

# Package ‘MALDIquant’

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**Suggests** knitr, testthat (>= 0.8)

**Description** A complete analysis pipeline for matrix-assisted laser desorption/ionization-time-of-flight (MALDI-TOF) and other two-dimensional mass spectrometry data. In addition to commonly used plotting and processing methods it includes distinctive features, namely baseline subtraction methods such as morphological filters (TopHat) or the statistics-sensitive non-linear iterative peak-clipping algorithm (SNIP), peak alignment using warping functions, handling of replicated measurements as well as allowing spectra with different resolutions.

**License** GPL (>= 3)

**URL** <https://strimmerlab.github.io/software/malDIquant/>  
<https://github.com/sgibb/MALDIquant/>

**BugReports** <https://github.com/sgibb/MALDIquant/issues/>

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## Description

MALDIquant provides a complete analysis pipeline for matrix-assisted laser desorption/ionization-time-of-flight (MALDI-TOF) and other two-dimensional mass spectrometry data.

In addition to commonly used plotting and processing methods it includes distinctive features, namely baseline subtraction methods such as morphological filters (TopHat) or the statistics-sensitive non-linear iterative peak-clipping algorithm (SNIP), peak alignment using warping functions, handling of replicated measurements as well as allowing spectra with different resolutions.

For a first overview see `vignette("MALDIquant-intro", package="MALDIquant")` and/or run `demo("MALDIquant")`.

## Details

Package: MALDIquant  
License: GPL (>= 3)  
URL: <https://strimmerlab.github.io/software/malDIquant/>

Main classes:

- `MassPeaks`: Represents a peak list of a single spectrum.
- `MassSpectrum`: Represents a single spectrum.

The accompanying website (see below) provides example R scripts to illustrate the functionality of this package, too.

## Author(s)

Sebastian Gibb

Maintainer: Sebastian Gibb <[mail@sebastiangibb.de](mailto:mail@sebastiangibb.de)>

## References

S. Gibb and K. Strimmer. 2012. MALDIquant: a versatile R package for the analysis of mass spectrometry data. *Bioinformatics* 28: 2270-2271. <https://academic.oup.com/bioinformatics/article-abstract/28/17/2270/246552>

Website: <https://strimmerlab.github.io/software/malDIquant/>

## See Also

- Introduction: `vignette("MALDIquant-intro", package="MALDIquant")`.
- Run demo files: `demo("MALDIquant")`.
- List all available manual pages: `library(help="MALDIquant")`.
- MALDIquant website: <https://strimmerlab.github.io/software/malDIquant/>.
- more MALDIquant examples and complete analyses: <https://github.com/sgibb/MALDIquantExamples/>.

---

 AbstractMassObject-class

*Class "AbstractMassObject"*


---

## Description

[AbstractMassObject](#) is an abstract (means pure virtual) class. It is the parent class of [MassSpectrum](#) and [MassPeaks](#). It shouldn't create or handle by the user because it is for internal use only.

## Derived classes

[MassPeaks](#), [MassSpectrum](#)

## Slots

**mass**: numeric, mass or mass-to-charge ratio

**intensity**: numeric, intensities for measured mass-to-charge ratios

**metaData**: list, some metadata to describe the spectrum

## Methods

[ **signature**(x = "AbstractMassObject", i = "numeric"): Extracts a range of an [AbstractMassObject](#) object and returns a new one.

**as.matrix** **signature**(x = "AbstractMassObject"): Converts an [AbstractMassObject](#) object to a matrix with 2 columns (mass, intensity).

**coordinates** **signature**(object = "AbstractMassObject"): Accessor function for coordinates stored in object generated from imaging mass spectrometry data.

**coordinates**<- **signature**(object = "AbstractMassObject", value = "numeric|matrix") Replacement function for coordinates used in imaging mass spectrometry datasets.

**intensity** **signature**(object = "AbstractMassObject"): Accessor function for slot intensity.

**intensity**<- **signature**(object = "AbstractMassObject", value = "numeric") Replacement function for slot intensity.

**isEmpty** **signature**(object = "AbstractMassObject"): Returns TRUE if length of intensity is 0 or all intensity values are 0.

**length** **signature**(x = "AbstractMassObject"): Returns length of slot intensity.

**lines** **signature**(x = "AbstractMassObject"): Extended function for adding [AbstractMassObject](#) object as a line to a specific plot. See [lines](#) for details.

**mass** **signature**(object = "AbstractMassObject"): Accessor function for slot mass.

**mass**<- **signature**(object = "AbstractMassObject", value = "numeric") Replacement function for slot mass.

**mz** **signature**(object = "AbstractMassObject"): Accessor function for slot mass.

**mz**<- **signature**(object = "AbstractMassObject", value = "numeric") Replacement function for slot mass.

**metaData** signature(object = "AbstractMassObject"): Accessor function for slot metaData.

**metaData<-** signature(object = "AbstractMassObject"): Replacement function for slot metaData.

**plot** signature(x = "AbstractMassObject", y = "missing"): Extended function for plotting an AbstractMassObject object. See [plot](#), [AbstractMassObject](#), [missing-method](#) for details.

**points** signature(x = "AbstractMassObject"): Extended function for adding [AbstractMassObject](#) object as points to a specific plot. See [points](#) for details.

**trim** signature(object = "AbstractMassObject", range = "numeric"): Trim an AbstractMassObject object. See [trim](#), [AbstractMassObject](#), [numeric-method](#) for details.

**transformIntensity** signature(object = "AbstractMassObject"): Transforms the intensities of an AbstractMassObject object. See [transformIntensity](#), [AbstractMassObject-method](#) for details.

### Author(s)

Sebastian Gibb <[mail@sebastiangibb.de](mailto:mail@sebastiangibb.de)>

### See Also

[MassPeaks](#), [MassSpectrum](#), [plot](#), [AbstractMassObject](#), [missing-method](#), [transformIntensity](#), [AbstractMassObject-trim](#), [AbstractMassObject-numeric-method](#)

Website: <https://strimmerlab.github.io/software/maldiquant/>

### Examples

```
## load package
library("MALDIquant")

## create example spectrum
s <- createMassSpectrum(mass=1:10, intensity=11:20,
                       metaData=list(name="Example Spectrum"))

## get intensity
intensity(s)

## get mass
mass(s)

## get metaData
metaData(s)

## replace metaData
metaData(s) <- list(name="Spectrum")

## trim spectrum
trim(s, c(2, 9))

## select a range
s[3:6]
```

---

alignSpectra                      *Align MassSpectrum objects.*

---

### Description

This function aligns a list of [MassSpectrum](#) objects (spectra alignment is also known as *warping/phase correction*).

### Usage

```
alignSpectra(spectra, halfWindowSize=20, noiseMethod="MAD", SNR=2,
             reference, tolerance=0.002, warpingMethod="lowess",
             allowNoMatches=FALSE, emptyNoMatches=FALSE, ...)
```

### Arguments

|                |  |
|----------------|--|
| spectra        | list, list of <a href="#">MassSpectrum</a> objects.  |
| halfWindowSize | numeric, half window size; see <a href="#">detectPeaks</a> .   |
| noiseMethod    | a noise estimation method; see <a href="#">detectPeaks</a> .   |
| SNR            | single numeric value. SNR is an abbreviation for <i>signal-to-noise-ratio</i> ; see <a href="#">detectPeaks</a> .  |
| reference      | <a href="#">MassPeaks</a> , reference object to which the samples (1) should be aligned. If missing <a href="#">referencePeaks</a> is used; see <a href="#">determineWarpingFunctions</a> .  |
| tolerance      | double, maximal relative deviation of a peak position (mass) to be considered as identical. Must be multiplied by $10^{-6}$ for ppm, e.g. use <code>tolerance=5e-6</code> for 5 ppm; see <a href="#">determineWarpingFunctions</a> . |
| warpingMethod  | used basic warping function; see <a href="#">determineWarpingFunctions</a> .   |
| allowNoMatches | logical, don't throw an error if an <a href="#">MassPeaks</a> object could not match to the reference; see <a href="#">determineWarpingFunctions</a> .   |
| emptyNoMatches | logical, if TRUE (default: FALSE) the intensity values of <a href="#">MassSpectrum</a> or <a href="#">MassPeaks</a> objects with missing (NA) warping functions are set to zero; see <a href="#">warpMassSpectra</a> .               |
| ...            | arguments to be passed to <a href="#">detectPeaks</a> , <a href="#">MassSpectrum-method</a> .  |

### Details

`alignSpectra` is a wrapper function around [detectPeaks](#), [determineWarpingFunctions](#) and [warpMassSpectra](#). Please call these functions manually if you need finer control (e.g. plotting of warping functions).

### Value

Returns a list of aligned [MassSpectrum](#) objects.

**Author(s)**

Sebastian Gibb <mail@sebastiangibb.de>

**See Also**

[detectPeaks](#), [determineWarpingFunctions](#), [referencePeaks](#), [warpMassSpectra](#), [MassSpectrum](#)  
demo("warping")

Website: <https://strimmerlab.github.io/software/maldiquant/>

**Examples**

```
## load package
library("MALDIquant")

## load example data
data("fiedler2009subset", package="MALDIquant")

## running typical workflow

## transform intensities
spectra <- transformIntensity(fiedler2009subset, method="sqrt")

## smooth spectra
spectra <- smoothIntensity(spectra, method="MovingAverage")

## baseline correction
spectra <- removeBaseline(spectra)

## align spectra
spectra <- alignSpectra(spectra)
```

---

averageMassSpectra     *Averages [MassSpectrum](#) objects.*

---

**Description**

This function averages [MassSpectrum](#) objects.

