Package 'BHTSpack'

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Type Package

Title Bayesian Multi-Plate High-Throughput Screening of Compounds

Version 0.6

Description Can be used for joint identification of candidate compound hits from multiple assays, in drug discovery. This package implements the framework of I. D. Shterev, D. B. Dunson, C. Chan and G. D. Sempowski. ``Bayesian Multi-Plate High-Throughput Screening of Compounds", Scientific Reports 8(1):9551, 2018. This project was funded by the Division of Allergy, Immunology, and Transplantation, National Institute of Allergy and Infectious Diseases, National Institutes of Health, Department of Health and Human Services, under contract No. HHSN272201400054C entitled ``Adjuvant Discovery For Vaccines Against West Nile Virus and Influenza", awarded to Duke University and lead by Drs. Herman Staats and Soman Abraham.

Depends R (>= 3.2.3), R2HTML (>= 2.3.2), xtable (>= 1.8-2)

VignetteBuilder knitr

Suggests knitr

License GPL-3

LazyLoad yes

NeedsCompilation yes

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`BHTSpack-package` *Bayesian Multi-Plate High-Throughput Screening of Compounds*

Description

Can be used for joint identification of candidate hits from multiple assays, in drug discovery. This package implements the framework of I. D. Shterev, D. B. Dunson, C. Chan and G. D. Sempowski. "Bayesian Multi Plate High Throughput Screening of Compounds", arXiv:1709.10041, September 2017. This project was funded by the Division of Allergy, Immunology, and Transplantation, National Institute of Allergy and Infectious Diseases, National Institutes of Health, Department of Health and Human Services, under contract No. HHSN272201400054C entitled "Adjuvant Discovery For Vaccines Against West Nile Virus and Influenza", awarded to Duke University and lead by Drs. Herman Staats and Soman Abraham.

Details

The DESCRIPTION file:

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Title:	Bayesian Multi-Plate High-Throughput Screening of Compounds
Version:	0.6
Authors@R:	c(person(c("Ivo", "D."), "Shterev", role = c("aut", "cre"), email = "i.shterev@gmail.com"), person(c("Davi
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VignetteBuilder:	knitr
Suggests:	knitr

License:	GPL-3
LazyLoad:	yes
Author:	Ivo D. Shterev [aut, cre], David B. Dunson [aut], Cliburn Chan [aut], Gregory D. Sempowski [aut]
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Author(s)

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References

I. D. Shterev, D. B. Dunson, C. Chan and G. D. Sempowski. "Bayesian Multi-Plate High-Throughput Screening of Compounds". *Scientific Reports*, 8(1):9551, 2018.

---

abfun *package internal function*

Description

package internal function

Usage

```
abfun(m, v)
```

Arguments

m	Description
v	Description

Examples

```
abfun(0.26, 10^-4)
```

`alpha.u` *package internal function*

Description

package internal function

Usage

```
alpha.u(nu, a0, b0, H)
```

Arguments

nu	Description
a0	Description
b0	Description
H	Description

Examples

```
M = 5
H = 10
a = 10^-6
b = 10^-6
nu = lapply(1:M, function(x){rbeta(H, a, b)})
alpha.u(nu, a, b, H)
```

b.u *package internal function*

Description

package internal function

Usage

```
b.u(hatpai)
```

Arguments

hatpai	Description

Examples

```
pai = 0.5
M = 10
H = 10
K = 5
n = 100

z = abs(rnorm(n))

sigma1 = abs(rnorm(K))
sigma0 = abs(rnorm(K))

mu1 = abs(rnorm(K))
mu0 = abs(rnorm(K))

hk0 = matrix(sample(K, M*H, replace=TRUE), M, H)
hk1 = matrix(sample(K, M*H, replace=TRUE), M, H)

nu.h0 = lapply(1:H, function(x){rbeta(1,5,5)})
nu.h1 = lapply(1:H, function(x){rbeta(1,5,5)})

ph0 = lapply(nu.h0, lambda.u)
ph1 = lapply(nu.h1, lambda.u)

b.u(hatpai.u(z, hk1, hk0, ph1, ph0, sigma1, sigma0, mu1, mu0, pai, H, n))
```

bhts *Bayesian High-Throughput Screening*

Description

This is the package main function.

Usage

```
bhts(Z, iters, H, K, mu00=NULL, mu10=NULL, a.alpha, b.alpha, a.tau, b.tau,
pnorm=FALSE, s=NULL, store=FALSE)
```

Arguments

Z	A list of compounds.
iters	Number of iterations to perform.
H	Number of local DP components.
K	Number of global DP components.
mu00	Activity level (mean) of non-hit compounds
mu10	Activity level (mean) of hit compounds
a.alpha	Gamma shape parameter specifying local DP concentration prior.
b.alpha	Gamma rate parameter specifying local DP concentration prior.
a.tau	Gamma shape parameter specifying global DP concentration prior.
b.tau	Gamma rate parameter specifying global DP concentration prior.
pnorm	Plate normalization. If $TRUE$, each plate is normalized to zero mean and unit variance, prior to analysis. Default is $FALSE$.
s	Random seed (for reproducibility purposes). Default is $NULL$.
store	If $TRUE$, all samples of certain latent variables are stored in the output object. Default is $FALSE$.

