s). Mutually exclusive with <code>rate</code> .
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details**Mathematical Details**

The probability density function (pdf) can be derived from the change of variables rule (since the distribution is logically equivalent to `tfb_log()(tfd_gamma(. .))`):

$$\text{pdf}(x; \alpha, \beta > 0) = \exp(x)^{(\alpha - 1)} \exp(-\exp(x) \beta) / Z + x$$

$$Z = \text{Gamma}(\alpha) \beta^{(-\alpha)}$$

where:

- `concentration` = α , $\alpha > 0$,
- `rate` = β , $\beta > 0$,
- Z is the normalizing constant of the corresponding Gamma distribution, and
- Gamma is the [gamma function](#).

The cumulative density function (cdf) is,

$$\text{cdf}(x; \alpha, \beta, x) = \text{GammaInc}(\alpha, \beta \exp(x)) / \text{Gamma}(\alpha)$$

where `GammaInc` is the **lower incomplete Gamma function**.

Distribution parameters are automatically broadcast in all functions. Samples of this distribution are reparameterized (pathwise differentiable). The derivatives are computed using the approach described in Figurnov et al., 2018.

Value

a distribution instance.

References

- Michael Figurnov, Shakir Mohamed, Andriy Mnih. **Implicit Reparameterization Gradients**. *arXiv preprint arXiv:1805.08498*, 2018.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

`tfd_exp_inverse_gamma` *ExpInverseGamma* distribution.

Description

The `ExpInverseGamma` distribution is defined over the real numbers such that $X \sim \text{ExpInverseGamma}(\cdot)$ $\Rightarrow \exp(X) \sim \text{InverseGamma}(\cdot)$. The distribution is logically equivalent to `tfd_log()(tfd_inverse_gamma(.))`, but can be sampled with much better precision.

Usage

```
tfd_exp_inverse_gamma(
  concentration,
  scale = NULL,
  log_scale = NULL,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "ExpGamma"
)
```

Arguments

concentration	Floating point tensor, the concentration params of the distribution(s). Must contain only positive values.
scale	Floating point tensor, the scale params of the distribution(s). Must contain only positive values. Mutually exclusive with log_scale.
log_scale	Floating point tensor, the natural logarithm of the scale params of the distribution(s). Mutually exclusive with scale.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details**Mathematical Details**

The probability density function (pdf) is very similar to ExpGamma,

$$\text{pdf}(x; \alpha, \beta > 0) = \frac{\exp(-x)^{\alpha-1} \exp(-\exp(-x) \beta)}{Z - x}$$

$$Z = \Gamma(\alpha) \beta^{-(\alpha)}$$

where:

- concentration = alpha,
- scale = beta,
- Z is the normalizing constant, and,
- Gamma is the **gamma function**.

The cumulative density function (cdf) is,

$$\text{cdf}(x; \alpha, \beta, x) = 1 - \Gamma\text{Inc}(\alpha, \beta \exp(-x)) / \Gamma(\alpha)$$

where `GammaInc` is the **upper incomplete Gamma function**.

Distribution parameters are automatically broadcast in all functions. Samples of this distribution are reparameterized (pathwise differentiable). The derivatives are computed using the approach described in Figurnov et al, 2018.

Value

a distribution instance.

References

- Michael Figurnov, Shakir Mohamed, Andriy Mnih. **Implicit Reparameterization Gradients**. *arXiv preprint arXiv:1805.08498*, 2018.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffusion_mixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

`tfd_exp_relaxed_one_hot_categorical`

ExpRelaxedOneHotCategorical distribution with temperature and logits.

Description

ExpRelaxedOneHotCategorical distribution with temperature and logits.

Usage

```
tfd_exp_relaxed_one_hot_categorical(  
    temperature,  
    logits = NULL,  
    probs = NULL,  
    validate_args = FALSE,  
    allow_nan_stats = TRUE,  
    name = "ExpRelaxedOneHotCategorical"  
)
```

Arguments

temperature	An 0-D Tensor, representing the temperature of a set of ExpRelaxedCategorical distributions. The temperature should be positive.
logits	An N-D Tensor, $N \geq 1$, representing the log probabilities of a set of ExpRelaxedCategorical distributions. The first $N - 1$ dimensions index into a batch of independent distributions and the last dimension represents a vector of logits for each class. Only one of logits or probs should be passed in.
probs	An N-D Tensor, $N \geq 1$, representing the probabilities of a set of ExpRelaxedCategorical distributions. The first $N - 1$ dimensions index into a batch of independent distributions and the last dimension represents a vector of probabilities for each class. Only one of logits or probs should be passed in.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Value

a distribution instance.

See Also

For usage examples see e.g. [tfd_sample\(\)](#), [tfd_log_prob\(\)](#), [tfd_mean\(\)](#).

tfd_finite_discrete *The finite discrete distribution.*

Description

The FiniteDiscrete distribution is parameterized by either probabilities or log-probabilities of a set of K possible outcomes, which is defined by a strictly ascending list of K values.

Usage

```
tfd_finite_discrete(
    outcomes,
    logits = NULL,
    probs = NULL,
    rtol = NULL,
    atol = NULL,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = "FiniteDiscrete"
)
```

Arguments

outcomes	A 1-D floating or integer Tensor, representing a list of possible outcomes in strictly ascending order.
logits	A floating N-D Tensor, $N \geq 1$, representing the log probabilities of a set of FiniteDiscrete distributions. The first $N - 1$ dimensions index into a batch of independent distributions and the last dimension represents a vector of logits for each discrete value. Only one of logits or probs should be passed in.
probs	A floating N-D Tensor, $N \geq 1$, representing the probabilities of a set of Finite-Discrete distributions. The first $N - 1$ dimensions index into a batch of independent distributions and the last dimension represents a vector of probabilities for each discrete value. Only one of logits or probs should be passed in.
rtol	Tensor with same dtype as outcomes. The relative tolerance for floating number comparison. Only effective when outcomes is a floating Tensor. Default is $10 * \text{eps}$.
atol	Tensor with same dtype as outcomes. The absolute tolerance for floating number comparison. Only effective when outcomes is a floating Tensor. Default is $10 * \text{eps}$.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.

name string prefixed to Ops created by this class.

Details

Note: `log_prob`, `prob`, `cdf`, `mode`, and `entropy` are differentiable with respect to logits or probs but not with respect to outcomes.

Mathematical Details

The probability mass function (pmf) is,

$$\text{pmf}(x; \text{pi}, \text{qi}) = \prod_j \text{pi}_j^{[x == \text{qi}_j]}$$

Value

a distribution instance.

See Also

For usage examples see e.g. [tfd_sample\(\)](#), [tfd_log_prob\(\)](#), [tfd_mean\(\)](#).

tfd_gamma

Gamma distribution

Description

The Gamma distribution is defined over positive real numbers using parameters concentration (aka "alpha") and rate (aka "beta").

Usage

```
tfd_gamma(
  concentration,
  rate,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "Gamma"
)
```

Arguments

concentration	Floating point tensor, the concentration params of the distribution(s). Must contain only positive values.
rate	Floating point tensor, the inverse scale params of the distribution(s). Must contain only positive values.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.

allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

Mathematical Details

The probability density function (pdf) is,

$$\text{pdf}(x; \alpha, \beta, x > 0) = x^{(\alpha - 1)} \exp(-x \beta) / Z$$

$$Z = \text{Gamma}(\alpha) \beta^{(-\alpha)}$$

where

- concentration = alpha, alpha > 0,
- rate = beta, beta > 0,
- Z is the normalizing constant, and,
- Gamma is the [gamma function](#).

The cumulative density function (cdf) is,

$$\text{cdf}(x; \alpha, \beta, x > 0) = \text{GammaInc}(\alpha, \beta x) / \text{Gamma}(\alpha)$$

where GammaInc is the [lower incomplete Gamma function](#). The parameters can be intuited via their relationship to mean and stddev,

$$\text{concentration} = \alpha = (\text{mean} / \text{stddev})^{**2}$$

$$\text{rate} = \beta = \text{mean} / \text{stddev}^{**2} = \text{concentration} / \text{mean}$$

Distribution parameters are automatically broadcast in all functions; see examples for details.

Warning: The samples of this distribution are always non-negative. However, the samples that are smaller than `np.finfo(dtype).tiny` are rounded to this value, so it appears more often than it should. This should only be noticeable when the concentration is very small, or the rate is very large. See note in `tf.random_gamma` docstring. Samples of this distribution are reparameterized (pathwise differentiable). The derivatives are computed using the approach described in the paper [Michael Figurnov, Shakir Mohamed, Andriy Mnih. Implicit Reparameterization Gradients, 2018](#)

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_gamma_gamma

Gamma-Gamma distribution

Description

Gamma-Gamma is a **compound distribution** defined over positive real numbers using parameters `concentration`, `mixing_concentration` and `mixing_rate`.

Usage

```
tfd_gamma_gamma(
  concentration,
  mixing_concentration,
  mixing_rate,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "GammaGamma"
)
```

Arguments

concentration	Floating point tensor, the concentration params of the distribution(s). Must contain only positive values.
mixing_concentration	Floating point tensor, the concentration params of the mixing Gamma distribution(s). Must contain only positive values.
mixing_rate	Floating point tensor, the rate params of the mixing Gamma distribution(s). Must contain only positive values.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

This distribution is also referred to as the beta of the second kind (B2), and can be useful for transaction value modeling, as in Fader and Hardi, 2013.

Mathematical Details

It is derived from the following Gamma-Gamma hierarchical model by integrating out the random variable beta.

$$\begin{aligned} \text{beta} &\sim \text{Gamma}(\alpha_0, \beta_0) \\ X \mid \text{beta} &\sim \text{Gamma}(\alpha, \text{beta}) \end{aligned}$$

where

- concentration = alpha
- mixing_concentration = alpha₀
- mixing_rate = beta₀

The probability density function (pdf) is

$$\begin{aligned} &x^{(\alpha - 1)} \\ \text{pdf}(x; \alpha, \alpha_0, \beta_0) &= Z * (x + \beta_0)^{(\alpha + \alpha_0)} \end{aligned}$$

where the normalizing constant $Z = \text{Beta}(\alpha, \alpha_0) * \beta_0^{(-\alpha_0)}$. Samples of this distribution are reparameterized as samples of the Gamma distribution are reparameterized using the technique described in (Figurnov et al., 2018).

@section References:

- Peter S. Fader, Bruce G. S. Hardi. *The Gamma-Gamma Model of Monetary Value. Technical Report, 2013.*
- Michael Figurnov, Shakir Mohamed, Andriy Mnih. *Implicit Reparameterization Gradients. arXiv preprint arXiv:1805.08498, 2018*

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

`tfd_gaussian_process` *Marginal distribution of a Gaussian process at finitely many points.*

Description

A Gaussian process (GP) is an indexed collection of random variables, any finite collection of which are jointly Gaussian. While this definition applies to finite index sets, it is typically implicit that the index set is infinite; in applications, it is often some finite dimensional real or complex vector space. In such cases, the GP may be thought of as a distribution over (real- or complex-valued) functions defined over the index set.

Usage

```
tfd_gaussian_process(  
    kernel,  
    index_points,  
    mean_fn = NULL,
```

```

    observation_noise_variance = 0,
    jitter = 1e-06,
    validate_args = FALSE,
    allow_nan_stats = FALSE,
    name = "GaussianProcess"
)

```

Arguments

kernel	PositiveSemidefiniteKernel-like instance representing the GP's covariance function.
index_points	float Tensor representing finite (batch of) vector(s) of points in the index set over which the GP is defined. Shape has the form [b1, ..., bB, e1, f1, ..., fF] where F is the number of feature dimensions and must equal kernel\$feature_ndims and e1 is the number (size) of index points in each batch (we denote it e1 to distinguish it from the number of inducing index points, denoted e2 below). Ultimately the GaussianProcess distribution corresponds to an e1-dimensional multivariate normal. The batch shape must be broadcastable with kernel\$batch_shape, the batch shape of inducing_index_points, and any batch dims yielded by mean_fn.
mean_fn	function that acts on index points to produce a (batch of) vector(s) of mean values at those index points. Takes a Tensor of shape [b1, ..., bB, f1, ..., fF] and returns a Tensor whose shape is (broadcastable with) [b1, ..., bB]. Default value: NULL implies constant zero function.
observation_noise_variance	float Tensor representing the variance of the noise in the Normal likelihood distribution of the model. May be batched, in which case the batch shape must be broadcastable with the shapes of all other batched parameters (kernel\$batch_shape, index_points, etc.). Default value: 0.
jitter	float scalar Tensor added to the diagonal of the covariance matrix to ensure positive definiteness of the covariance matrix. Default value: 1e-6.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

Just as Gaussian distributions are fully specified by their first and second moments, a Gaussian process can be completely specified by a mean and covariance function. Let S denote the index set and K the space in which each indexed random variable takes its values (again, often \mathbb{R} or \mathbb{C}). The mean function is then a map $m: S \rightarrow K$, and the covariance function, or kernel, is a positive-definite

function $k: (S \times S) \rightarrow K$. The properties of functions drawn from a GP are entirely dictated (up to translation) by the form of the kernel function.

This Distribution represents the marginal joint distribution over function values at a given finite collection of points $[x[1], \dots, x[N]]$ from the index set S . By definition, this marginal distribution is just a multivariate normal distribution, whose mean is given by the vector $[m(x[1]), \dots, m(x[N])]$ and whose covariance matrix is constructed from pairwise applications of the kernel function to the given inputs:

$$\begin{array}{cccc|c} | & k(x[1], x[1]) & k(x[1], x[2]) & \dots & k(x[1], x[N]) & | \\ | & k(x[2], x[1]) & k(x[2], x[2]) & \dots & k(x[2], x[N]) & | \\ | & \dots & \dots & \dots & \dots & | \\ | & k(x[N], x[1]) & k(x[N], x[2]) & \dots & k(x[N], x[N]) & | \end{array}$$

For this to be a valid covariance matrix, it must be symmetric and positive definite; hence the requirement that k be a positive definite function (which, by definition, says that the above procedure will yield PD matrices).

We also support the inclusion of zero-mean Gaussian noise in the model, via the `observation_noise_variance` parameter. This augments the generative model to

$$\begin{aligned} f &\sim \text{GP}(m, k) \\ (y[i] \mid f, x[i]) &\sim \text{Normal}(f(x[i]), s) \end{aligned}$$

where

- m is the mean function
- k is the covariance kernel function
- f is the function drawn from the GP
- $x[i]$ are the index points at which the function is observed
- $y[i]$ are the observed values at the index points
- s is the scale of the observation noise.

Note that this class represents an *unconditional* Gaussian process; it does not implement posterior inference conditional on observed function evaluations. This class is useful, for example, if one wishes to combine a GP prior with a non-conjugate likelihood using MCMC to sample from the posterior.

Mathematical Details

The probability density function (pdf) is a multivariate normal whose parameters are derived from the GP's properties:

$$\begin{aligned} \text{pdf}(x; \text{index_points}, \text{mean_fn}, \text{kernel}) &= \exp(-0.5 * y) / Z \\ K &= (\text{kernel.matrix}(\text{index_points}, \text{index_points}) + \\ &\quad (\text{observation_noise_variance} + \text{jitter}) * \text{eye}(N)) \\ y &= (x - \text{mean_fn}(\text{index_points}))^T @ K @ (x - \text{mean_fn}(\text{index_points})) \\ Z &= (2 * \text{pi})^{*.5 * N} |\text{det}(K)|^{*.5} \end{aligned}$$

where:

- `index_points` are points in the index set over which the GP is defined,
- `mean_fn` is a callable mapping the index set to the GP's mean values,
- `kernel` is `PositiveSemidefiniteKernel`-like and represents the covariance function of the GP,
- `observation_noise_variance` represents (optional) observation noise.
- `jitter` is added to the diagonal to ensure positive definiteness up to machine precision (otherwise Cholesky-decomposition is prone to failure),
- `eye(N)` is an N-by-N identity matrix.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

`tfd_gaussian_process_regression_model`

Posterior predictive distribution in a conjugate GP regression model.

Description

Posterior predictive distribution in a conjugate GP regression model.

Usage

```
tfd_gaussian_process_regression_model(
    kernel,
    index_points = NULL,
    observation_index_points = NULL,
    observations = NULL,
    observation_noise_variance = 0,
    predictive_noise_variance = NULL,
    mean_fn = NULL,
    jitter = 1e-06,
    validate_args = FALSE,
    allow_nan_stats = FALSE,
    name = "GaussianProcessRegressionModel"
)
```

Arguments

kernel	PositiveSemidefiniteKernel-like instance representing the GP's covariance function.
index_points	float Tensor representing finite (batch of) vector(s) of points in the index set over which the GP is defined. Shape has the form [b1, ..., bB, e1, f1, ..., fF] where F is the number of feature dimensions and must equal kernel\$feature_ndims and e1 is the number (size) of index points in each batch (we denote it e1 to distinguish it from the number of inducing index points, denoted e2 below). Ultimately the GaussianProcess distribution corresponds to an e1-dimensional multivariate normal. The batch shape must be broadcastable with kernel\$batch_shape, the batch shape of inducing_index_points, and any batch dims yielded by mean_fn.
observation_index_points	Tensor representing finite collection, or batch of collections, of points in the index set for which some data has been observed. Shape has the form [b1, ..., bB, e, f1, ..., fF] where F is the number of feature dimensions and must equal kernel\$feature_ndims, and e is the number (size) of index points in each batch. [b1, ..., bB, e] must be broadcastable with the shape of observations, and [b1, ..., bB] must be broadcastable with the shapes of all other batched parameters (kernel.batch_shape, index_points, etc). The default value is None, which corresponds to the empty set of observations, and simply results in the prior predictive model (a GP with noise of variance predictive_noise_variance).
observations	Tensor representing collection, or batch of collections, of observations corresponding to observation_index_points. Shape has the form [b1, ..., bB, e], which must be broadcastable with the batch and example shapes of observation_index_points. The batch shape [b1, ..., bB] must be broadcastable with the shapes of all other batched parameters (kernel.batch_shape, index_points, etc.). The default value is None, which corresponds to the empty set of observations, and simply results in the prior predictive model (a GP with noise of variance predictive_noise_variance).
observation_noise_variance	float Tensor representing the variance of the noise in the Normal likelihood

	distribution of the model. May be batched, in which case the batch shape must be broadcastable with the shapes of all other batched parameters (<code>kernel\$batch_shape</code> , <code>index_points</code> , etc.). Default value: <code>0</code> .
<code>predictive_noise_variance</code>	Tensor representing the variance in the posterior predictive model. If <code>None</code> , we simply re-use <code>observation_noise_variance</code> for the posterior predictive noise. If set explicitly, however, we use this value. This allows us, for example, to omit predictive noise variance (by setting this to zero) to obtain noiseless posterior predictions of function values, conditioned on noisy observations.
<code>mean_fn</code>	callable that acts on <code>index_points</code> to produce a collection, or batch of collections, of mean values at <code>index_points</code> . Takes a Tensor of shape <code>[b1, ..., bB, f1, ..., fF]</code> and returns a Tensor whose shape is broadcastable with <code>[b1, ..., bB]</code> . Default value: <code>None</code> implies the constant zero function.
<code>jitter</code>	float scalar Tensor added to the diagonal of the covariance matrix to ensure positive definiteness of the covariance matrix. Default value: <code>1e-6</code> .
<code>validate_args</code>	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
<code>allow_nan_stats</code>	Logical, default <code>TRUE</code> . When <code>TRUE</code> , statistics (e.g., mean, mode, variance) use the value <code>NaN</code> to indicate the result is undefined. When <code>FALSE</code> , an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`,

tfd_skellam(), tfd_spherical_uniform(), tfd_student_t_process(), tfd_student_t(), tfd_transformed_distribution(), tfd_triangular(), tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian(), tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(), tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(), tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(), tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()

tfd_generalized_normal

The Generalized Normal distribution.

Description

The Generalized Normal (or Generalized Gaussian) generalizes the Normal distribution with an additional shape parameter. It is parameterized by location `loc`, scale `scale` and shape power.

Usage

```
tfd_generalized_normal(
  loc,
  scale,
  power,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "GeneralizedNormal"
)
```

Arguments

<code>loc</code>	Floating point tensor; the means of the distribution(s).
<code>scale</code>	Floating point tensor; the scale of the distribution(s). Must contain only positive values.
<code>power</code>	Floating point tensor; the shape parameter of the distribution(s). Must contain only positive values. <code>loc</code> , <code>scale</code> and <code>power</code> must have compatible shapes for broadcasting.
<code>validate_args</code>	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
<code>allow_nan_stats</code>	Logical, default <code>TRUE</code> . When <code>TRUE</code> , statistics (e.g., mean, mode, variance) use the value <code>NaN</code> to indicate the result is undefined. When <code>FALSE</code> , an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Details

Mathematical details The probability density function (pdf) is,

$$\text{pdf}(x; \text{loc}, \text{scale}, \text{power}) = 1 / (2 * \text{scale} * \text{Gamma}(1 + 1 / \text{power})) * \exp(-(|x - \text{loc}| / \text{scale}) ^ \text{power})$$

where loc is the mean, scale is the scale, and, power is the shape parameter. If the power is above two, the distribution becomes platykurtic. A power equal to two results in a Normal distribution. A power smaller than two produces a leptokurtic (heavy-tailed) distribution. Mean and scale behave the same way as in the equivalent Normal distribution.

See https://en.wikipedia.org/w/index.php?title=Generalized_normal_distribution&oldid=954254464 for the definitions used here, including CDF, variance and entropy. See https://sccn.ucsd.edu/wiki/Generalized_Gaussian_Probability_Distribution for the sampling method used here.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_generalized_pareto

The Generalized Pareto distribution.

Description

The Generalized Pareto distributions are a family of continuous distributions on the reals. Special cases include Exponential (when $\text{loc} = 0$, $\text{concentration} = 0$), Pareto (when $\text{concentration} > 0$, $\text{loc} = \text{scale} / \text{concentration}$), and Uniform (when $\text{concentration} = -1$).

Usage

```
tfd_generalized_pareto(
  loc,
  scale,
  concentration,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = NULL
)
```

Arguments

loc	The location / shift of the distribution. GeneralizedPareto is a location-scale distribution. This parameter lower bounds the distribution's support. Must broadcast with scale, concentration. Floating point Tensor.
scale	The scale of the distribution. GeneralizedPareto is a location-scale distribution, so doubling the scale doubles a sample and halves the density. Strictly positive floating point Tensor. Must broadcast with loc, concentration.
concentration	The shape parameter of the distribution. The larger the magnitude, the more the distribution concentrates near loc (for $\text{concentration} \geq 0$) or near $\text{loc} - (\text{scale}/\text{concentration})$ (for $\text{concentration} < 0$). Floating point Tensor.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

This distribution is often used to model the tails of other distributions. As a member of the location-scale family, $X \sim \text{GeneralizedPareto}(\text{loc}=\text{loc}, \text{scale}=\text{scale}, \text{concentration}=\text{conc})$ maps to $Y \sim \text{GeneralizedPareto}(\text{loc}=0, \text{scale}=1, \text{concentration}=\text{conc})$ via $Y = (X - \text{loc}) / \text{scale}$.

For positive concentrations, the distribution is equivalent to a hierarchical Exponential-Gamma model with $X|rate \sim \text{Exponential}(rate)$ and $rate \sim \text{Gamma}(concentration=1 / concentration, scale=scale / concentration)$. In the following, `samps1` and `samps2` are identically distributed:

```
genp <- tfd_generalized_pareto(loc = 0, scale = scale, concentration = conc)
samps1 <- genp %>% tfd_sample(1000)
jd <- tfd_joint_distribution_named(
  list(
    rate = tfd_gamma(1 / genp$concentration, genp$scale / genp$concentration),
    x = function(rate) tfd_exponential(rate))
samps2 <- jd %>% tfd_sample(1000) %>% .$x
```

The support of the distribution is always lower bounded by `loc`. When `concentration < 0`, the support is also upper bounded by `loc + scale / abs(concentration)`.

Mathematical Details

The probability density function (pdf) is,

$$\text{pdf}(x; \mu, \sigma, \text{shp}, x > \mu) = (1 + \text{shp} * (x - \mu) / \sigma)^{*(-1 / \text{shp} - 1)} / \sigma$$

where:

- `concentration = shp`, any real value,
- `scale = sigma`, `sigma > 0`,
- `loc = mu`.

The cumulative density function (cdf) is,

$$\text{cdf}(x; \mu, \sigma, \text{shp}, x > \mu) = 1 - (1 + \text{shp} * (x - \mu) / \sigma)^{*(-1 / \text{shp})}$$

Distribution parameters are automatically broadcast in all functions; see examples for details. Samples of this distribution are reparameterized (pathwise differentiable).

Value

a distribution instance.

See Also

For usage examples see e.g. [tfd_sample\(\)](#), [tfd_log_prob\(\)](#), [tfd_mean\(\)](#).

tfd_geometric	<i>Geometric distribution</i>
---------------	-------------------------------

Description

The Geometric distribution is parameterized by p , the probability of a positive event. It represents the probability that in $k + 1$ Bernoulli trials, the first k trials failed, before seeing a success. The pmf of this distribution is:

Usage

```
tfd_geometric(
  logits = NULL,
  probs = NULL,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "Geometric"
)
```

Arguments

logits	Floating-point Tensor with shape $[B_1, \dots, B_b]$ where $b \geq 0$ indicates the number of batch dimensions. Each entry represents logits for the probability of success for independent Geometric distributions and must be in the range $(-\infty, \infty]$. Only one of <code>logits</code> or <code>probs</code> should be specified.
probs	Positive floating-point Tensor with shape $[B_1, \dots, B_b]$ where $b \geq 0$ indicates the number of batch dimensions. Each entry represents the probability of success for independent Geometric distributions and must be in the range $(0, 1]$. Only one of <code>logits</code> or <code>probs</code> should be specified.
validate_args	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
allow_nan_stats	Logical, default <code>TRUE</code> . When <code>TRUE</code> , statistics (e.g., mean, mode, variance) use the value <code>NaN</code> to indicate the result is undefined. When <code>FALSE</code> , an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

Mathematical Details

$$\text{pmf}(k; p) = (1 - p)^{k-1} * p$$

where:

- p is the success probability, $0 < p \leq 1$, and,
- k is a non-negative integer.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_gumbel

Scalar Gumbel distribution with location loc and scale parameters

Description

Mathematical details

Usage

```
tfd_gumbel(
  loc,
  scale,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "Gumbel"
)
```

Arguments

loc	Floating point tensor, the means of the distribution(s).
scale	Floating point tensor, the scales of the distribution(s). ‘scale‘ must contain only positive values.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic’s batch members are undefined.
name	name prefixed to Ops created by this class.

Details

The probability density function (pdf) of this distribution is,

$$\text{pdf}(x; \mu, \sigma) = \exp(-(x - \mu) / \sigma - \exp(-(x - \mu) / \sigma)) / \sigma$$

where loc = mu and scale = sigma.

The cumulative density function of this distribution is, $\text{cdf}(x; \mu, \sigma) = \exp(-\exp(-(x - \mu) / \sigma))$

The Gumbel distribution is a member of the **location-scale family**, i.e., it can be constructed as,

$$\begin{aligned} X &\sim \text{Gumbel}(\text{loc}=\theta, \text{scale}=1) \\ Y &= \text{loc} + \text{scale} * X \end{aligned}$$

Value

a distribution instance.

See Also

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`,

```
tfd_relaxed_one_hot_categorical(), tfd_sample_distribution(), tfd_sinh_arcsinh(),
tfd_skellam(), tfd_spherical_uniform(), tfd_student_t_process(), tfd_student_t(), tfd_transformed_distribution(),
tfd_triangular(), tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian(),
tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()
```

tfd_half_cauchy	<i>Half-Cauchy distribution</i>
-----------------	---------------------------------

Description

The half-Cauchy distribution is parameterized by a `loc` and a `scale` parameter. It represents the right half of the two symmetric halves in a [Cauchy distribution](#).

Usage

```
tfd_half_cauchy(
  loc,
  scale,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "HalfCauchy"
)
```

Arguments

<code>loc</code>	Floating-point Tensor; the location(s) of the distribution(s).
<code>scale</code>	Floating-point Tensor; the scale(s) of the distribution(s). Must contain only positive values.
<code>validate_args</code>	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
<code>allow_nan_stats</code>	Logical, default <code>TRUE</code> . When <code>TRUE</code> , statistics (e.g., mean, mode, variance) use the value <code>NaN</code> to indicate the result is undefined. When <code>FALSE</code> , an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Details

Mathematical Details

The probability density function (pdf) for the half-Cauchy distribution is given by

$$\text{pdf}(x; \text{loc}, \text{scale}) = 2 / (\pi \text{scale} (1 + z^2))$$

$$z = (x - \text{loc}) / \text{scale}$$

where `loc` is a scalar in R and `scale` is a positive scalar in R. The support of the distribution is given by the interval `[loc, infinity)`.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_half_normal	<i>Half-Normal distribution with scale scale</i>
-----------------	--

Description

Mathematical details

Usage

```
tfd_half_normal(
  scale,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "HalfNormal"
)
```

Arguments

scale	Floating point tensor; the scales of the distribution(s). Must contain only positive values.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

The half normal is a transformation of a centered normal distribution. If some random variable X has normal distribution,

$$X \sim \text{Normal}(0, \text{scale})$$

$$Y = |X|$$

Then Y will have half normal distribution. The probability density function (pdf) is:

$$\text{pdf}(x; \text{scale}, x > 0) = \sqrt{2} / (\text{scale} * \sqrt{\pi}) * \exp(- 1/2 * (x / \text{scale}) ** 2)$$

Where $\text{scale} = \text{sigma}$ is the standard deviation of the underlying normal distribution.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`,

```
tfd_relaxed_one_hot_categorical(), tfd_sample_distribution(), tfd_sinh_arcsinh(),
tfd_skellam(), tfd_spherical_uniform(), tfd_student_t_process(), tfd_student_t(), tfd_transformed_distribution(),
tfd_triangular(), tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian(),
tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()
```

```
tfd_hidden_markov_model
```

Hidden Markov model distribution

Description

The `HiddenMarkovModel` distribution implements a (batch of) hidden Markov models where the initial states, transition probabilities and observed states are all given by user-provided distributions.

Usage

```
tfd_hidden_markov_model(
  initial_distribution,
  transition_distribution,
  observation_distribution,
  num_steps,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "HiddenMarkovModel"
)
```

Arguments

<code>initial_distribution</code>	A Categorical-like instance. Determines probability of first hidden state in Markov chain. The number of categories must match the number of categories of <code>transition_distribution</code> as well as both the rightmost batch dimension of <code>transition_distribution</code> and the rightmost batch dimension of <code>observation_distribution</code> .
<code>transition_distribution</code>	A Categorical-like instance. The rightmost batch dimension indexes the probability distribution of each hidden state conditioned on the previous hidden state.
<code>observation_distribution</code>	A <code>tfp\$Distributions\$Distribution</code> -like instance. The rightmost batch dimension indexes the distribution of each observation conditioned on the corresponding hidden state.
<code>num_steps</code>	The number of steps taken in Markov chain. An integer.
<code>validate_args</code>	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .

<code>allow_nan_stats</code>	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Details

This model assumes that the transition matrices are fixed over time. In this model, there is a sequence of integer-valued hidden states: $z[0]$, $z[1]$, ..., $z[\text{num_steps} - 1]$ and a sequence of observed states: $x[0]$, ..., $x[\text{num_steps} - 1]$.

The distribution of $z[0]$ is given by `initial_distribution`. The conditional probability of $z[i + 1]$ given $z[i]$ is described by the batch of distributions in `transition_distribution`. For a batch of hidden Markov models, the coordinates before the rightmost one of the `transition_distribution` batch correspond to indices into the hidden Markov model batch. The rightmost coordinate of the batch is used to select which distribution $z[i + 1]$ is drawn from. The distributions corresponding to the probability of $z[i + 1]$ conditional on $z[i] == k$ is given by the elements of the batch whose rightmost coordinate is k .

Similarly, the conditional distribution of $z[i]$ given $x[i]$ is given by the batch of `observation_distribution`. When the rightmost coordinate of `observation_distribution` is k it gives the conditional probabilities of $x[i]$ given $z[i] == k$. The probability distribution associated with the `HiddenMarkovModel` distribution is the marginal distribution of $x[0], \dots, x[\text{num_steps} - 1]$.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`

tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
 tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
 tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
 tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()

tfd_horseshoe	<i>Horseshoe distribution</i>
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Description

The so-called 'horseshoe' distribution is a Cauchy-Normal scale mixture, proposed as a sparsity-inducing prior for Bayesian regression. It is symmetric around zero, has heavy (Cauchy-like) tails, so that large coefficients face relatively little shrinkage, but an infinitely tall spike at 0, which pushes small coefficients towards zero. It is parameterized by a positive scalar scale parameter: higher values yield a weaker sparsity-inducing effect.

Usage

```
tfd_horseshoe(  
  scale,  
  validate_args = FALSE,  
  allow_nan_stats = TRUE,  
  name = "Horseshoe"  
)
```

Arguments

scale	Floating point tensor; the scales of the distribution(s). Must contain only positive values.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

Mathematical details

The Horseshoe distribution is centered at zero, with scale parameter λ . It is defined by:

$$\text{horseshoe}(\text{scale} = \lambda) \sim \text{Normal}(0, \lambda * \sigma)$$

where $\sigma \sim \text{half_cauchy}(0, 1)$

Value

a distribution instance.

References

- [Carvalho, Polson, Scott. Handling Sparsity via the Horseshoe \(2008\).](#)
- [Barry, Parlange, Li. Approximation for the exponential integral \(2000\).](#)

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space_model()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_independent

Independent distribution from batch of distributions

Description

This distribution is useful for regarding a collection of independent, non-identical distributions as a single random variable. For example, the Independent distribution composed of a collection of Bernoulli distributions might define a distribution over an image (where each Bernoulli is a distribution over each pixel).

Usage

```
tfd_independent(
  distribution,
  reinterpreted_batch_ndims = NULL,
  validate_args = FALSE,
  name = paste0("Independent", distribution$name)
)
```

Arguments

distribution The base distribution instance to transform. Typically an instance of `Distribution`

reinterpreted_batch_ndims Scalar, integer number of rightmost batch dims which will be regarded as event dims. When `NULL` all but the first batch axis (batch axis 0) will be transferred to event dimensions (analogous to `tf.layers.flatten`).

validate_args Logical, default `FALSE`. When `TRUE` distribution parameters are checked for validity despite possibly degrading runtime performance. When `FALSE` invalid inputs may silently render incorrect outputs. Default value: `FALSE`.

name The name for ops managed by the distribution. Default value: `Independent + distribution.name`.

Details

More precisely, a collection of B (independent) E -variate random variables (rv) $\{X_1, \dots, X_B\}$, can be regarded as a $[B, E]$ -variate random variable (X_1, \dots, X_B) with probability $p(x_1, \dots, x_B) = p_1(x_1) * \dots * p_B(x_B)$ where $p_b(X_b)$ is the probability of the b -th rv. More generally B, E can be arbitrary shapes. Similarly, the `Independent` distribution specifies a distribution over $[B, E]$ -shaped events. It operates by reinterpreting the rightmost batch dims as part of the event dimensions. The `reinterpreted_batch_ndims` parameter controls the number of batch dims which are absorbed as event dims; `reinterpreted_batch_ndims <= len(batch_shape)`. For example, the `log_prob` function entails a `reduce_sum` over the rightmost `reinterpreted_batch_ndims` after calling the base distribution's `log_prob`. In other words, since the batch dimension(s) index independent distributions, the resultant multivariate will have independent components.

Mathematical Details

The probability function is,

```
prob(x; reinterpreted_batch_ndims) =
  tf.reduce_prod(dist.prob(x), axis=-1-range(reinterpreted_batch_ndims))
```

Value

a distribution instance.

See Also

For usage examples see e.g. [tfd_sample\(\)](#), [tfd_log_prob\(\)](#), [tfd_mean\(\)](#).

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space_model()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

<code>tfd_inverse_gamma</code>	<i>InverseGamma distribution</i>
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Description

The InverseGamma distribution is defined over positive real numbers using parameters concentration (aka "alpha") and scale (aka "beta").

Usage

```
tfd_inverse_gamma(
  concentration,
  scale,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "InverseGamma"
)
```

Arguments

`concentration` Floating point tensor, the concentration params of the distribution(s). Must contain only positive values.

scale	Floating point tensor, the scale params of the distribution(s). Must contain only positive values. This parameter was called rate before release 0.8.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

Mathematical Details

The probability density function (pdf) is,

$$\text{pdf}(x; \alpha, \beta, x > 0) = x^{-(\alpha + 1)} \exp(-\beta / x) / Z$$

$$Z = \Gamma(\alpha) \beta^{-\alpha}$$

where:

- concentration = alpha,
- scale = beta,
- Z is the normalizing constant, and,
- Gamma is the [gamma function](#).

The cumulative density function (cdf) is,

$$\text{cdf}(x; \alpha, \beta, x > 0) = \Gamma\text{Inc}(\alpha, \beta / x) / \Gamma(\alpha)$$

where `GammaInc` is the [upper incomplete Gamma function](https://en.wikipedia.org/wiki/Incomplete_gamma_function)

The parameters can be intuited via their relationship to mean and variance when these moments exist,

$$\text{mean} = \beta / (\alpha - 1) \text{ when } \alpha > 1 \quad \text{variance} = \beta^2 / (\alpha - 1)^2 / (\alpha - 2) \text{ when } \alpha > 2$$

i.e., under the same conditions:

$$\alpha = \text{mean}^2 / \text{variance} + 2 \quad \beta = \text{mean} * (\text{mean}^2 / \text{variance} + 1)$$

Distribution parameters are automatically broadcast in all functions; see examples for details.

Samples of this distribution are reparameterized (pathwise differentiable).