**Usage**

```
averageMassSpectra(l, labels, method=c("mean", "median", "sum"), ...)
```

**Arguments**

**l**                    list, list of [MassSpectrum](#) objects.

**labels**               list, list of [factors](#) (one for each [MassSpectrum](#) object) to do groupwise averaging.

method            used aggregation function.  
...                arguments to be passed to underlying functions (currently only `mc.cores` is supported).

### Details

The mass of the averaged `MassSpectrum` object will be the mass of the first non-empty `MassSpectrum` object (of each group).

### Value

Returns a single (no labels given) or a `list` (labels given) of averaged `MassSpectrum` objects.

### Author(s)

Sebastian Gibb <[mail@sebastiangibb.de](mailto:mail@sebastiangibb.de)>

### See Also

`MassSpectrum`, `mergeMassPeaks`

Website: <https://strimmerlab.github.io/software/maldiquant/>

### Examples

```
## load package
library("MALDIquant")

## create four MassSpectrum objects and add them to a list
s <- list(createMassSpectrum(mass=1:5, intensity=1:5),
          createMassSpectrum(mass=1:5, intensity=1:5),
          createMassSpectrum(mass=1:5, intensity=6:10),
          createMassSpectrum(mass=1:5, intensity=6:10))

## average all four MassSpectrum objects into a single new one
## by sum their intensities
## (no labels, returns only one new MassSpectrum object)
summedSpectra <- averageMassSpectra(s, method="sum")

## only average MassSpectrum objects in a group
## (e.g. useful for technical replicates)
## (two different labels, returns a list of two new MassPeaks objects)
groups <- factor(c("a", "a", "b", "b"), levels=c("a", "b"))
averagedSpectra <- averageMassSpectra(s, labels=groups, method="mean")
```



---

|          |  |
|----------|--|
| binPeaks | <i>Align Peaks into discrete bins.</i> |
|----------|--|

---

### Description

This function looks for similar peaks (mass) across [MassPeaks](#) objects and equalizes their mass.

### Usage

```
binPeaks(l, method=c("strict", "relaxed", "reference"), tolerance=0.002)
```

### Arguments

|           |  |
|-----------|--|
| l         | list, list of <a href="#">MassPeaks</a> objects.   |
| method    | bin creation rule. "strict" creates bins never containing two or more peaks of the same sample. "relaxed" allows multiple peaks of the same sample in one bin. "reference" generates bins around the mass values from the first <a href="#">MassPeaks</a> object in l. |
| tolerance | double, maximal relative deviation of a peak position (mass) to be considered as identical. Must be multiplied by $10^{-6}$ for ppm, e.g. use tolerance=5e-6 for 5 ppm.  |

### Details

The algorithm is based on the following workflow:

1. Put all mass in a sorted vector.
2. Calculate differences between each neighbor.
3. Divide the mass vector at the largest gap (largest difference) and form a left and a right bin.
4. Rerun step 3 for the left and/or the right bin if they don't fulfill the following criteria:
  - All peaks in a bin are near to the mean (method == "strict" or method == "relaxed") ( $\text{abs}(\text{mass} - \text{meanMass}) / \text{meanMass} < \text{tolerance}$ ) or the reference mass (method == "reference";  $\text{abs}(\text{mass} - \text{reference}) / \text{reference} < \text{tolerance}$ ).
  - method == "strict": The bin doesn't contain two or more peaks of the same sample.

method == "strict": The new peak positions (mass value) are the mean mass of a bin.

method == "relaxed": The new peak positions for the highest peaks of each sample in a bin are generated by the mean mass of this peaks. The lower peaks are not changed.

method == "reference": The new peak positions for the highest peaks of each sample in a bin are generated by the mass of peaks of the first [MassPeaks](#) object. Lower peaks are not changed.

### Value

Returns a [list](#) of mass adjusted [MassPeaks](#) objects.

**Author(s)**

Sebastian Gibb <mail@sebastiangibb.de>

**See Also**

[intensityMatrix](#), [MassPeaks](#)

Website: <https://strimmerlab.github.io/software/maldiquant/>

**Examples**

```
## load package
library("MALDIquant")

## create two MassPeaks objects
p <- list(createMassPeaks(mass=seq(100, 500, 100), intensity=1:5),
          createMassPeaks(mass=c(seq(100.2, 300.2, 100), 395), intensity=1:4))

binnedPeaks <- binPeaks(p, tolerance=0.002)

## compare result
iM1 <- intensityMatrix(p)
iM2 <- intensityMatrix(binnedPeaks)

all(dim(iM1) == c(2, 9)) # TRUE
all(dim(iM2) == c(2, 6)) # TRUE

show(iM2)

## increase tolerance
binnedPeaks <- binPeaks(p, tolerance=0.1)

iM3 <- intensityMatrix(binnedPeaks)

all(dim(iM3) == c(2, 5)) # TRUE

show(iM3)

## differences between "strict" and "relaxed"
p <- c(createMassPeaks(mass=c(1, 1.01, 3), intensity=c(2, 1, 1)),
       createMassPeaks(mass=c(0.99, 3), intensity=rep(1, 2)),
       createMassPeaks(mass=c(1.02, 3), intensity=rep(1, 2)))

intensityMatrix(binPeaks(p, method="strict", tolerance=0.05))
intensityMatrix(binPeaks(p, method="relaxed", tolerance=0.05))

## use a reference
ref <- createMassPeaks(mass=c(1, 3), intensity=rep(1, 2))
## include the reference
intensityMatrix(binPeaks(c(ref, p), method="reference", tolerance=0.05))
## drop the reference
intensityMatrix(binPeaks(c(ref, p), method="reference", tolerance=0.05)[-1])
```

---

`calibrateIntensity-methods`*Calibrates intensities of a MassSpectrum object.*

---

## Description

This function calibrates (normalize) intensities of `MassSpectrum` objects.

## Usage

```
## S4 method for signature 'MassSpectrum'  
calibrateIntensity(object,  
  method=c("TIC", "PQN", "median"), range, ...)  
## S4 method for signature 'list'  
calibrateIntensity(object,  
  method=c("TIC", "PQN", "median"), range, ...)
```

## Arguments

|                     |   |
|---------------------|---|
| <code>object</code> | <code>MassSpectrum</code> object or a list of <code>MassSpectrum</code> objects.  |
| <code>method</code> | the calibration method to be used. This should be one of "TIC", "PQN" or "median". See 'Details' section.   |
| <code>range</code>  | numeric of length 2, if given the scaling factor is calculated on the mass range from <code>range[1L]</code> to <code>range[2L]</code> and applied to the whole spectrum. |
| <code>...</code>    | arguments to be passed to other functions. Currently only <code>mc.cores</code> is supported if <code>object</code> is a list.  |

## Details

A number of different calibration methods are provided:

"TIC": The TIC (*Total Ion Current*) of a `MassSpectrum` object is set to one. If `range` is given the *TIC* is only calculated for the intensities in the specified mass range.

"PQN": The PQN (*Probabilistic Quotient Normalization*) is described in *Dieterle et al 2006*. `calibrateIntensity` uses the following algorithm:

1. Calibrate all spectra using the "TIC" calibration.
2. Calculate a median reference spectrum.
3. Calculate the quotients of all intensities of the spectra with those of the reference spectrum.
4. Calculate the median of these quotients for each spectrum.
5. Divide all intensities of each spectrum by its median of quotients.

"median": The median of intensities of a `MassSpectrum` object is set to one.

## Value

Returns a modified `MassSpectrum` object with calibrated intensities.

**Author(s)**

Sebastian Gibb <mail@sebastiangibb.de>

**References**

F. Dieterle, A. Ross, G. Schlotterbeck, and Hans Senn. 2006. Probabilistic quotient normalization as robust method to account for dilution of complex biological mixtures. Application in 1H NMR metabonomics. *Analytical Chemistry* 78(13): 4281-4290.

**See Also**

[MassSpectrum](#)

Website: <https://strimmerlab.github.io/software/maldiquant/>

**Examples**

```
## load package
library("MALDIquant")

## load example data
data("fiedler2009subset", package="MALDIquant")

## baseline correction
b <- removeBaseline(fiedler2009subset)

## calibrate intensity values
calibrateIntensity(b, method="TIC")

## calibrate intensity values using TIC for a specific mass range
calibrateIntensity(b, method="TIC", range=c(3000, 5000))
```

---

createMassPeaks      *Creates a MassPeaks object.*

---

**Description**

This function creates a [MassPeaks](#) object. Normally it shouldn't called by the user. Try [detectPeaks](#), [MassSpectrum-method](#) instead.

**Usage**

```
createMassPeaks(mass, intensity, snr=rep.int(NA_real_, length(intensity)),
                metaData=list())
```

**Arguments**

|           |   |
|-----------|---|
| mass      | vector, mass or mass-to-charge ratio.                   |
| intensity | vector, intensities for measured mass-to-charge ratios. |
| snr       | vector, signal-to-noise ratios for intensity values.    |
| metaData  | list, some metadata to describe the peaks.              |

**Value**

Returns a [MassPeaks](#) object.