Value

This function returns a list consisting of the following elements:

hatpai	A list of vectors of posterior probabilities, estimating the probability of a compound being a hit.
dat.store	If $store = TRUE$ (default is $FALSE$), the output contains a list of $iters \times K$ matrices of samples. Each matrix contains the samples of a separate latent variable. At each iteration, the following six variables are stored in a different row of their corrpsponding matrix, $(\lambda_1^{(0)}, \ldots, \lambda_K^{(0)})$, $(\lambda_1^{(1)}, \ldots, \lambda_K^{(1)})$, $(\mu_{01}, \ldots, \mu_{0K})$, $(\mu_{11}, \ldots, \mu_{1K})$, $(\sigma_{01}^2, \ldots, \sigma_{0K}^2)$ and $(\sigma_{11}^2, \ldots, \sigma_{1K}^2)$.

Examples

```
set.seed(1234)
Nmax = 100
M = 100
n = sample(Nmax, M, replace=TRUE)
Z = lapply(n, function(x){abs(rnorm(x))})
bhts(Z, iters=100, H=10, K=5, mu00=0, mu10=10, a.alpha=10, b.alpha=5, a.tau=10, b.tau=5)
```

---

bhts2HTML *Convert to HTML*

Description

This function creates an HTML file.

Usage

```
bhts2HTML(dat, dir, fname, title=NULL, bgcolor="#BBBBEE")
```

Arguments

dat	An object which is the output of $bhts()$.
dir	Directory in which to store the file.
fname	File name.
title	The title of the html file.
bgcolor	Color for the html background.

Examples

```
#See package vignette
```

---

data.create *Create Synthetic Data*

Description

This function generates synthetic compound data.

Usage

```
data.create(N, nr, nc, M, p, s=NULL, covrow=NULL, covcol=NULL, c=0.0001, mat=FALSE)
```

Arguments

N	Number of compounds per plate.
nr	Number of plate rows.
nc	Number of plate columns.
M	Number of plates.
p	Probability of a compound being a hit.
s	Random seed (for reproducibility perposes). Default is $NULL$.
covrow	Noise plate row-covariance matrix. Default is $NULL$.
covcol	Noise plate column-covariance matrix. Default is $NULL$.
c	Constant for scaling plate noise. Default is 0.0001.
mat	Specifies a matrix ($TRUE$) or a vector ($FALSE$) plate format. Default is $FALSE$.

Value

This function returns a list consisting of the following elements:

Z	A list of matrices ($mat = TRUE$) or vectors ($mat = FALSE$) of compounds.
B	A list of compound indicators specifying a hit (1) or a non-hit (0).
I	A list of compound indicators specifying the muxture component (from 1 to K).

Examples

```
#See package vignette
```

---

fdr.r *package internal function*

Description

package internal function

Usage

```
fdr.r(r, hatpai, fdr)
```

Arguments

r	Description
hatpai	Description
fdr	Description

Examples

```
#See package vignette
```

h.pr.u *package internal function*

Description

package internal function

Usage

```
h.pr.u(z, ih, mu, sigma, pk, K, H, n)
```

Arguments

z

ih

mu

sigma

pk

K

H

n

Examples

```
Nmax = 100
K = 5
H = 10
M = 20
n = sample(Nmax, M, replace=TRUE)

z = abs(rnorm(sum(n)))
mu = abs(rnorm(K))
sigma = 1/rgamma(n=K, shape=10, rate=10)

ih = sample(H,sum(n), replace=TRUE)

pk = lambda.u(rbeta(K, 1, 1))

h.pr.u(z, ih, mu, sigma, pk, K, H, n)
```

hatpai.u *package internal function*

Description

package internal function

Usage

```
hatpai.u(z, hk1, hk0, ph1, ph0, sigma1, sigma0, mu1, mu0, pai, H, n)
```

Arguments

z

hk1

hk0

ph1

ph0

sigma1

sigma0

mu1

mu0

pai

H

n

Value

value

Examples

```
pai = 0.5
M = 10
H = 10
K = 5
n = 100

z = abs(rnorm(n))

sigma1 = abs(rnorm(K))
sigma0 = abs(rnorm(K))

mu1 = abs(rnorm(K))
mu0 = abs(rnorm(K))
```

```
hk0 = matrix(sample(K, M*H, replace=TRUE), M, H)
hk1 = matrix(sample(K, M*H, replace=TRUE), M, H)

nu.h0 = lapply(1:H, function(x){rbeta(1,5,5)})
nu.h1 = lapply(1:H, function(x){rbeta(1,5,5)})

ph0 = lapply(nu.h0, lambda.u)
ph1 = lapply(nu.h1, lambda.u)

hatpai.u(z, hk1, hk0, ph1, ph0, sigma1, sigma0, mu1, mu0, pai, H, n)
```