The derivatives are computed using the approach described in the paper

[Michael Figurnov, Shakir Mohamed, Andriy Mnih. Implicit Reparameterization Gradients, 2018](<https://arxiv.org/abs/1802.02982>)

[gamma function]: `R:gamma%20function`

[upper incomplete Gamma function]: `R:upper%20incomplete%20Gamma%20function`

[Michael Figurnov, Shakir Mohamed, Andriy Mnih. Implicit Reparameterization Gradients, 2018]: `R:Michael`

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space_model()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

`tfd_inverse_gaussian` *Inverse Gaussian distribution*

Description

The **inverse Gaussian distribution** is parameterized by a `loc` and a concentration parameter. It's also known as the Wald distribution. Some, e.g., the Python `scipy` package, refer to the special case when `loc` is 1 as the Wald distribution.

Usage

```
tfd_inverse_gaussian(
    loc,
    concentration,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = "InverseGaussian"
)
```

Arguments

<code>loc</code>	Floating-point Tensor, the loc params. Must contain only positive values.
<code>concentration</code>	Floating-point Tensor, the concentration params. Must contain only positive values.
<code>validate_args</code>	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
<code>allow_nan_stats</code>	Logical, default <code>TRUE</code> . When <code>TRUE</code> , statistics (e.g., mean, mode, variance) use the value <code>NaN</code> to indicate the result is undefined. When <code>FALSE</code> , an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Details

The "inverse" in the name does not refer to the distribution associated to the multiplicative inverse of a random variable. Rather, the cumulant generating function of this distribution is the inverse to that of a Gaussian random variable.

Mathematical Details

The probability density function (pdf) is,

$$\text{pdf}(x; \mu, \lambda) = [\lambda / (2 \pi x^3)]^{0.5} \exp\{-\lambda(x - \mu)^2 / (2 \mu^2 x)\}$$

where

- `loc` = μ
- `concentration` = λ .

The support of the distribution is defined on $(0, \infty)$. Mapping to R and Python `scipy`'s parameterization:

- R: `statmod::invgauss`
- `mean` = `loc`
- `shape` = `concentration`
- `dispersion` = `1 / concentration`. Used only if `shape` is `NULL`.
- Python: `scipy.stats.invgauss`
- `mu` = `loc / concentration`
- `scale` = `concentration`

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space_model()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_johnson_s_u

Johnson's SU-distribution.

Description

This distribution has parameters: shape parameters skewness and tailweight, location loc, and scale.

Usage

```
tfd_johnson_s_u(
    skewness,
    tailweight,
    loc,
    scale,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = NULL
)
```


Arguments

skewness	Floating-point Tensor. Skewness of the distribution(s).
tailweight	Floating-point Tensor. Tail weight of the distribution(s). <code>tailweight</code> must contain only positive values.
loc	Floating-point Tensor. The mean(s) of the distribution(s).
scale	Floating-point Tensor. The scaling factor(s) for the distribution(s). Note that <code>scale</code> is not technically the standard deviation of this distribution but has semantics more similar to standard deviation than variance.
validate_args	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
allow_nan_stats	Logical, default <code>TRUE</code> . When <code>TRUE</code> , statistics (e.g., mean, mode, variance) use the value <code>NaN</code> to indicate the result is undefined. When <code>FALSE</code> , an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details**Mathematical details**

The probability density function (pdf) is,

$$\text{pdf}(x; s, t, \xi, \sigma) = \exp(-0.5 (s + t \operatorname{arcsinh}(y))^{**2}) / Z$$

where,

$$\begin{aligned} s &= \text{skewness} \\ t &= \text{tailweight} \\ y &= (x - \xi) / \sigma \\ Z &= \sigma \sqrt{2 \pi} \sqrt{1 + y^{**2}} / t \end{aligned}$$

where:

- `loc = xi`,
- `scale = sigma`, and,
- `Z` is the normalization constant. The JohnsonSU distribution is a member of the **location-scale family**, i.e., it can be constructed as,

$$\begin{aligned} X &\sim \text{JohnsonSU}(\text{skewness}, \text{tailweight}, \text{loc}=0, \text{scale}=1) \\ Y &= \text{loc} + \text{scale} * X \end{aligned}$$

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space_model()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

`tfd_joint_distribution_named`

Joint distribution parameterized by named distribution-making functions.

Description

This distribution enables both sampling and joint probability computation from a single model specification. A joint distribution is a collection of possibly interdependent distributions. Like `JointDistributionSequential`, `JointDistributionNamed` is parameterized by several distribution-making functions. Unlike `JointDistributionNamed`, each distribution-making function must have its own key. Additionally every distribution-making function's arguments must refer to only specified keys.

Usage

```
tfd_joint_distribution_named(model, validate_args = FALSE, name = NULL)
```

Arguments

model	named list of distribution-making functions each with required args corresponding only to other keys in the named list.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
name	The name for ops managed by the distribution. Default value: "JointDistributionNamed".

Details**Mathematical Details**

Internally JointDistributionNamed implements the chain rule of probability. That is, the probability function of a length-d vector x is,

$$p(x) = \text{prod}\{ p(x[i] \mid x[:i]) : i = 0, \dots, (d - 1) \}$$

The JointDistributionNamed is parameterized by a dict (or namedtuple) composed of either:

1. `tfp$distributions$Distribution`-like instances or,
2. functions which return a `tfp$distributions$Distribution`-like instance. The "conditioned on" elements are represented by the function's required arguments; every argument must correspond to a key in the named distribution-making functions. Distribution-makers which are directly a `Distribution`-like instance are allowed for convenience and semantically identical a zero argument function. When the maker takes no arguments it is preferable to directly provide the distribution instance.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space_model()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`.

```
tfd_normal(), tfd_one_hot_categorical(), tfd_pareto(), tfd_pixel_cnn(), tfd_poisson_log_normal_quadrature(),
tfd_poisson(), tfd_power_spherical(), tfd_probit_bernoulli(), tfd_quantized(), tfd_relaxed_bernoulli(),
tfd_relaxed_one_hot_categorical(), tfd_sample_distribution(), tfd_sinh_arcsinh(),
tfd_skellam(), tfd_spherical_uniform(), tfd_student_t_process(), tfd_student_t(), tfd_transformed_distribution(),
tfd_triangular(), tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian(),
tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()
```

```
tfd_joint_distribution_named_auto_batched
```

Joint distribution parameterized by named distribution-making functions.

Description

This class provides automatic vectorization and alternative semantics for `tfd_joint_distribution_named()`, which in many cases allows for simplifications in the model specification.

Usage

```
tfd_joint_distribution_named_auto_batched(
  model,
  batch_ndims = 0,
  use_vectorized_map = TRUE,
  validate_args = FALSE,
  name = NULL
)
```

Arguments

<code>model</code>	A generator that yields a sequence of <code>tfd\$Distribution</code> -like instances.
<code>batch_ndims</code>	integer Tensor number of batch dimensions. The <code>batch_shapes</code> of all component distributions must be such that the prefixes of length <code>batch_ndims</code> broadcast to a consistent joint batch shape. Default value: <code>0</code> .
<code>use_vectorized_map</code>	logical. Whether to use <code>tf\$vectorized_map</code> to automatically vectorize evaluation of the model. This allows the model specification to focus on drawing a single sample, which is often simpler, but some ops may not be supported. Default value: <code>TRUE</code> .
<code>validate_args</code>	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
<code>name</code>	name prefixed to Ops created by this class.

Details

Automatic vectorization

Auto-vectorized variants of `JointDistribution` allow the user to avoid explicitly annotating a model's vectorization semantics. When using manually-vectorized joint distributions, each operation in the model must account for the possibility of batch dimensions in `Distributions` and their samples. By contrast, auto-vectorized models need only describe a *single* sample from the joint distribution; any batch evaluation is automated using `tf$vectorized_map` as required. In many cases this allows for significant simplifications. For example, the following manually-vectorized `tfd_joint_distribution_named()` model:

```
model <- tfd_joint_distribution_sequential(
  list(
    x = tfd_normal(loc = 0, scale = tf$ones(3L)),
    y = tfd_normal(loc = 0, scale = 1),
    z = function(y, x) {
      tfd_normal(loc = x[reticulate::py_ellipsis(), 1:2] + y[reticulate::py_ellipsis(), tf$newaxis], s
    }
  )
)
```

can be written in auto-vectorized form as

```
model <- tfd_joint_distribution_sequential_auto_batched(
  list(
    x = tfd_normal(loc = 0, scale = tf$ones(3L)),
    y = tfd_normal(loc = 0, scale = 1),
    z = function(y, x) {tfd_normal(loc = x[1:2] + y, scale = 1)}
  )
)
```

in which we were able to avoid explicitly accounting for batch dimensions when indexing and slicing computed quantities in the third line. Note: auto-vectorization is still experimental and some TensorFlow ops may be unsupported. It can be disabled by setting `use_vectorized_map=FALSE`.

Alternative batch semantics This class also provides alternative semantics for specifying a batch of independent (non-identical) joint distributions. Instead of simply summing the `log_probs` of component distributions (which may have different shapes), it first reduces the component `log_probs` to ensure that `jd$log_prob(jd$sample())` always returns a scalar, unless `batch_ndims` is explicitly set to a nonzero value (in which case the result will have the corresponding tensor rank).

The essential changes are:

- An event of `JointDistributionNamedAutoBatched` is the list of tensors produced by `$sample()`; thus, the `event_shape` is the list containing the shapes of sampled tensors. These combine both the event and batch dimensions of the component distributions. By contrast, the event shape of a base `JointDistributions` does not include batch dimensions of component distributions.
- The `batch_shape` is a global property of the entire model, rather than a per-component property as in base `JointDistributions`. The global batch shape must be a prefix of the batch shapes of each component; the length of this prefix is specified by an optional argument `batch_ndims`. If `batch_ndims` is not specified, the model has batch shape `()`.#'

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_joint_distribution_sequential

Joint distribution parameterized by distribution-making functions

Description

This distribution enables both sampling and joint probability computation from a single model specification.

Usage

```
tfd_joint_distribution_sequential(model, validate_args = FALSE, name = NULL)
```

Arguments

model	list of either <code>tfp\$distributions\$Distribution</code> instances and/or functions which take the <code>k</code> previous distributions and returns a new <code>tfp\$distributions\$Distribution</code> instance.
-------	--

validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
name	name prefixed to Ops created by this class.

Details

A joint distribution is a collection of possibly interdependent distributions. Like `tf$keras$Sequential`, the `JointDistributionSequential` can be specified via a list of functions (each responsible for making a `tfp$distributions$Distribution`-like instance). Unlike `tf$keras$Sequential`, each function can depend on the output of all previous elements rather than only the immediately previous.

Mathematical Details

The `JointDistributionSequential` implements the chain rule of probability.

That is, the probability function of a length-d vector x is,

$$p(x) = \prod\{ p(x[i] \mid x[:i]) : i = 0, \dots, (d - 1) \}$$

The `JointDistributionSequential` is parameterized by a list comprised of either:

1. `tfp$distributions$Distribution`-like instances or,
2. callables which return a `tfp$distributions$Distribution`-like instance. Each list element implements the i -th *full conditional distribution*, $p(x[i] \mid x[:i])$. The "conditioned on" elements are represented by the callable's required arguments. Directly providing a `Distribution`-like instance is a convenience and is semantically identical to a zero argument callable. Denote the i -th callables non-default arguments as `args[i]`. Since the callable is the conditional manifest, $0 \leq \text{len}(\text{args}[i]) \leq i - 1$. When $\text{len}(\text{args}[i]) < i - 1$, the callable only depends on a subset of the previous distributions, specifically those at indexes: `range(i - 1, i - 1 - num_args[i], -1)`.

Name resolution: The names of `JointDistributionSequential` components are defined by explicit-name arguments passed to distributions (`tfd.Normal(0., 1., name='x')`) and/or by the argument names in distribution-making functions (`lambda x: tfd.Normal(x., 1.)`). Both approaches may be used in the same distribution, as long as they are consistent; referring to a single component by multiple names will raise a `ValueError`. Unnamed components will be assigned a dummy name.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`,

```
tfd_generalized_normal(), tfd_geometric(), tfd_gumbel(), tfd_half_cauchy(), tfd_half_normal(),
tfd_hidden_markov_model(), tfd_horseshoe(), tfd_independent(), tfd_inverse_gamma(),
tfd_inverse_gaussian(), tfd_johnson_s_u(), tfd_joint_distribution_named_auto_batched(),
tfd_joint_distribution_named(), tfd_joint_distribution_sequential_auto_batched(),
tfd_kumaraswamy(), tfd_laplace(), tfd_linear_gaussian_state_space_model(), tfd_lkj(),
tfd_log_logistic(), tfd_log_normal(), tfd_logistic(), tfd_mixture_same_family(), tfd_mixture(),
tfd_multinomial(), tfd_multivariate_normal_diag_plus_low_rank(), tfd_multivariate_normal_diag(),
tfd_multivariate_normal_full_covariance(), tfd_multivariate_normal_linear_operator(),
tfd_multivariate_normal_tri_l(), tfd_multivariate_student_t_linear_operator(), tfd_negative_binomial(),
tfd_normal(), tfd_one_hot_categorical(), tfd_pareto(), tfd_pixel_cnn(), tfd_poisson_log_normal_quadrature(),
tfd_poisson(), tfd_power_spherical(), tfd_probit_bernoulli(), tfd_quantized(), tfd_relaxed_bernoulli(),
tfd_relaxed_one_hot_categorical(), tfd_sample_distribution(), tfd_sinh_arcsinh(),
tfd_skellam(), tfd_spherical_uniform(), tfd_student_t_process(), tfd_student_t(), tfd_transformed_distribution(),
tfd_triangular(), tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian(),
tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()
```

```
tfd_joint_distribution_sequential_auto_batched
```

Joint distribution parameterized by distribution-making functions.

Description

This class provides automatic vectorization and alternative semantics for `tfd_joint_distribution_sequential()`, which in many cases allows for simplifications in the model specification.

Usage

```
tfd_joint_distribution_sequential_auto_batched(
  model,
  batch_ndims = 0,
  use_vectorized_map = TRUE,
  validate_args = FALSE,
  name = NULL
)
```

Arguments

<code>model</code>	A generator that yields a sequence of <code>tfd\$Distribution</code> -like instances.
<code>batch_ndims</code>	integer Tensor number of batch dimensions. The <code>batch_shapes</code> of all component distributions must be such that the prefixes of length <code>batch_ndims</code> broadcast to a consistent joint batch shape. Default value: <code>0</code> .

use_vectorized_map	logical. Whether to use <code>tf\$vectorized_map</code> to automatically vectorize evaluation of the model. This allows the model specification to focus on drawing a single sample, which is often simpler, but some ops may not be supported. Default value: TRUE.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
name	name prefixed to Ops created by this class.

Details

Automatic vectorization

Auto-vectorized variants of `JointDistribution` allow the user to avoid explicitly annotating a model's vectorization semantics. When using manually-vectorized joint distributions, each operation in the model must account for the possibility of batch dimensions in `Distributions` and their samples. By contrast, auto-vectorized models need only describe a *single* sample from the joint distribution; any batch evaluation is automated using `tf$vectorized_map` as required. In many cases this allows for significant simplifications. For example, the following manually-vectorized `tfd_joint_distribution_sequential()` model:

```
model <- tfd_joint_distribution_sequential(
  list(
    tfd_normal(loc = 0, scale = tf$ones(3L)),
    tfd_normal(loc = 0, scale = 1),
    function(y, x) {
      tfd_normal(loc = x[reticulate::py_ellipsis(), 1:2] + y[reticulate::py_ellipsis(), tf$newaxis], s
    }
  )
)
```

can be written in auto-vectorized form as

```
model <- tfd_joint_distribution_sequential_auto_batched(
  list(
    tfd_normal(loc = 0, scale = tf$ones(3L)),
    tfd_normal(loc = 0, scale = 1),
    function(y, x) {tfd_normal(loc = x[1:2] + y, scale = 1)}
  )
)
```

in which we were able to avoid explicitly accounting for batch dimensions when indexing and slicing computed quantities in the third line. Note: auto-vectorization is still experimental and some TensorFlow ops may be unsupported. It can be disabled by setting `use_vectorized_map=FALSE`.

Alternative batch semantics This class also provides alternative semantics for specifying a batch of independent (non-identical) joint distributions. Instead of simply summing the `log_probs` of component distributions (which may have different shapes), it first reduces the component `log_probs` to

ensure that `jd$log_prob(jd$sample())` always returns a scalar, unless `batch_ndims` is explicitly set to a nonzero value (in which case the result will have the corresponding tensor rank).

The essential changes are:

- An event of `JointDistributionSequentialAutoBatched` is the list of tensors produced by `$sample()`; thus, the `event_shape` is the list containing the shapes of sampled tensors. These combine both the event and batch dimensions of the component distributions. By contrast, the event shape of a base `JointDistributions` does not include batch dimensions of component distributions.
- The `batch_shape` is a global property of the entire model, rather than a per-component property as in base `JointDistributions`. The global batch shape must be a prefix of the batch shapes of each component; the length of this prefix is specified by an optional argument `batch_ndims`. If `batch_ndims` is not specified, the model has batch shape `()`.#

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space_model()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag()`, `tfd_multivariate_normal()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature_compound()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian_process()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_kl_divergence *Computes the Kullback–Leibler divergence.*

Description

Denote this distribution by p and the other distribution by q . Assuming p, q are absolutely continuous with respect to reference measure r , the KL divergence is defined as: $KL[p, q] = E_p[\log(p(X)/q(X))] = -\int_F p(x) \log q(x) dr(x) + \int_F p(x) \log p(x) dr(x) = H[p, q] - H[p]$ where F denotes the support of the random variable $X \sim p$, $H[\cdot, \cdot]$ denotes (Shannon) cross entropy, and $H[\cdot]$ denotes (Shannon) entropy.

Usage

```
tfd_kl_divergence(distribution, other, name = "kl_divergence")
```

Arguments

distribution	The distribution being used.
other	tfp\$distributions\$Distribution instance.
name	String prepended to names of ops created by this function.

Value

self\$type Tensor with shape $[B_1, \dots, B_n]$ representing n different calculations of the Kullback–Leibler divergence.

See Also

Other distribution_methods: [tfd_cdf\(\)](#), [tfd_covariance\(\)](#), [tfd_cross_entropy\(\)](#), [tfd_entropy\(\)](#), [tfd_log_cdf\(\)](#), [tfd_log_prob\(\)](#), [tfd_log_survival_function\(\)](#), [tfd_mean\(\)](#), [tfd_mode\(\)](#), [tfd_prob\(\)](#), [tfd_quantile\(\)](#), [tfd_sample\(\)](#), [tfd_stddev\(\)](#), [tfd_survival_function\(\)](#), [tfd_variance\(\)](#)

Examples

```
d1 <- tfd_normal(loc = c(1, 2), scale = c(1, 0.5))
d2 <- tfd_normal(loc = c(1.5, 2), scale = c(1, 0.5))
d1 %>% tfd_kl_divergence(d2)
```

tfd_kumaraswamy *Kumaraswamy distribution*

Description

The Kumaraswamy distribution is defined over the (0, 1) interval using parameters `concentration1` (aka "alpha") and `concentration0` (aka "beta"). It has a shape similar to the Beta distribution, but is easier to reparameterize.

Usage

```
tfd_kumaraswamy(
  concentration1 = 1,
  concentration0 = 1,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "Kumaraswamy"
)
```

Arguments

<code>concentration1</code>	Positive floating-point Tensor indicating mean number of successes; aka "alpha". Implies <code>self.dtype</code> and <code>self.batch_shape</code> , i.e., <code>concentration1.shape = [N1, N2, ..., Nm] = self.batch_shape</code> .
<code>concentration0</code>	Positive floating-point Tensor indicating mean number of failures; aka "beta". Otherwise has same semantics as <code>concentration1</code> .
<code>validate_args</code>	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
<code>allow_nan_stats</code>	Logical, default <code>TRUE</code> . When <code>TRUE</code> , statistics (e.g., mean, mode, variance) use the value <code>NaN</code> to indicate the result is undefined. When <code>FALSE</code> , an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Details

Mathematical Details

The probability density function (pdf) is,

$$\text{pdf}(x; \alpha, \beta) = \alpha * \beta * x^{(\alpha - 1)} * (1 - x^{\alpha})^{(\beta - 1)}$$

where:

- `concentration1 = alpha`,
- `concentration0 = beta`, Distribution parameters are automatically broadcast in all functions.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space_model()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag_plus_low_rank_linear_operator()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_laplace

Laplace distribution with location loc and scale parameters

Description

Mathematical details

Usage

```
tfd_laplace(
  loc,
  scale,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "Laplace"
)
```

Arguments

loc	Floating point tensor which characterizes the location (center) of the distribution.
scale	Positive floating point tensor which characterizes the spread of the distribution.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

The probability density function (pdf) of this distribution is,

$$\text{pdf}(x; \mu, \sigma) = \exp(-|x - \mu| / \sigma) / Z$$

$$Z = 2 \sigma$$

where $\text{loc} = \mu$, $\text{scale} = \sigma$, and Z is the normalization constant.

Note that the Laplace distribution can be thought of two exponential distributions spliced together "back-to-back." The Laplace distribution is a member of the **location-scale family**, i.e., it can be constructed as,

$$X \sim \text{Laplace}(\text{loc}=0, \text{scale}=1)$$

$$Y = \text{loc} + \text{scale} * X$$

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_linear_gaussian_state_space_model()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal()`.

```
tfd_multivariate_normal_full_covariance(), tfd_multivariate_normal_linear_operator(),
tfd_multivariate_normal_tri_l(), tfd_multivariate_student_t_linear_operator(), tfd_negative_binomial(),
tfd_normal(), tfd_one_hot_categorical(), tfd_pareto(), tfd_pixel_cnn(), tfd_poisson_log_normal_quadrature(),
tfd_poisson(), tfd_power_spherical(), tfd_probit_bernoulli(), tfd_quantized(), tfd_relaxed_bernoulli(),
tfd_relaxed_one_hot_categorical(), tfd_sample_distribution(), tfd_sinh_arcsinh(),
tfd_skellam(), tfd_spherical_uniform(), tfd_student_t_process(), tfd_student_t(), tfd_transformed_distribution(),
tfd_triangular(), tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian(),
tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()
```

```
tfd_linear_gaussian_state_space_model
```

Observation distribution from a linear Gaussian state space model

Description

The state space model, sometimes called a Kalman filter, posits a latent state vector z_t of dimension `latent_size` that evolves over time following linear Gaussian transitions, $z_{t+1} = F * z_t + N(b; Q)$ for transition matrix F , bias b and covariance matrix Q . At each timestep, we observe a noisy projection of the latent state $x_t = H * z_t + N(c; R)$. The transition and observation models may be fixed or may vary between timesteps.

Usage

```
tfd_linear_gaussian_state_space_model(
    num_timesteps,
    transition_matrix,
    transition_noise,
    observation_matrix,
    observation_noise,
    initial_state_prior,
    initial_step = 0L,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = "LinearGaussianStateSpaceModel"
)
```

Arguments

`num_timesteps` Integer Tensor total number of timesteps.

`transition_matrix`

A transition operator, represented by a Tensor or LinearOperator of shape `[latent_size, latent_size]`, or by a callable taking as argument a scalar integer Tensor t and returning a Tensor or LinearOperator representing the transition operator from latent state at time t to time $t + 1$.

<code>transition_noise</code>	An instance of <code>tfd\$MultivariateNormalLinearOperator</code> with event shape <code>[latent_size]</code> , representing the mean and covariance of the transition noise model, or a callable taking as argument a scalar integer Tensor <code>t</code> and returning such a distribution representing the noise in the transition from time <code>t</code> to time <code>t + 1</code> .
<code>observation_matrix</code>	An observation operator, represented by a Tensor or <code>LinearOperator</code> of shape <code>[observation_size, latent_size]</code> , or by a callable taking as argument a scalar integer Tensor <code>t</code> and returning a timestep-specific Tensor or <code>LinearOperator</code> .
<code>observation_noise</code>	An instance of <code>tfd.MultivariateNormalLinearOperator</code> with event shape <code>[observation_size]</code> , representing the mean and covariance of the observation noise model, or a callable taking as argument a scalar integer Tensor <code>t</code> and returning a timestep-specific noise model.
<code>initial_state_prior</code>	An instance of <code>MultivariateNormalLinearOperator</code> representing the prior distribution on latent states; must have event shape <code>[latent_size]</code> .
<code>initial_step</code>	optional integer specifying the time of the first modeled timestep. This is added as an offset when passing timesteps <code>t</code> to (optional) callables specifying timestep-specific transition and observation models.
<code>validate_args</code>	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
<code>allow_nan_stats</code>	Logical, default <code>TRUE</code> . When <code>TRUE</code> , statistics (e.g., mean, mode, variance) use the value <code>NaN</code> to indicate the result is undefined. When <code>FALSE</code> , an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	<code>name</code> prefixed to Ops created by this class.

Details

This Distribution represents the marginal distribution on observations, $p(x)$. The marginal `log_prob` is computed by Kalman filtering, and `sample` by an efficient forward recursion. Both operations require time linear in T , the total number of timesteps.

Shapes

The event shape is `[num_timesteps, observation_size]`, where `observation_size` is the dimension of each observation x_t . The observation and transition models must return consistent shapes. This implementation supports vectorized computation over a batch of models. All of the parameters (prior distribution, transition and observation operators and noise models) must have a consistent batch shape.

Time-varying processes

Any of the model-defining parameters (prior distribution, transition and observation operators and noise models) may be specified as a callable taking an integer timestep `t` and returning a time-dependent value. The dimensionality (`latent_size` and `observation_size`) must be the same at all timesteps.

Importantly, the timestep is passed as a Tensor, not a Python integer, so any conditional behavior must occur *inside* the TensorFlow graph. For example, suppose we want to use a different transition model on even days than odd days. It does *not* work to write

```
transition_matrix <- function(t) {
  if(t %% 2 == 0) even_day_matrix else odd_day_matrix
}
```

since the value of `t` is not fixed at graph-construction time. Instead we need to write

```
transition_matrix <- function(t) {
  tf$cond(tf$equal(tf$mod(t, 2), 0), function() even_day_matrix, function() odd_day_matrix)
}
```

so that TensorFlow can switch between operators appropriately at runtime.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature_compound()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian_process()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_lkj

*LKJ distribution on correlation matrices***Description**

This is a one-parameter of distributions on correlation matrices. The probability density is proportional to the determinant raised to the power of the parameter: $\text{pdf}(X; \eta) = Z(\eta) * \det(X) ** (\eta - 1)$, where $Z(\eta)$ is a normalization constant. The uniform distribution on correlation matrices is the special case $\eta = 1$.

Usage

```
tfd_lkj(
    dimension,
    concentration,
    input_output_cholesky = FALSE,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = "LKJ"
)
```

Arguments

<code>dimension</code>	integer. The dimension of the correlation matrices to sample.
<code>concentration</code>	float or double Tensor. The positive concentration parameter of the LKJ distributions. The pdf of a sample matrix X is proportional to $\det(X) ** (\text{concentration} - 1)$.
<code>input_output_cholesky</code>	Logical. If TRUE, functions whose input or output have the semantics of samples assume inputs are in Cholesky form and return outputs in Cholesky form. In particular, if this flag is TRUE, input to <code>log_prob</code> is presumed of Cholesky form and output from <code>sample</code> is of Cholesky form. Setting this argument to TRUE is purely a computational optimization and does not change the underlying distribution. Additionally, validation checks which are only defined on the multiplied-out form are omitted, even if <code>validate_args</code> is TRUE. Default value: FALSE (i.e., input/output does not have Cholesky semantics).
<code>validate_args</code>	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
<code>allow_nan_stats</code>	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Details

The distribution is named after Lewandowski, Kurowicka, and Joe, who gave a sampler for the distribution in Lewandowski, Kurowicka, Joe, 2009.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_logistic

Logistic distribution with location loc and scale parameters

Description

Mathematical details

Usage

```
tfd_logistic(
  loc,
  scale,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "Logistic"
)
```

Arguments

loc	Floating point tensor, the means of the distribution(s).
scale	Floating point tensor, the scales of the distribution(s). Must contain only positive values.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

The cumulative density function of this distribution is:

$$\text{cdf}(x; \mu, \sigma) = 1 / (1 + \exp(-(x - \mu) / \sigma))$$

where $\text{loc} = \mu$ and $\text{scale} = \sigma$.

The Logistic distribution is a member of the **location-scale family**, i.e., it can be constructed as,

$$\begin{aligned} X &\sim \text{Logistic}(\text{loc}=0, \text{scale}=1) \\ Y &= \text{loc} + \text{scale} * X \end{aligned}$$

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`,

```
tfd_generalized_normal(), tfd_geometric(), tfd_gumbel(), tfd_half_cauchy(), tfd_half_normal(),
tfd_hidden_markov_model(), tfd_horseshoe(), tfd_independent(), tfd_inverse_gamma(),
tfd_inverse_gaussian(), tfd_johnson_s_u(), tfd_joint_distribution_named_auto_batched(),
tfd_joint_distribution_named(), tfd_joint_distribution_sequential_auto_batched(),
tfd_joint_distribution_sequential(), tfd_kumaraswamy(), tfd_laplace(), tfd_linear_gaussian_state_space,
tfd_lkj(), tfd_log_logistic(), tfd_log_normal(), tfd_mixture_same_family(), tfd_mixture(),
tfd_multinomial(), tfd_multivariate_normal_diag_plus_low_rank(), tfd_multivariate_normal_diag(),
tfd_multivariate_normal_full_covariance(), tfd_multivariate_normal_linear_operator(),
tfd_multivariate_normal_tri_l(), tfd_multivariate_student_t_linear_operator(), tfd_negative_binomial(),
tfd_normal(), tfd_one_hot_categorical(), tfd_pareto(), tfd_pixel_cnn(), tfd_poisson_log_normal_quadrature,
tfd_poisson(), tfd_power_spherical(), tfd_probit_bernoulli(), tfd_quantized(), tfd_relaxed_bernoulli(),
tfd_relaxed_one_hot_categorical(), tfd_sample_distribution(), tfd_sinh_arcsinh(),
tfd_skellam(), tfd_spherical_uniform(), tfd_student_t_process(), tfd_student_t(), tfd_transformed_distribution,
tfd_triangular(), tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian,
tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()
```

tfd_logit_normal	<i>The Logit-Normal distribution</i>
------------------	--------------------------------------

Description

The Logit-Normal distribution models positive-valued random variables whose logit (i.e., sigmoid_inverse , i.e., $\log(p) - \log(1-p)$) is normally distributed with mean loc and standard deviation scale . It is constructed as the sigmoid transformation, (i.e., $1 / (1 + \exp(-x))$) of a Normal distribution.

Usage

```
tfd_logit_normal(
  loc,
  scale,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "LogitNormal"
)
```

Arguments

<code>loc</code>	Floating point tensor; the means of the distribution(s).
<code>scale</code>	floating point tensor; the stddevs of the distribution(s). Must contain only positive values.
<code>validate_args</code>	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.

allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Value

a distribution instance.

See Also

For usage examples see e.g. [tfd_sample\(\)](#), [tfd_log_prob\(\)](#), [tfd_mean\(\)](#).

tfd_log_cdf	<i>Log cumulative distribution function.</i>
-------------	--

Description

Given random variable X , the cumulative distribution function cdf is: $\text{tfd_log_cdf}(x) := \text{Log}[P[X \leq x]]$ Often, a numerical approximation can be used for $\text{tfd_log_cdf}(x)$ that yields a more accurate answer than simply taking the logarithm of the cdf when $x \ll -1$.

Usage

```
tfd_log_cdf(distribution, value, ...)
```

Arguments

distribution	The distribution being used.
value	float or double Tensor.
...	Additional parameters passed to Python.

Value

a Tensor of shape $\text{sample_shape}(x) + \text{self}\$batch_shape$ with values of type $\text{self}\$dtype$.

See Also

Other distribution_methods: [tfd_cdf\(\)](#), [tfd_covariance\(\)](#), [tfd_cross_entropy\(\)](#), [tfd_entropy\(\)](#), [tfd_kl_divergence\(\)](#), [tfd_log_prob\(\)](#), [tfd_log_survival_function\(\)](#), [tfd_mean\(\)](#), [tfd_mode\(\)](#), [tfd_prob\(\)](#), [tfd_quantile\(\)](#), [tfd_sample\(\)](#), [tfd_stddev\(\)](#), [tfd_survival_function\(\)](#), [tfd_variance\(\)](#)

Examples

```
d <- tfd_normal(loc = c(1, 2), scale = c(1, 0.5))
x <- d %>% tfd_sample()
d %>% tfd_log_cdf(x)
```

tfd_log_logistic	<i>The log-logistic distribution.</i>
------------------	---------------------------------------

Description

The LogLogistic distribution models positive-valued random variables whose logarithm is a logistic distribution with loc `loc` and scale `scale`. It is constructed as the exponential transformation of a Logistic distribution.

Usage

```
tfd_log_logistic(
  loc,
  scale,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "LogLogistic"
)
```

Arguments

<code>loc</code>	Floating-point Tensor; the loc of the underlying logistic distribution(s).
<code>scale</code>	Floating-point Tensor; the scale of the underlying logistic distribution(s).
<code>validate_args</code>	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
<code>allow_nan_stats</code>	Logical, default <code>TRUE</code> . When <code>TRUE</code> , statistics (e.g., mean, mode, variance) use the value <code>NaN</code> to indicate the result is undefined. When <code>FALSE</code> , an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Value

a distribution instance.

See Also

For usage examples see e.g. [tfd_sample\(\)](#), [tfd_log_prob\(\)](#), [tfd_mean\(\)](#).

Other distributions: [tfd_autoregressive\(\)](#), [tfd_batch_reshape\(\)](#), [tfd_bates\(\)](#), [tfd_bernoulli\(\)](#), [tfd_beta_binomial\(\)](#), [tfd_beta\(\)](#), [tfd_binomial\(\)](#), [tfd_categorical\(\)](#), [tfd_cauchy\(\)](#), [tfd_chi2\(\)](#), [tfd_chi\(\)](#), [tfd_cholesky_lkj\(\)](#), [tfd_continuous_bernoulli\(\)](#), [tfd_deterministic\(\)](#), [tfd_dirichlet_multinomial\(\)](#), [tfd_dirichlet\(\)](#), [tfd_empirical\(\)](#), [tfd_exp_gamma\(\)](#), [tfd_exp_inverse_gamma\(\)](#), [tfd_exponential\(\)](#), [tfd_gamma_gamma\(\)](#), [tfd_gamma\(\)](#), [tfd_gaussian_process_regression_model\(\)](#), [tfd_gaussian_process\(\)](#), [tfd_generalized_normal\(\)](#), [tfd_geometric\(\)](#), [tfd_gumbel\(\)](#), [tfd_half_cauchy\(\)](#), [tfd_half_normal\(\)](#), [tfd_hidden_markov_model\(\)](#), [tfd_horseshoe\(\)](#), [tfd_independent\(\)](#), [tfd_inverse_gamma\(\)](#),

```

tfd_inverse_gaussian(), tfd_johnson_s_u(), tfd_joint_distribution_named_auto_batched(),
tfd_joint_distribution_named(), tfd_joint_distribution_sequential_auto_batched(),
tfd_joint_distribution_sequential(), tfd_kumaraswamy(), tfd_laplace(), tfd_linear_gaussian_state_space,
tfd_lkj(), tfd_log_normal(), tfd_logistic(), tfd_mixture_same_family(), tfd_mixture(),
tfd_multinomial(), tfd_multivariate_normal_diag_plus_low_rank(), tfd_multivariate_normal_diag(),
tfd_multivariate_normal_full_covariance(), tfd_multivariate_normal_linear_operator(),
tfd_multivariate_normal_tri_l(), tfd_multivariate_student_t_linear_operator(), tfd_negative_binomial(),
tfd_normal(), tfd_one_hot_categorical(), tfd_pareto(), tfd_pixel_cnn(), tfd_poisson_log_normal_quadrature,
tfd_poisson(), tfd_power_spherical(), tfd_probit_bernoulli(), tfd_quantized(), tfd_relaxed_bernoulli(),
tfd_relaxed_one_hot_categorical(), tfd_sample_distribution(), tfd_sinh_arcsinh(),
tfd_skellam(), tfd_spherical_uniform(), tfd_student_t_process(), tfd_student_t(), tfd_transformed_distribution,
tfd_triangular(), tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian,
tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()

```

tfd_log_normal	<i>Log-normal distribution</i>
----------------	--------------------------------

Description

The LogNormal distribution models positive-valued random variables whose logarithm is normally distributed with mean `loc` and standard deviation `scale`. It is constructed as the exponential transformation of a Normal distribution.

The LogNormal distribution models positive-valued random variables whose logarithm is normally distributed with mean `loc` and standard deviation `scale`. It is constructed as the exponential transformation of a Normal distribution.

Usage

```

tfd_log_normal(
  loc,
  scale,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "LogNormal"
)

```

```

tfd_log_normal(
  loc,
  scale,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "LogNormal"
)

```


Arguments

<code>loc</code>	Floating-point Tensor; the means of the underlying Normal distribution(s).
<code>scale</code>	Floating-point Tensor; the stddevs of the underlying Normal distribution(s).
<code>validate_args</code>	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
<code>allow_nan_stats</code>	Logical, default <code>TRUE</code> . When <code>TRUE</code> , statistics (e.g., mean, mode, variance) use the value <code>NaN</code> to indicate the result is undefined. When <code>FALSE</code> , an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_log_prob	<i>Log probability density/mass function.</i>
--------------	---

Description

Log probability density/mass function.

Usage

```
tfd_log_prob(distribution, value, ...)
```

Arguments

distribution	The distribution being used.
value	float or double Tensor.
...	Additional parameters passed to Python.

Value

a Tensor of shape `sample_shape(x) + self$batch_shape` with values of type `self$dtype`.