**Author(s)**

Sebastian Gibb <[mail@sebastiangibb.de](mailto:mail@sebastiangibb.de)>

**See Also**

[detectPeaks](#), [MassSpectrum-method](#), [MassPeaks](#)

Website: <https://strimmerlab.github.io/software/maldiquant/>

**Examples**

```
## load package
library("MALDIquant")

## create a MassPeaks object by default constructor
s <- createMassPeaks(mass=1:100, intensity=rnorm(100)^2,
                    metaData=list(name="example peaks"))

## show some details
s
```

---

createMassSpectrum     *Creates a MassSpectrum object.*

---

**Description**

This function creates a [MassSpectrum](#) object.

**Usage**

```
createMassSpectrum(mass, intensity, metaData=list())
```

**Arguments**

|           |  |
|-----------|--|
| mass      | vector, mass or mass-to-charge ratio                   |
| intensity | vector, intensities for measured mass-to-charge ratios |
| metaData  | list, some metadata to describe the spectrum           |

**Value**

Returns a [MassSpectrum](#) object.

**Author(s)**

Sebastian Gibb <mail@sebastiangibb.de>

**See Also**

[MassSpectrum](#)

Website: <https://strimmerlab.github.io/software/maldiquant/>

**Examples**

```
## load package
library("MALDIquant")

## create a MassSpectrum object by default constructor
s <- createMassSpectrum(mass=1:100, intensity=rnorm(100)^2,
                       metaData=list(name="example spectrum"))

## show some details
s
```

---

detectPeaks-methods    *Detects peaks in a MassSpectrum object.*

---

**Description**

This method looks for peaks in mass spectrometry data (represented by a [MassSpectrum](#) object). A peak is a local maximum above a user defined noise threshold.

**Usage**

```
## S4 method for signature 'MassSpectrum'
detectPeaks(object,
            halfWindowSize=20, method=c("MAD", "SuperSmoother"), SNR=2,
            ...)
## S4 method for signature 'list'
detectPeaks(object, ...)
```

**Arguments**

**object**            [MassSpectrum](#) object or a list of [MassSpectrum](#) objects.

**halfWindowSize**   numeric, half window size.  
The resulting window reaches from `mass[currentIndex-halfWindowSize]` to `mass[currentIndex+halfWindowSize]`. A local maximum have to be the highest one in the given window to be recognized as peak.

method a noise estimation function; see [estimateNoise,MassSpectrum-method](#).  
SNR single numeric value. SNR is an abbreviation for *signal-to-noise-ratio*. A local maximum has to be higher than SNR\*noise to be recognize as peak.  
... arguments to be passed to [estimateNoise,MassSpectrum-method](#). If object is a list `mc.cores` is also supported.

### Value

Returns a [MassPeaks](#) object.

### Author(s)

Sebastian Gibb <[mail@sebastiangibb.de](mailto:mail@sebastiangibb.de)>

### See Also

[MassPeaks](#), [MassSpectrum](#), [estimateNoise,MassSpectrum-method](#)

`demo("peaks")`

Website: <https://strimmerlab.github.io/software/maldiquant/>

### Examples

```
## load package
library("MALDIquant")

## load example data
data("fiedler2009subset", package="MALDIquant")

## choose only the first mass spectrum
s <- fiedler2009subset[[1]]

## transform intensities
s <- transformIntensity(s, method="sqrt")

## smoothing spectrum
s <- smoothIntensity(s, method="MovingAverage")

## remove baseline
s <- removeBaseline(s)

## plot spectrum
plot(s)

## call peak detection
p <- detectPeaks(s)

## draw peaks on the plot
points(p)

## label 10 highest peaks
top10 <- intensity(p) %in% sort(intensity(p), decreasing=TRUE)[1:10]
```

```
labelPeaks(p, index=top10)
```

---

```
determineWarpingFunctions
```

*Determine warping functions of MassPeaks objects.*

---

### Description

This function determines a warping function for a list of [AbstractMassObject](#) objects (warping is also known as *phase correction/spectra alignment*).

### Usage

```
determineWarpingFunctions(l, reference, tolerance=0.002,
                          method=c("lowess", "linear", "quadratic", "cubic"),
                          allowNoMatches=FALSE,
                          plot=FALSE, plotInteractive=FALSE, ...)
```

### Arguments

|                 |  |
|-----------------|--|
| l               | list, list of <a href="#">MassPeaks</a> objects.   |
| reference       | <a href="#">MassPeaks</a> , reference object to which the samples (l) should be aligned. If missing <a href="#">referencePeaks</a> is used.  |
| tolerance       | double, maximal relative deviation of a peak position (mass) to be considered as identical. Must be multiplied by $10^{-6}$ for ppm, e.g. use <code>tolerance=5e-6</code> for 5 ppm. |
| method          | used basic warping function.   |
| allowNoMatches  | logical, don't throw an error if an <a href="#">MassPeaks</a> object could not match to the reference.   |
| plot            | logical, if TRUE a warping plot is drawn for each sample.  |
| plotInteractive | logical, if FALSE a non-interactive device (e.g. pdf) is used for warping plots.   |
| ...             | arguments to be passed to <code>warpingFunction</code>   |

### Details

`warpingFunction`: `determineWarpingFunctions` estimates a warping function to overcome the difference between mass in reference and in the current sample. To calculate the differences each reference peak would match with the highest sample peak in the nearer neighborhood (defined by mass of reference peak\*`tolerance`).

`allowNoMatches`: If `allowNoMatches` is TRUE a warning instead of an error is thrown if an [MassPeaks](#) object could not match to the reference. The returned list of warping functions will contain NA for this object (same index in the list). `plotInteractive`: If `plot` is TRUE a lot of output is created (each sample in l gets its own plot). That's why an non-interactive devices is recommended:



```
## create a device
pdf()
## calculate warping functions
w <- determineWarpingFunctions(p, plot=TRUE)
## close device
dev.off()
```

### Value

Returns a list of individual warping functions. The attribute `nmatch` contains the number of matches of each `MassPeaks` element in `l` against reference.

### Author(s)

Sebastian Gibb <mail@sebastiangibb.de>

### See Also

[referencePeaks](#), [warpMassPeaks](#), [warpMassSpectra](#), [MassPeaks](#)  
`demo("warping")`

Website: <https://strimmerlab.github.io/software/maldiquant/>

### Examples

```
## load package
library("MALDIquant")

## create a reference MassPeaks object
r <- createMassPeaks(mass=1:5, intensity=1:5)

## create test samples
p <- list(createMassPeaks(mass=((1:5)*1.01), intensity=1:5),
          createMassPeaks(mass=((1:5)*0.99), intensity=1:5))

## create an interactive device with 2 rows
par(mfrow=c(2, 1))
## calculate warping function
## (using a linear function as basic warping function)
## and show warping plot
w <- determineWarpingFunctions(p, tolerance=0.02, method="linear",
                              plot=TRUE, plotInteractive=TRUE)
par(mfrow=c(1, 1))

## access number of matches
attr(w, "nmatch")

## w contains the individual warping functions
warpedPeaks <- warpMassPeaks(p, w)

## compare results
```

```
all(mass(r) == mass(warpedPeaks[[1]])) # TRUE
all(mass(r) == mass(warpedPeaks[[2]])) # TRUE

## realistic example

## load example data
data("fiedler2009subset", package="MALDIquant")

## running typical workflow

## use only four spectra of the subset
spectra <- fiedler2009subset[1:4]

## transform intensities
spectra <- transformIntensity(spectra, method="sqrt")

## smooth spectra
spectra <- smoothIntensity(spectra, method="MovingAverage")

## baseline correction
spectra <- removeBaseline(spectra)

## detect peaks
peaks <- detectPeaks(spectra)

## create an interactive device with 2 rows
par(mfrow=c(4, 1))
## calculate warping functions (using LOWESS based basic function [default])
w <- determineWarpingFunctions(peaks, plot=TRUE, plotInteractive=TRUE)
par(mfrow=c(1, 1))

## realistic example with user defined reference/calibration peaks

## use the workflow above for fiedler2009subset

## create reference peaks
refPeaks <- createMassPeaks(mass=c(1207, 1264, 1351, 1466, 1616, 2769, 2932,
                                   3191, 3262, 4091, 4209, 5904, 7762, 9285),
                           intensity=rep(1, 14))

## create an interactive device with 2 rows
par(mfrow=c(4, 1))
## calculate warping functions (using a quadratic function as basic function)
w <- determineWarpingFunctions(peaks, reference=refPeaks, method="quadratic",
                              plot=TRUE, plotInteractive=TRUE)
par(mfrow=c(1, 1))
```

---

 estimateBaseline-methods

*Estimates the baseline of a MassSpectrum object.*


---

## Description

This method estimates the baseline of mass spectrometry data (represented by a [MassSpectrum](#) object).

## Usage

```
## S4 method for signature 'MassSpectrum'
estimateBaseline(object,
  method=c("SNIP", "TopHat", "ConvexHull", "median"),
  ...)
```

## Arguments

|        |   |
|--------|---|
| object | <a href="#">MassSpectrum</a> object   |
| method | used baseline estimation method, one of "SNIP", "TopHat", "ConvexHull" or "median". |
| ...    | arguments to be passed to method  |

## Details

"SNIP": This baseline estimation is based on the Statistics-sensitive Non-linear Iterative Peak-clipping algorithm (SNIP) described in Ryan et al 1988.