`ind.u` *package internal function*

Description

package internal function

Usage

```
ind.u(pr)
```

Arguments

```
pr
```

Examples

```
Nmax = 100
K = 5
H = 10
M = 20
n = sample(Nmax, M, replace=TRUE)

z = abs(rnorm(sum(n)))
mu = abs(rnorm(K))
sigma = 1/rgamma(n=K, shape=10, rate=10)

hk = sample(K,M*H, replace=TRUE)

ph = as.vector(sapply(1:M, function(x){lambda.u(rbeta(H, 1, 1))}))

ind.u(z.pr.u(z, hk, mu, sigma, ph, H, n))
```

lambda.u *package internal function*

Description

package internal function

Usage

```
lambda.u(nu)
```

Arguments

nu

Examples

```
H = 5
nu = rbeta(H, 1, 1)
lambda.u(nu)
```

lg.mu.sig *package internal function*

Description

package internal function

Usage

```
lg.mu.sig(m, v)
```

Arguments

m	Description
v	Description

Examples

```
#See package vignette
```

mu.k.u *package internal function*

Description

package internal function

Usage

```
mu.k.u(k, ik, z, sigma, mu0)
```

Arguments

k

ik

z

sigma

mu0

Value

Describe

Examples

```
K = 5
n = 100
z = abs(rnorm(n))
sigma = sapply(1:K, function(x){1/rgamma(n=1, shape=10, rate=10)})
mu0 = 0
ik = sample(K, n, replace=TRUE)
sapply(1:K, mu.k.u, ik, z, sigma, mu0)
```

nu.u *package internal function*

Description

package internal function

Usage

```
nu.u(ind, tau, H)
```

Arguments

ind

tau

H

Value

Describe

Examples

```
H = 5
n = 100
tau = rgamma(1, 1, 1)
ind = sample(H, n, replace=TRUE)
nu.u(ind, tau, H)
```

---

pai.u *package internal function*

Description

package internal function

Usage

```
pai.u(b, a.pai, b.pai)
```

Arguments

b

a.pai

b.pai

Value

Describe

Examples

```
n = 100
b = rbinom(n, 1, 0.5)
a.pai = 10^-6
b.pai = 10^-6
pai.u(b, a.pai, b.pai)
```

ptrace *Trace (ACF) Plots*

Description

This function outputs trace plots of certain latent variables.

Usage

```
ptrace(res, var, ndisc, nr, nc, type="trace")
```

Arguments

res	An output object from $bhts()$.
var	Variable for which to display convergence diagnostic plots. Current options are $mu0$ (displaying $\mu_{01}, \ldots, \mu_{0K}$), $mu1$ (displaying $\mu_{11}, \ldots, \mu_{1K}$), $sigma0$ (displaying $\sigma^2_{01}, \ldots, \sigma^2_{0K}$), $sigma1$ (displaying $\sigma^2_{11}, \ldots, \sigma^2_{1K}$), $pk0$ (displaying $\lambda^{(0)}_1, \ldots, \lambda^{(0)}_K$) and $pk1$ (displaying $\lambda^{(1)}_1, \ldots, \lambda^{(1)}_K$).
ndisc	Number of iterations for which to discard samples.
nr	Number of rows in the resulting composite plot.
nc	Number of columns in the resulting composite plot.
type	Type of convergence diagnostic. Currently implemented are trace plots (default $type = "trace"$) and ACF plots ($type = "acf"$)

Examples

```
#See package vignette
```

r.fdr *Significant Hits*

Description

This function determines significant hits, based on a specified expected FDR.

Usage

```
r.fdr(res, fdr=0.05)
```

Arguments

res	An output object from $bhts()$.
fdr	Expected FDR (default is 0.05).

Value

This function returns a list consisting of the following elements:

res	A data frame containing significant hits and their probabilities.
r	The computed significant hit probability threshold.

Examples

```
#See package vignette
```

---

sig.k.u *package internal function*

Description

package internal function

Usage

```
sig.k.u(k, ik, z, mu0, a0, b0)
```

Arguments

k
ik
z
mu0
a0
b0

Examples

```
K = 5
n = 100
z = abs(rnorm(n))
mu0 = 0
ik = sample(K, n, replace=TRUE)
a0 = 5
b0 = 5
sapply(1:K, sig.k.u, ik, z, mu0, a0, b0)
```

tau.u *package internal function*

Description

package internal function

Usage

```
tau.u(nu, a0, b0)
```

Arguments

nu	Description
a0	Description
b0	Description

Examples

```
K = 5
a = 10^-6
b = 10^-6
nu = rbeta(K, a, b)
tau.u(nu, a, b)
```

z.pr.u *package internal function*

Description

package internal function

Usage

```
z.pr.u(z, hk, mu, sigma, ph, H, n)
```

Arguments

z

hk

mu

sigma

ph

H

n

Examples

```
Nmax = 100
K = 5
H = 10
M = 20
n = sample(Nmax, M, replace=TRUE)

z = abs(rnorm(sum(n)))
mu = abs(rnorm(K))
sigma = 1/rgamma(n=K, shape=10, rate=10)

hk = sample(K,M*H, replace=TRUE)

ph = as.vector(sapply(1:M, function(x){lambda.u(rbeta(H, 1, 1))}))

z.pr.u(z, hk, mu, sigma, ph, H, n)
```

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