See Also

Other `distribution_methods`: [tfd_cdf\(\)](#), [tfd_covariance\(\)](#), [tfd_cross_entropy\(\)](#), [tfd_entropy\(\)](#), [tfd_kl_divergence\(\)](#), [tfd_log_cdf\(\)](#), [tfd_log_survival_function\(\)](#), [tfd_mean\(\)](#), [tfd_mode\(\)](#), [tfd_prob\(\)](#), [tfd_quantile\(\)](#), [tfd_sample\(\)](#), [tfd_stddev\(\)](#), [tfd_survival_function\(\)](#), [tfd_variance\(\)](#)

Examples

```
d <- tfd_normal(loc = c(1, 2), scale = c(1, 0.5))
x <- d %>% tfd_sample()
d %>% tfd_log_prob(x)
```

tfd_log_survival_function	<i>Log survival function.</i>
---------------------------	-------------------------------

Description

Given random variable X , the survival function is defined: `tfd_log_survival_function(x) = $\text{Log}[P[X > x]] = \text{Log}[1 - P[X \leq x]] = \text{Log}[1 - \text{cdf}(x)]$`

Usage

```
tfd_log_survival_function(distribution, value, ...)
```

Arguments

`distribution` The distribution being used.
`value` float or double Tensor.
`...` Additional parameters passed to Python.

Details

Typically, different numerical approximations can be used for the log survival function, which are more accurate than $1 - \text{cdf}(x)$ when $x \gg 1$.

Value

a Tensor of shape `sample_shape(x) + self$batch_shape` with values of type `self$dtype`.

See Also

Other `distribution_methods`: [tfd_cdf\(\)](#), [tfd_covariance\(\)](#), [tfd_cross_entropy\(\)](#), [tfd_entropy\(\)](#), [tfd_kl_divergence\(\)](#), [tfd_log_cdf\(\)](#), [tfd_log_prob\(\)](#), [tfd_mean\(\)](#), [tfd_mode\(\)](#), [tfd_prob\(\)](#), [tfd_quantile\(\)](#), [tfd_sample\(\)](#), [tfd_stddev\(\)](#), [tfd_survival_function\(\)](#), [tfd_variance\(\)](#)

Examples

```
d <- tfd_normal(loc = c(1, 2), scale = c(1, 0.5))
x <- d %>% tfd_sample()
d %>% tfd_log_survival_function(x)
```

tfd_mean	<i>Mean.</i>
----------	--------------

Description

Mean.

Usage

```
tfd_mean(distribution, ...)
```

Arguments

`distribution` The distribution being used.
`...` Additional parameters passed to Python.

Value

a Tensor of shape `sample_shape(x) + self$batch_shape` with values of type `self$dtype`.

See Also

Other `distribution_methods`: `tfd_cdf()`, `tfd_covariance()`, `tfd_cross_entropy()`, `tfd_entropy()`, `tfd_kl_divergence()`, `tfd_log_cdf()`, `tfd_log_prob()`, `tfd_log_survival_function()`, `tfd_mode()`, `tfd_prob()`, `tfd_quantile()`, `tfd_sample()`, `tfd_stddev()`, `tfd_survival_function()`, `tfd_variance()`

Examples

```
d <- tfd_normal(loc = c(1, 2), scale = c(1, 0.5))
d %>% tfd_mean()
```

tfd_mixture

Mixture distribution

Description

The Mixture object implements batched mixture distributions. The mixture model is defined by a Categorical distribution (the mixture) and a list of Distribution objects.

Usage

```
tfd_mixture(
  cat,
  components,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "Mixture"
)
```

Arguments

<code>cat</code>	A Categorical distribution instance, representing the probabilities of distributions.
<code>components</code>	A list or tuple of Distribution instances. Each instance must have the same type, be defined on the same domain, and have matching <code>event_shape</code> and <code>batch_shape</code> .
<code>validate_args</code>	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
<code>allow_nan_stats</code>	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Details

Methods supported include `tfd_log_prob`, `tfd_prob`, `tfd_mean`, `tfd_sample`, and `entropy_lower_bound`.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

`tfd_mixture_same_family`

Mixture (same-family) distribution

Description

The `MixtureSameFamily` distribution implements a (batch of) mixture distribution where all components are from different parameterizations of the same distribution type. It is parameterized by a Categorical "selecting distribution" (over k components) and a components distribution, i.e., a Distribution with a rightmost batch shape (equal to $[k]$) which indexes each (batch of) component.

Usage

```
tfd_mixture_same_family(
  mixture_distribution,
  components_distribution,
  reparameterize = FALSE,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "MixtureSameFamily"
)
```

Arguments

mixture_distribution tfp\$distributions\$Categorical-like instance. Manages the probability of selecting components. The number of categories must match the rightmost batch dimension of the `components_distribution`. Must have either scalar `batch_shape` or `batch_shape` matching `components_distribution$batch_shape[-1]`.

components_distribution tfp\$distributions\$Distribution-like instance. Right-most batch dimension indexes components.

reparameterize Logical, default FALSE. Whether to reparameterize samples of the distribution using implicit reparameterization gradients (Figurnov et al., 2018). The gradients for the mixture logits are equivalent to the ones described by (Graves, 2016). The gradients for the components parameters are also computed using implicit reparameterization (as opposed to ancestral sampling), meaning that all components are updated every step. Only works when: (1) `components_distribution` is fully reparameterized; (2) `components_distribution` is either a scalar distribution or fully factorized (tfd.Independent applied to a scalar distribution); (3) batch shape has a known rank. Experimental, may be slow and produce infs/NaNs.

validate_args Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.

allow_nan_stats Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.

name name prefixed to Ops created by this class.

Value

a distribution instance.

References

- Michael Figurnov, Shakir Mohamed and Andriy Mnih. Implicit reparameterization gradients. In *Neural Information Processing Systems*, 2018.

- Alex Graves. Stochastic Backpropagation through Mixture Density Distributions. *arXiv*, 2016.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag()`, `tfd_multivariate_normal()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature_compound()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian_process()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_mode

Mode.

Description

Mode.

Usage

```
tfd_mode(distribution, ...)
```

Arguments

`distribution` The distribution being used.
`...` Additional parameters passed to Python.

Value

a Tensor of shape `sample_shape(x) + self$batch_shape` with values of type `self$dtype`.

See Also

Other `distribution_methods`: `tfd_cdf()`, `tfd_covariance()`, `tfd_cross_entropy()`, `tfd_entropy()`, `tfd_kl_divergence()`, `tfd_log_cdf()`, `tfd_log_prob()`, `tfd_log_survival_function()`, `tfd_mean()`, `tfd_prob()`, `tfd_quantile()`, `tfd_sample()`, `tfd_stddev()`, `tfd_survival_function()`, `tfd_variance()`

Examples

```
d <- tfd_normal(loc = c(1, 2), scale = c(1, 0.5))
d %>% tfd_mode()
```

<code>tfd_multinomial</code>	<i>Multinomial distribution</i>
------------------------------	---------------------------------

Description

This Multinomial distribution is parameterized by `probs`, a (batch of) length- K prob (probability) vectors ($K > 1$) such that `tf.reduce_sum(probs, -1) = 1`, and a `total_count` number of trials, i.e., the number of trials per draw from the Multinomial. It is defined over a (batch of) length- K vector counts such that `tf$reduce_sum(counts, -1) = total_count`. The Multinomial is identically the Binomial distribution when $K = 2$.

Usage

```
tfd_multinomial(
  total_count,
  logits = NULL,
  probs = NULL,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "Multinomial"
)
```

Arguments

<code>total_count</code>	Non-negative floating point tensor with shape broadcastable to $[N_1, \dots, N_m]$ with $m \geq 0$. Defines this as a batch of $N_1 \times \dots \times N_m$ different Multinomial distributions. Its components should be equal to integer values.
<code>logits</code>	Floating point tensor representing unnormalized log-probabilities of a positive event with shape broadcastable to $[N_1, \dots, N_m, K]$ $m \geq 0$, and the same dtype as <code>total_count</code> . Defines this as a batch of $N_1 \times \dots \times N_m$ different K class Multinomial distributions. Only one of <code>logits</code> or <code>probs</code> should be passed in.

<code>probs</code>	Positive floating point tensor with shape broadcastable to $[N_1, \dots, N_m, K]$ $m \geq 0$ and same dtype as <code>total_count</code> . Defines this as a batch of $N_1 \times \dots \times N_m$ different K class Multinomial distributions. <code>probs</code> 's components in the last portion of its shape should sum to 1. Only one of <code>logits</code> or <code>probs</code> should be passed in.
<code>validate_args</code>	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
<code>allow_nan_stats</code>	Logical, default <code>TRUE</code> . When <code>TRUE</code> , statistics (e.g., mean, mode, variance) use the value <code>NaN</code> to indicate the result is undefined. When <code>FALSE</code> , an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Details

Mathematical Details

The Multinomial is a distribution over K -class counts, i.e., a length- K vector of non-negative integer counts $n = [n_0, \dots, n_{\{K-1\}}]$. The probability mass function (pmf) is,

$$\text{pmf}(n; \pi, N) = \frac{\prod_j (\pi_j)^{n_j}}{Z}$$

$$Z = \frac{\prod_j n_j!}{N!}$$

where:

- `probs = pi = [pi_0, ..., pi_{\{K-1\}}]`, $\pi_j > 0$, $\sum_j \pi_j = 1$,
- `total_count = N`, N a positive integer,
- Z is the normalization constant, and,
- $N!$ denotes N factorial.

Distribution parameters are automatically broadcast in all functions; see examples for details.

Pitfalls

The number of classes, K , must not exceed:

- the largest integer representable by `self.dtype`, i.e., $2^{(\text{mantissa_bits}+1)}$ (IEEE754),
- the maximum Tensor index, i.e., $2^{31}-1$.

Note: This condition is validated only when `validate_args = TRUE`.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_multivariate_normal_diag

Multivariate normal distribution on \mathbb{R}^k

Description

The Multivariate Normal distribution is defined over \mathbb{R}^k and parameterized by a (batch of) length- k loc vector (aka "mu") and a (batch of) $k \times k$ scale matrix; `covariance = scale @ scale`. Where `@` denotes matrix-multiplication.

Usage

```
tfd_multivariate_normal_diag(
    loc = NULL,
    scale_diag = NULL,
    scale_identity_multiplier = NULL,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = "MultivariateNormalDiag"
)
```

Arguments

<code>loc</code>	Floating-point Tensor. If this is set to NULL, <code>loc</code> is implicitly 0. When specified, may have shape $[B_1, \dots, B_b, k]$ where $b \geq 0$ and k is the event size.
<code>scale_diag</code>	Non-zero, floating-point Tensor representing a diagonal matrix added to scale. May have shape $[B_1, \dots, B_b, k]$, $b \geq 0$, and characterizes b -batches of $k \times k$ diagonal matrices added to scale. When both <code>scale_identity_multiplier</code> and <code>scale_diag</code> are NULL then scale is the Identity.
<code>scale_identity_multiplier</code>	Non-zero, floating-point Tensor representing a scaled-identity-matrix added to scale. May have shape $[B_1, \dots, B_b]$, $b \geq 0$, and characterizes b -batches of scaled $k \times k$ identity matrices added to scale. When both <code>scale_identity_multiplier</code> and <code>scale_diag</code> are NULL then scale is the Identity.
<code>validate_args</code>	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
<code>allow_nan_stats</code>	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Details**Mathematical Details**

The probability density function (pdf) is,

$$\begin{aligned} \text{pdf}(x; \text{loc}, \text{scale}) &= \exp(-0.5 \|y\|^2) / Z \\ y &= \text{inv}(\text{scale}) @ (x - \text{loc}) \\ Z &= (2 \pi)^{0.5 k} |\det(\text{scale})| \end{aligned}$$

where:

- `loc` is a vector in \mathbb{R}^k ,
- `scale` is a linear operator in $\mathbb{R}^{k \times k}$, $\text{cov} = \text{scale} @ \text{scale}^T$,
- Z denotes the normalization constant, and,
- $\|y\|^2$ denotes the squared Euclidean norm of y .

A (non-batch) scale matrix is:

$$\text{scale} = \text{diag}(\text{scale_diag} + \text{scale_identity_multiplier} * \text{ones}(k))$$

where:

- `scale_diag.shape = [k]`, and,
- `scale_identity_multiplier.shape = []`.

Additional leading dimensions (if any) will index batches.

If both `scale_diag` and `scale_identity_multiplier` are `NULL`, then `scale` is the Identity matrix. The MultivariateNormal distribution is a member of the **location-scale family**, i.e., it can be constructed as,

```
X ~ MultivariateNormal(loc=0, scale=1) # Identity scale, zero shift.
Y = scale @ X + loc
```

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature_compound()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian_process()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

`tfd_multivariate_normal_diag_plus_low_rank`

Multivariate normal distribution on \mathbb{R}^k

Description

The Multivariate Normal distribution is defined over \mathbb{R}^k and parameterized by a (batch of) length- k `loc` vector (aka "`mu`") and a (batch of) $k \times k$ `scale` matrix; `covariance = scale @ scale`. Where `@` denotes matrix-multiplication.

Usage

```
tfd_multivariate_normal_diag_plus_low_rank(
  loc = NULL,
  scale_diag = NULL,
  scale_identity_multiplier = NULL,
  scale_perturb_factor = NULL,
  scale_perturb_diag = NULL,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "MultivariateNormalDiagPlusLowRank"
)
```

Arguments

<code>loc</code>	Floating-point Tensor. If this is set to NULL, <code>loc</code> is implicitly 0. When specified, may have shape $[B_1, \dots, B_b, k]$ where $b \geq 0$ and k is the event size.
<code>scale_diag</code>	Non-zero, floating-point Tensor representing a diagonal matrix added to scale. May have shape $[B_1, \dots, B_b, k]$, $b \geq 0$, and characterizes b -batches of $k \times k$ diagonal matrices added to scale. When both <code>scale_identity_multiplier</code> and <code>scale_diag</code> are NULL then scale is the Identity.
<code>scale_identity_multiplier</code>	Non-zero, floating-point Tensor representing a scaled-identity-matrix added to scale. May have shape $[B_1, \dots, B_b]$, $b \geq 0$, and characterizes b -batches of scaled $k \times k$ identity matrices added to scale. When both <code>scale_identity_multiplier</code> and <code>scale_diag</code> are NULL then scale is the Identity.
<code>scale_perturb_factor</code>	Floating-point Tensor representing a rank- r perturbation added to scale. May have shape $[B_1, \dots, B_b, k, r]$, $b \geq 0$, and characterizes b -batches of rank- r updates to scale. When NULL, no rank- r update is added to scale.#'
<code>scale_perturb_diag</code>	Floating-point Tensor representing a diagonal matrix inside the rank- r perturbation added to scale. May have shape $[B_1, \dots, B_b, r]$, $b \geq 0$, and characterizes b -batches of $r \times r$ diagonal matrices inside the perturbation added to scale. When NULL, an identity matrix is used inside the perturbation. Can only be specified if <code>scale_perturb_factor</code> is also specified.
<code>validate_args</code>	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
<code>allow_nan_stats</code>	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Details

Mathematical Details

The probability density function (pdf) is,

$$\begin{aligned} \text{pdf}(x; \text{loc}, \text{scale}) &= \exp(-0.5 \|y\|^2) / Z \\ y &= \text{inv}(\text{scale}) @ (x - \text{loc}) \\ Z &= (2 \pi)^{0.5 k} |\det(\text{scale})| \end{aligned}$$

where:

- `loc` is a vector in \mathbb{R}^k ,
- `scale` is a linear operator in $\mathbb{R}^{k \times k}$, `cov = scale @ scale.T`,
- `Z` denotes the normalization constant, and,
- `\|y\|^2` denotes the squared Euclidean norm of `y`.

A (non-batch) scale matrix is:

$$\begin{aligned} \text{scale} &= \text{diag}(\text{scale_diag} + \text{scale_identity_multiplier} \text{ones}(k)) + \\ &\text{scale_perturb_factor} @ \text{diag}(\text{scale_perturb_diag}) @ \text{scale_perturb_factor.T} \end{aligned}$$

where:

- `scale_diag.shape = [k]`,
- `scale_identity_multiplier.shape = []`,
- `scale_perturb_factor.shape = [k, r]`, typically $k \gg r$, and,
- `scale_perturb_diag.shape = [r]`.

Additional leading dimensions (if any) will index batches. If both `scale_diag` and `scale_identity_multiplier` are `NULL`, then `scale` is the Identity matrix. The MultivariateNormal distribution is a member of the **location-scale family**, i.e., it can be constructed as,

$$\begin{aligned} X &\sim \text{MultivariateNormal}(\text{loc}=\theta, \text{scale}=1) \quad \# \text{ Identity scale, zero shift.} \\ Y &= \text{scale} @ X + \text{loc} \end{aligned}$$

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`.

```
tfd_lkj(), tfd_log_logistic(), tfd_log_normal(), tfd_logistic(), tfd_mixture_same_family(),
tfd_mixture(), tfd_multinomial(), tfd_multivariate_normal_diag(), tfd_multivariate_normal_full_covariance(),
tfd_multivariate_normal_linear_operator(), tfd_multivariate_normal_tri_l(), tfd_multivariate_student_t(),
tfd_negative_binomial(), tfd_normal(), tfd_one_hot_categorical(), tfd_pareto(), tfd_pixel_cnn(),
tfd_poisson_log_normal_quadrature_compound(), tfd_poisson(), tfd_power_spherical(),
tfd_probit_bernoulli(), tfd_quantized(), tfd_relaxed_bernoulli(), tfd_relaxed_one_hot_categorical(),
tfd_sample_distribution(), tfd_sinh_arcsinh(), tfd_skellam(), tfd_spherical_uniform(),
tfd_student_t_process(), tfd_student_t(), tfd_transformed_distribution(), tfd_triangular(),
tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian_process(),
tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()
```

```
tfd_multivariate_normal_full_covariance
```

Multivariate normal distribution on \mathbb{R}^k

Description

The Multivariate Normal distribution is defined over \mathbb{R}^k and parameterized by a (batch of) length- k loc vector (aka "mu") and a (batch of) $k \times k$ scale matrix; `covariance = scale @ scale.T` where `@` denotes matrix-multiplication.

Usage

```
tfd_multivariate_normal_full_covariance(
    loc = NULL,
    covariance_matrix = NULL,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = "MultivariateNormalFullCovariance"
)
```

Arguments

<code>loc</code>	Floating-point Tensor. If this is set to <code>NULL</code> , <code>loc</code> is implicitly \emptyset . When specified, may have shape <code>[B1, ..., Bb, k]</code> where <code>b >= 0</code> and <code>k</code> is the event size.
<code>covariance_matrix</code>	Floating-point, symmetric positive definite Tensor of same dtype as <code>loc</code> . The strict upper triangle of <code>covariance_matrix</code> is ignored, so if <code>covariance_matrix</code> is not symmetric no error will be raised (unless <code>validate_args</code> is <code>TRUE</code>). <code>covariance_matrix</code> has shape <code>[B1, ..., Bb, k, k]</code> where <code>b >= 0</code> and <code>k</code> is the event size.
<code>validate_args</code>	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .

allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

Mathematical Details

The probability density function (pdf) is,

$$\text{pdf}(x; \text{loc}, \text{scale}) = \exp(-0.5 \|y\|^2) / Z$$

$$y = \text{inv}(\text{scale}) @ (x - \text{loc})$$

$$Z = (2 \pi)^{0.5 k} |\det(\text{scale})|$$

where:

- loc is a vector in \mathbb{R}^k ,
- scale is a linear operator in $\mathbb{R}^{k \times k}$, $\text{cov} = \text{scale} @ \text{scale}.T$,
- Z denotes the normalization constant, and,
- $\|y\|^2$ denotes the squared Euclidean norm of y.

The MultivariateNormal distribution is a member of the **location-scale family**, i.e., it can be constructed as,

$$X \sim \text{MultivariateNormal}(\text{loc}=\theta, \text{scale}=1) \quad \# \text{ Identity scale, zero shift.}$$

$$Y = \text{scale} @ X + \text{loc}$$

The batch_shape is the broadcast shape between loc and covariance_matrix arguments. The event_shape is given by last dimension of the matrix implied by covariance_matrix. The last dimension of loc (if provided) must broadcast with this. A non-batch covariance_matrix matrix is a $k \times k$ symmetric positive definite matrix. In other words it is (real) symmetric with all eigenvalues strictly positive. Additional leading dimensions (if any) will index batches.

Value

a distribution instance.

See Also

For usage examples see e.g. [tfd_sample\(\)](#), [tfd_log_prob\(\)](#), [tfd_mean\(\)](#).

Other distributions: [tfd_autoregressive\(\)](#), [tfd_batch_reshape\(\)](#), [tfd_bates\(\)](#), [tfd_bernoulli\(\)](#), [tfd_beta_binomial\(\)](#), [tfd_beta\(\)](#), [tfd_binomial\(\)](#), [tfd_categorical\(\)](#), [tfd_cauchy\(\)](#), [tfd_chi2\(\)](#), [tfd_chi\(\)](#), [tfd_cholesky_lkj\(\)](#), [tfd_continuous_bernoulli\(\)](#), [tfd_deterministic\(\)](#), [tfd_dirichlet_multinomial\(\)](#), [tfd_dirichlet\(\)](#), [tfd_empirical\(\)](#), [tfd_exp_gamma\(\)](#), [tfd_exp_inverse_gamma\(\)](#), [tfd_exponential\(\)](#), [tfd_gamma_gamma\(\)](#), [tfd_gamma\(\)](#), [tfd_gaussian_process_regression_model\(\)](#), [tfd_gaussian_process\(\)](#), [tfd_generalized_normal\(\)](#), [tfd_geometric\(\)](#), [tfd_gumbel\(\)](#), [tfd_half_cauchy\(\)](#), [tfd_half_normal\(\)](#), [tfd_hidden_markov_model\(\)](#), [tfd_horseshoe\(\)](#), [tfd_independent\(\)](#), [tfd_inverse_gamma\(\)](#), [tfd_inverse_gaussian\(\)](#), [tfd_johnson_s_u\(\)](#), [tfd_joint_distribution_named_auto_batched\(\)](#),


```
tfd_joint_distribution_named(), tfd_joint_distribution_sequential_auto_batched(),
tfd_joint_distribution_sequential(), tfd_kumaraswamy(), tfd_laplace(), tfd_linear_gaussian_state_space,
tfd_lkj(), tfd_log_logistic(), tfd_log_normal(), tfd_logistic(), tfd_mixture_same_family(),
tfd_mixture(), tfd_multinomial(), tfd_multivariate_normal_diag_plus_low_rank(), tfd_multivariate_normal,
tfd_multivariate_normal_linear_operator(), tfd_multivariate_normal_tri_l(), tfd_multivariate_student_t,
tfd_negative_binomial(), tfd_normal(), tfd_one_hot_categorical(), tfd_pareto(), tfd_pixel_cnn(),
tfd_poisson_log_normal_quadrature_compound(), tfd_poisson(), tfd_power_spherical(),
tfd_probit_bernoulli(), tfd_quantized(), tfd_relaxed_bernoulli(), tfd_relaxed_one_hot_categorical(),
tfd_sample_distribution(), tfd_sinh_arcsinh(), tfd_skellam(), tfd_spherical_uniform(),
tfd_student_t_process(), tfd_student_t(), tfd_transformed_distribution(), tfd_triangular(),
tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian_process(),
tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()
```

tfd_multivariate_normal_linear_operator

The multivariate normal distribution on \mathbb{R}^k

Description

The Multivariate Normal distribution is defined over \mathbb{R}^k and parameterized by a (batch of) length- k loc vector (aka "mu") and a (batch of) $k \times k$ scale matrix; covariance = scale @ scale. Twhere @ denotes matrix-multiplication.

Usage

```
tfd_multivariate_normal_linear_operator(
  loc = NULL,
  scale = NULL,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "MultivariateNormalLinearOperator"
)
```

Arguments

loc	Floating-point Tensor. If this is set to NULL, loc is implicitly \emptyset . When specified, may have shape $[B_1, \dots, B_b, k]$ where $b \geq 0$ and k is the event size.
scale	Instance of LinearOperator with same dtype as loc and shape $[B_1, \dots, B_b, k, k]$.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.

name name prefixed to Ops created by this class.

Details

Mathematical Details

The probability density function (pdf) is,

$$\begin{aligned} \text{pdf}(x; \text{loc}, \text{scale}) &= \exp(-0.5 \|y\|^{**2}) / Z \\ y &= \text{inv}(\text{scale}) @ (x - \text{loc}) \\ Z &= (2 \text{ pi})^{**(\text{0.5 k})} |\det(\text{scale})| \end{aligned}$$

where:

- loc is a vector in \mathbb{R}^k ,
- scale is a linear operator in $\mathbb{R}^{k \times k}$, $\text{cov} = \text{scale} @ \text{scale.T}$,
- Z denotes the normalization constant, and,
- $\|y\|^{**2}$ denotes the squared Euclidean norm of y .

The MultivariateNormal distribution is a member of the **location-scale family**, i.e., it can be constructed as,

```
X ~ MultivariateNormal(loc=0, scale=1) # Identity scale, zero shift.
Y = scale @ X + loc
```

The `batch_shape` is the broadcast shape between `loc` and `scale` arguments. The `event_shape` is given by last dimension of the matrix implied by `scale`. The last dimension of `loc` (if provided) must broadcast with this. Recall that `covariance = scale @ scale.T`. Additional leading dimensions (if any) will index batches.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t()`.

```
tfd_negative_binomial(), tfd_normal(), tfd_one_hot_categorical(), tfd_pareto(), tfd_pixel_cnn(),
tfd_poisson_log_normal_quadrature_compound(), tfd_poisson(), tfd_power_spherical(),
tfd_probit_bernoulli(), tfd_quantized(), tfd_relaxed_bernoulli(), tfd_relaxed_one_hot_categorical(),
tfd_sample_distribution(), tfd_sinh_arcsinh(), tfd_skellam(), tfd_spherical_uniform(),
tfd_student_t_process(), tfd_student_t(), tfd_transformed_distribution(), tfd_triangular(),
tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian_process(),
tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()
```

```
tfd_multivariate_normal_tri_l
```

The multivariate normal distribution on \mathbb{R}^k

Description

The Multivariate Normal distribution is defined over \mathbb{R}^k and parameterized by a (batch of) length- k loc vector (aka "mu") and a (batch of) $k \times k$ scale matrix; covariance = scale @ scale. Twhere @ denotes matrix-multiplication.

Usage

```
tfd_multivariate_normal_tri_l(
    loc = NULL,
    scale_tril = NULL,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = "MultivariateNormalTriL"
)
```

Arguments

loc	Floating-point Tensor. If this is set to NULL, loc is implicitly \emptyset . When specified, may have shape $[B_1, \dots, B_b, k]$ where $b \geq 0$ and k is the event size.
scale_tril	Floating-point, lower-triangular Tensor with non-zero diagonal elements. scale_tril has shape $[B_1, \dots, B_b, k, k]$ where $b \geq 0$ and k is the event size.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details**Mathematical Details**

The probability density function (pdf) is,

$$\begin{aligned} \text{pdf}(x; \text{loc}, \text{scale}) &= \exp(-0.5 \|y\|^2) / Z \\ y &= \text{inv}(\text{scale}) @ (x - \text{loc}) \\ Z &= (2 \pi)^{0.5 k} |\det(\text{scale})| \end{aligned}$$

where:

- loc is a vector in \mathbb{R}^k ,
- scale is a linear operator in $\mathbb{R}^{k \times k}$, $\text{cov} = \text{scale} @ \text{scale} . T$,
- Z denotes the normalization constant, and,
- $\|y\|^2$ denotes the squared Euclidean norm of y .

A (non-batch) scale matrix is:

```
scale = scale_tril
```

where `scale_tril` is lower-triangular $k \times k$ matrix with non-zero diagonal, i.e., `tf$diag_part(scale_tril) != 0`. Additional leading dimensions (if any) will index batches.

The MultivariateNormal distribution is a member of the **location-scale family**, i.e., it can be constructed as,

```
X ~ MultivariateNormal(loc=0, scale=1) # Identity scale, zero shift.
Y = scale @ X + loc
```

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`,

```
tfd_one_hot_categorical(), tfd_pareto(), tfd_pixel_cnn(), tfd_poisson_log_normal_quadrature_compound(),
tfd_poisson(), tfd_power_spherical(), tfd_probit_bernoulli(), tfd_quantized(), tfd_relaxed_bernoulli(),
tfd_relaxed_one_hot_categorical(), tfd_sample_distribution(), tfd_sinh_arcsinh(),
tfd_skellam(), tfd_spherical_uniform(), tfd_student_t_process(), tfd_student_t(), tfd_transformed_distribution(),
tfd_triangular(), tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian(),
tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()
```

tfd_multivariate_student_t_linear_operator

Multivariate Student's t-distribution on \mathbb{R}^k

Description

Mathematical Details

Usage

```
tfd_multivariate_student_t_linear_operator(
    df,
    loc,
    scale,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = "MultivariateStudentTLinearOperator"
)
```

Arguments

df	A positive floating-point Tensor. Has shape $[B_1, \dots, B_b]$ where $b \geq 0$.
loc	Floating-point Tensor. Has shape $[B_1, \dots, B_b, k]$ where k is the event size.
scale	Instance of <code>LinearOperator</code> with a floating dtype and shape $[B_1, \dots, B_b, k, k]$.
validate_args	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
allow_nan_stats	Logical, default <code>TRUE</code> . When <code>TRUE</code> , statistics (e.g., mean, mode, variance) use the value <code>NaN</code> to indicate the result is undefined. When <code>FALSE</code> , an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

The probability density function (pdf) is,

$$\text{pdf}(x; \text{df}, \text{loc}, \text{Sigma}) = (1 + \|y\|^2 / \text{df})^{-(0.5 (\text{df} + k))} / Z$$

where,

$$y = \text{inv}(\text{Sigma}) (x - \text{loc})$$

$$Z = \text{abs}(\text{det}(\text{Sigma})) \sqrt{\text{df} \pi}^{**k} \text{Gamma}(0.5 \text{df}) / \text{Gamma}(0.5 (\text{df} + k))$$

where:

- df is a positive scalar.
- loc is a vector in \mathbb{R}^k ,
- Sigma is a positive definite shape matrix in $\mathbb{R}^{k \times k}$, parameterized as `scale @ scale.T` in this class,
- Z denotes the normalization constant, and,
- $\|y\|^2$ denotes the squared Euclidean norm of y.

The Multivariate Student's t-distribution distribution is a member of the **location-scale family**, i.e., it can be constructed as,

```
X ~ MultivariateT(loc=0, scale=1) # Identity scale, zero shift.
Y = scale @ X + loc
```

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature_compound()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`

tfd_triangular(), tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian(),
 tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
 tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
 tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
 tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()

tfd_negative_binomial *NegativeBinomial distribution*

Description

The NegativeBinomial distribution is related to the experiment of performing Bernoulli trials in sequence. Given a Bernoulli trial with probability p of success, the NegativeBinomial distribution represents the distribution over the number of successes s that occur until we observe f failures.

Usage

```
tfd_negative_binomial(
    total_count,
    logits = NULL,
    probs = NULL,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = "NegativeBinomial"
)
```

Arguments

total_count	Non-negative floating-point Tensor with shape broadcastable to $[B_1, \dots, B_b]$ with $b \geq 0$ and the same dtype as probs or logits. Defines this as a batch of $N_1 \times \dots \times N_m$ different Negative Binomial distributions. In practice, this represents the number of negative Bernoulli trials to stop at (the total_count of failures), but this is still a valid distribution when total_count is a non-integer.
logits	Floating-point Tensor with shape broadcastable to $[B_1, \dots, B_b]$ where $b \geq 0$ indicates the number of batch dimensions. Each entry represents logits for the probability of success for independent Negative Binomial distributions and must be in the open interval $(-\infty, \infty)$. Only one of logits or probs should be specified.
probs	Positive floating-point Tensor with shape broadcastable to $[B_1, \dots, B_b]$ where $b \geq 0$ indicates the number of batch dimensions. Each entry represents the probability of success for independent Negative Binomial distributions and must be in the open interval $(0, 1)$. Only one of logits or probs should be specified.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.

allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

The probability mass function (pmf) is,

$$\text{pmf}(s; f, p) = p^s (1 - p)^{f-s} / Z$$

$$Z = s! (f - s)! / (f - 1)!$$

where:

- total_count = f,
- probs = p,
- Z is the normalizing constant, and,
- n! is the factorial of n.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature_compound()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_normal	<i>Normal distribution with loc and scale parameters</i>
------------	--

Description

Mathematical details

Usage

```
tfd_normal(
  loc,
  scale,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "Normal"
)
```

Arguments

loc	Floating point tensor; the means of the distribution(s).
scale	floating point tensor; the stddevs of the distribution(s). Must contain only positive values.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

The probability density function (pdf) is,

$$\text{pdf}(x; \mu, \sigma) = \exp(-0.5 (x - \mu)^2 / \sigma^2) / Z$$

$$Z = (2 \pi \sigma^2)^{0.5}$$

where loc = mu is the mean, scale = sigma is the std. deviation, and, Z is the normalization constant. The Normal distribution is a member of the **location-scale family**, i.e., it can be constructed as,

$$X \sim \text{Normal}(\text{loc}=0, \text{scale}=1)$$

$$Y = \text{loc} + \text{scale} * X$$

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature_compound()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_one_hot_categorical

OneHotCategorical distribution

Description

The categorical distribution is parameterized by the log-probabilities of a set of classes. The difference between `OneHotCategorical` and `Categorical` distributions is that `OneHotCategorical` is a discrete distribution over one-hot bit vectors whereas `Categorical` is a discrete distribution over positive integers. `OneHotCategorical` is equivalent to `Categorical` except `Categorical` has `event_dim=()` while `OneHotCategorical` has `event_dim=K`, where `K` is the number of classes.

Usage

```
tfd_one_hot_categorical(
  logits = NULL,
```

```

    probs = NULL,
    dtype = tf$int32,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = "OneHotCategorical"
  )

```

Arguments

logits	An N-D Tensor, $N \geq 1$, representing the log probabilities of a set of Categorical distributions. The first $N - 1$ dimensions index into a batch of independent distributions and the last dimension represents a vector of logits for each class. Only one of logits or probs should be passed in.
probs	An N-D Tensor, $N \geq 1$, representing the probabilities of a set of Categorical distributions. The first $N - 1$ dimensions index into a batch of independent distributions and the last dimension represents a vector of probabilities for each class. Only one of logits or probs should be passed in.
dtype	The type of the event samples (default: int32).
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

This class provides methods to create indexed batches of OneHotCategorical distributions. If the provided logits or probs is rank 2 or higher, for every fixed set of leading dimensions, the last dimension represents one single OneHotCategorical distribution. When calling distribution functions (e.g. `dist.prob(x)`), logits and `x` are broadcast to the same shape (if possible). In all cases, the last dimension of logits, `x` represents single OneHotCategorical distributions.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`,

```
tfd_hidden_markov_model(), tfd_horseshoe(), tfd_independent(), tfd_inverse_gamma(),
tfd_inverse_gaussian(), tfd_johnson_s_u(), tfd_joint_distribution_named_auto_batched(),
tfd_joint_distribution_named(), tfd_joint_distribution_sequential_auto_batched(),
tfd_joint_distribution_sequential(), tfd_kumaraswamy(), tfd_laplace(), tfd_linear_gaussian_state_space,
tfd_lkj(), tfd_log_logistic(), tfd_log_normal(), tfd_logistic(), tfd_mixture_same_family(),
tfd_mixture(), tfd_multinomial(), tfd_multivariate_normal_diag_plus_low_rank(), tfd_multivariate_normal,
tfd_multivariate_normal_full_covariance(), tfd_multivariate_normal_linear_operator(),
tfd_multivariate_normal_tri_l(), tfd_multivariate_student_t_linear_operator(), tfd_negative_binomial(),
tfd_normal(), tfd_pareto(), tfd_pixel_cnn(), tfd_poisson_log_normal_quadrature_compound(),
tfd_poisson(), tfd_power_spherical(), tfd_probit_bernoulli(), tfd_quantized(), tfd_relaxed_bernoulli(),
tfd_relaxed_one_hot_categorical(), tfd_sample_distribution(), tfd_sinh_arcsinh(),
tfd_skellam(), tfd_spherical_uniform(), tfd_student_t_process(), tfd_student_t(), tfd_transformed_distribution,
tfd_triangular(), tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian,
tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()
```

tfd_pareto

Pareto distribution

Description

The Pareto distribution is parameterized by a scale and a concentration parameter.

Usage

```
tfd_pareto(
  concentration,
  scale = 1,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "Pareto"
)
```

Arguments

concentration	Floating point tensor. Must contain only positive values.
scale	Floating point tensor, equivalent to mode. scale also restricts the domain of this distribution to be in [scale, inf). Must contain only positive values. Default value: 1.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.

allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

Mathematical Details

The probability density function (pdf) is,

$$\text{pdf}(x; \text{alpha}, \text{scale}, x \geq \text{scale}) = \text{alpha} * \text{scale} ** \text{alpha} / x ** (\text{alpha} + 1)$$

where `concentration = alpha`.

Note that `scale` acts as a scaling parameter, since
 $\text{Pareto}(c, \text{scale}).\text{pdf}(x) == \text{Pareto}(c, 1).\text{pdf}(x / \text{scale})$.
 The support of the distribution is defined on `[scale, infinity)`.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature_compound()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffusion_mixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_pert

Modified PERT distribution for modeling expert predictions.

Description

The PERT distribution is a loc-scale family of Beta distributions fit onto a real interval between low and high values set by the user, along with a peak to indicate the expert's most frequent prediction, and temperature to control how sharp the peak is.

Usage

```
tfd_pert(
  low,
  peak,
  high,
  temperature = 4,
  validate_args = FALSE,
  allow_nan_stats = FALSE,
  name = "Pert"
)
```

Arguments

low	lower bound
peak	most frequent value
high	upper bound
temperature	controls the shape of the distribution
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

The distribution is similar to a **Triangular distribution** (i.e. `tfd.Triangular`) but with a smooth peak.

Mathematical Details

In terms of a Beta distribution, PERT can be expressed as

$$\text{PERT} \sim \text{loc} + \text{scale} * \text{Beta}(\text{concentration1}, \text{concentration0})$$

where

```
loc = low
scale = high - low
concentration1 = 1 + temperature * (peak - low)/(high - low)
concentration0 = 1 + temperature * (high - peak)/(high - low)
temperature > 0
```

The support is [low, high]. The peak must fit in that interval: $low < peak < high$. The temperature is a positive parameter that controls the shape of the distribution. Higher values yield a sharper peak. The standard PERT distribution is obtained when $temperature = 4$.