The algorithm based on the following equation:

$$y_i(k) = \min\left\{y_i, \frac{(y_{i-k} + y_{i+k})}{2}\right\}$$

It has two additional arguments namely `iterations` and `decreasing`. `iterations` controls the window size ( $k$ ; similar to `halfWindowSize` in "TopHat", "Median") of the algorithm. The resulting window reaches from `mass[cur_index-iterations]` to `mass[cur_index+iterations]`. `decreasing`: In Morhac 2009 a decreasing clipping window is suggested to get a smoother baseline. For `decreasing = TRUE` (`decreasing = FALSE`)  $k=iterations$  is decreased (increased) by one until zero (`iterations`) is reached. The default setting is `decreasing = TRUE`.

"TopHat": This algorithm applies a moving minimum (erosion filter) and subsequently a moving maximum (dilation filter) filter on the intensity values. The implementation is based on van Herk 1996. It has an additional `halfWindowSize` argument determining the half size of the moving window for the TopHat filter. The resulting window reaches from `mass[cur_index-halfWindowSize]` to `mass[cur_index+halfWindowSize]`.

"ConvexHull": The baseline estimation is based on a convex hull constructed below the spectrum.

"median": This baseline estimation uses a moving median. It is based on `runmed`. The additional argument `halfWindowSize` corresponds to the `k` argument in `runmed` ( $k = 2 * \text{halfWindowSize} + 1$ ) and controls the half size of the moving window. The resulting window reaches from `mass[cur_index-halfWindowSize]` to `mass[cur_index+halfWindowSize]`.

### Value

Returns a two column matrix (first column: mass, second column: intensity) of the estimated baseline.

### Author(s)

Sebastian Gibb <[mail@sebastiangibb.de](mailto:mail@sebastiangibb.de)>

### References

"SNIP":

C.G. Ryan, E. Clayton, W.L. Griffin, S.H. Sie, and D.R. Cousens. 1988. Snip, a statistics-sensitive background treatment for the quantitative analysis of pixe spectra in geoscience applications. Nuclear Instruments and Methods in Physics Research Section B: Beam Interactions with Materials and Atoms, 34(3): 396-402.

M. Morhac. 2009. An algorithm for determination of peak regions and baseline elimination in spectroscopic data. Nuclear Instruments and Methods in Physics Research Section A: Accelerators, Spectrometers, Detectors and Associated Equipment, 600(2), 478-487.

"TopHat":

M. van Herk. 1992. A Fast Algorithm for Local Minimum and Maximum Filters on Rectangular and Octagonal Kernels. Pattern Recognition Letters 13.7: 517-521.

J. Y. Gil and M. Werman. 1996. Computing 2-Dimensional Min, Median and Max Filters. IEEE Transactions: 504-507.

"ConvexHull":

Andrew, A. M. 1979. Another efficient algorithm for convex hulls in two dimensions. Information Processing Letters, 9(5), 216-219.

### See Also

[MassSpectrum](#), [removeBaseline](#), [MassSpectrum-method](#)

`demo("baseline")`

Website: <https://strimmerlab.github.io/software/maldiquant/>

### Examples

```
## load package
library("MALDIquant")

## load example data
data("fiedler2009subset", package="MALDIquant")
```

```
## choose only the first mass spectrum
s <- fiedler2009subset[[1]]

## SNIP
plot(s)

## estimate baseline
b <- estimateBaseline(s, method="SNIP", iterations=100)

## draw baseline on the plot
lines(b, col="red")

## TopHat
plot(s)

## estimate baseline (try different parameters)
b1 <- estimateBaseline(s, method="TopHat", halfWindowSize=75)
b2 <- estimateBaseline(s, method="TopHat", halfWindowSize=150)

## draw baselines on the plot
lines(b1, col=2)
lines(b2, col=3)

## draw legend
legend(x="topright", lwd=1, legend=paste0("halfWindowSize=", c(75, 150)),
      col=c(2, 3))

## ConvexHull
plot(s)

## estimate baseline
b <- estimateBaseline(s, method="ConvexHull")

## draw baseline on the plot
lines(b, col="red")

## Median
plot(s)

## estimate baseline
b <- estimateBaseline(s, method="median")

## draw baseline on the plot
lines(b, col="red")
```

## Description

This method estimates the noise of mass spectrometry data (represented by a [MassSpectrum](#) object).

## Usage

```
## S4 method for signature 'MassSpectrum'  
estimateNoise(object,  
  method=c("MAD", "SuperSmoother"),  
  ...)
```

## Arguments

|        |  |
|--------|--|
| object | <a href="#">MassSpectrum</a> object                            |
| method | used noise estimation method, one of "MAD" or "SuperSmoother". |
| ...    | arguments to be passed to method.                              |

## Details

"MAD": This function estimates the noise of mass spectrometry data by calculating the median absolute deviation, see also [mad](#).

"SuperSmoother": This function estimates the noise of mass spectrometry data using Friedman's Super Smoother. Please refer [supsmu](#) for details and additional arguments.

## Value

Returns a two column matrix (first column: mass, second column: intensity) of the estimated noise.

## Author(s)

Sebastian Gibb <[mail@sebastiangibb.de](mailto:mail@sebastiangibb.de)>

## See Also

[MassSpectrum](#), [detectPeaks](#), [MassSpectrum-method](#), [mad](#), [supsmu](#)  
Website: <https://strimmerlab.github.io/software/maldiquant/>

## Examples

```
## load package  
library("MALDIquant")  
  
## load example data  
data("fiedler2009subset", package="MALDIquant")  
  
## choose only the first mass spectrum  
s <- fiedler2009subset[[1]]  
  
## transform intensities
```

```
s <- transformIntensity(s, method="sqrt")

## remove baseline
s <- removeBaseline(s)

## plot spectrum
plot(s)

## estimate noise
nm <- estimateNoise(s, method="MAD")
nss <- estimateNoise(s, method="SuperSmoother")

## draw noise on the plot
lines(nm, col=2)
lines(nss, col=4)

## draw legend
legend(x="topright", lwd=1, legend=c("MAD", "SuperSmoother"),
      col=c(2, 4))
```

---

fiedler2009subset      *Example Mass Spectra (raw)*

---

## Description

This dataset contains 16 example mass spectra. It is used to demonstrate the usage of [MALDIquant](#)-package.

## Usage

```
data(fiedler2009subset)
```

## Format

A list containing 16 [MassSpectrum](#)-class objects.

## Details

The dataset is a subset of data used in *Fiedler et al 2009*.

It contains spectra of 8 different patients (each one has 2 technical replicates).

| list_index | laboratory | patient_id | sex    | age | type    |
|------------|------------|------------|--------|-----|---------|
| 1          | Leipzig    | LC77       | male   | 37  | control |
| 2          | Leipzig    | LC77       | male   | 37  | control |
| 3          | Leipzig    | LC213      | female | 51  | control |
| 4          | Leipzig    | LC213      | female | 51  | control |
| 5          | Leipzig    | LT178      | male   | 58  | cancer  |
| 6          | Leipzig    | LT178      | male   | 58  | cancer  |
| 7          | Leipzig    | LT157      | male   | 60  | cancer  |
| 8          | Leipzig    | LT157      | male   | 60  | cancer  |

|    |            |       |        |    |         |
|----|------------|-------|--------|----|---------|
| 9  | Heidelberg | HC49  | male   | 43 | control |
| 10 | Heidelberg | HC49  | male   | 43 | control |
| 11 | Heidelberg | HC54  | female | 71 | control |
| 12 | Heidelberg | HC54  | female | 71 | control |
| 13 | Heidelberg | HT151 | male   | 53 | cancer  |
| 14 | Heidelberg | HT151 | male   | 53 | cancer  |
| 15 | Heidelberg | HT429 | female | 58 | cancer  |
| 16 | Heidelberg | HT429 | female | 58 | cancer  |

## References

G.M. Fiedler, A.B. Leichtle, J. Kase, S. Baumann, U. Ceglarek, K. Felix, T. Conrad, H. Witzigmann, A. Weimann, C. Schütte, J. Hauss, M. Büchler and J. Thiery  
 “Serum Peptidome Profiling Revealed Platelet Factor 4 as a Potential Discriminating Peptide Associated with Pancreatic Cancer”  
 Clinical Cancer Research, 11(15): 3812-3819, 2009  
 ISSN 1557-3265; doi:10.1158/1078-0432.CCR-08-2701  
<https://clincancerres.aacrjournals.org/content/15/11/3812>

## See Also

[MassSpectrum-class](#)

Website: <https://strimmerlab.github.io/software/maldiquant/>

---

|             |                                     |
|-------------|-------------------------------------|
| filterPeaks | <i>Removes less frequent peaks.</i> |
|-------------|-------------------------------------|

---

## Description

This function removes infrequently occurring peaks in a list of [MassPeaks](#) objects.

## Usage

```
filterPeaks(l, minFrequency, minNumber, labels, mergeWhitelists=FALSE)
```

## Arguments

|              |   |
|--------------|---|
| l            | list, list of <a href="#">MassPeaks</a> objects.  |
| minFrequency | double, remove all peaks which occur in less than minFrequency*length(l) <a href="#">MassPeaks</a> objects. It is a relative threshold.   |
| minNumber    | double, remove all peaks which occur in less than minNumber <a href="#">MassPeaks</a> objects. It is an absolute threshold.   |
| labels       | factor, (one for each <a href="#">MassPeaks</a> object) to do groupwise filtering. The <i>levels</i> of the factor label define the groups. If not specified a single group is assumed. |



mergeWhitelists

logical, if FALSE the filtering criteria are applied groupwise. If TRUE peaks that survive the filtering in one group (level of labels) these peaks are also kept in other groups even if their frequencies are below minFrequency.