Value

a distribution instance.

See Also

For usage examples see e.g. [tfd_sample\(\)](#), [tfd_log_prob\(\)](#), [tfd_mean\(\)](#).

tfd_pixel_cnn

The Pixel CNN++ distribution

Description

Pixel CNN++ (Salimans et al., 2017) models a distribution over image data, parameterized by a neural network. It builds on Pixel CNN and Conditional Pixel CNN, as originally proposed by (van den Oord et al., 2016). The model expresses the joint distribution over pixels as the product of conditional distributions: $p(x|h) = \prod_{i=0, \dots, d} p(x[i] | x[0:i], h)$, in which $p(x[i] | x[0:i], h)$ is the probability of the i -th pixel conditional on the pixels that preceded it in raster order (color channels in RGB order, then left to right, then top to bottom). h is optional additional data on which to condition the image distribution, such as class labels or VAE embeddings. The Pixel CNN++ network enforces the dependency structure among pixels by applying a mask to the kernels of the convolutional layers that ensures that the values for each pixel depend only on other pixels up and to the left. Pixel values are modeled with a mixture of quantized logistic distributions, which can take on a set of distinct integer values (e.g. between 0 and 255 for an 8-bit image). Color intensity v of each pixel is modeled as: $v \sim \sum_{i=0, \dots, k} q[i] * \text{quantized_logistic}(\text{loc}[i], \text{scale}[i])$, in which k is the number of mixture components and the $q[i]$ are the Categorical probabilities over the components.

Usage

```
tfd_pixel_cnn(
  image_shape,
  conditional_shape = NULL,
  num_resnet = 5,
  num_hierarchies = 3,
```

```

num_filters = 160,
num_logistic_mix = 10,
receptive_field_dims = c(3, 3),
dropout_p = 0.5,
resnet_activation = "concat_elu",
use_weight_norm = TRUE,
use_data_init = TRUE,
high = 255,
low = 0,
dtype = tf$float32,
name = "PixelCNN"
)

```

Arguments

<code>image_shape</code>	3D TensorShape or tuple for the [height, width, channels] dimensions of the image.
<code>conditional_shape</code>	TensorShape or tuple for the shape of the conditional input, or NULL if there is no conditional input.
<code>num_resnet</code>	integer, the number of layers (shown in Figure 2 of https://arxiv.org/abs/1606.05328) within each highest-level block of Figure 2 of https://pdfs.semanticscholar.org/9e90/6792f67cbdda7b7777b69284a81044857656.pdf .
<code>num_hierarchies</code>	integer, the number of highest-level blocks (separated by expansions/contractions of dimensions in Figure 2 of https://pdfs.semanticscholar.org/9e90/6792f67cbdda7b7777b69284a81044857656.pdf).
<code>num_filters</code>	integer, the number of convolutional filters.
<code>num_logistic_mix</code>	integer, number of components in the logistic mixture distribution.
<code>receptive_field_dims</code>	tuple, height and width in pixels of the receptive field of the convolutional layers above and to the left of a given pixel. The width (second element of the tuple) should be odd. Figure 1 (middle) of https://arxiv.org/abs/1606.05328 shows a receptive field of (3, 5) (the row containing the current pixel is included in the height). The default of (3, 3) was used to produce the results in https://pdfs.semanticscholar.org/9e90/6792f67cbdda7b7777b69284a81044857656.pdf .
<code>dropout_p</code>	float, the dropout probability. Should be between 0 and 1.
<code>resnet_activation</code>	string, the type of activation to use in the resnet blocks. May be 'concat_elu', 'elu', or 'relu'.
<code>use_weight_norm</code>	logical, if TRUE then use weight normalization (works only in Eager mode).
<code>use_data_init</code>	logical, if TRUE then use data-dependent initialization (has no effect if <code>use_weight_norm</code> is FALSE).
<code>high</code>	integer, the maximum value of the input data (255 for an 8-bit image).
<code>low</code>	integer, the minimum value of the input data.
<code>dtype</code>	Data type of the Distribution.
<code>name</code>	string, the name of the Distribution.

Value

a distribution instance.

References

- Tim Salimans, Andrej Karpathy, Xi Chen, and Diederik P. Kingma. PixelCNN++: Improving the PixelCNN with Discretized Logistic Mixture Likelihood and Other Modifications. In *International Conference on Learning Representations, 2017*.
- Aaron van den Oord, Nal Kalchbrenner, Oriol Vinyals, Lasse Espeholt, Alex Graves, and Koray Kavukcuoglu. Conditional Image Generation with PixelCNN Decoders. In *Neural Information Processing Systems, 2016*.
- Aaron van den Oord, Nal Kalchbrenner, and Koray Kavukcuoglu. Pixel Recurrent Neural Networks. In *International Conference on Machine Learning, 2016*.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_poisson_log_normal_quadrature_compound()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

Description

The Plackett-Luce distribution is defined over permutations of fixed length. It is parameterized by a positive score vector of same length. This class provides methods to create indexed batches of PlackettLuce distributions. If the provided scores is rank 2 or higher, for every fixed set of leading dimensions, the last dimension represents one single PlackettLuce distribution. When calling distribution functions (e.g. `dist.log_prob(x)`), scores and `x` are broadcast to the same shape (if possible). In all cases, the last dimension of scores, `x` represents single PlackettLuce distributions.

Usage

```
tfd_plackett_luce(
    scores,
    dtype = tf.int32,
    validate_args = False,
    allow_nan_stats = False,
    name = "PlackettLuce"
)
```

Arguments

<code>scores</code>	An N-D Tensor, $N \geq 1$, representing the scores of a set of elements to be permuted. The first $N - 1$ dimensions index into a batch of independent distributions and the last dimension represents a vector of scores for the elements.
<code>dtype</code>	The type of the event samples (default: <code>int32</code>).
<code>validate_args</code>	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
<code>allow_nan_stats</code>	Logical, default <code>TRUE</code> . When <code>TRUE</code> , statistics (e.g., mean, mode, variance) use the value <code>NaN</code> to indicate the result is undefined. When <code>FALSE</code> , an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Details**Mathematical Details**

The Plackett-Luce is a distribution over permutation vectors p of length k where the permutation p is an arbitrary ordering of k indices $\{0, 1, \dots, k-1\}$.

The probability mass function (pmf) is,

$$\begin{aligned} \text{pmf}(p; s) &= \prod_i s_{\{p_i\}} / (Z - Z_i) \\ Z &= \sum_{\{j=0\}^{k-1}} s_j \\ Z_i &= \sum_{\{j=0\}^{i-1}} s_{\{p_j\}} \text{ for } i > 0 \text{ and } 0 \text{ for } i=0 \end{aligned}$$

where scores = $s = [s_0, \dots, s_{\{k-1\}}]$, $s_i \geq 0$.

Samples from Plackett-Luce distribution are generated sequentially as follows.

```

Initialize normalization `N_0 = Z`
For `i` in `{0, 1, ..., k-1}`
  1. Sample  $i$ -th element of permutation
     `p_i ~ Categorical(probs=[s_0/N_i, ..., s_{k-1}/N_i])`
  2. Update normalization
     `N_{i+1} = N_i - s_{p_i}`
  3. Mask out sampled index for subsequent rounds
     `s_{p_i} = 0`
Return p

```

Alternately, an equivalent way to sample from this distribution is to sort Gumbel perturbed log-scores (Aditya et al. 2019)

```

p = argsort(log s + g) ~ PlackettLuce(s)
g = [g_0, ..., g_{k-1}], g_i ~ Gumbel(0, 1)

```

Value

a distribution instance.

References

- Aditya Grover, Eric Wang, Aaron Zweig, Stefano Ermon. Stochastic Optimization of Sorting Networks via Continuous Relaxations. ICLR 2019.

See Also

For usage examples see e.g. [tfd_sample\(\)](#), [tfd_log_prob\(\)](#), [tfd_mean\(\)](#).

tfd_poisson

Poisson distribution

Description

The Poisson distribution is parameterized by an event rate parameter.

Usage

```

tfd_poisson(
  rate = NULL,
  log_rate = NULL,
  interpolate_nondiscrete = TRUE,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "Poisson"
)

```

Arguments

rate	Floating point tensor, the rate parameter. rate must be positive. Must specify exactly one of rate and log_rate.
log_rate	Floating point tensor, the log of the rate parameter. Must specify exactly one of rate and log_rate.
interpolate_nondiscrete	Logical. When FALSE, log_prob returns -inf (and prob returns 0) for non-integer inputs. When TRUE, log_prob evaluates the continuous function $k * \log_rate - \lgamma(k+1) - rate$, which matches the Poisson pmf at integer arguments k (note that this function is not itself a normalized probability log-density). Default value: TRUE.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details**Mathematical Details**

The probability mass function (pmf) is,

$$\text{pmf}(k; \text{lambda}, k \geq 0) = (\text{lambda}^k / k!) / Z$$

$$Z = \exp(\text{lambda}).$$

where rate = lambda and Z is the normalizing constant.

Value

a distribution instance.

See Also

For usage examples see e.g. [tfd_sample\(\)](#), [tfd_log_prob\(\)](#), [tfd_mean\(\)](#).

Other distributions: [tfd_autoregressive\(\)](#), [tfd_batch_reshape\(\)](#), [tfd_bates\(\)](#), [tfd_bernoulli\(\)](#), [tfd_beta_binomial\(\)](#), [tfd_beta\(\)](#), [tfd_binomial\(\)](#), [tfd_categorical\(\)](#), [tfd_cauchy\(\)](#), [tfd_chi2\(\)](#), [tfd_chi\(\)](#), [tfd_cholesky_lkj\(\)](#), [tfd_continuous_bernoulli\(\)](#), [tfd_deterministic\(\)](#), [tfd_dirichlet_multinomial\(\)](#), [tfd_dirichlet\(\)](#), [tfd_empirical\(\)](#), [tfd_exp_gamma\(\)](#), [tfd_exp_inverse_gamma\(\)](#), [tfd_exponential\(\)](#), [tfd_gamma_gamma\(\)](#), [tfd_gamma\(\)](#), [tfd_gaussian_process_regression_model\(\)](#), [tfd_gaussian_process\(\)](#), [tfd_generalized_normal\(\)](#), [tfd_geometric\(\)](#), [tfd_gumbel\(\)](#), [tfd_half_cauchy\(\)](#), [tfd_half_normal\(\)](#), [tfd_hidden_markov_model\(\)](#), [tfd_horseshoe\(\)](#), [tfd_independent\(\)](#), [tfd_inverse_gamma\(\)](#), [tfd_inverse_gaussian\(\)](#), [tfd_johnson_s_u\(\)](#), [tfd_joint_distribution_named_auto_batched\(\)](#), [tfd_joint_distribution_named\(\)](#), [tfd_joint_distribution_sequential_auto_batched\(\)](#), [tfd_joint_distribution_sequential\(\)](#), [tfd_kumaraswamy\(\)](#), [tfd_laplace\(\)](#), [tfd_linear_gaussian_state_space\(\)](#).

```
tfd_lkj(), tfd_log_logistic(), tfd_log_normal(), tfd_logistic(), tfd_mixture_same_family(),
tfd_mixture(), tfd_multinomial(), tfd_multivariate_normal_diag_plus_low_rank(), tfd_multivariate_normal(),
tfd_multivariate_normal_full_covariance(), tfd_multivariate_normal_linear_operator(),
tfd_multivariate_normal_tri_l(), tfd_multivariate_student_t_linear_operator(), tfd_negative_binomial(),
tfd_normal(), tfd_one_hot_categorical(), tfd_pareto(), tfd_pixel_cnn(), tfd_poisson_log_normal_quadrature_compound(),
tfd_power_spherical(), tfd_probit_bernoulli(), tfd_quantized(), tfd_relaxed_bernoulli(),
tfd_relaxed_one_hot_categorical(), tfd_sample_distribution(), tfd_sinh_arcsinh(),
tfd_skellam(), tfd_spherical_uniform(), tfd_student_t_process(), tfd_student_t(), tfd_transformed_distribution(),
tfd_triangular(), tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian(),
tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()
```

```
tfd_poisson_log_normal_quadrature_compound
```

```
PoissonLogNormalQuadratureCompound distribution
```

Description

The `PoissonLogNormalQuadratureCompound` is an approximation to a Poisson-LogNormal **compound distribution**, i.e.,

$$p(k|loc, scale) = \int_{d=0}^{\infty} \text{LogNormal}(l | loc, scale) \text{Poisson}(k | l) \text{prob}[d] \text{Poisson}(k | \lambda(\text{grid}[d])) : d=0, \dots, \text{deg}-1$$

Usage

```
tfd_poisson_log_normal_quadrature_compound(
  loc,
  scale,
  quadrature_size = 8,
  quadrature_fn = tfp$distributions$quadrature_scheme_lognormal_quantiles,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "PoissonLogNormalQuadratureCompound"
)
```

Arguments

<code>loc</code>	float-like (batch of) scalar Tensor; the location parameter of the LogNormal prior.
<code>scale</code>	float-like (batch of) scalar Tensor; the scale parameter of the LogNormal prior.
<code>quadrature_size</code>	integer scalar representing the number of quadrature points.

quadrature_fn	Function taking loc, scale, quadrature_size, validate_args and returning tuple(grid,probs) representing the LogNormal grid and corresponding normalized weight. Default value: quadrature_scheme_lognormal_quantiles.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

By default, the grid is chosen as quantiles of the LogNormal distribution parameterized by loc, scale and the prob vector is $[1. / \text{quadrature_size}] * \text{quadrature_size}$.

In the non-approximation case, a draw from the LogNormal prior represents the Poisson rate parameter. Unfortunately, the non-approximate distribution lacks an analytical probability density function (pdf). Therefore the PoissonLogNormalQuadratureCompound class implements an approximation based on **quadrature**. Note: although the PoissonLogNormalQuadratureCompound is approximately the Poisson-LogNormal compound distribution, it is itself a valid distribution. Viz., it possesses a sample, log_prob, mean, variance, etc. which are all mutually consistent.

Mathematical Details

The PoissonLogNormalQuadratureCompound approximates a Poisson-LogNormal **compound distribution**. Using variable-substitution and **numerical quadrature** (default: based on LogNormal quantiles) we can redefine the distribution to be a parameter-less convex combination of deg different Poisson samples. That is, defined over positive integers, this distribution is parameterized by a (batch of) loc and scale scalars.

The probability density function (pdf) is,

$$\text{pdf}(k \mid \text{loc}, \text{scale}, \text{deg}) = \text{sum}\{ \text{prob}[d] \text{Poisson}(k \mid \text{lambda}=\text{exp}(\text{grid}[d])) : d=0, \dots, \text{deg}-1 \}$$

Note: probs returned by (optional) quadrature_fn are presumed to be either a length-quadrature_size vector or a batch of vectors in 1-to-1 correspondence with the returned grid. (I.e., broadcasting is only partially supported.)

Value

a distribution instance.

See Also

For usage examples see e.g. [tfd_sample\(\)](#), [tfd_log_prob\(\)](#), [tfd_mean\(\)](#).

Other distributions: [tfd_autoregressive\(\)](#), [tfd_batch_reshape\(\)](#), [tfd_bates\(\)](#), [tfd_bernoulli\(\)](#), [tfd_beta_binomial\(\)](#), [tfd_beta\(\)](#), [tfd_binomial\(\)](#), [tfd_categorical\(\)](#), [tfd_cauchy\(\)](#), [tfd_chi2\(\)](#), [tfd_chi\(\)](#), [tfd_cholesky_lkj\(\)](#), [tfd_continuous_bernoulli\(\)](#), [tfd_deterministic\(\)](#), [tfd_dirichlet_multinomial\(\)](#), [tfd_dirichlet\(\)](#), [tfd_empirical\(\)](#), [tfd_exp_gamma\(\)](#), [tfd_exp_inverse_gamma\(\)](#), [tfd_exponential\(\)](#),

```
tfd_gamma_gamma(), tfd_gamma(), tfd_gaussian_process_regression_model(), tfd_gaussian_process(),
tfd_generalized_normal(), tfd_geometric(), tfd_gumbel(), tfd_half_cauchy(), tfd_half_normal(),
tfd_hidden_markov_model(), tfd_horseshoe(), tfd_independent(), tfd_inverse_gamma(),
tfd_inverse_gaussian(), tfd_johnson_s_u(), tfd_joint_distribution_named_auto_batched(),
tfd_joint_distribution_named(), tfd_joint_distribution_sequential_auto_batched(),
tfd_joint_distribution_sequential(), tfd_kumaraswamy(), tfd_laplace(), tfd_linear_gaussian_state_space(),
tfd_lkj(), tfd_log_logistic(), tfd_log_normal(), tfd_logistic(), tfd_mixture_same_family(),
tfd_mixture(), tfd_multinomial(), tfd_multivariate_normal_diag_plus_low_rank(), tfd_multivariate_normal(),
tfd_multivariate_normal_full_covariance(), tfd_multivariate_normal_linear_operator(),
tfd_multivariate_normal_tri_l(), tfd_multivariate_student_t_linear_operator(), tfd_negative_binomial(),
tfd_normal(), tfd_one_hot_categorical(), tfd_pareto(), tfd_pixel_cnn(), tfd_poisson(),
tfd_power_spherical(), tfd_probit_bernoulli(), tfd_quantized(), tfd_relaxed_bernoulli(),
tfd_relaxed_one_hot_categorical(), tfd_sample_distribution(), tfd_sinh_arcsinh(),
tfd_skellam(), tfd_spherical_uniform(), tfd_student_t_process(), tfd_student_t(), tfd_transformed_distribution(),
tfd_triangular(), tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian(),
tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()
```

tfd_power_spherical *The Power Spherical distribution over unit vectors on S^{n-1} .*

Description

The Power Spherical distribution is a distribution over vectors on the unit hypersphere S^{n-1} embedded in n dimensions (\mathbb{R}^n). It serves as an alternative to the von Mises-Fisher distribution with a simpler (faster) `log_prob` calculation, as well as a reparameterizable sampler. In contrast, the Power Spherical distribution does have `-mean_direction` as a point with zero density (and hence a neighborhood around that having arbitrarily small density), in contrast with the von Mises-Fisher distribution which has non-zero density everywhere. NOTE: `mean_direction` is not in general the mean of the distribution. For spherical distributions, the mean is generally not in the support of the distribution.

Usage

```
tfd_power_spherical(
    mean_direction,
    concentration,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = "PowerSpherical"
)
```

Arguments

`mean_direction` Floating-point Tensor with shape $[B_1, \dots, B_n, N]$. A unit vector indicating the mode of the distribution, or the unit-normalized direction of the mean.

concentration	Floating-point Tensor having batch shape [B1, ... Bn] broadcastable with mean_direction. The level of concentration of samples around the mean_direction. concentration=0 indicates a uniform distribution over the unit hypersphere, and concentration=+inf indicates a Deterministic distribution (delta function) at mean_direction.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

Mathematical details

The probability density function (pdf) is,

$$\text{pdf}(x; \mu, \text{kappa}) = C(\text{kappa}) (1 + \mu^T x)^{-k}$$

where,

$$C(\text{kappa}) = 2^{-(a+b)} \pi^b \Gamma(a) / \Gamma(a+b)$$

$$a = (n-1) / 2 + k$$

$$b = (n-1) / 2.$$

where

- mean_direction = mu; a unit vector in R^k ,
- concentration = kappa; scalar real ≥ 0 , concentration of samples around mean_direction, where 0 pertains to the uniform distribution on the hypersphere, and ∞ indicates a delta function at mean_direction.

Value

a distribution instance.

See Also

For usage examples see e.g. [tfd_sample\(\)](#), [tfd_log_prob\(\)](#), [tfd_mean\(\)](#).

Other distributions: [tfd_autoregressive\(\)](#), [tfd_batch_reshape\(\)](#), [tfd_bates\(\)](#), [tfd_bernoulli\(\)](#), [tfd_beta_binomial\(\)](#), [tfd_beta\(\)](#), [tfd_binomial\(\)](#), [tfd_categorical\(\)](#), [tfd_cauchy\(\)](#), [tfd_chi2\(\)](#), [tfd_chi\(\)](#), [tfd_cholesky_lkj\(\)](#), [tfd_continuous_bernoulli\(\)](#), [tfd_deterministic\(\)](#), [tfd_dirichlet_multinomial\(\)](#), [tfd_dirichlet\(\)](#), [tfd_empirical\(\)](#), [tfd_exp_gamma\(\)](#), [tfd_exp_inverse_gamma\(\)](#), [tfd_exponential\(\)](#), [tfd_gamma_gamma\(\)](#), [tfd_gamma\(\)](#), [tfd_gaussian_process_regression_model\(\)](#), [tfd_gaussian_process\(\)](#), [tfd_generalized_normal\(\)](#), [tfd_geometric\(\)](#), [tfd_gumbel\(\)](#), [tfd_half_cauchy\(\)](#), [tfd_half_normal\(\)](#), [tfd_hidden_markov_model\(\)](#), [tfd_horseshoe\(\)](#), [tfd_independent\(\)](#), [tfd_inverse_gamma\(\)](#), [tfd_inverse_gaussian\(\)](#), [tfd_johnson_s_u\(\)](#), [tfd_joint_distribution_named_auto_batched\(\)](#), [tfd_joint_distribution_named\(\)](#), [tfd_joint_distribution_sequential_auto_batched\(\)](#), [tfd_joint_distribution_sequential\(\)](#), [tfd_kumaraswamy\(\)](#), [tfd_laplace\(\)](#), [tfd_linear_gaussian_state_space\(\)](#).

tfd_lkj(), tfd_log_logistic(), tfd_log_normal(), tfd_logistic(), tfd_mixture_same_family(),
 tfd_mixture(), tfd_multinomial(), tfd_multivariate_normal_diag_plus_low_rank(), tfd_multivariate_normal(),
 tfd_multivariate_normal_full_covariance(), tfd_multivariate_normal_linear_operator(),
 tfd_multivariate_normal_tri_l(), tfd_multivariate_student_t_linear_operator(), tfd_negative_binomial(),
 tfd_normal(), tfd_one_hot_categorical(), tfd_pareto(), tfd_pixel_cnn(), tfd_poisson_log_normal_quadrature(),
 tfd_poisson(), tfd_probit_bernoulli(), tfd_quantized(), tfd_relaxed_bernoulli(), tfd_relaxed_one_hot_categorical(),
 tfd_sample_distribution(), tfd_sinh_arcsinh(), tfd_skellam(), tfd_spherical_uniform(),
 tfd_student_t_process(), tfd_student_t(), tfd_transformed_distribution(), tfd_triangular(),
 tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian_process(),
 tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
 tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
 tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
 tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()

tfd_prob

Probability density/mass function.

Description

Probability density/mass function.

Usage

```
tfd_prob(distribution, value, ...)
```

Arguments

distribution	The distribution being used.
value	float or double Tensor.
...	Additional parameters passed to Python.

Value

a Tensor of shape `sample_shape(x) + self$batch_shape` with values of type `self$dtype`.

See Also

Other `distribution_methods`: [tfd_cdf\(\)](#), [tfd_covariance\(\)](#), [tfd_cross_entropy\(\)](#), [tfd_entropy\(\)](#),
[tfd_kl_divergence\(\)](#), [tfd_log_cdf\(\)](#), [tfd_log_prob\(\)](#), [tfd_log_survival_function\(\)](#), [tfd_mean\(\)](#),
[tfd_mode\(\)](#), [tfd_quantile\(\)](#), [tfd_sample\(\)](#), [tfd_stddev\(\)](#), [tfd_survival_function\(\)](#), [tfd_variance\(\)](#)

Examples

```
d <- tfd_normal(loc = c(1, 2), scale = c(1, 0.5))
x <- d %>% tfd_sample()
d %>% tfd_prob(x)
```

tfd_probit_bernoulli *ProbitBernoulli distribution.*

Description

The ProbitBernoulli distribution with probs parameter, i.e., the probability of a 1 outcome (vs a 0 outcome). Unlike a regular Bernoulli distribution, which uses the logistic (aka 'sigmoid') function to go from the un-constrained parameters to probabilities, this distribution uses the CDF of the **standard normal distribution**:

$$p(x=1; probits) = 0.5 * (1 + \text{erf}(probits / \text{sqrt}(2)))$$

$$p(x=0; probits) = 1 - p(x=1; probits)$$

Where erf is the **error function**. A typical application of this distribution is in **probit regression**.

Usage

```
tfd_probit_bernoulli(
    probits = NULL,
    probs = NULL,
    dtype = tf$int32,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = "ProbitBernoulli"
)
```

Arguments

probits	An N-D Tensor representing the probit-odds of a 1 event. Each entry in the Tensor parameterizes an independent ProbitBernoulli distribution where the probability of an event is <code>normal_cdf(probits)</code> . Only one of <code>probits</code> or <code>probs</code> should be passed in.
probs	An N-D Tensor representing the probability of a 1 event. Each entry in the Tensor parameterizes an independent ProbitBernoulli distribution. Only one of <code>probits</code> or <code>probs</code> should be passed in.
dtype	The type of the event samples. Default: <code>int32</code> .
validate_args	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
allow_nan_stats	Logical, default <code>TRUE</code> . When <code>TRUE</code> , statistics (e.g., mean, mode, variance) use the value <code>NaN</code> to indicate the result is undefined. When <code>FALSE</code> , an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian_process()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_quantile

Quantile function. Aka "inverse cdf" or "percent point function".

Description

Given random variable X and p in $[0, 1]$, the quantile is: `tfd_quantile(p) := x` such that $P[X \leq x] = p$

Usage

```
tfd_quantile(distribution, value, ...)
```

Arguments

<code>distribution</code>	The distribution being used.
<code>value</code>	float or double Tensor.
<code>...</code>	Additional parameters passed to Python.

Value

a Tensor of shape `sample_shape(x) + self$batch_shape` with values of type `self$dtype`.

See Also

Other `distribution_methods`: `tfd_cdf()`, `tfd_covariance()`, `tfd_cross_entropy()`, `tfd_entropy()`, `tfd_kl_divergence()`, `tfd_log_cdf()`, `tfd_log_prob()`, `tfd_log_survival_function()`, `tfd_mean()`, `tfd_mode()`, `tfd_prob()`, `tfd_sample()`, `tfd_stddev()`, `tfd_survival_function()`, `tfd_variance()`

Examples

```
d <- tfd_normal(loc = c(1, 2), scale = c(1, 0.5))
d %>% tfd_quantile(0.5)
```

tfd_quantized

Distribution representing the quantization $Y = \text{ceiling}(X)$

Description

Definition in Terms of Sampling

Usage

```
tfd_quantized(
  distribution,
  low = NULL,
  high = NULL,
  validate_args = FALSE,
  name = "QuantizedDistribution"
)
```

Arguments

<code>distribution</code>	The base distribution class to transform. Typically an instance of <code>Distribution</code> .
<code>low</code>	Tensor with same <code>dtype</code> as this distribution and shape able to be added to samples. Should be a whole number. Default <code>NULL</code> . If provided, base distribution's <code>prob</code> should be defined at <code>low</code> .
<code>high</code>	Tensor with same <code>dtype</code> as this distribution and shape able to be added to samples. Should be a whole number. Default <code>NULL</code> . If provided, base distribution's <code>prob</code> should be defined at <code>high - 1</code> . <code>high</code> must be strictly greater than <code>low</code> .
<code>validate_args</code>	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
<code>name</code>	name prefixed to Ops created by this class.

Details

1. Draw X
2. Set $Y \leftarrow \text{ceiling}(X)$
3. If $Y < \text{low}$, reset $Y \leftarrow \text{low}$
4. If $Y > \text{high}$, reset $Y \leftarrow \text{high}$
5. Return Y

Definition in Terms of the Probability Mass Function

Given scalar random variable X , we define a discrete random variable Y supported on the integers as follows:

$$\begin{aligned}
 P[Y = j] &:= P[X \leq \text{low}], & \text{if } j == \text{low}, \\
 &:= P[X > \text{high} - 1], & j == \text{high}, \\
 &:= 0, & \text{if } j < \text{low} \text{ or } j > \text{high}, \\
 &:= P[j - 1 < X \leq j], & \text{all other } j.
 \end{aligned}$$

Conceptually, without cutoffs, the quantization process partitions the real line \mathbb{R} into half open intervals, and identifies an integer j with the right endpoints:

$$\begin{array}{cccccccc}
 \mathbb{R} = \dots & (-2, -1] & (-1, 0] & (0, 1] & (1, 2] & (2, 3] & (3, 4] & \dots \\
 j = \dots & -1 & 0 & 1 & 2 & 3 & 4 & \dots
 \end{array}$$

$P[Y = j]$ is the mass of X within the j th interval. If $\text{low} = 0$, and $\text{high} = 2$, then the intervals are redrawn and j is re-assigned:

$$\begin{array}{cccc}
 \mathbb{R} = (-\text{infy}, 0] & (0, 1] & (1, \text{infy}) \\
 j = & 0 & 1 & 2
 \end{array}$$

$P[Y = j]$ is still the mass of X within the j th interval.

@section References:

- [Tim Salimans, Andrej Karpathy, Xi Chen, and Diederik P. Kingma. PixelCNN++: Improving the PixelCNN with discretized logistic mixture likelihood and other modifications. International Conference on Learning Representations_, 2017.](#)
- [Aaron van den Oord et al. Parallel WaveNet: Fast High-Fidelity Speech Synthesis. arXiv preprint arXiv:1711.10433, 2017.](#)

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`,

```
tfd_gamma_gamma(), tfd_gamma(), tfd_gaussian_process_regression_model(), tfd_gaussian_process(),
tfd_generalized_normal(), tfd_geometric(), tfd_gumbel(), tfd_half_cauchy(), tfd_half_normal(),
tfd_hidden_markov_model(), tfd_horseshoe(), tfd_independent(), tfd_inverse_gamma(),
tfd_inverse_gaussian(), tfd_johnson_s_u(), tfd_joint_distribution_named_auto_batched(),
tfd_joint_distribution_named(), tfd_joint_distribution_sequential_auto_batched(),
tfd_joint_distribution_sequential(), tfd_kumaraswamy(), tfd_laplace(), tfd_linear_gaussian_state_space,
tfd_lkj(), tfd_log_logistic(), tfd_log_normal(), tfd_logistic(), tfd_mixture_same_family(),
tfd_mixture(), tfd_multinomial(), tfd_multivariate_normal_diag_plus_low_rank(), tfd_multivariate_normal,
tfd_multivariate_normal_full_covariance(), tfd_multivariate_normal_linear_operator(),
tfd_multivariate_normal_tri_l(), tfd_multivariate_student_t_linear_operator(), tfd_negative_binomial(),
tfd_normal(), tfd_one_hot_categorical(), tfd_pareto(), tfd_pixel_cnn(), tfd_poisson_log_normal_quadrature,
tfd_poisson(), tfd_power_spherical(), tfd_probit_bernoulli(), tfd_relaxed_bernoulli(),
tfd_relaxed_one_hot_categorical(), tfd_sample_distribution(), tfd_sinh_arcsinh(),
tfd_skellam(), tfd_spherical_uniform(), tfd_student_t_process(), tfd_student_t(), tfd_transformed_distribution,
tfd_triangular(), tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian,
tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()
```

tfd_relaxed_bernoulli *RelaxedBernoulli distribution with temperature and logits parameters*

Description

The RelaxedBernoulli is a distribution over the unit interval (0,1), which continuously approximates a Bernoulli. The degree of approximation is controlled by a temperature: as the temperature goes to 0 the RelaxedBernoulli becomes discrete with a distribution described by the logits or probs parameters, as the temperature goes to infinity the RelaxedBernoulli becomes the constant distribution that is identically 0.5.

Usage

```
tfd_relaxed_bernoulli(
    temperature,
    logits = NULL,
    probs = NULL,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = "RelaxedBernoulli"
)
```

Arguments

`temperature` An 0-D Tensor, representing the temperature of a set of RelaxedBernoulli distributions. The temperature should be positive.

logits	An N-D Tensor representing the log-odds of a positive event. Each entry in the Tensor parametrizes an independent RelaxedBernoulli distribution where the probability of an event is <code>sigmoid(logits)</code> . Only one of logits or probs should be passed in.
probs	An N-D Tensor representing the probability of a positive event. Each entry in the Tensor parameterizes an independent Bernoulli distribution. Only one of logits or probs should be passed in.
validate_args	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
allow_nan_stats	Logical, default <code>TRUE</code> . When <code>TRUE</code> , statistics (e.g., mean, mode, variance) use the value <code>NaN</code> to indicate the result is undefined. When <code>FALSE</code> , an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

The RelaxedBernoulli distribution is a reparameterized continuous distribution that is the binary special case of the RelaxedOneHotCategorical distribution (Maddison et al., 2016; Jang et al., 2016). For details on the binary special case see the appendix of Maddison et al. (2016) where it is referred to as BinConcrete. If you use this distribution, please cite both papers.

Some care needs to be taken for loss functions that depend on the log-probability of RelaxedBernoullis, because computing log-probabilities of the RelaxedBernoulli can suffer from underflow issues. In many case loss functions such as these are invariant under invertible transformations of the random variables. The KL divergence, found in the variational autoencoder loss, is an example. Because RelaxedBernoullis are sampled by a Logistic random variable followed by a `tf.sigmoid` op, one solution is to treat the Logistic as the random variable and `tf.sigmoid` as downstream. The KL divergences of two Logistics, which are always followed by a `tf.sigmoid` op, is equivalent to evaluating KL divergences of RelaxedBernoulli samples. See Maddison et al., 2016 for more details where this distribution is called the BinConcrete. An alternative approach is to evaluate Bernoulli log probability or KL directly on relaxed samples, as done in Jang et al., 2016. In this case, guarantees on the loss are usually violated. For instance, using a Bernoulli KL in a relaxed ELBO is no longer a lower bound on the log marginal probability of the observation. Thus care and early stopping are important.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`,

```
tfd_gamma_gamma(), tfd_gamma(), tfd_gaussian_process_regression_model(), tfd_gaussian_process(),
tfd_generalized_normal(), tfd_geometric(), tfd_gumbel(), tfd_half_cauchy(), tfd_half_normal(),
tfd_hidden_markov_model(), tfd_horseshoe(), tfd_independent(), tfd_inverse_gamma(),
tfd_inverse_gaussian(), tfd_johnson_s_u(), tfd_joint_distribution_named_auto_batched(),
tfd_joint_distribution_named(), tfd_joint_distribution_sequential_auto_batched(),
tfd_joint_distribution_sequential(), tfd_kumaraswamy(), tfd_laplace(), tfd_linear_gaussian_state_space,
tfd_lkj(), tfd_log_logistic(), tfd_log_normal(), tfd_logistic(), tfd_mixture_same_family(),
tfd_mixture(), tfd_multinomial(), tfd_multivariate_normal_diag_plus_low_rank(), tfd_multivariate_normal,
tfd_multivariate_normal_full_covariance(), tfd_multivariate_normal_linear_operator(),
tfd_multivariate_normal_tri_l(), tfd_multivariate_student_t_linear_operator(), tfd_negative_binomial(),
tfd_normal(), tfd_one_hot_categorical(), tfd_pareto(), tfd_pixel_cnn(), tfd_poisson_log_normal_quadrature,
tfd_poisson(), tfd_power_spherical(), tfd_probit_bernoulli(), tfd_quantized(), tfd_relaxed_one_hot_categ,
tfd_sample_distribution(), tfd_sinh_arcsinh(), tfd_skellam(), tfd_spherical_uniform(),
tfd_student_t_process(), tfd_student_t(), tfd_transformed_distribution(), tfd_triangular(),
tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian_process(),
tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()
```

tfd_relaxed_one_hot_categorical

RelaxedOneHotCategorical distribution with temperature and logits

Description

The `RelaxedOneHotCategorical` is a distribution over random probability vectors, vectors of positive real values that sum to one, which continuously approximates a `OneHotCategorical`. The degree of approximation is controlled by a temperature: as the temperature goes to 0 the `RelaxedOneHotCategorical` becomes discrete with a distribution described by the logits or probs parameters, as the temperature goes to infinity the `RelaxedOneHotCategorical` becomes the constant distribution that is identically the constant vector of $(1/\text{event_size}, \dots, 1/\text{event_size})$. The `RelaxedOneHotCategorical` distribution was concurrently introduced as the Gumbel-Softmax (Jang et al., 2016) and Concrete (Maddison et al., 2016) distributions for use as a reparameterized continuous approximation to the `Categorical` one-hot distribution. If you use this distribution, please cite both papers.

Usage

```
tfd_relaxed_one_hot_categorical(
  temperature,
  logits = NULL,
  probs = NULL,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "RelaxedOneHotCategorical"
)
```


Arguments

temperature	An 0-D Tensor, representing the temperature of a set of RelaxedOneHotCategorical distributions. The temperature should be positive.
logits	An N-D Tensor, $N \geq 1$, representing the log probabilities of a set of RelaxedOneHotCategorical distributions. The first $N - 1$ dimensions index into a batch of independent distributions and the last dimension represents a vector of logits for each class. Only one of logits or probs should be passed in.
probs	An N-D Tensor, $N \geq 1$, representing the probabilities of a set of RelaxedOneHotCategorical distributions. The first $N - 1$ dimensions index into a batch of independent distributions and the last dimension represents a vector of probabilities for each class. Only one of logits or probs should be passed in.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Value

a distribution instance.

References

- Eric Jang, Shixiang Gu, and Ben Poole. Categorical Reparameterization with Gumbel-Softmax. 2016.
- Chris J. Maddison, Andriy Mnih, and Yee Whye Teh. The Concrete Distribution: A Continuous Relaxation of Discrete Random Variables. 2016.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`,

```
tfd_multivariate_normal_tri_l(), tfd_multivariate_student_t_linear_operator(), tfd_negative_binomial(),
tfd_normal(), tfd_one_hot_categorical(), tfd_pareto(), tfd_pixel_cnn(), tfd_poisson_log_normal_quadrature(),
tfd_poisson(), tfd_power_spherical(), tfd_probit_bernoulli(), tfd_quantized(), tfd_relaxed_bernoulli(),
tfd_sample_distribution(), tfd_sinh_arcsinh(), tfd_skellam(), tfd_spherical_uniform(),
tfd_student_t_process(), tfd_student_t(), tfd_transformed_distribution(), tfd_triangular(),
tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian_process(),
tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()
```

tfd_sample

Generate samples of the specified shape.

Description

Note that a call to `tfd_sample()` without arguments will generate a single sample.

Usage

```
tfd_sample(distribution, sample_shape = list(), ...)
```

Arguments

<code>distribution</code>	The distribution being used.
<code>sample_shape</code>	0D or 1D int32 Tensor. Shape of the generated samples.
<code>...</code>	Additional parameters passed to Python.

Value

a Tensor with prepended dimensions `sample_shape`.

See Also

Other `distribution_methods`: [tfd_cdf\(\)](#), [tfd_covariance\(\)](#), [tfd_cross_entropy\(\)](#), [tfd_entropy\(\)](#), [tfd_kl_divergence\(\)](#), [tfd_log_cdf\(\)](#), [tfd_log_prob\(\)](#), [tfd_log_survival_function\(\)](#), [tfd_mean\(\)](#), [tfd_mode\(\)](#), [tfd_prob\(\)](#), [tfd_quantile\(\)](#), [tfd_stddev\(\)](#), [tfd_survival_function\(\)](#), [tfd_variance\(\)](#)

Examples

```
d <- tfd_normal(loc = c(1, 2), scale = c(1, 0.5))
d %>% tfd_sample()
```

tfd_sample_distribution

Sample distribution via independent draws.

Description

This distribution is useful for reducing over a collection of independent, identical draws. It is otherwise identical to the input distribution.

Usage

```
tfd_sample_distribution(
  distribution,
  sample_shape = list(),
  validate_args = FALSE,
  name = NULL
)
```

Arguments

distribution	The base distribution instance to transform. Typically an instance of <code>Distribution</code> .
sample_shape	integer scalar or vector <code>Tensor</code> representing the shape of a single sample.
validate_args	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
name	The name for ops managed by the distribution. Default value: <code>NULL</code> (i.e., <code>'Sample' + distribution\$name</code>).

Details

Mathematical Details The probability function is,

$$p(x) = \prod\{ p(x[i]) : i = 0, \dots, (n - 1) \}$$

Value

a distribution instance.

See Also

For usage examples see e.g. [tfd_sample\(\)](#), [tfd_log_prob\(\)](#), [tfd_mean\(\)](#).

Other distributions: [tfd_autoregressive\(\)](#), [tfd_batch_reshape\(\)](#), [tfd_bates\(\)](#), [tfd_bernoulli\(\)](#), [tfd_beta_binomial\(\)](#), [tfd_beta\(\)](#), [tfd_binomial\(\)](#), [tfd_categorical\(\)](#), [tfd_cauchy\(\)](#), [tfd_chi2\(\)](#), [tfd_chi\(\)](#), [tfd_cholesky_lkj\(\)](#), [tfd_continuous_bernoulli\(\)](#), [tfd_deterministic\(\)](#), [tfd_dirichlet_multinomial\(\)](#), [tfd_dirichlet\(\)](#), [tfd_empirical\(\)](#), [tfd_exp_gamma\(\)](#), [tfd_exp_inverse_gamma\(\)](#), [tfd_exponential\(\)](#), [tfd_gamma_gamma\(\)](#), [tfd_gamma\(\)](#), [tfd_gaussian_process_regression_model\(\)](#), [tfd_gaussian_process\(\)](#),

```
tfd_generalized_normal(), tfd_geometric(), tfd_gumbel(), tfd_half_cauchy(), tfd_half_normal(),
tfd_hidden_markov_model(), tfd_horseshoe(), tfd_independent(), tfd_inverse_gamma(),
tfd_inverse_gaussian(), tfd_johnson_s_u(), tfd_joint_distribution_named_auto_batched(),
tfd_joint_distribution_named(), tfd_joint_distribution_sequential_auto_batched(),
tfd_joint_distribution_sequential(), tfd_kumaraswamy(), tfd_laplace(), tfd_linear_gaussian_state_space,
tfd_lkj(), tfd_log_logistic(), tfd_log_normal(), tfd_logistic(), tfd_mixture_same_family(),
tfd_mixture(), tfd_multinomial(), tfd_multivariate_normal_diag_plus_low_rank(), tfd_multivariate_normal,
tfd_multivariate_normal_full_covariance(), tfd_multivariate_normal_linear_operator(),
tfd_multivariate_normal_tri_l(), tfd_multivariate_student_t_linear_operator(), tfd_negative_binomial(),
tfd_normal(), tfd_one_hot_categorical(), tfd_pareto(), tfd_pixel_cnn(), tfd_poisson_log_normal_quadrature,
tfd_poisson(), tfd_power_spherical(), tfd_probit_bernoulli(), tfd_quantized(), tfd_relaxed_bernoulli(),
tfd_relaxed_one_hot_categorical(), tfd_sinh_arcsinh(), tfd_skellam(), tfd_spherical_uniform(),
tfd_student_t_process(), tfd_student_t(), tfd_transformed_distribution(), tfd_triangular(),
tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian_process(),
tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()
```

tfd_sinh_arcsinh	<i>The SinhArcsinh transformation of a distribution on (-inf, inf)</i>
------------------	--

Description

This distribution models a random variable, making use of a SinhArcsinh transformation (which has adjustable tailweight and skew), a rescaling, and a shift. The SinhArcsinh transformation of the Normal is described in great depth in [Sinh-arcsinh distributions](#). Here we use a slightly different parameterization, in terms of tailweight and skewness. Additionally we allow for distributions other than Normal, and control over scale as well as a "shift" parameter loc.

Usage

```
tfd_sinh_arcsinh(
    loc,
    scale,
    skewness = NULL,
    tailweight = NULL,
    distribution = NULL,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = "SinhArcsinh"
)
```

Arguments

loc	Floating-point Tensor.
scale	Tensor of same dtype as loc.

skewness	Skewness parameter. Default is 0.0 (no skew).
tailweight	Tailweight parameter. Default is 1.0 (unchanged tailweight)
distribution	<code>tf\$Distributions\$Distribution</code> -like instance. Distribution that is transformed to produce this distribution. Default is <code>tfd_normal(0,1)</code> . Must be a scalar-batch, scalar-event distribution. Typically <code>distribution\$reparameterization_type = FULLY_REPARAMETERIZED</code> or it is a function of non-trainable parameters. WARNING: If you backprop through a <code>SinhArcsinh</code> sample and <code>distribution</code> is not <code>FULLY_REPARAMETERIZED</code> yet is a function of trainable variables, then the gradient will be incorrect!
validate_args	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
allow_nan_stats	Logical, default <code>TRUE</code> . When <code>TRUE</code> , statistics (e.g., mean, mode, variance) use the value <code>NaN</code> to indicate the result is undefined. When <code>FALSE</code> , an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

Mathematical Details

Given random variable Z , we define the `SinhArcsinh` transformation of Z , Y , parameterized by $(\text{loc}, \text{scale}, \text{skewness}, \text{tailweight})$, via the relation:

$$\begin{aligned}
 Y &:= \text{loc} + \text{scale} * F(Z) * (2 / F_0(2)) \\
 F(Z) &:= \text{Sinh}(\text{Arcsinh}(Z) + \text{skewness} * \text{tailweight}) \\
 F_0(Z) &:= \text{Sinh}(\text{Arcsinh}(Z) * \text{tailweight})
 \end{aligned}$$

This distribution is similar to the location-scale transformation $L(Z) := \text{loc} + \text{scale} * Z$ in the following ways:

- If $\text{skewness} = 0$ and $\text{tailweight} = 1$ (the defaults), $F(Z) = Z$, and then $Y = L(Z)$ exactly.
- loc is used in both to shift the result by a constant factor.
- The multiplication of scale by $2 / F_0(2)$ ensures that if $\text{skewness} = 0$ $P[Y - \text{loc} \leq 2 * \text{scale}] = P[L(Z) - \text{loc} \leq 2 * \text{scale}]$. Thus it can be said that the weights in the tails of Y and $L(Z)$ beyond $\text{loc} + 2 * \text{scale}$ are the same.

This distribution is different than $\text{loc} + \text{scale} * Z$ due to the reshaping done by F :

- Positive (negative) skewness leads to positive (negative) skew.
- positive skew means, the mode of $F(Z)$ is "tilted" to the right.
- positive skew means positive values of $F(Z)$ become more likely, and negative values become less likely.
- Larger (smaller) tailweight leads to fatter (thinner) tails.
- Fatter tails mean larger values of $|F(Z)|$ become more likely.

- `tailweight < 1` leads to a distribution that is "flat" around $Y = \text{loc}$, and a very steep drop-off in the tails.
- `tailweight > 1` leads to a distribution more peaked at the mode with heavier tails.

To see the argument about the tails, note that for $|Z| \gg 1$ and $|Z| \gg (|\text{skewness}| * \text{tailweight})^{**\text{tailweight}}$, we have $Y \approx 0.5 Z^{**\text{tailweight}} e^{*(\text{sign}(Z) \text{skewness} * \text{tailweight})}$.

To see the argument regarding multiplying scale by $2 / F_{-0}(2)$,

$$\begin{aligned} P[(Y - \text{loc}) / \text{scale} \leq 2] &= P[F(Z) * (2 / F_{-0}(2)) \leq 2] \\ &= P[F(Z) \leq F_{-0}(2)] \\ &= P[Z \leq 2] \quad (\text{if } F = F_{-0}). \end{aligned}$$

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian_process()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_skellam	<i>Skellam distribution.</i>
-------------	------------------------------

Description

The Skellam distribution is parameterized by two rate parameters, `rate1` and `rate2`. Its samples are defined as:

```
x ~ Poisson(rate1)
y ~ Poisson(rate2)
z = x - y
z ~ Skellam(rate1, rate2)
```

where the samples `x` and `y` are assumed to be independent.

Usage

```
tfd_skellam(
  rate1 = NULL,
  rate2 = NULL,
  log_rate1 = NULL,
  log_rate2 = NULL,
  force_probs_to_zero_outside_support = FALSE,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "Skellam"
)
```

Arguments

<code>rate1</code>	Floating point tensor, the first rate parameter. <code>rate1</code> must be positive. Must specify exactly one of <code>rate1</code> and <code>log_rate1</code>
<code>rate2</code>	Floating point tensor, the second rate parameter. <code>rate</code> must be positive. Must specify exactly one of <code>rate2</code> and <code>log_rate2</code> .
<code>log_rate1</code>	Floating point tensor, the log of the first rate parameter. Must specify exactly one of <code>rate1</code> and <code>log_rate1</code> .
<code>log_rate2</code>	Floating point tensor, the log of the second rate parameter. Must specify exactly one of <code>rate2</code> and <code>log_rate2</code> .
<code>force_probs_to_zero_outside_support</code>	logical. When <code>TRUE</code> , <code>log_prob</code> returns <code>-inf</code> (and <code>prob</code> returns <code>0</code>) for non-integer inputs. When <code>FALSE</code> , <code>log_prob</code> evaluates the Skellam pmf as a continuous function (note that this function is not itself a normalized probability log-density). Default value: <code>FALSE</code> .
<code>validate_args</code>	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .

<code>allow_nan_stats</code>	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Details

Mathematical Details The probability mass function (pmf) is,

$$\text{pmf}(k; l1, l2) = (l1 / l2) ** (k / 2) * I_k(2 * \text{sqrt}(l1 * l2)) / Z$$

$$Z = \exp(l1 + l2).$$

where $rate1 = l1$, $rate2 = l2$, Z is the normalizing constant and I_k is the modified bessel function of the first kind.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag_plus_low_rank_linear_operator()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_spherical_uniform *The uniform distribution over unit vectors on S^{n-1} .*

Description

The uniform distribution on the unit hypersphere S^{n-1} embedded in n dimensions (R^n).

Usage

```
tfd_spherical_uniform(
    dimension,
    batch_shape = list(),
    dtype = tf$float32,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = "SphericalUniform"
)
```

Arguments

dimension	integer. The dimension of the embedded space where the sphere resides.
batch_shape	Positive integer-like vector-shaped Tensor representing the new shape of the batch dimensions. Default value: [].
dtype	dtype of the generated samples.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

Mathematical details

The probability density function (pdf) is,

$$\text{pdf}(x; n) = 1. / A(n)$$

where,

$$A(n) = 2 * \pi^{n / 2} / \text{Gamma}(n / 2),$$

Gamma being the Gamma function.

where $n = \text{dimension}$; corresponds to S^{n-1} embedded in R^n .

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_stddev

Standard deviation.

Description

Standard deviation is defined as, $\text{stddev} = \sqrt{E[(X - E[X])^2]}$ where X is the random variable associated with this distribution, E denotes expectation, and $\text{Var}[\text{shape}] = \text{batch_shape} + \text{event_shape}$.

Usage

```
tfd_stddev(distribution, ...)
```

Arguments

`distribution` The distribution being used.
`...` Additional parameters passed to Python.

Value

a Tensor of shape `sample_shape(x) + self$batch_shape` with values of type `self$dtype`.

See Also

Other `distribution_methods`: [tfd_cdf\(\)](#), [tfd_covariance\(\)](#), [tfd_cross_entropy\(\)](#), [tfd_entropy\(\)](#), [tfd_kl_divergence\(\)](#), [tfd_log_cdf\(\)](#), [tfd_log_prob\(\)](#), [tfd_log_survival_function\(\)](#), [tfd_mean\(\)](#), [tfd_mode\(\)](#), [tfd_prob\(\)](#), [tfd_quantile\(\)](#), [tfd_sample\(\)](#), [tfd_survival_function\(\)](#), [tfd_variance\(\)](#)

Examples

```
d <- tfd_normal(loc = c(1, 2), scale = c(1, 0.5))
d %>% tfd_stddev()
```

<code>tfd_student_t</code>	<i>Student's t-distribution</i>
----------------------------	---------------------------------

Description

This distribution has parameters: degree of freedom `df`, location `loc`, and `scale`.

Usage

```
tfd_student_t(
  df,
  loc,
  scale,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "StudentT"
)
```

Arguments

<code>df</code>	Floating-point Tensor. The degrees of freedom of the distribution(s). <code>df</code> must contain only positive values.
<code>loc</code>	Floating-point Tensor. The mean(s) of the distribution(s).
<code>scale</code>	Floating-point Tensor. The scaling factor(s) for the distribution(s). Note that <code>scale</code> is not technically the standard deviation of this distribution but has semantics more similar to standard deviation than variance.
<code>validate_args</code>	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .

allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

Mathematical details

The probability density function (pdf) is,

$$\text{pdf}(x; \text{df}, \text{mu}, \text{sigma}) = (1 + y^{*2} / \text{df})^{*(-0.5 (\text{df} + 1))} / Z$$

where,

$$y = (x - \text{mu}) / \text{sigma}$$

$$Z = \text{abs}(\text{sigma}) \sqrt{\text{df} \pi} \text{Gamma}(0.5 \text{df}) / \text{Gamma}(0.5 (\text{df} + 1))$$

where:

- loc = mu,
- scale = sigma, and,
- Z is the normalization constant, and,
- Gamma is the [gamma function](#). The StudentT distribution is a member of the [location-scale family](#), i.e., it can be constructed as,

$$X \sim \text{StudentT}(\text{df}, \text{loc}=0, \text{scale}=1)$$

$$Y = \text{loc} + \text{scale} * X$$

Notice that scale has semantics more similar to standard deviation than variance. However it is not actually the std. deviation; the Student's t-distribution std. dev. is $\text{scale} \sqrt{\text{df} / (\text{df} - 2)}$ when $\text{df} > 2$.

Samples of this distribution are reparameterized (pathwise differentiable). The derivatives are computed using the approach described in the paper [Michael Figurnov, Shakir Mohamed, Andriy Mnih. Implicit Reparameterization Gradients, 2018](#)

Value

a distribution instance.

See Also

For usage examples see e.g. [tfd_sample\(\)](#), [tfd_log_prob\(\)](#), [tfd_mean\(\)](#).

Other distributions: [tfd_autoregressive\(\)](#), [tfd_batch_reshape\(\)](#), [tfd_bates\(\)](#), [tfd_bernoulli\(\)](#), [tfd_beta_binomial\(\)](#), [tfd_beta\(\)](#), [tfd_binomial\(\)](#), [tfd_categorical\(\)](#), [tfd_cauchy\(\)](#), [tfd_chi2\(\)](#), [tfd_chi\(\)](#), [tfd_cholesky_lkj\(\)](#), [tfd_continuous_bernoulli\(\)](#), [tfd_deterministic\(\)](#), [tfd_dirichlet_multinomial\(\)](#), [tfd_dirichlet\(\)](#), [tfd_empirical\(\)](#), [tfd_exp_gamma\(\)](#), [tfd_exp_inverse_gamma\(\)](#), [tfd_exponential\(\)](#), [tfd_gamma_gamma\(\)](#), [tfd_gamma\(\)](#), [tfd_gaussian_process_regression_model\(\)](#), [tfd_gaussian_process\(\)](#), [tfd_generalized_normal\(\)](#), [tfd_geometric\(\)](#), [tfd_gumbel\(\)](#), [tfd_half_cauchy\(\)](#), [tfd_half_normal\(\)](#),

```
tfd_hidden_markov_model(), tfd_horseshoe(), tfd_independent(), tfd_inverse_gamma(),
tfd_inverse_gaussian(), tfd_johnson_s_u(), tfd_joint_distribution_named_auto_batched(),
tfd_joint_distribution_named(), tfd_joint_distribution_sequential_auto_batched(),
tfd_joint_distribution_sequential(), tfd_kumaraswamy(), tfd_laplace(), tfd_linear_gaussian_state_space(),
tfd_lkj(), tfd_log_logistic(), tfd_log_normal(), tfd_logistic(), tfd_mixture_same_family(),
tfd_mixture(), tfd_multinomial(), tfd_multivariate_normal_diag_plus_low_rank(), tfd_multivariate_normal_diag_plus_low_rank(),
tfd_multivariate_normal_full_covariance(), tfd_multivariate_normal_linear_operator(),
tfd_multivariate_normal_tri_l(), tfd_multivariate_student_t_linear_operator(), tfd_negative_binomial(),
tfd_normal(), tfd_one_hot_categorical(), tfd_pareto(), tfd_pixel_cnn(), tfd_poisson_log_normal_quadrature(),
tfd_poisson(), tfd_power_spherical(), tfd_probit_bernoulli(), tfd_quantized(), tfd_relaxed_bernoulli(),
tfd_relaxed_one_hot_categorical(), tfd_sample_distribution(), tfd_sinh_arcsinh(),
tfd_skellam(), tfd_spherical_uniform(), tfd_student_t_process(), tfd_transformed_distribution(),
tfd_triangular(), tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian(),
tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()
```

tfd_student_t_process *Marginal distribution of a Student's T process at finitely many points*

Description

A Student's T process (TP) is an indexed collection of random variables, any finite collection of which are jointly Multivariate Student's T. While this definition applies to finite index sets, it is typically implicit that the index set is infinite; in applications, it is often some finite dimensional real or complex vector space. In such cases, the TP may be thought of as a distribution over (real- or complex-valued) functions defined over the index set.

Usage

```
tfd_student_t_process(
  df,
  kernel,
  index_points,
  mean_fn = NULL,
  jitter = 1e-06,
  validate_args = FALSE,
  allow_nan_stats = FALSE,
  name = "StudentTProcess"
)
```

Arguments

df Positive Floating-point Tensor representing the degrees of freedom. Must be greater than 2.

kernel	PositiveSemidefiniteKernel-like instance representing the TP's covariance function.
index_points	float Tensor representing finite (batch of) vector(s) of points in the index set over which the TP is defined. Shape has the form [b1, ..., bB, e, f1, ..., fF] where F is the number of feature dimensions and must equal kernel.feature_ndims and e is the number (size) of index points in each batch. Ultimately this distribution corresponds to a e-dimensional multivariate Student's T. The batch shape must be broadcastable with kernel.batch_shape and any batch dims yielded by mean_fn.
mean_fn	Function that acts on index_points to produce a (batch of) vector(s) of mean values at index_points. Takes a Tensor of shape [b1, ..., bB, f1, ..., fF] and returns a Tensor whose shape is broadcastable with [b1, ..., bB]. Default value: NULL implies constant zero function.
jitter	float scalar Tensor added to the diagonal of the covariance matrix to ensure positive definiteness of the covariance matrix. Default value: 1e-6.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

Just as Student's T distributions are fully specified by their degrees of freedom, location and scale, a Student's T process can be completely specified by a degrees of freedom parameter, mean function and covariance function.

Let S denote the index set and K the space in which each indexed random variable takes its values (again, often \mathbb{R} or \mathbb{C}). The mean function is then a map $m: S \rightarrow K$, and the covariance function, or kernel, is a positive-definite function $k: (S \times S) \rightarrow K$. The properties of functions drawn from a TP are entirely dictated (up to translation) by the form of the kernel function.

This Distribution represents the marginal joint distribution over function values at a given finite collection of points $[x[1], \dots, x[N]]$ from the index set S . By definition, this marginal distribution is just a multivariate Student's T distribution, whose mean is given by the vector $[m(x[1]), \dots, m(x[N])]$ and whose covariance matrix is constructed from pairwise applications of the kernel function to the given inputs:

$$\begin{array}{cccc|}
 | & k(x[1], x[1]) & k(x[1], x[2]) & \dots & k(x[1], x[N]) & | \\
 | & k(x[2], x[1]) & k(x[2], x[2]) & \dots & k(x[2], x[N]) & | \\
 | & \dots & \dots & \dots & \dots & | \\
 | & k(x[N], x[1]) & k(x[N], x[2]) & \dots & k(x[N], x[N]) & |
 \end{array}$$

For this to be a valid covariance matrix, it must be symmetric and positive definite; hence the requirement that k be a positive definite function (which, by definition, says that the above procedure

will yield PD matrices). Note also we use a parameterization as suggested in Shat et al. (2014), which requires df to be greater than 2. This allows for the covariance for any finite dimensional marginal of the TP (a multivariate Student's T distribution) to just be the PD matrix generated by the kernel.

Mathematical Details

The probability density function (pdf) is a multivariate Student's T whose parameters are derived from the TP's properties:

```
pdf(x; df, index_points, mean_fn, kernel) = MultivariateStudentT(df, loc, K)
K = (df - 2) / df * (kernel.matrix(index_points, index_points) + jitter * eye(N))
loc = (x - mean_fn(index_points))^T @ K @ (x - mean_fn(index_points))
```

where:

- df is the degrees of freedom parameter for the TP.
- $index_points$ are points in the index set over which the TP is defined,
- $mean_fn$ is a callable mapping the index set to the TP's mean values,
- $kernel$ is `PositiveSemidefiniteKernel`-like and represents the covariance function of the TP,
- $jitter$ is added to the diagonal to ensure positive definiteness up to machine precision (otherwise Cholesky-decomposition is prone to failure),
- $eye(N)$ is an N -by- N identity matrix.

Value

a distribution instance.

References

- [Amar Shah, Andrew Gordon Wilson, and Zoubin Ghahramani. Student-t Processes as Alternatives to Gaussian Processes. In *Artificial Intelligence and Statistics, 2014*.](#)

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal()`.

```
tfd_multivariate_normal_full_covariance(), tfd_multivariate_normal_linear_operator(),
tfd_multivariate_normal_tri_l(), tfd_multivariate_student_t_linear_operator(), tfd_negative_binomial(),
tfd_normal(), tfd_one_hot_categorical(), tfd_pareto(), tfd_pixel_cnn(), tfd_poisson_log_normal_quadrature(),
tfd_poisson(), tfd_power_spherical(), tfd_probit_bernoulli(), tfd_quantized(), tfd_relaxed_bernoulli(),
tfd_relaxed_one_hot_categorical(), tfd_sample_distribution(), tfd_sinh_arcsinh(),
tfd_skellam(), tfd_spherical_uniform(), tfd_student_t(), tfd_transformed_distribution(),
tfd_triangular(), tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian(),
tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()
```

tfd_survival_function *Survival function.*

Description

Given random variable X , the survival function is defined: $\text{tfd_survival_function}(x) = P[X > x] = 1 - P[X \leq x] = 1 - \text{cdf}(x)$.

Usage

```
tfd_survival_function(distribution, value, ...)
```

Arguments

distribution	The distribution being used.
value	float or double Tensor.
...	Additional parameters passed to Python.

Value

a Tensor of shape $\text{sample_shape}(x) + \text{self}\batch_shape with values of type $\text{self}\$\text{dtype}$.

See Also

Other `distribution_methods`: [tfd_cdf\(\)](#), [tfd_covariance\(\)](#), [tfd_cross_entropy\(\)](#), [tfd_entropy\(\)](#), [tfd_kl_divergence\(\)](#), [tfd_log_cdf\(\)](#), [tfd_log_prob\(\)](#), [tfd_log_survival_function\(\)](#), [tfd_mean\(\)](#), [tfd_mode\(\)](#), [tfd_prob\(\)](#), [tfd_quantile\(\)](#), [tfd_sample\(\)](#), [tfd_stddev\(\)](#), [tfd_variance\(\)](#)

Examples

```
d <- tfd_normal(loc = c(1, 2), scale = c(1, 0.5))
x <- d %>% tfd_sample()
d %>% tfd_survival_function(x)
```

tfd_transformed_distribution
A Transformed Distribution

Description

A TransformedDistribution models $p(y)$ given a base distribution $p(x)$, and a deterministic, invertible, differentiable transform, $Y = g(X)$. The transform is typically an instance of the Bijector class and the base distribution is typically an instance of the Distribution class.

Usage

```
tfd_transformed_distribution(
    distribution,
    bijector,
    batch_shape = NULL,
    event_shape = NULL,
    kwargs_split_fn = NULL,
    validate_args = FALSE,
    parameters = NULL,
    name = NULL
)
```

Arguments

distribution	The base distribution instance to transform. Typically an instance of Distribution.
bijector	The object responsible for calculating the transformation. Typically an instance of Bijector.
batch_shape	integer vector Tensor which overrides distribution batch_shape; valid only if distribution.is_scalar_batch().
event_shape	integer vector Tensor which overrides distribution event_shape; valid only if distribution.is_scalar_event().
kwargs_split_fn	Python callable which takes a kwargs dict and returns a tuple of kwargs dicts for each of the distribution and bijector parameters respectively. Default value: <code>_default_kwargs_split_fn</code> (i.e., <code>lambda kwargs: (kwargs.get('distribution_kwargs', {}), kwargs)</code>).
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
parameters	Locals dict captured by subclass constructor, to be used for copy/slice re-instantiation operations.
name	The name for ops managed by the distribution. Default value: <code>bijector.name + distribution.name</code> .

Details

A Bijector is expected to implement the following functions:

- forward,
- inverse,
- inverse_log_det_jacobian.

The semantics of these functions are outlined in the Bijector documentation.

We now describe how a TransformedDistribution alters the input/outputs of a Distribution associated with a random variable (rv) X . Write $\text{cdf}(Y=y)$ for an absolutely continuous cumulative distribution function of random variable Y ; write the probability density function $\text{pdf}(Y=y) := d^k / (dy_1, \dots, dy_k) \text{cdf}(Y=y)$ for its derivative wrt to Y evaluated at y . Assume that $Y = g(X)$ where g is a deterministic diffeomorphism, i.e., a non-random, continuous, differentiable, and invertible function. Write the inverse of g as $X = g^{-1}(Y)$ and $(J \circ g)(x)$ for the Jacobian of g evaluated at x .

A TransformedDistribution implements the following operations:

- `sample` Mathematically: $Y = g(X)$ Programmatically: `bijector.forward(distribution.sample(...))`
- `log_prob` Mathematically: $(\log \circ \text{pdf})(Y=y) = (\log \circ \text{pdf} \circ g^{-1})(y) + (\log \circ \text{abs} \circ \text{det} \circ J \circ g^{-1})(y)$ Programmatically: `(distribution.log_prob(bijector.inverse(y)) + bijector.inverse_log_det_jacob...`
- `log_cdf` Mathematically: $(\log \circ \text{cdf})(Y=y) = (\log \circ \text{cdf} \circ g^{-1})(y)$ Programmatically: `distribution.log_cdf(bijector.inverse(x))`
- and similarly for: `cdf`, `prob`, `log_survival_function`, `survival_function`.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`,

```
tfd_skellam(), tfd_spherical_uniform(), tfd_student_t_process(), tfd_student_t(), tfd_triangular(),
tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian_process(),
tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()
```

tfd_triangular	<i>Triangular distribution with low, high and peak parameters</i>
----------------	---

Description

The parameters low, high and peak must be shaped in a way that supports broadcasting (e.g., high - low is a valid operation).

Usage

```
tfd_triangular(
  low = 0,
  high = 1,
  peak = 0.5,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "Triangular"
)
```

Arguments

low	Floating point tensor, lower boundary of the output interval. Must have low < high. Default value: 0.
high	Floating point tensor, upper boundary of the output interval. Must have low < high. Default value: 1.
peak	Floating point tensor, mode of the output interval. Must have low <= peak and peak <= high. Default value: 0.5.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag_plus_low_rank_linear_operator()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian_process()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

`tfd_truncated_cauchy` *The Truncated Cauchy distribution.*

Description

The truncated Cauchy is a Cauchy distribution bounded between `low` and `high` (the pdf is 0 outside these bounds and renormalized). Samples from this distribution are differentiable with respect to `loc` and `scale`, but not with respect to the bounds `low` and `high`.

Usage

```
tfd_truncated_cauchy(
    loc,
    scale,
    low,
    high,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = "TruncatedCauchy"
)
```

Arguments

loc	Floating point tensor; the modes of the corresponding non-truncated Cauchy distribution(s).
scale	Floating point tensor; the scales of the distribution(s). Must contain only positive values.
low	float Tensor representing lower bound of the distribution's support. Must be such that low < high.
high	float Tensor representing upper bound of the distribution's support. Must be such that low < high.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details**Mathematical Details**

The probability density function (pdf) of this distribution is:

$$\text{pdf}(x; \text{loc}, \text{scale}, \text{low}, \text{high}) = \begin{cases} 1 / (\pi * \text{scale} * (1 + z^{**2}) * A) & \text{for } \text{low} \leq x \leq \text{high} \\ 0 & \text{otherwise} \end{cases}$$

where

$$z = (x - \text{loc}) / \text{scale}$$

$$A = \text{CauchyCDF}((\text{high} - \text{loc}) / \text{scale}) - \text{CauchyCDF}((\text{low} - \text{loc}) / \text{scale})$$

where CauchyCDF is the cumulative density function of the Cauchy distribution with 0 mean and unit variance. This is a scalar distribution so the event shape is always scalar and the dimensions of the parameters define the batch_shape.

Value

a distribution instance.

See Also

For usage examples see e.g. [tfd_sample\(\)](#), [tfd_log_prob\(\)](#), [tfd_mean\(\)](#).

Other distributions: [tfd_autoregressive\(\)](#), [tfd_batch_reshape\(\)](#), [tfd_bates\(\)](#), [tfd_bernoulli\(\)](#), [tfd_beta_binomial\(\)](#), [tfd_beta\(\)](#), [tfd_binomial\(\)](#), [tfd_categorical\(\)](#), [tfd_cauchy\(\)](#), [tfd_chi2\(\)](#), [tfd_chi\(\)](#), [tfd_cholesky_lkj\(\)](#), [tfd_continuous_bernoulli\(\)](#), [tfd_deterministic\(\)](#), [tfd_dirichlet_multinomial\(\)](#), [tfd_dirichlet\(\)](#), [tfd_empirical\(\)](#), [tfd_exp_gamma\(\)](#), [tfd_exp_inverse_gamma\(\)](#), [tfd_exponential\(\)](#), [tfd_gamma_gamma\(\)](#), [tfd_gamma\(\)](#), [tfd_gaussian_process_regression_model\(\)](#), [tfd_gaussian_process\(\)](#),

```

tfd_generalized_normal(), tfd_geometric(), tfd_gumbel(), tfd_half_cauchy(), tfd_half_normal(),
tfd_hidden_markov_model(), tfd_horseshoe(), tfd_independent(), tfd_inverse_gamma(),
tfd_inverse_gaussian(), tfd_johnson_s_u(), tfd_joint_distribution_named_auto_batched(),
tfd_joint_distribution_named(), tfd_joint_distribution_sequential_auto_batched(),
tfd_joint_distribution_sequential(), tfd_kumaraswamy(), tfd_laplace(), tfd_linear_gaussian_state_space,
tfd_lkj(), tfd_log_logistic(), tfd_log_normal(), tfd_logistic(), tfd_mixture_same_family(),
tfd_mixture(), tfd_multinomial(), tfd_multivariate_normal_diag_plus_low_rank(), tfd_multivariate_normal,
tfd_multivariate_normal_full_covariance(), tfd_multivariate_normal_linear_operator(),
tfd_multivariate_normal_tri_l(), tfd_multivariate_student_t_linear_operator(), tfd_negative_binomial(),
tfd_normal(), tfd_one_hot_categorical(), tfd_pareto(), tfd_pixel_cnn(), tfd_poisson_log_normal_quadrature,
tfd_poisson(), tfd_power_spherical(), tfd_probit_bernoulli(), tfd_quantized(), tfd_relaxed_bernoulli(),
tfd_relaxed_one_hot_categorical(), tfd_sample_distribution(), tfd_sinh_arcsinh(),
tfd_skellam(), tfd_spherical_uniform(), tfd_student_t_process(), tfd_student_t(), tfd_transformed_distribution,
tfd_triangular(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian_process(),
tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()

```

tfd_truncated_normal *Truncated Normal distribution*

Description

The truncated normal is a normal distribution bounded between low and high (the pdf is 0 outside these bounds and renormalized). Samples from this distribution are differentiable with respect to loc, scale as well as the bounds, low and high, i.e., this implementation is fully reparameterizable. For more details, see [here](#).

Usage

```

tfd_truncated_normal(
  loc,
  scale,
  low,
  high,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "TruncatedNormal"
)

```

Arguments

loc	Floating point tensor; the means of the distribution(s).
scale	floating point tensor; the stddevs of the distribution(s). Must contain only positive values.

low	float Tensor representing lower bound of the distribution's support. Must be such that low < high.
high	float Tensor representing upper bound of the distribution's support. Must be such that low < high.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

Mathematical Details

The probability density function (pdf) of this distribution is:

$$\text{pdf}(x; \text{loc}, \text{scale}, \text{low}, \text{high}) = \begin{cases} (2 \pi)^{-0.5} \exp(-0.5 y^2) / (\text{scale} * z) & \text{for } \text{low} \leq x \leq \text{high} \\ 0 & \text{otherwise} \end{cases}$$

$$y = (x - \text{loc}) / \text{scale}$$

$$z = \text{NormalCDF}((\text{high} - \text{loc}) / \text{scale}) - \text{NormalCDF}((\text{low} - \text{loc}) / \text{scale})$$

where:

- NormalCDF is the cumulative density function of the Normal distribution with 0 mean and unit variance.

This is a scalar distribution so the event shape is always scalar and the dimensions of the parameters defined the batch_shape.

Value

a distribution instance.

See Also

For usage examples see e.g. [tfd_sample\(\)](#), [tfd_log_prob\(\)](#), [tfd_mean\(\)](#).

Other distributions: [tfd_autoregressive\(\)](#), [tfd_batch_reshape\(\)](#), [tfd_bates\(\)](#), [tfd_bernoulli\(\)](#), [tfd_beta_binomial\(\)](#), [tfd_beta\(\)](#), [tfd_binomial\(\)](#), [tfd_categorical\(\)](#), [tfd_cauchy\(\)](#), [tfd_chi2\(\)](#), [tfd_chi\(\)](#), [tfd_cholesky_lkj\(\)](#), [tfd_continuous_bernoulli\(\)](#), [tfd_deterministic\(\)](#), [tfd_dirichlet_multinomial\(\)](#), [tfd_dirichlet\(\)](#), [tfd_empirical\(\)](#), [tfd_exp_gamma\(\)](#), [tfd_exp_inverse_gamma\(\)](#), [tfd_exponential\(\)](#), [tfd_gamma_gamma\(\)](#), [tfd_gamma\(\)](#), [tfd_gaussian_process_regression_model\(\)](#), [tfd_gaussian_process\(\)](#), [tfd_generalized_normal\(\)](#), [tfd_geometric\(\)](#), [tfd_gumbel\(\)](#), [tfd_half_cauchy\(\)](#), [tfd_half_normal\(\)](#), [tfd_hidden_markov_model\(\)](#), [tfd_horseshoe\(\)](#), [tfd_independent\(\)](#), [tfd_inverse_gamma\(\)](#), [tfd_inverse_gaussian\(\)](#), [tfd_johnson_s_u\(\)](#), [tfd_joint_distribution_named_auto_batched\(\)](#), [tfd_joint_distribution_named\(\)](#), [tfd_joint_distribution_sequential_auto_batched\(\)](#),

tfd_joint_distribution_sequential(), tfd_kumaraswamy(), tfd_laplace(), tfd_linear_gaussian_state_space,
 tfd_lkj(), tfd_log_logistic(), tfd_log_normal(), tfd_logistic(), tfd_mixture_same_family(),
 tfd_mixture(), tfd_multinomial(), tfd_multivariate_normal_diag_plus_low_rank(), tfd_multivariate_normal,
 tfd_multivariate_normal_full_covariance(), tfd_multivariate_normal_linear_operator(),
 tfd_multivariate_normal_tri_l(), tfd_multivariate_student_t_linear_operator(), tfd_negative_binomial(),
 tfd_normal(), tfd_one_hot_categorical(), tfd_pareto(), tfd_pixel_cnn(), tfd_poisson_log_normal_quadrature,
 tfd_poisson(), tfd_power_spherical(), tfd_probit_bernoulli(), tfd_quantized(), tfd_relaxed_bernoulli(),
 tfd_relaxed_one_hot_categorical(), tfd_sample_distribution(), tfd_sinh_arcsinh(),
 tfd_skellam(), tfd_spherical_uniform(), tfd_student_t_process(), tfd_student_t(), tfd_transformed_distribution,
 tfd_triangular(), tfd_truncated_cauchy(), tfd_uniform(), tfd_variational_gaussian_process(),
 tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
 tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
 tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
 tfd_wishart_tri_l(), tfd_wishart(), tfd_zipf()

tfd_uniform

Uniform distribution with low and high parameters

Description

Mathematical Details

Usage

```
tfd_uniform(  
  low = 0,  
  high = 1,  
  validate_args = FALSE,  
  allow_nan_stats = TRUE,  
  name = "Uniform"  
)
```

Arguments

low	Floating point tensor, lower boundary of the output interval. Must have low < high.
high	Floating point tensor, upper boundary of the output interval. Must have low < high.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

The probability density function (pdf) is,

$$\text{pdf}(x; a, b) = I[a \leq x < b] / Z$$

$$Z = b - a$$

where

- low = a,
- high = b,
- Z is the normalizing constant, and
- $I[\text{predicate}]$ is the **indicator function** for predicate.