### Details

For mergeWhitelists=FALSE the filtering uses a separate peak whitelist for each group specified by labels, and is done independently in each group. For mergeWhitelists=TRUE the peak whitelists are combined, which means that peaks that occur frequently in at least one group are also kept in all other groups.

If both minFrequency and minNumber arguments are specified the more stringent threshold is used.

### Value

Returns a [list](#) of filtered [MassPeaks](#) objects.

### Author(s)

Sebastian Gibb <[mail@sebastiangibb.de](mailto:mail@sebastiangibb.de)>

### See Also

[intensityMatrix](#), [MassPeaks](#)

Website: <https://strimmerlab.github.io/software/maldiquant/>

### Examples

```
## load package
library("MALDIquant")

## create four MassPeaks objects and add them to the list
p <- list(createMassPeaks(mass=1:2, intensity=1:2),
          createMassPeaks(mass=1:3, intensity=1:3),
          createMassPeaks(mass=1:4, intensity=1:4),
          createMassPeaks(mass=1:5, intensity=1:5))

## only keep peaks which occur in all MassPeaks objects
filteredPeaks <- filterPeaks(p, minFrequency=1)

## compare result
intensities <- intensityMatrix(filteredPeaks)

## peaks at mass 3,4,5 are removed
all(dim(intensities) == c(4, 2)) # TRUE
all(intensities[,1] == 1)      # TRUE
all(intensities[,2] == 2)      # TRUE

## only keep peaks which occur in all MassPeaks objects in a group
## (e.g. useful for technical replicates)
groups <- factor(c("a", "a", "b", "b"), levels=c("a", "b"))
```

```

filteredPeaks <- filterPeaks(p, minFrequency=1, labels=groups)

## peaks at mass 3 were removed in group "a"
filteredPeaks[groups == "a"]

## peaks at mass 5 were removed in group "b"
filteredPeaks[groups == "b"]

## only keep peaks which occur at least twice in a group
groups <- factor(c("a", "a", "b", "b", "b"), levels=c("a", "b"))
filteredPeaks <- filterPeaks(c(p, p[[3]]), minNumber=2, labels=groups)

## peaks at mass 3 were removed in group "a"
filteredPeaks[groups == "a"]

## peaks at mass 5 were removed in group "b"
filteredPeaks[groups == "b"]

## apply different minFrequency arguments to each group
groups <- factor(c("a", "a", "b", "b", "b"), levels=c("a", "b"))
filteredPeaks <- filterPeaks(c(p, p[[3]]), minFrequency=c(1, 2/3), labels=groups)
intensityMatrix(filteredPeaks)
#      1 2 3 4
#[1,] 1 2 NA NA
#[2,] 1 2 NA NA
#[3,] 1 2 3 4
#[4,] 1 2 3 4
#[4,] 1 2 3 4

## demonstrate the use of mergeWhitelists
groups <- factor(c("a", "a", "b", "b"), levels=c("a", "b"))

## default behaviour
filteredPeaks <- filterPeaks(p, minNumber=2, labels=groups)
intensityMatrix(filteredPeaks)
#      1 2 3 4
#[1,] 1 2 NA NA
#[2,] 1 2 NA NA
#[3,] 1 2 3 4
#[4,] 1 2 3 4

## use mergeWhitelists=TRUE to keep peaks of group "a" that match all filtering
## criteria in group "b"
## (please note that mass == 3 is not removed in the second MassPeaks object)
filteredPeaks <- filterPeaks(p, minNumber=2, labels=groups,
                             mergeWhitelists=TRUE)
intensityMatrix(filteredPeaks)
#      1 2 3 4
#[1,] 1 2 NA NA
#[2,] 1 2 3 NA
#[3,] 1 2 3 4
#[4,] 1 2 3 4

```



```
## find empty objects (there should not be any one)
findEmptyMassObjects(peakList)

## add an empty MassPeaks object to the list
peakList[[3]] <- createMassPeaks(mass=double(), intensity=double(),
                                metaData=list(name="empty MassPeaks object"))

## look for empty objects (isEmptyIdx == 3)
(isEmptyIdx <- findEmptyMassObjects(peakList))

## to remove all empty MassObjects from a list
length(peakList) # 3
peakList <- removeEmptyMassObjects(peakList)
length(peakList) # 2; WARNING: all indices could changed
```

---

|                 |  |
|-----------------|--|
| intensityMatrix | <i>Converts a list of MassPeaks objects into a matrix.</i> |
|-----------------|--|

---

## Description

This function converts a [list](#) of [MassPeaks](#) objects into a [matrix](#).

## Usage

```
intensityMatrix(peaks, spectra)
```

## Arguments

|         |   |
|---------|---|
| peaks   | list, list of <a href="#">MassPeaks</a> objects.  |
| spectra | list, list of <a href="#">MassSpectrum</a> objects. If a peak is missing the corresponding intensity value of the spectrum is used. If spectra is missing NA is used instead. |

## Details

peaks have to be binned by [binPeaks](#) before calling [intensityMatrix](#).

## Value

Returns a [matrix](#) containing intensities of all [MassPeaks](#) objects of peaks and interpolated intensity values for missing peaks if spectra was given or NA otherwise.

The [matrix](#) has `length(peaks)` rows (one row for each sample) and `length(unique(mass))` columns. There is an additional attribute "mass" that stores the mass values.

## Author(s)

Sebastian Gibb <[mail@sebastiangibb.de](mailto:mail@sebastiangibb.de)>

**See Also**

[binPeaks](#), [MassPeaks](#), [MassSpectrum](#)

Website: <https://strimmerlab.github.io/software/maldiquant/>

**Examples**

```
## load package
library("MALDIquant")

## create example MassPeaks objects
p <- list(createMassPeaks(mass=1:4,
                        intensity=11:14,
                        metaData=list(name="test mass peaks 1")),
          createMassPeaks(mass=2:5,
                        intensity=22:25,
                        metaData=list(name="test mass peaks 2")))

## converts MassPeaks objects into a matrix
intensityMatrix(p)

## realistic example

## load example data
data("fiedler2009subset", package="MALDIquant")

## transform intensities
s <- transformIntensity(fiedler2009subset, method="sqrt")

## smoothing spectrum
s <- smoothIntensity(s, method="MovingAverage")

## remove baseline
s <- removeBaseline(s)

## call peak detection
p <- detectPeaks(s)

## bin peaks
p <- binPeaks(p)

## convert MassPeaks objects into a matrix with missing intensity
## values
intensityMatrix(p)

## convert MassPeaks and MassSpectrum objects into a matrix without
## missing intensity values
intensityMatrix(p, s)
```

---

isMassSpectrum      *Tests for MassSpectrum or MassPeaks object.*

---

### Description

These functions test for a [MassSpectrum](#) or [MassPeaks](#) object.

### Usage

```
isMassSpectrum(x)
```

```
isMassPeaks(x)
```

### Arguments

x                    object to be tested.

### Value

Returns [TRUE](#) or [FALSE](#) depending on whether its argument is an [MassSpectrum](#) or [MassPeaks](#) object.

### Author(s)

Sebastian Gibb <[mail@sebastiangibb.de](mailto:mail@sebastiangibb.de)>

### See Also

[MassPeaks](#), [MassSpectrum](#), [AbstractMassObject](#)

Website: <https://strimmerlab.github.io/software/maldiquant/>

### Examples

```
## load package
library("MALDIquant")

## create a MassPeaks object
peaks <- createMassPeaks(mass=1:100, intensity=1:100,
                        metaData=list(name="example 1"))

## test
isMassPeaks(peaks)      # returns TRUE
isMassSpectrum(peaks)   # returns FALSE
isMassPeaks(double())   # returns FALSE
```

---

isMassSpectrumList      *Tests a list of MassSpectrum or MassPeaks objects.*

---

## Description

These functions test a [list](#) whether containing [MassSpectrum](#) or [MassSpectrum](#) objects.

## Usage

```
isMassSpectrumList(x)
```

```
isMassPeaksList(x)
```

## Arguments

x                      object to be tested.

## Value

Returns [TRUE](#) or [FALSE](#) depending on whether its argument is a [list](#) of [MassSpectrum](#) or [MassPeaks](#) objects.

## Author(s)

Sebastian Gibb <[mail@sebastiangibb.de](mailto:mail@sebastiangibb.de)>

## See Also

[MassPeaks](#), [MassSpectrum](#), [AbstractMassObject](#)

Website: <https://strimmerlab.github.io/software/maldiquant/>

## Examples

```
## load package
library("MALDIquant")

## create list
p <- list()

## test list
isMassPeaksList(p) # returns FALSE

## create two MassPeaks objects and add them to the list
p <- createMassPeaks(mass=1:100, intensity=1:100,
                    metaData=list(name="example 1"))
p <- createMassPeaks(mass=1:100, intensity=1:100,
                    metaData=list(name="example 2"))

## test list
```

```
isMassPeaksList(p) # returns TRUE
isMassSpectrumList(p) # returns FALSE
```

---

labelPeaks-methods      *Draws peak labels to plot.*

---

## Description

[labelPeaks](#) draws the corresponding mass values on top of the peaks stored in a [MassPeaks](#) object to a plot.