The parameters low and high must be shaped in a way that supports broadcasting (e.g., high - low is a valid operation).

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_variational_gaussian_process()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_variance	<i>Variance.</i>
--------------	------------------

Description

Variance is defined as, $\text{Var} = E[(X - E[X])**2]$ where X is the random variable associated with this distribution, E denotes expectation, and $\text{Var}\$shape = \text{batch_shape} + \text{event_shape}$.

Usage

```
tfd_variance(distribution, ...)
```

Arguments

distribution	The distribution being used.
...	Additional parameters passed to Python.

Value

a Tensor of shape $\text{sample_shape}(x) + \text{self}\$batch_shape$ with values of type $\text{self}\$dtype$.

See Also

Other `distribution_methods`: [tfd_cdf\(\)](#), [tfd_covariance\(\)](#), [tfd_cross_entropy\(\)](#), [tfd_entropy\(\)](#), [tfd_kl_divergence\(\)](#), [tfd_log_cdf\(\)](#), [tfd_log_prob\(\)](#), [tfd_log_survival_function\(\)](#), [tfd_mean\(\)](#), [tfd_mode\(\)](#), [tfd_prob\(\)](#), [tfd_quantile\(\)](#), [tfd_sample\(\)](#), [tfd_stddev\(\)](#), [tfd_survival_function\(\)](#)

Examples

```
d <- tfd_normal(loc = c(1, 2), scale = c(1, 0.5))
d %>% tfd_variance()
```

tfd_variational_gaussian_process

Posterior predictive of a variational Gaussian process

Description

This distribution implements the variational Gaussian process (VGP), as described in Titsias (2009) and Hensman (2013). The VGP is an inducing point-based approximation of an exact GP posterior. Ultimately, this Distribution class represents a marginal distribution over function values at a collection of `index_points`. It is parameterized by

- a kernel function,
- a mean function,
- the (scalar) observation noise variance of the normal likelihood,
- a set of index points,
- a set of inducing index points, and
- the parameters of the (full-rank, Gaussian) variational posterior distribution over function values at the inducing points, conditional on some observations.

Usage

```
tfd_variational_gaussian_process(
    kernel,
    index_points,
    inducing_index_points,
    variational_inducing_observations_loc,
    variational_inducing_observations_scale,
    mean_fn = NULL,
    observation_noise_variance = 0,
    predictive_noise_variance = 0,
    jitter = 1e-06,
    validate_args = FALSE,
    allow_nan_stats = FALSE,
    name = "VariationalGaussianProcess"
)
```

Arguments

<code>kernel</code>	PositiveSemidefiniteKernel-like instance representing the GP's covariance function.
<code>index_points</code>	float Tensor representing finite (batch of) vector(s) of points in the index set over which the VGP is defined. Shape has the form <code>[b1, ..., bB, e1, f1, ..., fF]</code> where <code>F</code> is the number of feature dimensions and must equal <code>kernel\$feature_ndims</code> and <code>e1</code> is the number (size) of index points in each batch (we denote it <code>e1</code> to distinguish it from the numer of inducing index points, denoted <code>e2</code> below). Ultimately the <code>VariationalGaussianProcess</code> distribution corresponds to an <code>e1</code> -dimensional multivariate normal. The batch shape must be broadcastable with <code>kernel\$batch_shape</code> , the batch shape of <code>inducing_index_points</code> , and any batch dims yielded by <code>mean_fn</code> .

<code>inducing_index_points</code>	float Tensor of locations of inducing points in the index set. Shape has the form <code>[b1, ..., bB, e2, f1, ..., fF]</code> , just like <code>index_points</code> . The batch shape components needn't be identical to those of <code>index_points</code> , but must be broadcast compatible with them.
<code>variational_inducing_observations_loc</code>	float Tensor; the mean of the (full-rank Gaussian) variational posterior over function values at the inducing points, conditional on observed data. Shape has the form <code>[b1, ..., bB, e2]</code> , where <code>b1, ..., bB</code> is broadcast compatible with other parameters' batch shapes, and <code>e2</code> is the number of inducing points.
<code>variational_inducing_observations_scale</code>	float Tensor; the scale matrix of the (full-rank Gaussian) variational posterior over function values at the inducing points, conditional on observed data. Shape has the form <code>[b1, ..., bB, e2, e2]</code> , where <code>b1, ..., bB</code> is broadcast compatible with other parameters and <code>e2</code> is the number of inducing points.
<code>mean_fn</code>	function that acts on index points to produce a (batch of) vector(s) of mean values at those index points. Takes a Tensor of shape <code>[b1, ..., bB, f1, ..., fF]</code> and returns a Tensor whose shape is (broadcastable with) <code>[b1, ..., bB]</code> . Default value: NULL implies constant zero function.
<code>observation_noise_variance</code>	float Tensor representing the variance of the noise in the Normal likelihood distribution of the model. May be batched, in which case the batch shape must be broadcastable with the shapes of all other batched parameters (<code>kernel\$batch_shape</code> , <code>index_points</code> , etc.). Default value: \emptyset .
<code>predictive_noise_variance</code>	float Tensor representing additional variance in the posterior predictive model. If NULL, we simply re-use <code>observation_noise_variance</code> for the posterior predictive noise. If set explicitly, however, we use the given value. This allows us, for example, to omit predictive noise variance (by setting this to zero) to obtain noiseless posterior predictions of function values, conditioned on noisy observations.
<code>jitter</code>	float scalar Tensor added to the diagonal of the covariance matrix to ensure positive definiteness of the covariance matrix. Default value: $1e-6$.
<code>validate_args</code>	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
<code>allow_nan_stats</code>	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Details

A VGP is "trained" by selecting any kernel parameters, the locations of the inducing index points, and the variational parameters. Titsias (2009) and Hensman (2013) describe a variational lower bound on the marginal log likelihood of observed data, which this class offers through the `variational_loss`

method (this is the negative lower bound, for convenience when plugging into a TF Optimizer's minimize function). Training may be done in minibatches.

Titsias (2009) describes a closed form for the optimal variational parameters, in the case of sufficiently small observational data (ie, small enough to fit in memory but big enough to warrant approximating the GP posterior). A method to compute these optimal parameters in terms of the full observational data set is provided as a staticmethod, `optimal_variational_posterior`. It returns a `MultivariateNormalLinearOperator` instance with optimal location and scale parameters.

Mathematical Details

Notation We will in general be concerned about three collections of index points, and it'll be good to give them names:

- $x[1], \dots, x[N]$: observation index points – locations of our observed data.
- $z[1], \dots, z[M]$: inducing index points – locations of the "summarizing" inducing points
- $t[1], \dots, t[P]$: predictive index points – locations where we are making posterior predictions based on observations and the variational parameters.

To lighten notation, we'll use X, Z, T to denote the above collections. Similarly, we'll denote by $f(X)$ the collection of function values at each of the $x[i]$, and by Y , the collection of (noisy) observed data at each $x[i]$. We'll denote kernel matrices generated from pairs of index points as K_{tt}, K_{xt}, K_{tz} , etc, e.g.,

$$K_{tz} = \begin{array}{cccc|c} | & k(t[1], z[1]) & k(t[1], z[2]) & \dots & k(t[1], z[M]) & | \\ | & k(t[2], z[1]) & k(t[2], z[2]) & \dots & k(t[2], z[M]) & | \\ | & \dots & \dots & \dots & \dots & | \\ | & k(t[P], z[1]) & k(t[P], z[2]) & \dots & k(t[P], z[M]) & | \end{array}$$

Preliminaries A Gaussian process is an indexed collection of random variables, any finite collection of which are jointly Gaussian. Typically, the index set is some finite-dimensional, real vector space, and indeed we make this assumption in what follows. The GP may then be thought of as a distribution over functions on the index set. Samples from the GP are functions *on the whole index set*; these can't be represented in finite compute memory, so one typically works with the marginals at a finite collection of index points. The properties of the GP are entirely determined by its mean function m and covariance function k . The generative process, assuming a mean-zero normal likelihood with stddev σ , is

$$f \sim \text{GP}(m, k) \\ Y | f(X) \sim \text{Normal}(f(X), \sigma), \quad i = 1, \dots, N$$

In finite terms (ie, marginalizing out all but a finite number of $f(X)$, σ), we can write

$$f(X) \sim \text{MVN}(\text{loc}=m(X), \text{cov}=K_{xx}) \\ Y | f(X) \sim \text{Normal}(f(X), \sigma), \quad i = 1, \dots, N$$

Posterior inference is possible in analytical closed form but becomes intractible as data sizes get large. See Rasmussen (2006) for details.

The VGP

The VGP is an inducing point-based approximation of an exact GP posterior, where two approximating assumptions have been made:

1. function values at non-inducing points are mutually independent conditioned on function values at the inducing points,
2. the (expensive) posterior over function values at inducing points conditional on observations is replaced with an arbitrary (learnable) full-rank Gaussian distribution,

$$q(f(Z)) = \text{MVN}(\text{loc}=m, \text{scale}=S),$$

where m and S are parameters to be chosen by optimizing an evidence lower bound (ELBO). The posterior predictive distribution becomes

$$q(f(T)) = \int df(Z) p(f(T) | f(Z)) q(f(Z)) = \text{MVN}(\text{loc} = A @ m, \text{scale} = B^{(1/2)})$$

where

$$A = K_{tz} @ K_{zz}^{-1}$$

$$B = K_{tt} - A @ (K_{zz} - S S^T) A^T$$

The approximate posterior predictive distribution $q(f(T))$ is what the `VariationalGaussianProcess` class represents.

Model selection in this framework entails choosing the kernel parameters, inducing point locations, and variational parameters. We do this by optimizing a variational lower bound on the marginal log likelihood of observed data. The lower bound takes the following form (see Titsias (2009) and Hensman (2013) for details on the derivation):

$$L(Z, m, S, Y) = \text{MVN}(\text{loc}=(K_{zx} @ K_{zz}^{-1}) @ m, \text{scale_diag}=\text{sigma}).\text{log_prob}(Y) -$$

$$(\text{Tr}(K_{xx} - K_{zx} @ K_{zz}^{-1} @ K_{xz}) +$$

$$\text{Tr}(S @ S^T @ K_{zz}^{-1} @ K_{zx} @ K_{xz} @ K_{zz}^{-1})) / (2 * \text{sigma}^2) -$$

$$\text{KL}(q(f(Z)) || p(f(Z)))$$

where in the final KL term, $p(f(Z))$ is the GP prior on inducing point function values. This variational lower bound can be computed on minibatches of the full data set (X, Y) . A method to compute the *negative* variational lower bound is implemented as `VariationalGaussianProcess$variational_loss`.

Optimal variational parameters

As described in Titsias (2009), a closed form optimum for the variational location and scale parameters, m and S , can be computed when the observational data are not prohibitively voluminous. The `optimal_variational_posterior` function computes the optimal variational posterior distribution over inducing point function values in terms of the GP parameters (mean and kernel functions), inducing point locations, observation index points, and observations. Note that the inducing index point locations must still be optimized even when these parameters are known functions of the inducing index points. The optimal parameters are computed as follows:

$$C = \text{sigma}^{-2} (K_{zz} + K_{zx} @ K_{xz})^{-1}$$

optimal Gaussian covariance: $K_{zz} @ C @ K_{zz}$

optimal Gaussian location: $\text{sigma}^{-2} K_{zz} @ C @ K_{zx} @ Y$

Value

a distribution instance.

References

- Titsias, M. "Variational Model Selection for Sparse Gaussian Process Regression", 2009.
- Hensman, J., Lawrence, N. "Gaussian Processes for Big Data", 2013.
- Carl Rasmussen, Chris Williams. Gaussian Processes For Machine Learning, 2006.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_vector_deterministic

Vector Deterministic Distribution

Description

The VectorDeterministic distribution is parameterized by a batch point loc in \mathbb{R}^k . The distribution is supported at this point only, and corresponds to a random variable that is constant, equal to loc .

Usage

```
tfd_vector_deterministic(
  loc,
  atol = NULL,
```

```

    rtol = NULL,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = "VectorDeterministic"
)

```

Arguments

<code>loc</code>	Numeric Tensor of shape $[B_1, \dots, B_b, k]$, with $b \geq 0, k \geq 0$ The point (or batch of points) on which this distribution is supported.
<code>atol</code>	Non-negative Tensor of same dtype as <code>loc</code> and broadcastable shape. The absolute tolerance for comparing closeness to <code>loc</code> . Default is 0.
<code>rtol</code>	Non-negative Tensor of same dtype as <code>loc</code> and broadcastable shape. The relative tolerance for comparing closeness to <code>loc</code> . Default is 0.
<code>validate_args</code>	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
<code>allow_nan_stats</code>	Logical, default <code>TRUE</code> . When <code>TRUE</code> , statistics (e.g., mean, mode, variance) use the value <code>NaN</code> to indicate the result is undefined. When <code>FALSE</code> , an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Details

See [Degenerate rv.](#)

Value

a distribution instance.

See Also

For usage examples see e.g. [tfd_sample\(\)](#), [tfd_log_prob\(\)](#), [tfd_mean\(\)](#).

tfd_vector_diffeomixture

VectorDiffeomixture distribution

Description

A vector diffeomixture (VDM) is a distribution parameterized by a convex combination of K component `loc` vectors, `loc[k]`, $k = 0, \dots, K-1$, and K scale matrices `scale[k]`, $k = 0, \dots, K-1$. It approximates the following **compound distribution** $p(x) = \int p(x | z) p(z) dz$, where z is in the K -simplex, and $p(x | z) := p(x | \text{loc}=\sum_k z[k] \text{loc}[k], \text{scale}=\sum_k z[k] \text{scale}[k])$

Usage

```
tfd_vector_diffeomixture(
  mix_loc,
  temperature,
  distribution,
  loc = NULL,
  scale = NULL,
  quadrature_size = 8,
  quadrature_fn = tfp$distributions$quadrature_scheme_softmaxnormal_quantiles,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "VectorDiffeomixture"
)
```

Arguments

<code>mix_loc</code>	float-like Tensor with shape $[b_1, \dots, b_B, K-1]$. In terms of samples, larger <code>mix_loc[... ,k]</code> \implies Z is more likely to put more weight on its kth component.
<code>temperature</code>	float-like Tensor. Broadcastable with <code>mix_loc</code> . In terms of samples, smaller temperature means one component is more likely to dominate. I.e., smaller temperature makes the VDM look more like a standard mixture of K components.
<code>distribution</code>	<code>tfp\$distributions\$Distribution</code> -like instance. Distribution from which <code>d</code> iid samples are used as input to the selected affine transformation. Must be a scalar-batch, scalar-event distribution. Typically <code>distribution\$reparameterization_type = FULLY_REPARAMETERIZED</code> or it is a function of non-trainable parameters. WARNING: If you backprop through a <code>VectorDiffeomixture</code> sample and the <code>distribution</code> is not <code>FULLY_REPARAMETERIZED</code> yet is a function of trainable variables, then the gradient will be incorrect!
<code>loc</code>	Length-K list of float-type Tensors. The k-th element represents the shift used for the k-th affine transformation. If the k-th item is <code>NULL</code> , <code>loc</code> is implicitly \emptyset . When specified, must have shape $[B_1, \dots, B_b, d]$ where $b \geq 0$ and <code>d</code> is the event size.
<code>scale</code>	Length-K list of <code>LinearOperators</code> . Each should be positive-definite and operate on a <code>d</code> -dimensional vector space. The k-th element represents the scale used for the k-th affine transformation. <code>LinearOperators</code> must have shape $[B_1, \dots, B_b, d, d]$, $b \geq 0$, i.e., characterizes <code>b</code> -batches of <code>d x d</code> matrices
<code>quadrature_size</code>	integer scalar representing number of quadrature points. Larger <code>quadrature_size</code> means $q_N(x)$ better approximates $p(x)$.
<code>quadrature_fn</code>	Function taking <code>normal_loc</code> , <code>normal_scale</code> , <code>quadrature_size</code> , <code>validate_args</code> and returning <code>tuple(grid, probs)</code> representing the <code>SoftmaxNormal</code> grid and corresponding normalized weight. Default value: <code>quadrature_scheme_softmaxnor</code>
<code>validate_args</code>	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .

allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

The integral $\int p(x | z) p(z) dz$ is approximated with a quadrature scheme adapted to the mixture density $p(z)$. The N quadrature points $z_{\{N, n\}}$ and weights $w_{\{N, n\}}$ (which are non-negative and sum to 1) are chosen such that $q_N(x) := \sum_{n=1}^N w_{\{n, N\}} p(x | z_{\{N, n\}}) \rightarrow p(x)$ as $N \rightarrow \infty$.

Since $q_N(x)$ is in fact a mixture (of N points), we may sample from q_N exactly. It is important to note that the VDM is *defined* as q_N above, and *not* $p(x)$. Therefore, sampling and pdf may be implemented as exact (up to floating point error) methods.

A common choice for the conditional $p(x | z)$ is a multivariate Normal. The implemented marginal $p(z)$ is the `SoftmaxNormal`, which is a $K-1$ dimensional Normal transformed by a `SoftmaxCentered` bijector, making it a density on the K -simplex. That is, $Z = \text{SoftmaxCentered}(X)$, $X = \text{Normal}(\text{mix_loc} / \text{temperature}, 1 / \text{temperature})$

The default quadrature scheme chooses $z_{\{N, n\}}$ as N midpoints of the quantiles of $p(z)$ (generalized quantiles if $K > 2$). See Dillon and Langmore (2018) for more details.

About Vector distributions in TensorFlow.

The `VectorDiffeomixture` is a non-standard distribution that has properties particularly useful in **variational Bayesian methods**. Conditioned on a draw from the `SoftmaxNormal`, $X|z$ is a vector whose components are linear combinations of affine transformations, thus is itself an affine transformation.

Note: The marginals $X_{1|v}$, ..., $X_{d|v}$ are *not* generally identical to some parameterization of distribution. This is due to the fact that the sum of draws from distribution are not generally itself the same distribution.

About Diffeomixtures and reparameterization.

The `VectorDiffeomixture` is designed to be reparameterized, i.e., its parameters are only used to transform samples from a distribution which has no trainable parameters. This property is important because backprop stops at sources of stochasticity. That is, as long as the parameters are used *after* the underlying source of stochasticity, the computed gradient is accurate. Reparametrization means that we can use gradient-descent (via backprop) to optimize Monte-Carlo objectives. Such objectives are a finite-sample approximation of an expectation and arise throughout scientific computing.

WARNING: If you backprop through a `VectorDiffeomixture` sample and the "base" distribution is both: not `FULLY_REPARAMETERIZED` and a function of trainable variables, then the gradient is not guaranteed correct!

Value

a distribution instance.

References

- Joshua Dillon and Ian Langmore. Quadrature Compound: An approximating family of distributions. *arXiv preprint arXiv:1801.03080*, 2018.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_vector_exponential_diag

The vectorization of the Exponential distribution on \mathbb{R}^k

Description

The vector exponential distribution is defined over a subset of \mathbb{R}^k , and parameterized by a (batch of) length- k loc vector and a (batch of) $k \times k$ scale matrix: $\text{covariance} = \text{scale} @ \text{scale.T}$, where $@$ denotes matrix-multiplication.

Usage

```
tfd_vector_exponential_diag(
    loc = NULL,
    scale_diag = NULL,
```

```

    scale_identity_multiplier = NULL,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = "VectorExponentialDiag"
  )

```

Arguments

<code>loc</code>	Floating-point Tensor. If this is set to NULL, <code>loc</code> is implicitly 0. When specified, may have shape $[B_1, \dots, B_b, k]$ where $b \geq 0$ and k is the event size.
<code>scale_diag</code>	Non-zero, floating-point Tensor representing a diagonal matrix added to <code>scale</code> . May have shape $[B_1, \dots, B_b, k]$, $b \geq 0$, and characterizes b -batches of $k \times k$ diagonal matrices added to <code>scale</code> . When both <code>scale_identity_multiplier</code> and <code>scale_diag</code> are NULL then <code>scale</code> is the Identity.
<code>scale_identity_multiplier</code>	Non-zero, floating-point Tensor representing a scaled-identity-matrix added to <code>scale</code> . May have shape $[B_1, \dots, B_b]$, $b \geq 0$, and characterizes b -batches of scaled $k \times k$ identity matrices added to <code>scale</code> . When both <code>scale_identity_multiplier</code> and <code>scale_diag</code> are NULL then <code>scale</code> is the Identity.
<code>validate_args</code>	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
<code>allow_nan_stats</code>	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Details

Mathematical Details The probability density function (pdf) is defined over the image of the `scale` matrix + `loc`, applied to the positive half-space: $\text{Supp} = \{\text{loc} + \text{scale} @ x : x \in \mathbb{R}^k, x_1 > 0, \dots, x_k > 0\}$. On this set,

$$\begin{aligned} \text{pdf}(y; \text{loc}, \text{scale}) &= \exp(-\|x\|_1) / Z, \quad \text{for } y \text{ in Supp} \\ x &= \text{inv}(\text{scale}) @ (y - \text{loc}), \\ Z &= |\det(\text{scale})|, \end{aligned}$$

where:

- `loc` is a vector in \mathbb{R}^k ,
- `scale` is a linear operator in $\mathbb{R}^{k \times k}$, $\text{cov} = \text{scale} @ \text{scale} . T$,
- Z denotes the normalization constant, and,
- $\|x\|_1$ denotes the l_1 norm of x , $\sum_i |x_i|$. The VectorExponential distribution is a member of the **location-scale family**, i.e., it can be constructed as,

$$\begin{aligned} X &= (X_1, \dots, X_k), \quad \text{each } X_i \sim \text{Exponential}(\text{rate}=1) \\ Y &= (Y_1, \dots, Y_k) = \text{scale} @ X + \text{loc} \end{aligned}$$

About `VectorExponential` and `Vector` distributions in TensorFlow.

The `VectorExponential` is a non-standard distribution that has useful properties. The marginals Y_1, \dots, Y_k are *not* Exponential random variables, due to the fact that the sum of Exponential random variables is not Exponential. Instead, Y is a vector whose components are linear combinations of Exponential random variables. Thus, Y lives in the vector space generated by vectors of Exponential distributions. This allows the user to decide the mean and covariance (by setting `loc` and `scale`), while preserving some properties of the Exponential distribution. In particular, the tails of Y_i will be (up to polynomial factors) exponentially decaying. To see this last statement, note that the pdf of Y_i is the convolution of the pdf of k independent Exponential random variables. One can then show by induction that distributions with exponential (up to polynomial factors) tails are closed under convolution.

The `batch_shape` is the broadcast shape between `loc` and `scale` arguments. The `event_shape` is given by last dimension of the matrix implied by `scale`. The last dimension of `loc` (if provided) must broadcast with this. Recall that `covariance = 2 * scale @ scale.T`. Additional leading dimensions (if any) will index batches. If both `scale_diag` and `scale_identity_multiplier` are `NULL`, then `scale` is the Identity matrix.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_vector_exponential_linear_operator

The vectorization of the Exponential distribution on \mathbb{R}^k

Description

The vector exponential distribution is defined over a subset of \mathbb{R}^k , and parameterized by a (batch of) length- k loc vector and a (batch of) $k \times k$ scale matrix: $\text{covariance} = \text{scale} @ \text{scale.T}$, where $@$ denotes matrix-multiplication.

Usage

```
tfd_vector_exponential_linear_operator(
    loc = NULL,
    scale = NULL,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = "VectorExponentialLinearOperator"
)
```

Arguments

loc	Floating point tensor; the means of the distribution(s).
scale	Instance of LinearOperator with same dtype as loc and shape $[B_1, \dots, B_b, k, k]$.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

Mathematical Details The probability density function (pdf) is

$$\text{pdf}(y; \text{loc}, \text{scale}) = \exp(-\|x\|_1) / Z, \quad \text{for } y \text{ in } S(\text{loc}, \text{scale}),$$

$$x = \text{inv}(\text{scale}) @ (y - \text{loc}),$$

$$Z = |\det(\text{scale})|,$$

where:

- loc is a vector in \mathbb{R}^k ,
- scale is a linear operator in $\mathbb{R}^{\{k \times k\}}$, $\text{cov} = \text{scale} @ \text{scale.T}$,
- $S = \{\text{loc} + \text{scale} @ x : x \text{ in } \mathbb{R}^k, x_1 > 0, \dots, x_k > 0\}$, is an image of the positive half-space,

- $\|x\|_1$ denotes the l1 norm of x , $\sum_i |x_i|$,
- Z denotes the normalization constant.

The VectorExponential distribution is a member of the **location-scale family**, i.e., it can be constructed as,

$$X = (X_1, \dots, X_k), \text{ each } X_i \sim \text{Exponential}(\text{rate}=1)$$

$$Y = (Y_1, \dots, Y_k) = \text{scale} @ X + \text{loc}$$

About VectorExponential and Vector distributions in TensorFlow.

The VectorExponential is a non-standard distribution that has useful properties. The marginals Y_1, \dots, Y_k are *not* Exponential random variables, due to the fact that the sum of Exponential random variables is not Exponential. Instead, Y is a vector whose components are linear combinations of Exponential random variables. Thus, Y lives in the vector space generated by vectors of Exponential distributions. This allows the user to decide the mean and covariance (by setting loc and scale), while preserving some properties of the Exponential distribution. In particular, the tails of Y_i will be (up to polynomial factors) exponentially decaying. To see this last statement, note that the pdf of Y_i is the convolution of the pdf of k independent Exponential random variables. One can then show by induction that distributions with exponential (up to polynomial factors) tails are closed under convolution.

The batch_shape is the broadcast shape between loc and scale arguments. The event_shape is given by last dimension of the matrix implied by scale . The last dimension of loc (if provided) must broadcast with this. Recall that $\text{covariance} = 2 * \text{scale} @ \text{scale}^T$. Additional leading dimensions (if any) will index batches.

#' @param loc Floating-point Tensor. If this is set to NULL, loc is implicitly 0. When specified, may have shape $[B_1, \dots, B_b, k]$ where $b \geq 0$ and k is the event size.

Value

a distribution instance.

See Also

For usage examples see e.g. [tfd_sample\(\)](#), [tfd_log_prob\(\)](#), [tfd_mean\(\)](#).

Other distributions: [tfd_autoregressive\(\)](#), [tfd_batch_reshape\(\)](#), [tfd_bates\(\)](#), [tfd_bernoulli\(\)](#), [tfd_beta_binomial\(\)](#), [tfd_beta\(\)](#), [tfd_binomial\(\)](#), [tfd_categorical\(\)](#), [tfd_cauchy\(\)](#), [tfd_chi2\(\)](#), [tfd_chi\(\)](#), [tfd_cholesky_lkj\(\)](#), [tfd_continuous_bernoulli\(\)](#), [tfd_deterministic\(\)](#), [tfd_dirichlet_multinomial\(\)](#), [tfd_dirichlet\(\)](#), [tfd_empirical\(\)](#), [tfd_exp_gamma\(\)](#), [tfd_exp_inverse_gamma\(\)](#), [tfd_exponential\(\)](#), [tfd_gamma_gamma\(\)](#), [tfd_gamma\(\)](#), [tfd_gaussian_process_regression_model\(\)](#), [tfd_gaussian_process\(\)](#), [tfd_generalized_normal\(\)](#), [tfd_geometric\(\)](#), [tfd_gumbel\(\)](#), [tfd_half_cauchy\(\)](#), [tfd_half_normal\(\)](#), [tfd_hidden_markov_model\(\)](#), [tfd_horseshoe\(\)](#), [tfd_independent\(\)](#), [tfd_inverse_gamma\(\)](#), [tfd_inverse_gaussian\(\)](#), [tfd_johnson_s_u\(\)](#), [tfd_joint_distribution_named_auto_batched\(\)](#), [tfd_joint_distribution_named\(\)](#), [tfd_joint_distribution_sequential_auto_batched\(\)](#), [tfd_joint_distribution_sequential\(\)](#), [tfd_kumaraswamy\(\)](#), [tfd_laplace\(\)](#), [tfd_linear_gaussian_state_space\(\)](#), [tfd_lkj\(\)](#), [tfd_log_logistic\(\)](#), [tfd_log_normal\(\)](#), [tfd_logistic\(\)](#), [tfd_mixture_same_family\(\)](#), [tfd_mixture\(\)](#), [tfd_multinomial\(\)](#), [tfd_multivariate_normal_diag_plus_low_rank\(\)](#), [tfd_multivariate_normal_full_covariance\(\)](#), [tfd_multivariate_normal_linear_operator\(\)](#), [tfd_multivariate_normal_tri_l\(\)](#), [tfd_multivariate_student_t_linear_operator\(\)](#), [tfd_negative_binomial\(\)](#)

```
tfd_normal(), tfd_one_hot_categorical(), tfd_pareto(), tfd_pixel_cnn(), tfd_poisson_log_normal_quadrature(),
tfd_poisson(), tfd_power_spherical(), tfd_probit_bernoulli(), tfd_quantized(), tfd_relaxed_bernoulli(),
tfd_relaxed_one_hot_categorical(), tfd_sample_distribution(), tfd_sinh_arcsinh(),
tfd_skellam(), tfd_spherical_uniform(), tfd_student_t_process(), tfd_student_t(), tfd_transformed_distribution(),
tfd_triangular(), tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian(),
tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_laplace_diag(),
tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(), tfd_von_mises_fisher(),
tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(), tfd_wishart_tri_l(),
tfd_wishart(), tfd_zipf()
```

tfd_vector_laplace_diag

The vectorization of the Laplace distribution on R^k

Description

The vector laplace distribution is defined over R^k , and parameterized by a (batch of) length- k loc vector (the means) and a (batch of) $k \times k$ scale matrix: covariance = $2 * \text{scale} @ \text{scale}^T$, where $@$ denotes matrix-multiplication.

Usage

```
tfd_vector_laplace_diag(
  loc = NULL,
  scale_diag = NULL,
  scale_identity_multiplier = NULL,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "VectorLaplaceDiag"
)
```

Arguments

loc	Floating-point Tensor. If this is set to NULL, loc is implicitly 0. When specified, may have shape $[B_1, \dots, B_b, k]$ where $b \geq 0$ and k is the event size.
scale_diag	Non-zero, floating-point Tensor representing a diagonal matrix added to scale. May have shape $[B_1, \dots, B_b, k]$, $b \geq 0$, and characterizes b -batches of $k \times k$ diagonal matrices added to scale. When both <code>scale_identity_multiplier</code> and <code>scale_diag</code> are NULL then scale is the Identity.
scale_identity_multiplier	Non-zero, floating-point Tensor representing a scaled-identity-matrix added to scale. May have shape $[B_1, \dots, B_b]$, $b \geq 0$, and characterizes b -batches of scaled $k \times k$ identity matrices added to scale. When both <code>scale_identity_multiplier</code> and <code>scale_diag</code> are NULL then scale is the Identity.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.

<code>allow_nan_stats</code>	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Details

Mathematical Details The probability density function (pdf) is,

$$\begin{aligned} \text{pdf}(x; \text{loc}, \text{scale}) &= \exp(-\|y\|_1) / Z \\ y &= \text{inv}(\text{scale}) @ (x - \text{loc}) \\ Z &= 2^{**k} |\det(\text{scale})| \end{aligned}$$

where:

- `loc` is a vector in \mathbb{R}^k ,
- `scale` is a linear operator in $\mathbb{R}^{k \times k}$, `cov = scale @ scale.T`,
- `Z` denotes the normalization constant, and,
- `\|y\|_1` denotes the l1 norm of `y`, `'sum_i |y_i|`.

A (non-batch) scale matrix is:

$$\text{scale} = \text{diag}(\text{scale_diag} + \text{scale_identity_multiplier} * \text{ones}(k))$$

where:

- `scale_diag.shape = [k]`, and,
- `scale_identity_multiplier.shape = []`. Additional leading dimensions (if any) will index batches. If both `scale_diag` and `scale_identity_multiplier` are NULL, then `scale` is the Identity matrix.

About VectorLaplace and Vector distributions in TensorFlow

The VectorLaplace is a non-standard distribution that has useful properties. The marginals Y_1, \dots, Y_k are *not* Laplace random variables, due to the fact that the sum of Laplace random variables is not Laplace. Instead, Y is a vector whose components are linear combinations of Laplace random variables. Thus, Y lives in the vector space generated by vectors of Laplace distributions. This allows the user to decide the mean and covariance (by setting `loc` and `scale`), while preserving some properties of the Laplace distribution. In particular, the tails of Y_i will be (up to polynomial factors) exponentially decaying. To see this last statement, note that the pdf of Y_i is the convolution of the pdf of k independent Laplace random variables. One can then show by induction that distributions with exponential (up to polynomial factors) tails are closed under convolution.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag_plus_low_rank_batched()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

`tfd_vector_laplace_linear_operator`

The vectorization of the Laplace distribution on \mathbb{R}^k

Description

The vector laplace distribution is defined over \mathbb{R}^k , and parameterized by a (batch of) length- k loc vector (the means) and a (batch of) $k \times k$ scale matrix: `covariance = 2 * scale @ scale.T`, where `@` denotes matrix-multiplication.

Usage

```
tfd_vector_laplace_linear_operator(
    loc = NULL,
    scale = NULL,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = "VectorLaplaceLinearOperator"
)
```

Arguments

<code>loc</code>	Floating-point Tensor. If this is set to NULL, <code>loc</code> is implicitly 0. When specified, may have shape $[B_1, \dots, B_b, k]$ where $b \geq 0$ and k is the event size.
<code>scale</code>	Instance of <code>LinearOperator</code> with same dtype as <code>loc</code> and shape $[B_1, \dots, B_b, k, k]$.
<code>validate_args</code>	Logical, default <code>FALSE</code> . When <code>TRUE</code> distribution parameters are checked for validity despite possibly degrading runtime performance. When <code>FALSE</code> invalid inputs may silently render incorrect outputs. Default value: <code>FALSE</code> .
<code>allow_nan_stats</code>	Logical, default <code>TRUE</code> . When <code>TRUE</code> , statistics (e.g., mean, mode, variance) use the value <code>NaN</code> to indicate the result is undefined. When <code>FALSE</code> , an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Details

Mathematical Details The probability density function (pdf) is,

$$\begin{aligned} \text{pdf}(x; \text{loc}, \text{scale}) &= \exp(-\|y\|_1) / Z, \\ y &= \text{inv}(\text{scale}) @ (x - \text{loc}), \\ Z &= 2^{**k} |\det(\text{scale})|, \end{aligned}$$

where:

- `loc` is a vector in \mathbb{R}^k ,
- `scale` is a linear operator in $\mathbb{R}^{k \times k}$, $\text{cov} = \text{scale} @ \text{scale} . T$,
- Z denotes the normalization constant, and,
- $\|y\|_1$ denotes the l_1 norm of y , $\sum_i |y_{i1}|$.

The `VectorLaplace` distribution is a member of the **location-scale family**, i.e., it can be constructed as,

$$\begin{aligned} X &= (X_1, \dots, X_k), \text{ each } X_i \sim \text{Laplace}(\text{loc}=0, \text{scale}=1) \\ Y &= (Y_1, \dots, Y_k) = \text{scale} @ X + \text{loc} \end{aligned}$$

About `VectorLaplace` and `Vector` distributions in TensorFlow

The `VectorLaplace` is a non-standard distribution that has useful properties. The marginals Y_1, \dots, Y_k are *not* Laplace random variables, due to the fact that the sum of Laplace random variables is not Laplace. Instead, Y is a vector whose components are linear combinations of Laplace random variables. Thus, Y lives in the vector space generated by vectors of Laplace distributions. This allows the user to decide the mean and covariance (by setting `loc` and `scale`), while preserving some properties of the Laplace distribution. In particular, the tails of Y_i will be (up to polynomial factors) exponentially decaying. To see this last statement, note that the pdf of Y_i is the convolution of the pdf of k independent Laplace random variables. One can then show by induction that distributions with exponential (up to polynomial factors) tails are closed under convolution.

The `batch_shape` is the broadcast shape between `loc` and `scale` arguments. The `event_shape` is given by last dimension of the matrix implied by `scale`. The last dimension of `loc` (if provided) must broadcast with this. Recall that $\text{covariance} = 2 * \text{scale} @ \text{scale} . T$. Additional leading dimensions (if any) will index batches.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_vector_sinh_arcsinh_diag

The (diagonal) SinhArcsinh transformation of a distribution on \mathbb{R}^k

Description

This distribution models a random vector $Y = (Y_1, \dots, Y_k)$, making use of a SinhArcsinh transformation (which has adjustable tailweight and skew), a rescaling, and a shift. The SinhArcsinh transformation of the Normal is described in great depth in [Sinh-arcsinh distributions](#). Here we use a slightly different parameterization, in terms of tailweight and skewness. Additionally we allow for distributions other than Normal, and control over scale as well as a "shift" parameter `loc`.

Usage

```
tfd_vector_sinh_arcsinh_diag(
  loc = NULL,
```

```

    scale_diag = NULL,
    scale_identity_multiplier = NULL,
    skewness = NULL,
    tailweight = NULL,
    distribution = NULL,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = "VectorSinhArcsinhDiag"
)

```

Arguments

<code>loc</code>	Floating-point Tensor. If this is set to NULL, loc is implicitly 0. When specified, may have shape $[B_1, \dots, B_b, k]$ where $b \geq 0$ and k is the event size.
<code>scale_diag</code>	Non-zero, floating-point Tensor representing a diagonal matrix added to scale. May have shape $[B_1, \dots, B_b, k]$, $b \geq 0$, and characterizes b -batches of $k \times k$ diagonal matrices added to scale. When both <code>scale_identity_multiplier</code> and <code>scale_diag</code> are NULL then scale is the Identity.
<code>scale_identity_multiplier</code>	Non-zero, floating-point Tensor representing a scale-identity-matrix added to scale. May have shape $[B_1, \dots, B_b]$, $b \geq 0$, and characterizes b -batches of scale $k \times k$ identity matrices added to scale. When both <code>scale_identity_multiplier</code> and <code>scale_diag</code> are NULL then scale is the Identity.
<code>skewness</code>	Skewness parameter. floating-point Tensor with shape broadcastable with <code>event_shape</code> .
<code>tailweight</code>	Tailweight parameter. floating-point Tensor with shape broadcastable with <code>event_shape</code> .
<code>distribution</code>	<code>tf\$Distributions\$Distribution</code> -like instance. Distribution from which k iid samples are used as input to transformation F . Default is <code>tfd_normal(loc = 0, scale = 1)</code> . Must be a scalar-batch, scalar-event distribution. Typically <code>distribution\$reparameterization_type = FULLY_REPARAMETERIZED</code> or it is a function of non-trainable parameters. WARNING: If you backprop through a <code>VectorSinhArcsinhDiag</code> sample and <code>distribution</code> is not <code>FULLY_REPARAMETERIZED</code> yet is a function of trainable variables, then the gradient will be incorrect!
<code>validate_args</code>	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
<code>allow_nan_stats</code>	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
<code>name</code>	name prefixed to Ops created by this class.

Details

Mathematical Details

Given iid random vector $Z = (Z_1, \dots, Z_k)$, we define the `VectorSinhArcsinhDiag` transformation of Z , Y , parameterized by $(loc, scale, skewness, tailweight)$, via the relation (with $@$ denoting matrix multiplication):

```

Y := loc + scale @ F(Z) * (2 / F_0(2))
F(Z) := Sinh( (Arcsinh(Z) + skewness) * tailweight )
F_0(Z) := Sinh( Arcsinh(Z) * tailweight )

```

This distribution is similar to the location-scale transformation $L(Z) := \text{loc} + \text{scale} @ Z$ in the following ways:

- If skewness = 0 and tailweight = 1 (the defaults), $F(Z) = Z$, and then $Y = L(Z)$ exactly.
- loc is used in both to shift the result by a constant factor.
- The multiplication of scale by $2 / F_0(2)$ ensures that if skewness = 0 $P[Y - \text{loc} \leq 2 * \text{scale}] = P[L(Z) - \text{loc} \leq 2 * \text{scale}]$. Thus it can be said that the weights in the tails of Y and $L(Z)$ beyond $\text{loc} + 2 * \text{scale}$ are the same. This distribution is different than $\text{loc} + \text{scale} @ Z$ due to the reshaping done by F :
 - Positive (negative) skewness leads to positive (negative) skew.
 - positive skew means, the mode of $F(Z)$ is "tilted" to the right.
 - positive skew means positive values of $F(Z)$ become more likely, and negative values become less likely.
 - Larger (smaller) tailweight leads to fatter (thinner) tails.
 - Fatter tails mean larger values of $|F(Z)|$ become more likely.
 - tailweight < 1 leads to a distribution that is "flat" around $Y = \text{loc}$, and a very steep drop-off in the tails.
 - tailweight > 1 leads to a distribution more peaked at the mode with heavier tails. To see the argument about the tails, note that for $|Z| \gg 1$ and $|Z| \gg (|\text{skewness}| * \text{tailweight})^{**\text{tailweight}}$, we have $Y \approx 0.5 Z^{**\text{tailweight}} e^{*(\text{sign}(Z) \text{skewness} * \text{tailweight})}$. To see the argument regarding multiplying scale by $2 / F_0(2)$,

$$\begin{aligned}
 P[(Y - \text{loc}) / \text{scale} \leq 2] &= P[F(Z) * (2 / F_0(2)) \leq 2] \\
 &= P[F(Z) \leq F_0(2)] \\
 &= P[Z \leq 2] \quad (\text{if } F = F_0).
 \end{aligned}$$

Value

a distribution instance.

See Also

For usage examples see e.g. [tfd_sample\(\)](#), [tfd_log_prob\(\)](#), [tfd_mean\(\)](#).

Other distributions: [tfd_autoregressive\(\)](#), [tfd_batch_reshape\(\)](#), [tfd_bates\(\)](#), [tfd_bernoulli\(\)](#), [tfd_beta_binomial\(\)](#), [tfd_beta\(\)](#), [tfd_binomial\(\)](#), [tfd_categorical\(\)](#), [tfd_cauchy\(\)](#), [tfd_chi2\(\)](#), [tfd_chi\(\)](#), [tfd_cholesky_lkj\(\)](#), [tfd_continuous_bernoulli\(\)](#), [tfd_deterministic\(\)](#), [tfd_dirichlet_multinomial\(\)](#), [tfd_dirichlet\(\)](#), [tfd_empirical\(\)](#), [tfd_exp_gamma\(\)](#), [tfd_exp_inverse_gamma\(\)](#), [tfd_exponential\(\)](#), [tfd_gamma_gamma\(\)](#), [tfd_gamma\(\)](#), [tfd_gaussian_process_regression_model\(\)](#), [tfd_gaussian_process\(\)](#), [tfd_generalized_normal\(\)](#), [tfd_geometric\(\)](#), [tfd_gumbel\(\)](#), [tfd_half_cauchy\(\)](#), [tfd_half_normal\(\)](#), [tfd_hidden_markov_model\(\)](#), [tfd_horseshoe\(\)](#), [tfd_independent\(\)](#), [tfd_inverse_gamma\(\)](#), [tfd_inverse_gaussian\(\)](#), [tfd_johnson_s_u\(\)](#), [tfd_joint_distribution_named_auto_batched\(\)](#), [tfd_joint_distribution_named\(\)](#), [tfd_joint_distribution_sequential_auto_batched\(\)](#), [tfd_joint_distribution_sequential\(\)](#), [tfd_kumaraswamy\(\)](#), [tfd_laplace\(\)](#), [tfd_linear_gaussian_state_space\(\)](#), [tfd_lkj\(\)](#), [tfd_log_logistic\(\)](#), [tfd_log_normal\(\)](#), [tfd_logistic\(\)](#), [tfd_mixture_same_family\(\)](#),

```
tfd_mixture(), tfd_multinomial(), tfd_multivariate_normal_diag_plus_low_rank(), tfd_multivariate_normal_diag_plus_low_rank_linear_operator(),
tfd_multivariate_normal_full_covariance(), tfd_multivariate_normal_linear_operator(),
tfd_multivariate_normal_tri_l(), tfd_multivariate_student_t_linear_operator(), tfd_negative_binomial(),
tfd_normal(), tfd_one_hot_categorical(), tfd_pareto(), tfd_pixel_cnn(), tfd_poisson_log_normal_quadrature(),
tfd_poisson(), tfd_power_spherical(), tfd_probit_bernoulli(), tfd_quantized(), tfd_relaxed_bernoulli(),
tfd_relaxed_one_hot_categorical(), tfd_sample_distribution(), tfd_sinh_arcsinh(),
tfd_skellam(), tfd_spherical_uniform(), tfd_student_t_process(), tfd_student_t(), tfd_transformed_distribution(),
tfd_triangular(), tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian(),
tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_von_mises_fisher(),
tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(), tfd_wishart_tri_l(),
tfd_wishart(), tfd_zipf()
```

tfd_von_mises

The von Mises distribution over angles

Description

The von Mises distribution is a univariate directional distribution. Similarly to Normal distribution, it is a maximum entropy distribution. The samples of this distribution are angles, measured in radians. They are 2π -periodic: $x = 0$ and $x = 2\pi$ are equivalent. This means that the density is also 2π -periodic. The generated samples, however, are guaranteed to be in $[-\pi, \pi]$ range. When concentration = 0, this distribution becomes a Uniform distribution on the $[-\pi, \pi]$ domain.

Usage

```
tfd_von_mises(
  loc,
  concentration,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "VonMises"
)
```

Arguments

loc	Floating point tensor, the circular means of the distribution(s).
concentration	Floating point tensor, the level of concentration of the distribution(s) around loc. Must take non-negative values. concentration = 0 defines a Uniform distribution, while concentration = +inf indicates a Deterministic distribution at loc.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

The von Mises distribution is a special case of von Mises-Fisher distribution for $n=2$. However, the TFP's VonMisesFisher implementation represents the samples and location as (x, y) points on a circle, while VonMises represents them as scalar angles.

Mathematical details The probability density function (pdf) of this distribution is,

$$\text{pdf}(x; \text{loc}, \text{concentration}) = \frac{\exp(\text{concentration} \cos(x - \text{loc}))}{Z}$$

$$Z = 2 * \pi * I_0(\text{concentration})$$

where:

- $I_0(\text{concentration})$ is the modified Bessel function of order zero;
- loc the circular mean of the distribution, a scalar. It can take arbitrary values, but it is 2π -periodic: loc and $\text{loc} + 2\pi$ result in the same distribution.
- $\text{concentration} \geq 0$ parameter is the concentration parameter. When $\text{concentration} = 0$, this distribution becomes a Uniform distribution on $[-\pi, \pi)$.

The parameters loc and concentration must be shaped in a way that supports broadcasting (e.g. $\text{loc} + \text{concentration}$ is a valid operation).

Value

a distribution instance.

See Also

For usage examples see e.g. [tfd_sample\(\)](#), [tfd_log_prob\(\)](#), [tfd_mean\(\)](#).

Other distributions: [tfd_autoregressive\(\)](#), [tfd_batch_reshape\(\)](#), [tfd_bates\(\)](#), [tfd_bernoulli\(\)](#), [tfd_beta_binomial\(\)](#), [tfd_beta\(\)](#), [tfd_binomial\(\)](#), [tfd_categorical\(\)](#), [tfd_cauchy\(\)](#), [tfd_chi2\(\)](#), [tfd_chi\(\)](#), [tfd_cholesky_lkj\(\)](#), [tfd_continuous_bernoulli\(\)](#), [tfd_deterministic\(\)](#), [tfd_dirichlet_multinomial\(\)](#), [tfd_dirichlet\(\)](#), [tfd_empirical\(\)](#), [tfd_exp_gamma\(\)](#), [tfd_exp_inverse_gamma\(\)](#), [tfd_exponential\(\)](#), [tfd_gamma_gamma\(\)](#), [tfd_gamma\(\)](#), [tfd_gaussian_process_regression_model\(\)](#), [tfd_gaussian_process\(\)](#), [tfd_generalized_normal\(\)](#), [tfd_geometric\(\)](#), [tfd_gumbel\(\)](#), [tfd_half_cauchy\(\)](#), [tfd_half_normal\(\)](#), [tfd_hidden_markov_model\(\)](#), [tfd_horseshoe\(\)](#), [tfd_independent\(\)](#), [tfd_inverse_gamma\(\)](#), [tfd_inverse_gaussian\(\)](#), [tfd_johnson_s_u\(\)](#), [tfd_joint_distribution_named_auto_batched\(\)](#), [tfd_joint_distribution_named\(\)](#), [tfd_joint_distribution_sequential_auto_batched\(\)](#), [tfd_joint_distribution_sequential\(\)](#), [tfd_kumaraswamy\(\)](#), [tfd_laplace\(\)](#), [tfd_linear_gaussian_state_space\(\)](#), [tfd_lkj\(\)](#), [tfd_log_logistic\(\)](#), [tfd_log_normal\(\)](#), [tfd_logistic\(\)](#), [tfd_mixture_same_family\(\)](#), [tfd_mixture\(\)](#), [tfd_multinomial\(\)](#), [tfd_multivariate_normal_diag_plus_low_rank\(\)](#), [tfd_multivariate_normal_diag\(\)](#), [tfd_multivariate_normal_full_covariance\(\)](#), [tfd_multivariate_normal_linear_operator\(\)](#), [tfd_multivariate_normal_tri_l\(\)](#), [tfd_multivariate_student_t_linear_operator\(\)](#), [tfd_negative_binomial\(\)](#), [tfd_normal\(\)](#), [tfd_one_hot_categorical\(\)](#), [tfd_pareto\(\)](#), [tfd_pixel_cnn\(\)](#), [tfd_poisson_log_normal_quadrature\(\)](#), [tfd_poisson\(\)](#), [tfd_power_spherical\(\)](#), [tfd_probit_bernoulli\(\)](#), [tfd_quantized\(\)](#), [tfd_relaxed_bernoulli\(\)](#), [tfd_relaxed_one_hot_categorical\(\)](#), [tfd_sample_distribution\(\)](#), [tfd_sinh_arcsinh\(\)](#), [tfd_skellam\(\)](#), [tfd_spherical_uniform\(\)](#), [tfd_student_t_process\(\)](#), [tfd_student_t\(\)](#), [tfd_transformed_distribution\(\)](#), [tfd_triangular\(\)](#), [tfd_truncated_cauchy\(\)](#), [tfd_truncated_normal\(\)](#), [tfd_uniform\(\)](#), [tfd_variational_gaussian\(\)](#), [tfd_vector_diffeomixture\(\)](#), [tfd_vector_exponential_diag\(\)](#), [tfd_vector_exponential_linear_operator\(\)](#), [tfd_vector_laplace_diag\(\)](#), [tfd_vector_laplace_linear_operator\(\)](#), [tfd_vector_sinh_arcsinh_diag\(\)](#),

[tfd_von_mises_fisher\(\)](#), [tfd_weibull\(\)](#), [tfd_wishart_linear_operator\(\)](#), [tfd_wishart_tri_l\(\)](#), [tfd_wishart\(\)](#), [tfd_zipf\(\)](#)

tfd_von_mises_fisher *The von Mises-Fisher distribution over unit vectors on S^{n-1}*

Description

The von Mises-Fisher distribution is a directional distribution over vectors on the unit hypersphere S^{n-1} embedded in n dimensions (\mathbb{R}^n).

Usage

```
tfd_von_mises_fisher(
    mean_direction,
    concentration,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = "VonMisesFisher"
)
```

Arguments

mean_direction	Floating-point Tensor with shape $[B_1, \dots, B_n, D]$. A unit vector indicating the mode of the distribution, or the unit-normalized direction of the mean. (This is <i>not</i> in general the mean of the distribution; the mean is not generally in the support of the distribution.) NOTE: D is currently restricted to ≤ 5 .
concentration	Floating-point Tensor having batch shape $[B_1, \dots, B_n]$ broadcastable with mean_direction. The level of concentration of samples around the mean_direction. concentration=0 indicates a uniform distribution over the unit hypersphere, and concentration=+inf indicates a Deterministic distribution (delta function) at mean_direction.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details

Mathematical details The probability density function (pdf) is,

$$\text{pdf}(x; \mu, \kappa) = C(\kappa) \exp(\kappa * \mu^T x)$$

where,

$$C(\kappa) = (2 \pi)^{-n/2} \kappa^{n/2-1} / I_{n/2-1}(\kappa),$$

$I_v(z)$ being the modified Bessel function of the first kind of order v

where:

- `mean_direction` = μ ; a unit vector in \mathbb{R}^k ,
- `concentration` = κ ; scalar real ≥ 0 , concentration of samples around `mean_direction`, where 0 pertains to the uniform distribution on the hypersphere, and `inf` indicates a delta function at `mean_direction`.

NOTE: Currently only `n` in 2, 3, 4, 5 are supported. For `n=5` some numerical instability can occur for low concentrations ($<.01$).

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

Description

The probability density function (pdf) of this distribution is,

$$\text{pdf}(x; \text{lambda}, k) = k / \text{lambda} * (x / \text{lambda}) ** (k - 1) * \exp(-(x / \text{lambda}) ** k)$$

where concentration = k and scale = lambda. The cumulative density function of this distribution is,

$$\text{cdf}(x; \text{lambda}, k) = 1 - \exp(-(x / \text{lambda}) ** k)$$

The Weibull distribution includes the Exponential and Rayleigh distributions as special cases:

$$\text{Exponential}(\text{rate}) = \text{Weibull}(\text{concentration}=1., 1. / \text{rate})$$

$$\text{Rayleigh}(\text{scale}) = \text{Weibull}(\text{concentration}=2., \text{sqrt}(2.) * \text{scale})$$

Usage

```
tfd_weibull(
    concentration,
    scale,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = "Weibull"
)
```

Arguments

concentration	Positive Float-type Tensor, the concentration param of the distribution. Must contain only positive values.
scale	Positive Float-type Tensor, the scale param of the distribution. Must contain only positive values.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag_plus_low_rank_linear_operator()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

tfd_wishart

The matrix Wishart distribution on positive definite matrices

Description

This distribution is defined by a scalar number of degrees of freedom `df` and an instance of `LinearOperator`, which provides matrix-free access to a symmetric positive definite operator, which defines the scale matrix.

Usage

```
tfd_wishart(
    df,
    scale = NULL,
    scale_tril = NULL,
    input_output_cholesky = FALSE,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = "Wishart"
)
```

Arguments

df	float or double tensor, the degrees of freedom of the distribution(s). df must be greater than or equal to k.
scale	float or double Tensor. The symmetric positive definite scale matrix of the distribution. Exactly one of scale and 'scale_tril must be passed.
scale_tril	float or double Tensor. The Cholesky factorization of the symmetric positive definite scale matrix of the distribution. Exactly one of scale and 'scale_tril must be passed.
input_output_cholesky	Logical. If TRUE, functions whose input or output have the semantics of samples assume inputs are in Cholesky form and return outputs in Cholesky form. In particular, if this flag is TRUE, input to log_prob is presumed of Cholesky form and output from sample, mean, and mode are of Cholesky form. Setting this argument to TRUE is purely a computational optimization and does not change the underlying distribution; for instance, mean returns the Cholesky of the mean, not the mean of Cholesky factors. The variance and stddev methods are unaffected by this flag. Default value: FALSE (i.e., input/output does not have Cholesky semantics).
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details**Mathematical Details**

The probability density function (pdf) is,

$$\text{pdf}(X; \text{df}, \text{scale}) = \frac{\det(X)^{0.5(\text{df}-k-1)} \exp(-0.5 \text{tr}[\text{inv}(\text{scale}) X])}{Z}$$

$$Z = 2^{0.5 \text{df} k} |\det(\text{scale})|^{0.5 \text{df}} \Gamma_k(0.5 \text{df})$$

where:

- df >= k denotes the degrees of freedom,
- scale is a symmetric, positive definite, k x k matrix,
- Z is the normalizing constant, and,
- Gamma_k is the **multivariate Gamma function**.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart_tri_l()`, `tfd_zipf()`

tfd_wishart_linear_operator

The matrix Wishart distribution on positive definite matrices

Description

This distribution is defined by a scalar number of degrees of freedom `df` and an instance of `LinearOperator`, which provides matrix-free access to a symmetric positive definite operator, which defines the scale matrix.

Usage

```
tfd_wishart_linear_operator(
    df,
    scale,
    input_output_cholesky = FALSE,
    validate_args = FALSE,
    allow_nan_stats = TRUE,
    name = "WishartLinearOperator"
)
```

Arguments

df	float or double tensor, the degrees of freedom of the distribution(s). df must be greater than or equal to k.
scale	float or double instance of LinearOperator.
input_output_cholesky	Logical. If TRUE, functions whose input or output have the semantics of samples assume inputs are in Cholesky form and return outputs in Cholesky form. In particular, if this flag is TRUE, input to log_prob is presumed of Cholesky form and output from sample, mean, and mode are of Cholesky form. Setting this argument to TRUE is purely a computational optimization and does not change the underlying distribution; for instance, mean returns the Cholesky of the mean, not the mean of Cholesky factors. The variance and stddev methods are unaffected by this flag. Default value: FALSE (i.e., input/output does not have Cholesky semantics).
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details**Mathematical Details**

The probability density function (pdf) is,

$$\text{pdf}(X; \text{df}, \text{scale}) = \frac{\det(X)^{0.5(\text{df}-k-1)} \exp(-0.5 \text{tr}[\text{inv}(\text{scale}) X])}{Z}$$

$$Z = 2^{0.5 \text{df} k} |\det(\text{scale})|^{0.5 \text{df}} \Gamma_k(0.5 \text{df})$$

where:

- df >= k denotes the degrees of freedom,
- scale is a symmetric, positive definite, k x k matrix,
- Z is the normalizing constant, and,
- Gamma_k is the **multivariate Gamma function**.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_tri_l()`, `tfd_wishart()`, `tfd_zipf()`

`tfd_wishart_tri_l`

The matrix Wishart distribution parameterized with Cholesky factors.

Description

This distribution is defined by a scalar degrees of freedom `df` and a scale matrix, expressed as a lower triangular Cholesky factor.

Usage

```
tfd_wishart_tri_l(
  df,
  scale_tril,
  input_output_cholesky = FALSE,
  validate_args = FALSE,
  allow_nan_stats = TRUE,
  name = "WishartTriL"
)
```


Arguments

df	float or double tensor, the degrees of freedom of the distribution(s). df must be greater than or equal to k.
scale_tril	float or double Tensor. The Cholesky factorization of the symmetric positive definite scale matrix of the distribution.
input_output_cholesky	Logical. If TRUE, functions whose input or output have the semantics of samples assume inputs are in Cholesky form and return outputs in Cholesky form. In particular, if this flag is TRUE, input to log_prob is presumed of Cholesky form and output from sample, mean, and mode are of Cholesky form. Setting this argument to TRUE is purely a computational optimization and does not change the underlying distribution; for instance, mean returns the Cholesky of the mean, not the mean of Cholesky factors. The variance and stddev methods are unaffected by this flag. Default value: FALSE (i.e., input/output does not have Cholesky semantics).
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Logical, default TRUE. When TRUE, statistics (e.g., mean, mode, variance) use the value NaN to indicate the result is undefined. When FALSE, an exception is raised if one or more of the statistic's batch members are undefined.
name	name prefixed to Ops created by this class.

Details**Mathematical Details**

The probability density function (pdf) is,

$$\text{pdf}(X; \text{df}, \text{scale}) = \frac{\det(X)^{0.5(\text{df}-k-1)} \exp(-0.5 \text{tr}[\text{inv}(\text{scale}) X])}{Z} \\ Z = 2^{0.5 \text{df} k} |\det(\text{scale})|^{0.5 \text{df}} \text{Gamma}_k(0.5 \text{df})$$

where:

- df \geq k denotes the degrees of freedom,
- scale is a symmetric, positive definite, k x k matrix,
- Z is the normalizing constant, and,
- Gamma_k is the **multivariate Gamma function**.

Value

a distribution instance.

See Also

For usage examples see e.g. `tfd_sample()`, `tfd_log_prob()`, `tfd_mean()`.

Other distributions: `tfd_autoregressive()`, `tfd_batch_reshape()`, `tfd_bates()`, `tfd_bernoulli()`, `tfd_beta_binomial()`, `tfd_beta()`, `tfd_binomial()`, `tfd_categorical()`, `tfd_cauchy()`, `tfd_chi2()`, `tfd_chi()`, `tfd_cholesky_lkj()`, `tfd_continuous_bernoulli()`, `tfd_deterministic()`, `tfd_dirichlet_multinomial()`, `tfd_dirichlet()`, `tfd_empirical()`, `tfd_exp_gamma()`, `tfd_exp_inverse_gamma()`, `tfd_exponential()`, `tfd_gamma_gamma()`, `tfd_gamma()`, `tfd_gaussian_process_regression_model()`, `tfd_gaussian_process()`, `tfd_generalized_normal()`, `tfd_geometric()`, `tfd_gumbel()`, `tfd_half_cauchy()`, `tfd_half_normal()`, `tfd_hidden_markov_model()`, `tfd_horseshoe()`, `tfd_independent()`, `tfd_inverse_gamma()`, `tfd_inverse_gaussian()`, `tfd_johnson_s_u()`, `tfd_joint_distribution_named_auto_batched()`, `tfd_joint_distribution_named()`, `tfd_joint_distribution_sequential_auto_batched()`, `tfd_joint_distribution_sequential()`, `tfd_kumaraswamy()`, `tfd_laplace()`, `tfd_linear_gaussian_state_space()`, `tfd_lkj()`, `tfd_log_logistic()`, `tfd_log_normal()`, `tfd_logistic()`, `tfd_mixture_same_family()`, `tfd_mixture()`, `tfd_multinomial()`, `tfd_multivariate_normal_diag_plus_low_rank()`, `tfd_multivariate_normal_diag()`, `tfd_multivariate_normal_full_covariance()`, `tfd_multivariate_normal_linear_operator()`, `tfd_multivariate_normal_tri_l()`, `tfd_multivariate_student_t_linear_operator()`, `tfd_negative_binomial()`, `tfd_normal()`, `tfd_one_hot_categorical()`, `tfd_pareto()`, `tfd_pixel_cnn()`, `tfd_poisson_log_normal_quadrature()`, `tfd_poisson()`, `tfd_power_spherical()`, `tfd_probit_bernoulli()`, `tfd_quantized()`, `tfd_relaxed_bernoulli()`, `tfd_relaxed_one_hot_categorical()`, `tfd_sample_distribution()`, `tfd_sinh_arcsinh()`, `tfd_skellam()`, `tfd_spherical_uniform()`, `tfd_student_t_process()`, `tfd_student_t()`, `tfd_transformed_distribution()`, `tfd_triangular()`, `tfd_truncated_cauchy()`, `tfd_truncated_normal()`, `tfd_uniform()`, `tfd_variational_gaussian()`, `tfd_vector_diffeomixture()`, `tfd_vector_exponential_diag()`, `tfd_vector_exponential_linear_operator()`, `tfd_vector_laplace_diag()`, `tfd_vector_laplace_linear_operator()`, `tfd_vector_sinh_arcsinh_diag()`, `tfd_von_mises_fisher()`, `tfd_von_mises()`, `tfd_weibull()`, `tfd_wishart_linear_operator()`, `tfd_wishart()`, `tfd_zipf()`

tfd_zipf

Zipf distribution

Description

The Zipf distribution is parameterized by a power parameter.

Usage

```
tfd_zipf(
  power,
  dtype = tf$int32,
  interpolate_nondiscrete = TRUE,
  sample_maximum_iterations = 100,
  validate_args = FALSE,
  allow_nan_stats = FALSE,
  name = "Zipf"
)
```

Arguments

power	Float like Tensor representing the power parameter. Must be strictly greater than 1.
dtype	The dtype of Tensor returned by sample. Default value: tf\$int32.
interpolate_nondiscrete	Logical. When FALSE, log_prob returns -inf (and prob returns 0) for non-integer inputs. When TRUE, log_prob evaluates the continuous function $-\text{power} \log(k) - \log(\zeta(\text{power}))$, which matches the Zipf pmf at integer arguments k (note that this function is not itself a normalized probability log-density). Default value: TRUE.
sample_maximum_iterations	Maximum number of iterations of allowable iterations in sample. When validate_args=TRUE, samples which fail to reach convergence (subject to this cap) are masked out with self\$dtype\$min or nan depending on self\$dtype\$is_integer. Default value: 100.
validate_args	Logical, default FALSE. When TRUE distribution parameters are checked for validity despite possibly degrading runtime performance. When FALSE invalid inputs may silently render incorrect outputs. Default value: FALSE.
allow_nan_stats	Default value: FALSE.
name	name prefixed to Ops created by this class.

Details

Mathematical Details The probability mass function (pmf) is,

$$\text{pmf}(k; \alpha, k \geq 0) = (k^{-\alpha}) / Z$$

$$Z = \zeta(\alpha).$$

where power = alpha and Z is the normalization constant. zeta is the [Riemann zeta function](#). Note that gradients with respect to the power parameter are not supported in the current implementation.

Value

a distribution instance.

See Also

For usage examples see e.g. [tfd_sample\(\)](#), [tfd_log_prob\(\)](#), [tfd_mean\(\)](#).

Other distributions: [tfd_autoregressive\(\)](#), [tfd_batch_reshape\(\)](#), [tfd_bates\(\)](#), [tfd_bernoulli\(\)](#), [tfd_beta_binomial\(\)](#), [tfd_beta\(\)](#), [tfd_binomial\(\)](#), [tfd_categorical\(\)](#), [tfd_cauchy\(\)](#), [tfd_chi2\(\)](#), [tfd_chi\(\)](#), [tfd_cholesky_lkj\(\)](#), [tfd_continuous_bernoulli\(\)](#), [tfd_deterministic\(\)](#), [tfd_dirichlet_multinomial\(\)](#), [tfd_dirichlet\(\)](#), [tfd_empirical\(\)](#), [tfd_exp_gamma\(\)](#), [tfd_exp_inverse_gamma\(\)](#), [tfd_exponential\(\)](#), [tfd_gamma_gamma\(\)](#), [tfd_gamma\(\)](#), [tfd_gaussian_process_regression_model\(\)](#), [tfd_gaussian_process\(\)](#), [tfd_generalized_normal\(\)](#), [tfd_geometric\(\)](#), [tfd_gumbel\(\)](#), [tfd_half_cauchy\(\)](#), [tfd_half_normal\(\)](#), [tfd_hidden_markov_model\(\)](#), [tfd_horseshoe\(\)](#), [tfd_independent\(\)](#), [tfd_inverse_gamma\(\)](#), [tfd_inverse_gaussian\(\)](#), [tfd_johnson_s_u\(\)](#), [tfd_joint_distribution_named_auto_batched\(\)](#),

```

tfd_joint_distribution_named(), tfd_joint_distribution_sequential_auto_batched(),
tfd_joint_distribution_sequential(), tfd_kumaraswamy(), tfd_laplace(), tfd_linear_gaussian_state_space,
tfd_lkj(), tfd_log_logistic(), tfd_log_normal(), tfd_logistic(), tfd_mixture_same_family(),
tfd_mixture(), tfd_multinomial(), tfd_multivariate_normal_diag_plus_low_rank(), tfd_multivariate_normal,
tfd_multivariate_normal_full_covariance(), tfd_multivariate_normal_linear_operator(),
tfd_multivariate_normal_tri_l(), tfd_multivariate_student_t_linear_operator(), tfd_negative_binomial(),
tfd_normal(), tfd_one_hot_categorical(), tfd_pareto(), tfd_pixel_cnn(), tfd_poisson_log_normal_quadrature,
tfd_poisson(), tfd_power_spherical(), tfd_probit_bernoulli(), tfd_quantized(), tfd_relaxed_bernoulli(),
tfd_relaxed_one_hot_categorical(), tfd_sample_distribution(), tfd_sinh_arcsinh(),
tfd_skellam(), tfd_spherical_uniform(), tfd_student_t_process(), tfd_student_t(), tfd_transformed_distribution,
tfd_triangular(), tfd_truncated_cauchy(), tfd_truncated_normal(), tfd_uniform(), tfd_variational_gaussian,
tfd_vector_diffeomixture(), tfd_vector_exponential_diag(), tfd_vector_exponential_linear_operator(),
tfd_vector_laplace_diag(), tfd_vector_laplace_linear_operator(), tfd_vector_sinh_arcsinh_diag(),
tfd_von_mises_fisher(), tfd_von_mises(), tfd_weibull(), tfd_wishart_linear_operator(),
tfd_wishart_tri_l(), tfd_wishart()

```

tfp	<i>Handle to the tensorflow_probability module</i>
-----	--

Description

Handle to the tensorflow_probability module

Usage

```
tfp
```

Format

An object of class `python.builtin.module` (inherits from `python.builtin.object`) of length 0.

Value

Module(tensorflow_probability)

tfp_version	<i>TensorFlow Probability Version</i>
-------------	---------------------------------------

Description

TensorFlow Probability Version

Usage

```
tfp_version()
```

Value

the Python TFP version

vi_amari_alpha	<i>The Amari-alpha Csiszar-function in log-space</i>
----------------	--

Description

A Csiszar-function is a member of $F = \{ f: \mathbb{R}_+ \rightarrow \mathbb{R} : f \text{ convex} \}$.

Usage

```
vi_amari_alpha(logu, alpha = 1, self_normalized = FALSE, name = NULL)
```

Arguments

logu	float-like Tensor representing $\log(u)$ from above.
alpha	float-like scalar.
self_normalized	logical indicating whether $f'(u=1)=0$. When $f'(u=1)=0$ the implied Csiszar f-Divergence remains non-negative even when p, q are unnormalized measures.
name	name prefixed to Ops created by this function.

Details

When `self_normalized = TRUE`, the Amari-alpha Csiszar-function is:

$$f(u) = \begin{cases} -\log(u) + (u - 1), & \alpha = 0 \\ u \log(u) - (u - 1), & \alpha = 1 \\ ((u^\alpha - 1) - \alpha (u - 1) / (\alpha (\alpha - 1))), & \text{otherwise} \end{cases}$$

When `self_normalized = FALSE` the $(u - 1)$ terms are omitted.

Warning: when $\alpha \neq 0$ and/or `self_normalized = True` this function makes non-log-space calculations and may therefore be numerically unstable for $|\log u| \gg 0$.

Value

amari_alpha_of_u float-like Tensor of the Csiszar-function evaluated at $u = \exp(\log u)$.

References

- A. Cichocki and S. Amari. "Families of Alpha-Beta-and GammaDivergences: Flexible and Robust Measures of Similarities." *Entropy*, vol. 12, no. 6, pp. 1532-1568, 2010.

See Also

Other vi-functions: [vi_arithmetic_geometric\(\)](#), [vi_chi_square\(\)](#), [vi_csiszar_vimco\(\)](#), [vi_dual_csiszar_function\(\)](#), [vi_fit_surrogate_posterior\(\)](#), [vi_jeffreys\(\)](#), [vi_jensen_shannon\(\)](#), [vi_kl_forward\(\)](#), [vi_kl_reverse\(\)](#), [vi_log1p_abs\(\)](#), [vi_modified_gan\(\)](#), [vi_monte_carlo_variational_loss\(\)](#), [vi_pearson\(\)](#), [vi_squared_hellinger\(\)](#), [vi_symmetrized_csiszar_function\(\)](#)

vi_arithmetic_geometric

The Arithmetic-Geometric Csiszar-function in log-space

Description

A Csiszar-function is a member of $F = \{ f: \mathbb{R}_+ \rightarrow \mathbb{R} : f \text{ convex} \}$.

Usage

```
vi_arithmetic_geometric(logu, self_normalized = FALSE, name = NULL)
```

Arguments

logu	float-like Tensor representing $\log(u)$ from above.
self_normalized	logical indicating whether $f(u=1)=0$. When $f(u=1)=0$ the implied Csiszar f-Divergence remains non-negative even when p, q are unnormalized measures.
name	name prefixed to Ops created by this function.

Details

When `self_normalized = True` the Arithmetic-Geometric Csiszar-function is:

$$f(u) = (1 + u) \log\left(\frac{1 + u}{\sqrt{u}}\right) - (1 + u) \log(2)$$

When `self_normalized = False` the $(1 + u) \log(2)$ term is omitted.

Observe that as an f-Divergence, this Csiszar-function implies:

$$D_f[p, q] = KL[m, p] + KL[m, q]$$

$$m(x) = 0.5 p(x) + 0.5 q(x)$$

In a sense, this divergence is the "reverse" of the Jensen-Shannon f-Divergence. This Csiszar-function induces a symmetric f-Divergence, i.e., $D_f[p, q] = D_f[q, p]$.

Warning: when `self_normalized = True` this function makes non-log-space calculations and may therefore be numerically unstable for $\log u \gg 0$.

Value

arithmetic_geometric_of_u: float-like Tensor of the Csiszar-function evaluated at $u = \exp(\log u)$.

See Also

Other vi-functions: [vi_amari_alpha\(\)](#), [vi_chi_square\(\)](#), [vi_csiszar_vimco\(\)](#), [vi_dual_csiszar_function\(\)](#), [vi_fit_surrogate_posterior\(\)](#), [vi_jeffreys\(\)](#), [vi_jensen_shannon\(\)](#), [vi_kl_forward\(\)](#), [vi_kl_reverse\(\)](#), [vi_log1p_abs\(\)](#), [vi_modified_gan\(\)](#), [vi_monte_carlo_variational_loss\(\)](#), [vi_pearson\(\)](#), [vi_squared_hellinger\(\)](#), [vi_symmetrized_csiszar_function\(\)](#)

 vi_chi_square

The chi-square Csiszar-function in log-space

Description

A Csiszar-function is a member of $F = \{ f: \mathbb{R}_+ \text{ to } \mathbb{R} : f \text{ convex} \}$.

Usage

```
vi_chi_square(logu, name = NULL)
```

Arguments

logu float-like Tensor representing $\log(u)$ from above.
 name name prefixed to Ops created by this function.

Details

The Chi-square Csiszar-function is:

$$f(u) = u^{**2} - 1$$

Warning: this function makes non-log-space calculations and may therefore be numerically unstable for $\text{llogul} \gg 0$.

Value

chi_square_of_u: float-like Tensor of the Csiszar-function evaluated at $u = \exp(\text{logu})$.

See Also

Other vi-functions: [vi_amari_alpha\(\)](#), [vi_arithmetic_geometric\(\)](#), [vi_csiszar_vimco\(\)](#), [vi_dual_csiszar_function\(\)](#), [vi_fit_surrogate_posterior\(\)](#), [vi_jeffreys\(\)](#), [vi_jensen_shannon\(\)](#), [vi_kl_forward\(\)](#), [vi_kl_reverse\(\)](#), [vi_log1p_abs\(\)](#), [vi_modified_gan\(\)](#), [vi_monte_carlo_variational_loss\(\)](#), [vi_pearson\(\)](#), [vi_squared_hellinger\(\)](#), [vi_symmetrized_csiszar_function\(\)](#)

vi_csiszar_vimco	<i>Use VIMCO to lower the variance of the gradient of csiszar_function(Avg(logu))</i>
------------------	---

Description

This function generalizes VIMCO (Mnih and Rezende, 2016) to Csiszar f-Divergences.