## Usage

```
## S4 method for signature 'MassPeaks'
labelPeaks(object,
  index,
  mass,
  labels,
  digits=3, underline=TRUE,
  verticalOffset=abs(diff(par("usr")[3:4]))*0.01,
  absoluteVerticalPos,
  adj=c(0.5, 0), cex=0.7, srt=0,
  avoidOverlap=FALSE,
  arrowLength=0, arrowLwd=0.5, arrowCol=1,
  ...)
```

## Arguments

|                                 |   |
|---------------------------------|---|
| object                          | <a href="#">MassPeaks</a> object.   |
| index                           | integer/logical, indices of peaks to label.   |
| mass                            | numeric, mass of peaks to label.  |
| labels                          | character, use labels instead of mass values as peak label.   |
| digits                          | integer, number of decimal places.  |
| underline                       | logical, underline peak values?   |
| verticalOffset                  | numeric, move label vertically (relative to peak height).   |
| absoluteVerticalPos             | numeric, absolute y value for the label. If missing verticalOffset is used.   |
| adj                             | numeric, adjust text to the left, center, right and top, center, bottom; see <a href="#">text</a> .                   |
| cex                             | numeric, font size, see <a href="#">par</a> .   |
| srt                             | numeric, the label rotation in degrees.   |
| avoidOverlap                    | logical, try to find label coordinates to avoid overlap.  |
| arrowLength, arrowLwd, arrowCol | arrow parameters, possible vectors. NA values in arrowCol cause the arrow to be omitted, see <a href="#">arrows</a> . |
| ...                             | arguments to be passed to <a href="#">text</a> .  |



## Details

Please note that `avoidOverlap = TRUE` is just supported for `srt %% 90 == 0` (means `srt` has to be a multiple of 90 degree).

## Author(s)

Sebastian Gibb

## See Also

[MassPeaks](#), [plot](#), [AbstractMassObject](#), [missing-method](#)

Website: <https://strimmerlab.github.io/software/maldiquant/>

## Examples

```
## load package
library("MALDIquant")

## create a MassPeaks object from scratch
p <- createMassPeaks(mass=1:20, intensity=sample(x=100:10000, size=20),
                    metaData=list(name="example"))

## plot peaks
plot(p)

## label the first 5 peaks
labelPeaks(p, index=1:5)

## label all peaks in mass range 15 to 20
labelPeaks(p, mass=15:20, underline=FALSE)

## label highest peaks (top 5)
top5 <- intensity(p) %in% sort(intensity(p), decreasing=TRUE)[1:5]
labelPeaks(p, index=top5, col="red")

## real example
data("fiedler2009subset")

## a simplified preprocessing
r <- removeBaseline(fiedler2009subset[[1]])
p <- detectPeaks(r)
plot(p)

## label highest peaks (top 10) and avoid label overlap
top10 <- sort(intensity(p), decreasing=TRUE, index.return=TRUE)$ix[1:10]
labelPeaks(p, index=top10, avoidOverlap=TRUE, digits=1)

## use own labels and rotate by 90 degree
plot(p)
labelPeaks(p, index=top10, labels=paste("TOP", 1:10), underline=FALSE,
```

```
srt=90, adj=c(0, 0.5), col=2)
```

---

MALDIquant-parallel *Parallel Support in Package MALDIquant*

---

## Description

**MALDIquant** offers multi-core support using `mclapply` and `mcmapply`. This approach is limited to unix-based platforms.

Please note that not all functions benefit from parallelisation. Often the overhead to create/copy objects outrun the time saving of parallel runs. This is true for functions that are very fast to compute (e.g. `sqrt`-transformation). That's why the default value for the `mc.cores` argument in all functions is 1L. It depends on the size of the dataset which step (often only `removeBaseline` and `detectPeaks`) benefits from parallelisation.

In general it is faster to encapsulate the complete workflow into a function and parallelise it using `mclapply` instead of using the `mc.cores` argument of each method. The reason is the reduced overhead for object management (only one split/combine is needed instead of doing these operations in each function again and again).

## Details

The following functions/methods support the `mc.cores` argument:

- `trim,list,numeric-method`
- `transformIntensity,list-method`
- `smoothIntensity,list-method`
- `removeBaseline,list-method`
- `calibrateIntensity,list-method`
- `detectPeaks,list-method`
- `alignSpectra`
- `averageMassSpectra`
- `mergeMassPeaks`

## See Also

`mclapply`, `mcmapply`

## Examples

```
## Not run:  
## load package  
library("MALDIquant")  
  
## load example data  
data("fiedler2009subset", package="MALDIquant")  
  
## run single-core baseline correction  
print(system.time(  

```

```

    b1 <- removeBaseline(fiedler2009subset, method="SNIP")
  ))

  if(.Platform$OS.type == "unix") {
    ## run multi-core baseline correction
    print(system.time(
      b2 <- removeBaseline(fiedler2009subset, method="SNIP", mc.cores=2)
    ))
    stopifnot(all.equal(b1, b2))
  }

  ## parallelise complete workflow
  workflow <- function(spectra, cores) {
    s <- transformIntensity(spectra, method="sqrt", mc.cores=cores)
    s <- smoothIntensity(s, method="SavitzkyGolay", halfWindowSize=10,
      mc.cores=cores)
    s <- removeBaseline(s, method="SNIP", iterations=100, mc.cores=cores)
    s <- calibrateIntensity(s, method="TIC", mc.cores=cores)
    detectPeaks(s, method="MAD", halfWindowSize=20, SNR=2, mc.cores=cores)
  }

  if(.Platform$OS.type == "unix") {
    ## parallelise the complete workflow is often faster because the overhead is
    ## reduced
    print(system.time(
      p1 <- unlist(parallel::mclapply(fiedler2009subset,
        function(x)workflow(list(x), cores=1),
        mc.cores=2), use.names=FALSE)
    ))
    print(system.time(
      p2 <- workflow(fiedler2009subset, cores=2)
    ))
    stopifnot(all.equal(p1, p2))
  }

  ## End(Not run)

```

---

 MassPeaks-class

 Class "MassPeaks"
 

---

## Description

[MassPeaks](#) represents extracted peaks of a single spectrum of a MALDI-TOF mass spectrometry measurement.

## Objects from the Class

[createMassPeaks](#): Creates a [MassPeaks](#) object.

**Extends**

Class [AbstractMassObject](#), directly.

**Slots**

snr: vector, signal-to-noise ratio

**Methods**

**labelPeaks** signature( $x = \text{"MassPeaks"}$ ): Draws peak labels to plot. See [labelPeaks, MassPeaks-method](#) for details.

**monoisotopicPeaks** signature( $x = \text{"MassPeaks"}$ ): Finds monoisotopic peaks in peak lists. See [monoisotopicPeaks, MassPeaks-method](#) for details.

**Author(s)**

Sebastian Gibb <[mail@sebastiangibb.de](mailto:mail@sebastiangibb.de)>

**See Also**

[createMassPeaks, detectPeaks, MassSpectrum-method, labelPeaks, MassPeaks-method, AbstractMassObject](#)

Website: <https://strimmerlab.github.io/software/maldiquant/>

---

MassSpectrum-class      *Class "MassSpectrum"*

---

**Description**

[MassSpectrum](#) represents a single spectrum of a MALDI-TOF mass spectrometry measurement. It provides an easy framework for doing some preprocessing steps like peak detection, baseline correction and much more.

**Objects from the Class**

[createMassSpectrum](#): Creates a [MassSpectrum](#) object.

**Extends**

Class [AbstractMassObject](#), directly.

## Methods

- calibrateIntensity** signature(x = "MassSpectrum"): Calibrates the intensity of a `MassSpectrum` object. See [calibrateIntensity, MassSpectrum-method](#) for details.
- detectPeaks** signature(x = "MassSpectrum"): Look for local maxima and estimate noise to extract peaks out of a `MassSpectrum` object. See [detectPeaks, MassSpectrum-method](#) for details.
- estimateBaseline** signature(x = "MassSpectrum"): Estimates the baseline of a `MassSpectrum` object. See [estimateBaseline, MassSpectrum-method](#) for details.
- estimateNoise** signature(x = "MassSpectrum"): Estimates the noise of a `MassSpectrum` object. See [estimateNoise, MassSpectrum-method](#) for details.
- isRegular** signature(object = "MassSpectrum"): Returns FALSE if the frequency of mass values with irregular intervals is greater than threshold (because object was measured in *centroid* mode or some intensity values were filtered).
- removeBaseline** signature(x = "MassSpectrum"): Estimates and removes the baseline of a `MassSpectrum` object. See [removeBaseline, MassSpectrum-method](#) for details.
- smoothIntensity** signature(object = "MassSpectrum"): Smooths the intensities of an `MassSpectrum` object. See [smoothIntensity, MassSpectrum-method](#) for details.
- totalIonCurrent** signature(object = "MassSpectrum"): Accessor function for Total Ion Current (TIC, area under the curve).

## Author(s)

Sebastian Gibb <[mail@sebastiangibb.de](mailto:mail@sebastiangibb.de)>

## See Also

[createMassSpectrum](#), [calibrateIntensity, MassSpectrum-method](#), [detectPeaks, MassSpectrum-method](#), [estimateBaseline, MassSpectrum-method](#), [estimateNoise, MassSpectrum-method](#), [removeBaseline, MassSpectrum-method](#), [smoothIntensity, MassSpectrum-method](#), [AbstractMassObject](#)

Website: <https://strimmerlab.github.io/software/maldiquant/>

## Examples

```
## load package
library("MALDIquant")

## create a MassSpectrum object by default constructor
s <- createMassSpectrum(mass=1:100, intensity=rnorm(100)^2,
                       metaData=list(name="example"))

## show some details
s

## plot spectrum
plot(s)

## get TIC
```

```
totalIonCurrent(s)

## modify intensity and metaData
intensity(s)[1:50] <- 0
metaData(s) <- list(name="modified example")

## plot again
plot(s)
```

---

match.closest

*Relaxed Value Matching*

---

## Description

`match.closest` returns a vector of the positions of (first) matches its first arguments in its second. In contrast to the similar `match` it just accept numeric arguments but has an additional tolerance argument that allows relaxed matching.

## Usage

```
match.closest(x, table, tolerance = Inf, nomatch = NA_integer_)
```

## Arguments

|                        |  |
|------------------------|--|
| <code>x</code>         | numeric, the values to be matched.   |
| <code>table</code>     | numeric, the values to be matched against. In contrast to <code>match</code> table has to be sorted in increasing order.   |
| <code>tolerance</code> | numeric, accepted tolerance. Use <code>Inf</code> to match without restrictions. Could be of length one or the same length as <code>table</code> .                                     |
| <code>nomatch</code>   | numeric, if the difference between the value in <code>x</code> and <code>table</code> is larger than <code>tolerance</code> <code>nomatch</code> is returned. Has to be of length one. |

## Value

An integer vector of the same length as `x` giving the closest position in `table` of the first match or `nomatch` if there is no match.

## See Also

[match](#)

## Examples

```
library("MALDIquant")
match.closest(c(1.1, 1.4, 9.8), 1:10)
# [1] 1 1 10
match.closest(c(1.1, 1.4, 9.8), 1:10, tolerance=0.25)
# [1] 1 NA 10
match.closest(c(1.1, 1.4, 9.8), 1:10, tolerance=0.25, nomatch=0)
# [1] 1 0 10

## this function is most useful if you want to subset an intensityMatrix
## by a few (reference) peaks

## create an example intensityMatrix
im <- matrix(1:10, nrow=2, dimnames=list(NULL, 1:5))
attr(im, "mass") <- 1:5
im
#      1 2 3 4 5
# [1,] 1 3 5 7 9
# [2,] 2 4 6 8 10
# attr(,"mass")
# [1] 1 2 3 4 5

## reference peaks
ref <- c(2.2, 4.8)

im[, match.closest(ref, attr(im, "mass"), tolerance=0.25, nomatch=0)]
#      2 5
# [1,] 3 9
# [2,] 4 10
```

---

|                |   |
|----------------|---|
| mergeMassPeaks | Merges <a href="#">MassPeaks</a> objects. |
|----------------|---|

---

## Description

This function merges [MassPeaks](#) objects.

## Usage

```
mergeMassPeaks(l, labels, method=c("mean", "median", "sum"), ignore.na=TRUE,
...)
```

## Arguments

|        |   |
|--------|---|
| l      | list, list of <a href="#">MassPeaks</a> objects.  |
| labels | list, list of <a href="#">factor</a> s (one for each <a href="#">MassPeaks</a> object) to do groupwise merging. |
| method | used merge method.  |

ignore.na      Should NA (positions where a peak is missing) ignored (ignore.na=TRUE) or treated as zero (ignore.na=FALSE)?

...            arguments to be passed to underlying functions (currently only mc.cores is supported).

### Value

Returns a single (no labels given) or a [list](#) (labels given) of merged [MassPeaks](#) objects.

### Author(s)

Sebastian Gibb <mail@sebastiangibb.de>

### See Also

[MassPeaks](#), [averageMassSpectra](#)

Website: <https://strimmerlab.github.io/software/maldiquant/>

### Examples

```
## load package
library("MALDIquant")

## create four MassPeaks objects and add them to the list
p <- list(createMassPeaks(mass=1:2, intensity=1:2),
          createMassPeaks(mass=1:3, intensity=1:3),
          createMassPeaks(mass=1:4, intensity=1:4),
          createMassPeaks(mass=1:5, intensity=1:5))

## merge all four MassPeaks objects into a single new one
## by sum their intensities
## (no labels, returns only one new MassPeaks object)
mergedPeaks <- mergeMassPeaks(p, method="sum")

## only merge MassPeaks objects in a group
## (two different labels, returns a list of two new MassPeaks objects)
groups <- factor(c("a", "a", "b", "b"), levels=c("a", "b"))
mergedPeaks <- mergeMassPeaks(p, labels=groups, method="mean")

## the same, but treat NA as zero
mergedPeaks <- mergeMassPeaks(p, labels=groups, method="mean", ignore.na=FALSE)
```

---

monoisotopicPeaks-methods

*Finds monoisotopic peaks in a MassPeaks object.*

---

### Description

This method looks for monoisotopic peaks in peak list data (represented by a [MassPeaks](#) objects). It is based on the poisson model for isotopic patterns described in Breen et al 2000.



**Usage**

```
## S4 method for signature 'MassPeaks'  
monoisotopicPeaks(object,  
  minCor=0.95, tolerance=1e-4, distance=1.00235, size=3L:10L)  
## S4 method for signature 'list'  
monoisotopicPeaks(object, ...)
```

**Arguments**

|           |  |
|-----------|--|
| object    | <a href="#">MassPeaks</a> object or a list of <a href="#">MassPeaks</a> objects.   |
| minCor    | double, minimal correlation between the peak pattern generated by the model and the experimental peaks in the <a href="#">MassPeaks</a> object to be recognized as isotopic pattern.   |
| tolerance | double, maximal relative deviation of peaks position (mass) to be considered as isotopic distance ( $\text{abs}((\text{mass}[i]+\text{distance})-\text{mass}[i+1])/\text{mass}[i]) < \text{tolerance}$ ).  |
| distance  | double, distance between two consecutive peaks in an isotopic pattern (default value taken from Park et al 2008). Could contain more than one value, e.g. $\text{distance}=(1:3)^{-1}$ to find isotopic patterns for multiple charged patterns (e.g. 1+, 2+, and 3+). Please note that the order matters here if there is a monoisotopic peak for charge state 1 and 3 it would be reported as charge 1 for $\text{distance}=(1:3)^{-1}$ and as 3 for $\text{distance}=(3:1)^{-1}$ respectively. |
| size      | double, size (length) of isotopic pattern, longer patterns are preferred over shorter ones.  |
| ...       | arguments to be passed to <a href="#">monoisotopicPeaks, MassPeaks-method</a> . If object is a list <code>mc.cores</code> is also supported.   |

**Value**

Returns a [MassPeaks](#) object with monoisotopic peaks only.

**Author(s)**

Sebastian Gibb <[mail@sebastiangibb.de](mailto:mail@sebastiangibb.de)>

**References**

K. Park, J.Y. Yoon, S. Lee, E. Paek, H. Park, H.J. Jung, and S.W. Lee. 2008. Isotopic peak intensity ratio based algorithm for determination of isotopic clusters and monoisotopic masses of polypeptides from high-resolution mass spectrometric data. *Analytical Chemistry*, 80: 7294-7303.

E.J. Breen, F.G. Hopwood, K.L. Williams, and M.R. Wilkins. 2000. Automatic poisson peak harvesting for high throughput protein identification. *Electrophoresis* 21: 2243-2251.

**See Also**

[MassPeaks](#), [detectPeaks](#), [MassSpectrum-method](#)

Website: <https://strimmerlab.github.io/software/malDIquant/>

**Examples**

```
## load package
library("MALDIquant")

## create example peaks
p <- createMassPeaks(mass=995:1005,
                    intensity=c(100, 10, 30, 10, 40, # noise
                               550, 330, 110, 10, # isotopic pattern
                               5, 15)) # more noise

m <- monoisotopicPeaks(p)
as.matrix(m)

## plot the peaks and mark the monoisotopic one
plot(p)
points(m, col=2, pch=4)
```

---

|           |  |
|-----------|--|
| msiSlices | <i>Turn a list of <code>AbstractMassObjects</code> into a mass spectrometry imaging slice.</i> |
|-----------|--|

---

**Description**

This function turns a mass spectrometry imaging dataset represented by a list of `AbstractMassObject` objects into an `intensityMatrix` for each slice (stored in an array).

**Usage**

```
msiSlices(x, center, tolerance, method=c("sum", "mean", "median"), adjust=TRUE)
```

**Arguments**

|                        |   |
|------------------------|---|
| <code>x</code>         | a list of <code>MassSpectrum/ MassPeaks</code> objects.   |
| <code>center</code>    | double, the center mas value of each slice.   |
| <code>tolerance</code> | double, specifies the thickness of the slices ( <code>center + c(-tolerance, tolerance)</code> ).   |
| <code>method</code>    | used aggregation function.  |
| <code>adjust</code>    | logical, if TRUE the lowest coordinates of the mass spectrometry imaging dataset are set to <code>c(x=1, y=1)</code> to avoid NA values at the borders. |

**Details**

Each `MassSpectrum/MassPeaks` object in `x` must contain a list named `imaging` with an element `pos` that stores the `x` and `y` value of the spectrum, e.g.:

```
> metaData(spectra[[1]])$imaging$pos
x y
1 5
```

**Value**

Returns an [array](#) of three dimensions. The first and second dimensions contains the x and y coordinates of the image. The third dimension represents the index of the center of each slice. There are two additional attributes, namely "center" and "tolerance" which store the original center and tolerance information.