Usage

```
vi_csiszar_vimco(
  f,
  p_log_prob,
  q,
  num_draws,
  num_batch_draws = 1,
  seed = NULL,
  name = NULL
)
```

Arguments

f	function representing a Csiszar-function in log-space.
p_log_prob	function representing the natural-log of the probability under distribution p. (In variational inference p is the joint distribution.)
q	tfd\$Distribution-like instance; must implement: sample(n, seed), and log_prob(x). (In variational inference q is the approximate posterior distribution.)
num_draws	Integer scalar number of draws used to approximate the f-Divergence expectation.
num_batch_draws	Integer scalar number of draws used to approximate the f-Divergence expectation.
seed	integer seed for q\$sample.
name	String prefixed to Ops created by this function.

Details

Note: if q.reparameterization_type = tfd.FULLY_REPARAMETERIZED, consider using monte_carlo_csiszar_f_diverg

The VIMCO loss is:

```
vimco = f(Avg{logu[i] : i=0,...,m-1})
where,
logu[i] = log( p(x, h[i]) / q(h[i] | x) )
h[i] iid~ q(H | x)
```

Interestingly, the VIMCO gradient is not the naive gradient of vimco. Rather, it is characterized by:

grad[vimco] - variance_reducing_term

where,

variance_reducing_term = Sum{ grad[log q(h[i] | x)] * (vimco - f(log Avg{h[j;i] : j=0,...,m-1})) #' : i=0
h[j;i] = u[j] for j!=i, GeometricAverage{ u[k] : k!=i} for j==i

(We omitted stop_gradient for brevity. See implementation for more details.) The Avg{h[j;i] : j} term is a kind of "swap-out average" where the i-th element has been replaced by the leave-i-out Geometric-average.

This implementation prefers numerical precision over efficiency, i.e., $O(\text{num_draws} * \text{num_batch_draws} * \text{prod}(\text{batch_shape}) * \text{prod}(\text{event_shape}))$. (The constant may be fairly large, perhaps around 12.)

Value

vimco The Csiszar f-Divergence generalized VIMCO objective

References

- [Andriy Mnih and Danilo Rezende. Variational Inference for Monte Carlo objectives. In *International Conference on Machine Learning*, 2016.](#)

See Also

Other vi-functions: [vi_amari_alpha\(\)](#), [vi_arithmetic_geometric\(\)](#), [vi_chi_square\(\)](#), [vi_dual_csiszar_function\(\)](#), [vi_fit_surrogate_posterior\(\)](#), [vi_jeffreys\(\)](#), [vi_jensen_shannon\(\)](#), [vi_kl_forward\(\)](#), [vi_kl_reverse\(\)](#), [vi_log1p_abs\(\)](#), [vi_modified_gan\(\)](#), [vi_monte_carlo_variational_loss\(\)](#), [vi_pearson\(\)](#), [vi_squared_hellinger\(\)](#), [vi_symmetrized_csiszar_function\(\)](#)

vi_dual_csiszar_function

Calculates the dual Csiszar-function in log-space

Description

A Csiszar-function is a member of $F = \{ f: \mathbb{R}_+ \text{ to } \mathbb{R} : f \text{ convex} \}$.

Usage

```
vi_dual_csiszar_function(logu, csiszar_function, name = NULL)
```

Arguments

logu float-like Tensor representing log(u) from above.
 csiszar_function function representing a Csiszar-function over log-domain.
 name name prefixed to Ops created by this function.

Details

The Csiszar-dual is defined as:

$$f^{*}(u) = u f(1 / u)$$

where f is some other Csiszar-function. For example, the dual of `kl_reverse` is `kl_forward`, i.e.,

$$\begin{aligned} f(u) &= -\log(u) \\ f^{*}(u) &= u f(1 / u) = -u \log(1 / u) = u \log(u) \end{aligned}$$

The dual of the dual is the original function:

$$f^{**}(u) = \{u f(1/u)\}^{*}(u) = u (1/u) f(1/(1/u)) = f(u)$$

Warning: this function makes non-log-space calculations and may therefore be numerically unstable for $|\log u| \gg 0$.

Value

`dual_f_of_u` float-like Tensor of the result of calculating the dual of f at $u = \exp(\log u)$.

See Also

Other vi-functions: [vi_amari_alpha\(\)](#), [vi_arithmetic_geometric\(\)](#), [vi_chi_square\(\)](#), [vi_csiszar_vimco\(\)](#), [vi_fit_surrogate_posterior\(\)](#), [vi_jeffreys\(\)](#), [vi_jensen_shannon\(\)](#), [vi_kl_forward\(\)](#), [vi_kl_reverse\(\)](#), [vi_log1p_abs\(\)](#), [vi_modified_gan\(\)](#), [vi_monte_carlo_variational_loss\(\)](#), [vi_pearson\(\)](#), [vi_squared_hellinger\(\)](#), [vi_symmetrized_csiszar_function\(\)](#)

`vi_fit_surrogate_posterior`

Fit a surrogate posterior to a target (unnormalized) log density

Description

The default behavior constructs and minimizes the negative variational evidence lower bound (ELBO), given by `q_samples <- surrogate_posterior$sample(num_draws)` `elbo_loss <- -tf$reduce_mean(target_log_prob_fn(q_samples) surrogate_posterior$log_prob(q_samples))`

Usage

```
vi_fit_surrogate_posterior(
  target_log_prob_fn,
  surrogate_posterior,
  optimizer,
  num_steps,
  convergence_criterion = NULL,
  trace_fn = tfp$vi$optimization$`_trace_loss`,
```

```

    variational_loss_fn = NULL,
    discrepancy_fn = tfp$vi$kl_reverse,
    sample_size = 1,
    importance_sample_size = 1,
    trainable_variables = NULL,
    jit_compile = NULL,
    seed = NULL,
    name = "fit_surrogate_posterior"
  )

```

Arguments

`target_log_prob_fn`

function that takes a set of Tensor arguments and returns a Tensor log-density. Given `q_sample <- surrogate_posterior$sample(sample_size)`, this will be (in Python) called as `target_log_prob_fn(q_sample)` if `q_sample` is a list or a tuple, `target_log_prob_fn(**q_sample)` if `q_sample` is a dictionary, or `target_log_prob_fn(q_sample)` if `q_sample` is a Tensor. It should support batched evaluation, i.e., should return a result of shape `[sample_size]`.

`surrogate_posterior`

A `tfp$distributions$Distribution` instance defining a variational posterior (could be a `tfp$distributions$JointDistribution`). Crucially, the distribution's `log_prob` and (if reparameterized) `sample` methods must directly invoke all ops that generate gradients to the underlying variables. One way to ensure this is to use `tfp$util$DeferredTensor` to represent any parameters defined as transformations of unconstrained variables, so that the transformations execute at runtime instead of at distribution creation.

`optimizer`

Optimizer instance to use. This may be a TF1-style `tf$train$Optimizer`, TF2-style `tf$optimizers$Optimizer`, or any Python-compatible object that implements `optimizer$apply_gradients(grads_and_vars)`.

`num_steps`

integer number of steps to run the optimizer.

`convergence_criterion`

Optional instance of `tfp$optimizer$convergence_criteria$ConvergenceCriterion` representing a criterion for detecting convergence. If `NULL`, the optimization will run for `num_steps` steps, otherwise, it will run for at *most* `num_steps` steps, as determined by the provided criterion. Default value: `NULL`.

`trace_fn`

function with signature `state = trace_fn(loss, grads, variables)`, where `state` may be a Tensor or nested structure of Tensors. The state values are accumulated (by `tf$scan`) and returned. The default `trace_fn` simply returns the loss, but in general can depend on the gradients and variables (if `trainable_variables` is not `NULL` then `variables==trainable_variables`; otherwise it is the list of all variables accessed during execution of `loss_fn()`), as well as any other quantities captured in the closure of `trace_fn`, for example, statistics of a variational distribution. Default value: `function(loss, grads, variables) loss`.

`variational_loss_fn`

function with signature `loss <- variational_loss_fn(target_log_prob_fn, surrogate_posterior, ...)` defining a variational loss function. The default is a Monte Carlo approximation

	to the standard evidence lower bound (ELBO), equivalent to minimizing the 'reverse' $KL[q p]$ divergence between the surrogate q and true posterior p . Default value: <code>functools.partial(tfp.vi.monte_carlo_variational_loss, discrepancy_fn=tf.vi.monte_carlo_variational_loss.discrepancy_fn)</code> .
<code>discrepancy_fn</code>	A function of Python callable representing a Csiszar f function in log-space. See the docs for <code>tf.vi.monte_carlo_variational_loss</code> for examples. This argument is ignored if a <code>variational_loss_fn</code> is explicitly specified. Default value: <code>tf.vi.kl_reverse</code> .
<code>sample_size</code>	integer number of Monte Carlo samples to use in estimating the variational divergence. Larger values may stabilize the optimization, but at higher cost per step in time and memory. Default value: 1.
<code>importance_sample_size</code>	An integer number of terms used to define an importance-weighted divergence. If <code>importance_sample_size > 1</code> , then the surrogate posterior is optimized to function as an importance-sampling proposal distribution. In this case, posterior expectations should be approximated by importance sampling, as demonstrated in the example below. This argument is ignored if a <code>variational_loss_fn</code> is explicitly specified. Default value: 1.
<code>trainable_variables</code>	Optional list of <code>tf.Variable</code> instances to optimize with respect to. If NULL, defaults to the set of all variables accessed during the computation of the variational bound, i.e., those defining <code>surrogate_posterior</code> and the model <code>target_log_prob_fn</code> . Default value: NULL.
<code>jit_compile</code>	If TRUE, compiles the loss function and gradient update using XLA. XLA performs compiler optimizations, such as fusion, and attempts to emit more efficient code. This may drastically improve the performance. See the docs for <code>tf.function</code> . Default value: NULL.
<code>seed</code>	integer to seed the random number generator.
<code>name</code>	name prefixed to ops created by this function. Default value: 'fit_surrogate_posterior'.

Details

This corresponds to minimizing the 'reverse' Kullback-Liebler divergence ($KL[q||p]$) between the variational distribution and the unnormalized `target_log_prob_fn`, and defines a lower bound on the marginal log likelihood, $\log p(x) \geq -\text{elbo_loss}$.

More generally, this function supports fitting variational distributions that minimize any **Csiszar f -divergence**.

Value

results Tensor or nested structure of Tensors, according to the return type of `result_fn`. Each Tensor has an added leading dimension of size `num_steps`, packing the trajectory of the result over the course of the optimization.

See Also

Other vi-functions: `vi_amari_alpha()`, `vi_arithmetic_geometric()`, `vi_chi_square()`, `vi_csiszar_vimco()`, `vi_dual_csiszar_function()`, `vi_jeffreys()`, `vi_jensen_shannon()`, `vi_kl_forward()`, `vi_kl_reverse()`,

[vi_log1p_abs\(\)](#), [vi_modified_gan\(\)](#), [vi_monte_carlo_variational_loss\(\)](#), [vi_pearson\(\)](#),
[vi_squared_hellinger\(\)](#), [vi_symmetrized_csiszar_function\(\)](#)

vi_jeffreys

The Jeffreys Csiszar-function in log-space

Description

A Csiszar-function is a member of $F = \{ f: \mathbb{R}_+ \text{ to } \mathbb{R} : f \text{ convex} \}$.

Usage

```
vi_jeffreys(logu, name = NULL)
```

Arguments

logu	float-like Tensor representing $\log(u)$ from above.
name	name prefixed to Ops created by this function.

Details

The Jeffreys Csiszar-function is:

$$\begin{aligned} f(u) &= 0.5 (u \log(u) - \log(u)) \\ &= 0.5 \text{ kl_forward} + 0.5 \text{ kl_reverse} \\ &= \text{symmetrized_csiszar_function}(\text{kl_reverse}) \\ &= \text{symmetrized_csiszar_function}(\text{kl_forward}) \end{aligned}$$

This Csiszar-function induces a symmetric f-Divergence, i.e., $D_f[p, q] = D_f[q, p]$.

Warning: this function makes non-log-space calculations and may therefore be numerically unstable for $|\log u| \gg 0$.

Value

jeffreys_of_u: float-like Tensor of the Csiszar-function evaluated at $u = \exp(\log u)$.

See Also

Other vi-functions: [vi_amari_alpha\(\)](#), [vi_arithmetic_geometric\(\)](#), [vi_chi_square\(\)](#), [vi_csiszar_vimco\(\)](#),
[vi_dual_csiszar_function\(\)](#), [vi_fit_surrogate_posterior\(\)](#), [vi_jensen_shannon\(\)](#), [vi_kl_forward\(\)](#),
[vi_kl_reverse\(\)](#), [vi_log1p_abs\(\)](#), [vi_modified_gan\(\)](#), [vi_monte_carlo_variational_loss\(\)](#),
[vi_pearson\(\)](#), [vi_squared_hellinger\(\)](#), [vi_symmetrized_csiszar_function\(\)](#)

vi_jensen_shannon *The Jensen-Shannon Csiszar-function in log-space*

Description

A Csiszar-function is a member of $F = \{ f:R_+ \text{ to } R : f \text{ convex} \}$.

Usage

```
vi_jensen_shannon(logu, self_normalized = FALSE, name = NULL)
```

Arguments

logu	float-like Tensor representing $\log(u)$ from above.
self_normalized	logical indicating whether $f(u=1)=0$. When $f(u=1)=0$ the implied Csiszar f-Divergence remains non-negative even when p, q are unnormalized measures.
name	name prefixed to Ops created by this function.

Details

When `self_normalized = True`, the Jensen-Shannon Csiszar-function is:

$$f(u) = u \log(u) - (1 + u) \log(1 + u) + (u + 1) \log(2)$$

When `self_normalized = False` the $(u + 1) \log(2)$ term is omitted.

Observe that as an f-Divergence, this Csiszar-function implies:

$$D_f[p, q] = KL[p, m] + KL[q, m]$$

$$m(x) = 0.5 p(x) + 0.5 q(x)$$

In a sense, this divergence is the "reverse" of the Arithmetic-Geometric f-Divergence.

This Csiszar-function induces a symmetric f-Divergence, i.e., $D_f[p, q] = D_f[q, p]$.

Warning: this function makes non-log-space calculations and may therefore be numerically unstable for $\log u \gg 0$.

Value

jensen_shannon_of_u, float-like Tensor of the Csiszar-function evaluated at $u = \exp(\log u)$.

References

- Lin, J. "Divergence measures based on the Shannon entropy." IEEE Trans. Inf. Th., 37, 145-151, 1991.

See Also

Other vi-functions: [vi_amari_alpha\(\)](#), [vi_arithmetic_geometric\(\)](#), [vi_chi_square\(\)](#), [vi_csiszar_vimco\(\)](#), [vi_dual_csiszar_function\(\)](#), [vi_fit_surrogate_posterior\(\)](#), [vi_jeffreys\(\)](#), [vi_kl_forward\(\)](#), [vi_kl_reverse\(\)](#), [vi_log1p_abs\(\)](#), [vi_modified_gan\(\)](#), [vi_monte_carlo_variational_loss\(\)](#), [vi_pearson\(\)](#), [vi_squared_hellinger\(\)](#), [vi_symmetrized_csiszar_function\(\)](#)

vi_kl_forward

*The forward Kullback-Leibler Csiszar-function in log-space***Description**

A Csiszar-function is a member of $F = \{ f: \mathbb{R}_+ \rightarrow \mathbb{R} : f \text{ convex} \}$.

Usage

```
vi_kl_forward(logu, self_normalized = FALSE, name = NULL)
```

Arguments

logu	float-like Tensor representing $\log(u)$ from above.
self_normalized	logical indicating whether $f(u=1)=0$. When $f(u=1)=0$ the implied Csiszar f-Divergence remains non-negative even when p, q are unnormalized measures.
name	name prefixed to Ops created by this function.

Details

When `self_normalized = TRUE`, the KL-reverse Csiszar-function is $f(u) = u \log(u) - (u - 1)$. When `self_normalized = FALSE` the $(u - 1)$ term is omitted. Observe that as an f-Divergence, this Csiszar-function implies: $D_f[p, q] = KL[q, p]$

The KL is "forward" because in maximum likelihood we think of minimizing q as in $KL[p, q]$.

Warning: when `self_normalized = TRUE` this function makes non-log-space calculations and may therefore be numerically unstable for $\log u \gg 0$.

Value

kl_forward_of_u: float-like Tensor of the Csiszar-function evaluated at $u = \exp(\log u)$.

See Also

Other vi-functions: [vi_amari_alpha\(\)](#), [vi_arithmetic_geometric\(\)](#), [vi_chi_square\(\)](#), [vi_csiszar_vimco\(\)](#), [vi_dual_csiszar_function\(\)](#), [vi_fit_surrogate_posterior\(\)](#), [vi_jeffreys\(\)](#), [vi_jensen_shannon\(\)](#), [vi_kl_reverse\(\)](#), [vi_log1p_abs\(\)](#), [vi_modified_gan\(\)](#), [vi_monte_carlo_variational_loss\(\)](#), [vi_pearson\(\)](#), [vi_squared_hellinger\(\)](#), [vi_symmetrized_csiszar_function\(\)](#)

 vi_kl_reverse

The reverse Kullback-Leibler Csiszar-function in log-space

Description

A Csiszar-function is a member of $F = \{ f: \mathbb{R}_+ \text{ to } \mathbb{R} : f \text{ convex} \}$.

Usage

```
vi_kl_reverse(logu, self_normalized = FALSE, name = NULL)
```

Arguments

logu	float-like Tensor representing $\log(u)$ from above.
self_normalized	logical indicating whether $f'(u=1)=0$. When $f'(u=1)=0$ the implied Csiszar f-Divergence remains non-negative even when p, q are unnormalized measures.
name	name prefixed to Ops created by this function.

Details

When `self_normalized = TRUE`, the KL-reverse Csiszar-function is $f(u) = -\log(u) + (u - 1)$. When `self_normalized = FALSE` the $(u - 1)$ term is omitted. Observe that as an f-Divergence, this Csiszar-function implies: $D_f[p, q] = KL[q, p]$

The KL is "reverse" because in maximum likelihood we think of minimizing q as in $KL[p, q]$.

Warning: when `self_normalized = TRUE` this function makes non-log-space calculations and may therefore be numerically unstable for $\log u \gg 0$.

Value

kl_reverse_of_u float-like Tensor of the Csiszar-function evaluated at $u = \exp(\log u)$.

See Also

Other vi-functions: [vi_amar_alpha\(\)](#), [vi_arithmetic_geometric\(\)](#), [vi_chi_square\(\)](#), [vi_csiszar_vimco\(\)](#), [vi_dual_csiszar_function\(\)](#), [vi_fit_surrogate_posterior\(\)](#), [vi_jeffreys\(\)](#), [vi_jensen_shannon\(\)](#), [vi_kl_forward\(\)](#), [vi_log1p_abs\(\)](#), [vi_modified_gan\(\)](#), [vi_monte_carlo_variational_loss\(\)](#), [vi_pearson\(\)](#), [vi_squared_hellinger\(\)](#), [vi_symmetrized_csiszar_function\(\)](#)

vi_log1p_abs

The log1p-abs Csiszar-function in log-space

Description

A Csiszar-function is a member of $F = \{ f: \mathbb{R}_+ \text{ to } \mathbb{R} : f \text{ convex} \}$.

Usage

```
vi_log1p_abs(logu, name = NULL)
```

Arguments

logu float-like Tensor representing $\log(u)$ from above.
name name prefixed to Ops created by this function.

Details

The Log1p-Abs Csiszar-function is:

$$f(u) = u^{**}(\text{sign}(u-1)) - 1$$

This function is so-named because it was invented from the following recipe. Choose a convex function g such that $g(0)=0$ and solve for f :

$$\log(1 + f(u)) = g(\log(u)).$$

\Leftrightarrow

$$f(u) = \exp(g(\log(u))) - 1$$

That is, the graph is identically g when y-axis is log1p-domain and x-axis is log-domain.

Warning: this function makes non-log-space calculations and may therefore be numerically unstable for $|\log u| \gg 0$.

Value

log1p_abs_of_u: float-like Tensor of the Csiszar-function evaluated at $u = \exp(\text{logu})$.

See Also

Other vi-functions: [vi_amari_alpha\(\)](#), [vi_arithmetic_geometric\(\)](#), [vi_chi_square\(\)](#), [vi_csiszar_vimco\(\)](#), [vi_dual_csiszar_function\(\)](#), [vi_fit_surrogate_posterior\(\)](#), [vi_jeffreys\(\)](#), [vi_jensen_shannon\(\)](#), [vi_kl_forward\(\)](#), [vi_kl_reverse\(\)](#), [vi_modified_gan\(\)](#), [vi_monte_carlo_variational_loss\(\)](#), [vi_pearson\(\)](#), [vi_squared_hellinger\(\)](#), [vi_symmetrized_csiszar_function\(\)](#)

vi_modified_gan	<i>The Modified-GAN Csiszar-function in log-space</i>
-----------------	---

Description

A Csiszar-function is a member of $F = \{ f: \mathbb{R}_+ \text{ to } \mathbb{R} : f \text{ convex} \}$.

Usage

```
vi_modified_gan(logu, self_normalized = FALSE, name = NULL)
```

Arguments

logu	float-like Tensor representing $\log(u)$ from above.
self_normalized	logical indicating whether $f(u=1)=0$. When $f(u=1)=0$ the implied Csiszar f-Divergence remains non-negative even when p, q are unnormalized measures.
name	name prefixed to Ops created by this function.

Details

When `self_normalized = True` the modified-GAN (Generative/Adversarial Network) Csiszar-function is:

$$f(u) = \log(1 + u) - \log(u) + 0.5 (u - 1)$$

When `self_normalized = False` the $0.5 (u - 1)$ is omitted.

The unmodified GAN Csiszar-function is identical to Jensen-Shannon (with `self_normalized = False`).

Warning: this function makes non-log-space calculations and may therefore be numerically unstable for $\log u \gg 0$.

Value

jensen_shannon_of_u, float-like Tensor of the Csiszar-function evaluated at $u = \exp(\log u)$.

See Also

Other vi-functions: [vi_amari_alpha\(\)](#), [vi_arithmetic_geometric\(\)](#), [vi_chi_square\(\)](#), [vi_csiszar_vimco\(\)](#), [vi_dual_csiszar_function\(\)](#), [vi_fit_surrogate_posterior\(\)](#), [vi_jeffreys\(\)](#), [vi_jensen_shannon\(\)](#), [vi_kl_forward\(\)](#), [vi_kl_reverse\(\)](#), [vi_log1p_abs\(\)](#), [vi_monte_carlo_variational_loss\(\)](#), [vi_pearson\(\)](#), [vi_squared_hellinger\(\)](#), [vi_symmetrized_csiszar_function\(\)](#)

 vi_monte_carlo_variational_loss

Monte-Carlo approximation of an f -Divergence variational loss

Description

Variational losses measure the divergence between an unnormalized target distribution p (provided via `target_log_prob_fn`) and a surrogate distribution q (provided as `surrogate_posterior`). When the target distribution is an unnormalized posterior from conditioning a model on data, minimizing the loss with respect to the parameters of `surrogate_posterior` performs approximate posterior inference.

Usage

```
vi_monte_carlo_variational_loss(
  target_log_prob_fn,
  surrogate_posterior,
  sample_size = 1L,
  importance_sample_size = 1L,
  discrepancy_fn = vi_kl_reverse,
  use_reparametrization = NULL,
  seed = NULL,
  name = NULL
)
```

Arguments

`target_log_prob_fn`

function that takes a set of Tensor arguments and returns a Tensor log-density. Given `q_sample <- surrogate_posterior$sample(sample_size)`, this will be (in Python) called as `target_log_prob_fn(q_sample)` if `q_sample` is a list or a tuple, `target_log_prob_fn(**q_sample)` if `q_sample` is a dictionary, or `target_log_prob_fn(q_sample)` if `q_sample` is a Tensor. It should support batched evaluation, i.e., should return a result of shape `[sample_size]`.

`surrogate_posterior`

A `tfp$distributions$Distribution` instance defining a variational posterior (could be a `tfp$distributions$JointDistribution`). Crucially, the distribution's `log_prob` and (if reparameterized) `sample` methods must directly invoke all ops that generate gradients to the underlying variables. One way to ensure this is to use `tfp$util$DeferredTensor` to represent any parameters defined as transformations of unconstrained variables, so that the transformations execute at runtime instead of at distribution creation.

`sample_size`

integer number of Monte Carlo samples to use in estimating the variational divergence. Larger values may stabilize the optimization, but at higher cost per step in time and memory. Default value: 1.

`importance_sample_size` integer number of terms used to define an importance-weighted divergence. If `importance_sample_size > 1`, then the `surrogate_posterior` is optimized to function as an importance-sampling proposal distribution. In this case it often makes sense to use importance sampling to approximate posterior expectations (see `tfp.vi.fit_surrogate_posterior` for an example). Default value: 1.

`discrepancy_fn` function representing a Csiszar f function in in log-space. That is, `discrepancy_fn(log(u)) = f(u)`, where f is convex in u . Default value: `vi_kl_reverse`.

`use_reparameterization` logical. When NULL (the default), automatically set to: `surrogate_posterior.reparameterization == tfp$distributions$FULLY_REPARAMETERIZED`. When TRUE uses the standard Monte-Carlo average. When FALSE uses the score-gradient trick. (See above for details.) When FALSE, consider using `csiszar_vimco`.

`seed` integer seed for `surrogate_posterior$sample`.

`name` name prefixed to Ops created by this function.

Details

This function defines divergences of the form $E_q[\text{discrepancy_fn}(\log p(z) - \log q(z))]$, sometimes known as **f-divergences**.

In the special case `discrepancy_fn(logu) == -logu` (the default `vi_kl_reverse`), this is the reverse Kullback-Liebler divergence $KL[q||p]$, whose negation applied to an unnormalized p is the widely-used evidence lower bound (ELBO). Other cases of interest available under `tfp$vi` include the forward $KL[p||q]$ (given by `vi_kl_forward(logu) == exp(logu) * logu`), total variation distance, Amari alpha-divergences, and more.

Csiszar f-divergences

A Csiszar function f is a convex function from \mathbb{R}^+ (the positive reals) to \mathbb{R} . The Csiszar f -Divergence is given by:

$$D_f[p(X), q(X)] := E_{\{q(X)\}}[f(p(X) / q(X))] \\ \approx m^{*-1} \sum_j^m f(p(x_j) / q(x_j)), \\ \text{where } x_j \sim \text{iid } q(X)$$

For example, $f = \lambda u: -\log(u)$ recovers $KL[q||p]$, while $f = \lambda u: u * \log(u)$ recovers the forward $KL[p||q]$. These and other functions are available in `tfp$vi`.

Tricks: Reparameterization and Score-Gradient

When q is "reparameterized", i.e., a diffeomorphic transformation of a parameterless distribution (e.g., $\text{Normal}(Y; m, s) \Leftrightarrow Y = sX + m, X \sim \text{Normal}(0,1)$), we can swap gradient and expectation, i.e., $\text{grad}[\text{Avg}\{s_i : i=1\dots n\}] = \text{Avg}\{\text{grad}[s_i] : i=1\dots n\}$ where $S_n = \text{Avg}\{s_i\}$ and $s_i = f(x_i), x_i \sim \text{iid } q(X)$.

However, if q is not reparameterized, TensorFlow's gradient will be incorrect since the chain-rule stops at samples of unreparameterized distributions. In this circumstance using the Score-Gradient trick results in an unbiased gradient, i.e.,

$$\text{grad}[E_q[f(X)]] \\ = \text{grad}[\int dx q(x) f(x)]$$

```

= int dx grad[ q(x) f(x) ]
= int dx [ q'(x) f(x) + q(x) f'(x) ]
= int dx q(x) [q'(x) / q(x) f(x) + f'(x) ]
= int dx q(x) grad[ f(x) q(x) / stop_grad[q(x)] ]
= E_q[ grad[ f(x) q(x) / stop_grad[q(x)] ] ]

```

Unless `q.reparameterization_type != tfd.FULLY_REPARAMETERIZED` it is usually preferable to set `use_reparameterization = True`.

Example Application: The Csiszar f-Divergence is a useful framework for variational inference. I.e., observe that,

$$\begin{aligned}
 f(p(x)) &= f(E_{\{q(Z | x)\}}[p(x, Z) / q(Z | x)]) \\
 &\leq E_{\{q(Z | x)\}}[f(p(x, Z) / q(Z | x))] \\
 &:= D_f[p(x, Z), q(Z | x)]
 \end{aligned}$$

The inequality follows from the fact that the "perspective" of f , i.e., $(s, t) \mapsto t f(s / t)$, is convex in (s, t) when s/t in $\text{domain}(f)$ and t is a real. Since the above framework includes the popular Evidence Lower Bound (ELBO) as a special case, i.e., $f(u) = -\log(u)$, we call this framework "Evidence Divergence Bound Optimization" (EDBO).

Value

`monte_carlo_variational_loss` float-like Tensor Monte Carlo approximation of the Csiszar f-Divergence.

References

- Ali, Syed Mumtaz, and Samuel D. Silvey. "A general class of coefficients of divergence of one distribution from another." *Journal of the Royal Statistical Society: Series B (Methodological)* 28.1 (1966): 131-142.

See Also

Other vi-functions: [vi_amari_alpha\(\)](#), [vi_arithmetic_geometric\(\)](#), [vi_chi_square\(\)](#), [vi_csiszar_vimco\(\)](#), [vi_dual_csiszar_function\(\)](#), [vi_fit_surrogate_posterior\(\)](#), [vi_jeffreys\(\)](#), [vi_jensen_shannon\(\)](#), [vi_kl_forward\(\)](#), [vi_kl_reverse\(\)](#), [vi_log1p_abs\(\)](#), [vi_modified_gan\(\)](#), [vi_pearson\(\)](#), [vi_squared_hellinger\(\)](#), [vi_symmetrized_csiszar_function\(\)](#)

vi_pearson

The Pearson Csiszar-function in log-space

Description

A Csiszar-function is a member of $F = \{ f: \mathbb{R}_+ \rightarrow \mathbb{R} : f \text{ convex} \}$.

Usage

```
vi_pearson(logu, name = NULL)
```

Arguments

logu float-like Tensor representing $\log(u)$ from above.
 name name prefixed to Ops created by this function.

Details

The Pearson Csiszar-function is:

$$f(u) = (u - 1)**2$$

Warning: this function makes non-log-space calculations and may therefore be numerically unstable for $|\log u| \gg 0$.

Value

pearson_of_u: float-like Tensor of the Csiszar-function evaluated at $u = \exp(\log u)$.

See Also

Other vi-functions: [vi_amari_alpha\(\)](#), [vi_arithmetic_geometric\(\)](#), [vi_chi_square\(\)](#), [vi_csiszar_vimco\(\)](#), [vi_dual_csiszar_function\(\)](#), [vi_fit_surrogate_posterior\(\)](#), [vi_jeffreys\(\)](#), [vi_jensen_shannon\(\)](#), [vi_kl_forward\(\)](#), [vi_kl_reverse\(\)](#), [vi_log1p_abs\(\)](#), [vi_modified_gan\(\)](#), [vi_monte_carlo_variational_loss\(\)](#), [vi_squared_hellinger\(\)](#), [vi_symmetrized_csiszar_function\(\)](#)

vi_squared_hellinger *The Squared-Hellinger Csiszar-function in log-space*

Description

A Csiszar-function is a member of $F = \{ f: \mathbb{R}_+ \text{ to } \mathbb{R} : f \text{ convex} \}$.

Usage

```
vi_squared_hellinger(logu, name = NULL)
```

Arguments

logu float-like Tensor representing $\log(u)$ from above.
 name name prefixed to Ops created by this function.

Details

The Squared-Hellinger Csiszar-function is:

$$f(u) = (\sqrt{u} - 1)**2$$

This Csiszar-function induces a symmetric f-Divergence, i.e., $D_f[p, q] = D_f[q, p]$.

Warning: this function makes non-log-space calculations and may therefore be numerically unstable for $|\log u| \gg 0$.

Value

Squared-Hellinger_of_u: float-like Tensor of the Csiszar-function evaluated at $u = \exp(\log u)$.

See Also

Other vi-functions: [vi_amari_alpha\(\)](#), [vi_arithmetic_geometric\(\)](#), [vi_chi_square\(\)](#), [vi_csiszar_vimco\(\)](#), [vi_dual_csiszar_function\(\)](#), [vi_fit_surrogate_posterior\(\)](#), [vi_jeffreys\(\)](#), [vi_jensen_shannon\(\)](#), [vi_kl_forward\(\)](#), [vi_kl_reverse\(\)](#), [vi_log1p_abs\(\)](#), [vi_modified_gan\(\)](#), [vi_monte_carlo_variational_loss\(\)](#), [vi_pearson\(\)](#), [vi_symmetrized_csiszar_function\(\)](#)

vi_symmetrized_csiszar_function

Symmetrizes a Csiszar-function in log-space

Description

A Csiszar-function is a member of $F = \{ f: \mathbb{R}_+ \rightarrow \mathbb{R} : f \text{ convex} \}$.

Usage

```
vi_symmetrized_csiszar_function(logu, csiszar_function, name = NULL)
```

Arguments

logu float-like Tensor representing $\log(u)$ from above.
 csiszar_function function representing a Csiszar-function over log-domain.
 name name prefixed to Ops created by this function.

Details

The symmetrized Csiszar-function is defined as:

$$f_g(u) = 0.5 g(u) + 0.5 u g(1 / u)$$

where g is some other Csiszar-function. We say the function is "symmetrized" because:

$$D_{\{f_g\}}[p, q] = D_{\{f_g\}}[q, p]$$

for all $p \ll \gg q$ (i.e., $\text{support}(p) = \text{support}(q)$).

There exists alternatives for symmetrizing a Csiszar-function. For example,

$$f_g(u) = \max(f(u), f^*(u)),$$

where f^* is the dual Csiszar-function, also implies a symmetric f -Divergence.

Example: When either of the following functions are symmetrized, we obtain the Jensen-Shannon Csiszar-function, i.e.,

$$g(u) = -\log(u) - (1 + u) \log((1 + u) / 2) + u - 1$$

$$h(u) = \log(4) + 2 u \log(u / (1 + u))$$

implies,

$$f_g(u) = f_h(u) = u \log(u) - (1 + u) \log((1 + u) / 2)$$

$$= \text{jensen_shannon}(\log(u)).$$

Warning: this function makes non-log-space calculations and may therefore be numerically unstable for $\lvert \log u \rvert \gg 0$.

Value

symmetrized_g_of_u: float-like Tensor of the result of applying the symmetrization of g evaluated at $u = \exp(\log u)$.

See Also

Other vi-functions: [vi_amari_alpha\(\)](#), [vi_arithmetic_geometric\(\)](#), [vi_chi_square\(\)](#), [vi_csiszar_vimco\(\)](#), [vi_dual_csiszar_function\(\)](#), [vi_fit_surrogate_posterior\(\)](#), [vi_jeffreys\(\)](#), [vi_jensen_shannon\(\)](#), [vi_kl_forward\(\)](#), [vi_kl_reverse\(\)](#), [vi_log1p_abs\(\)](#), [vi_modified_gan\(\)](#), [vi_monte_carlo_variational_loss\(\)](#), [vi_pearson\(\)](#), [vi_squared_hellinger\(\)](#)

vi_total_variation	<i>The Total Variation Csiszar-function in log-space</i>
--------------------	--

Description

A Csiszar-function is a member of $F = \{ f: \mathbb{R}_+ \rightarrow \mathbb{R} : f \text{ convex} \}$.

Usage

```
vi_total_variation(logu, name = NULL)
```

Arguments

logu	float-like Tensor representing $\log(u)$ from above.
name	name prefixed to Ops created by this function.

Details

The Total-Variation Csiszar-function is:

$$f(u) = 0.5 |u - 1|$$

Warning: this function makes non-log-space calculations and may therefore be numerically unstable for $\lvert \log u \rvert \gg 0$.

Value

total_variation_of_u: float-like Tensor of the Csiszar-function evaluated at $u = \exp(\text{logu})$.

See Also

Other vi-functions#: [vi_t_power\(\)](#), [vi_triangular\(\)](#)

vi_triangular	<i>The Triangular Csiszar-function in log-space</i>
---------------	---

Description

The Triangular Csiszar-function is:

Usage

```
vi_triangular(logu, name = NULL)
```

Arguments

logu	float-like Tensor representing $\log(u)$ from above.
name	name prefixed to Ops created by this function.

Details

$$f(u) = (u - 1)**2 / (1 + u)$$

Warning: this function makes non-log-space calculations and may therefore be numerically unstable for $|\text{llogu}| \gg 0$.

Value

triangular_of_u: float-like Tensor of the Csiszar-function evaluated at $u = \exp(\text{logu})$.

See Also

Other vi-functions#: [vi_t_power\(\)](#), [vi_total_variation\(\)](#)

 vi_t_power

The T-Power Csiszar-function in log-space

Description

A Csiszar-function is a member of $F = \{ f: \mathbb{R}_+ \text{ to } \mathbb{R} : f \text{ convex} \}$.

Usage

```
vi_t_power(logu, t, self_normalized = FALSE, name = NULL)
```

Arguments

logu	float-like Tensor representing $\log(u)$ from above.
t	Tensor of same dtype as logu and broadcastable shape.
self_normalized	logical indicating whether $f(u=1)=0$. When $f(u=1)=0$ the implied Csiszar f-Divergence remains non-negative even when p, q are unnormalized measures.
name	name prefixed to Ops created by this function.

Details

When `self_normalized = True` the T-Power Csiszar-function is:

$$f(u) = s [u^{*t} - 1 - t(u - 1)]$$

$$s = \begin{cases} -1 & 0 < t < 1 \\ +1 & \text{otherwise} \end{cases}$$

When `self_normalized = False` the $-t(u - 1)$ term is omitted.

This is similar to the `amari_alpha` Csiszar-function, with the associated divergence being the same up to factors depending only on `t`.

Warning: when `self_normalized = True` this function makes non-log-space calculations and may therefore be numerically unstable for $\log u \gg 0$.

Value

`t_power_of_u`: float-like Tensor of the Csiszar-function evaluated at $u = \exp(\log u)$.

See Also

Other vi-functions#: [vi_total_variation\(\)](#), [vi_triangular\(\)](#)

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