**Author(s)**

Sebastian Gibb <mail@sebastiangibb.de>

**See Also**

[AbstractMassObject](#), [MassSpectrum](#), [MassPeaks](#), [coordinates](#), [AbstractMassObject-method](#), [plotMsiSlice](#), [list-method](#)

Please find real examples on:

Website: <https://strimmerlab.github.io/software/maldiquant/>

Vignette: <https://github.com/sgibb/MALDIquantExamples/raw/master/inst/doc/nyakas2013.pdf>

Shiny: <https://github.com/sgibb/ims-shiny/>

**Examples**

```
## load package
library("MALDIquant")

## load example data
data("fiedler2009subset", package="MALDIquant")

## please note: this is NOT a MSI data set
## we just add some coordinates for demonstration
coordinates(fiedler2009subset) <- cbind(x=rep(1:4, 2), y=rep(1:2, each=4))

slices <- msiSlices(fiedler2009subset, center=c(5864.49, 8936.97),
                   tolerance=0.25)

slices
```

---

plot-methods

*Plots an AbstractMassObject object.*

---

**Description**

This is an overloaded method to allow plotting of an [AbstractMassObject](#) object.

## Usage

```
## S4 method for signature 'AbstractMassObject,missing'
plot(x, col="black",
     xlab=expression(italic(m/z)), ylab="intensity",
     type=ifelse(isMassPeaks(x), "h", "l"),
     xlim=c(ifelse(length(x@mass), min(x@mass, na.rm=TRUE), 0),
            ifelse(length(x@mass), max(x@mass, na.rm=TRUE), 1)),
     ylim=c(0, ifelse(length(x@intensity), max(x@intensity, na.rm=TRUE), 1)),
     main=x@metaData$name, sub=x@metaData$file,
     cex.sub=0.75, col.sub="#808080", ...)
```

## Arguments

|         |   |
|---------|---|
| x       | MassSpectrum object.  |
| col     | line colour, see <a href="#">par</a> .                                |
| xlab    | title for the x-axis, see <a href="#">title</a> .                     |
| ylab    | title for the y-axis, see <a href="#">title</a> .                     |
| type    | type of plot: see <a href="#">plot</a> .                              |
| xlim    | the x limits (x1, x2) of the plot, see <a href="#">plot.default</a> . |
| ylim    | the y limits (y1, y2) of the plot, see <a href="#">plot.default</a> . |
| main    | title for the plot, see <a href="#">title</a> .                       |
| sub     | sub title for the plot, see <a href="#">title</a> .                   |
| cex.sub | sub title font size, see <a href="#">par</a> .                        |
| col.sub | sub title color, see <a href="#">par</a> .                            |
| ...     | arguments to be passed to <a href="#">plot</a> .                      |

## Author(s)

Sebastian Gibb <[mail@sebastiangibb.de](mailto:mail@sebastiangibb.de)>

## See Also

[AbstractMassObject](#)

Website: <https://strimmerlab.github.io/software/maldiquant/>

## Examples

```
## load package
library("MALDIquant")

## create a MassSpectrum object by default constructor
s <- createMassSpectrum(mass=1:100, intensity=rnorm(100)^2,
                       metaData=list(name="example"))

## show some details
s
```

```
## plot spectrum
plot(s)
```

---

plotMsiSlice-methods *Plots a Mass Spectrometry Imaging dataset.*

---

## Description

This function allows to plot a slice of a mass spectrometry imaging dataset represented by a list of [AbstractMassObject](#) objects or an array or a matrix.

## Usage

```
## S4 method for signature 'list'
plotMsiSlice(x, center, tolerance,
  colRamp=colorRamp(c("black", "blue", "green", "yellow", "red")),
  interpolate=FALSE, legend=TRUE, alignLabels=FALSE, combine=FALSE,
  ...)
## S4 method for signature 'array'
plotMsiSlice(x,
  colRamp=colorRamp(c("black", "blue", "green", "yellow", "red")),
  interpolate=FALSE, legend=TRUE, alignLabels=FALSE, combine=FALSE,
  plotInteractive=FALSE, ...)
## S4 method for signature 'matrix'
plotMsiSlice(x,
  colRamp=colorRamp(c("black", "blue", "green", "yellow", "red")),
  interpolate=FALSE, scale=TRUE, legend=scale, ...)
```

## Arguments

|             |   |
|-------------|---|
| x           | The mass spectrometry imaging dataset. It could be a list of <a href="#">MassSpectrum/MassPeaks</a> objects or an array (e.g. generated by <a href="#">msiSlices</a> ) or a matrix.                 |
| center      | double, if x is a list of <a href="#">MassSpectrum/MassPeaks</a> objects this argument represent the <i>center</i> mass value of the slices, see <a href="#">msiSlices</a> for details.             |
| tolerance   | double, if center is given tolerance specifies the thickness of the slices (center + c(-tolerance, tolerance)), see <a href="#">msiSlices</a> for details.  |
| colRamp     | colours as colorRamp function, see <a href="#">colorRamp</a> for details. If combine=TRUE multiple colour functions must be applied as list with an length that equals the number of given centers. |
| interpolate | logical, use linear interpolation when drawing the image, see <a href="#">rasterImage</a> for details.  |
| scale       | logical, if TRUE all values are divided by the maximal value of the slice to get values between 0 and 1.  |

|                 |   |
|-----------------|---|
| legend          | logical, if TRUE a reference color gradient is plotted on the right hand side of the plot. The upper color represents the highest value in the slice and the lower color the lowest value respectively. The legend is disabled if scale=FALSE.  |
| alignLabels     | logical, if combine=TRUE and alignLabels=TRUE the center positions below the legend are aligned on the right margin otherwise the aligned to their corresponding gradient.  |
| combine         | logical, if TRUE multiple centers are plotted in one image. Therefore it would be necessary to apply a list of colRamp functions (one function for each center). The intensity values for each center of each pixel are compared against each other and the highest scaled intensity determines the center (and the corresponding colRamp). |
| plotInteractive | logical, if the slice array contains multiple centers, combine=FALSE and an interactive plotting device is used a warning is thrown and only the first center would be plotted. Use plotInteractive=TRUE to overwrite this behaviour and to plot multiple centers on an interactive device.   |
| ...             | arguments to be passed to <code>plot</code> , e.g. <code>main</code> .  |

### Details

Each `MassSpectrum/MassPeaks` object in `x` must contain a list named `imaging` with an element `pos` that stores the `x` and `y` value of the spectrum, e.g.:

```
> metaData(spectra[[1]])$imaging$pos
x y
1 5
```

### Author(s)

Sebastian Gibb <[mail@sebastiangibb.de](mailto:mail@sebastiangibb.de)>

### See Also

[AbstractMassObject](#), [MassSpectrum](#), [MassPeaks](#), [coordinates](#), [AbstractMassObject-method](#), [msiSlices](#), [plot](#), [MassSpectrum](#), [missing-method](#)

Please find real examples on:

Website: <https://strimmerlab.github.io/software/maldiquant/>

Vignette: <https://github.com/sgibb/MALDIquantExamples/raw/master/inst/doc/nyakas2013.pdf>

Shiny: <https://github.com/sgibb/ims-shiny/>

### Examples

```
## load package
library("MALDIquant")

## load example data
```

```
data("fiedler2009subset", package="MALDIquant")

## please note: this is NOT a MSI data set
## we just add some coordinates for demonstration
coordinates(fiedler2009subset) <- cbind(x=rep(1:4, 2), y=rep(1:2, each=4))

plotMsiSlice(fiedler2009subset, center=8936.97, tolerance=0.25)

plotMsiSlice(fiedler2009subset, center=c(5864.49, 8936.97), tolerance=0.25,
             combine=TRUE,
             colRamp=list(colorRamp(c("#000000", "#FF00FF")),
                          colorRamp(c("#000000", "#00FF00"))))
```

---

referencePeaks                      *Creates a reference [MassPeaks](#) object.*

---

## Description

This function creates a reference [MassPeaks](#) object (also called *Anchor Peaks*) from a list of [MassPeaks](#) objects.

Generally it is a combination of [binPeaks](#) and [filterPeaks](#)

## Usage

```
referencePeaks(l, method=c("strict", "relaxed"), minFrequency=0.9,
              tolerance=0.002)
```

## Arguments

|              |  |
|--------------|--|
| l            | list, list of <a href="#">MassPeaks</a> objects.   |
| method       | bin creation rule (see <a href="#">binPeaks</a> ).   |
| minFrequency | double, remove all peaks which occur in less than $\text{minFrequency} \times \text{length}(l)$ <a href="#">MassPeaks</a> objects.   |
| tolerance    | double, maximal relative deviation of a peak position (mass) to be considered as identical. Must be multiplied by $10^{-6}$ for ppm, e.g. use $\text{tolerance}=5e-6$ for 5 ppm. |

## Value

Returns a new [MassPeaks](#) objects.

The *intensity* slot of the returned [MassPeaks](#) represents the frequency of this mass position in all samples.

## Author(s)

Sebastian Gibb <mail@sebastiangibb.de>

## See Also

[binPeaks](#), [filterPeaks](#), [MassPeaks](#)

Website: <https://strimmerlab.github.io/software/maldiquant/>

## Examples

```
## load package
library("MALDIquant")

## create four MassPeaks objects and add them to the list
p<- list(createMassPeaks(mass=1:2, intensity=1:2),
         createMassPeaks(mass=1:3, intensity=1:3),
         createMassPeaks(mass=1:4, intensity=1:4),
         createMassPeaks(mass=1:5, intensity=1:5))

## only use peaks which occur in all MassPeaks objects as reference peaks
refPeaks <- referencePeaks(p, minFrequency=1)

mass(refPeaks)      # 1:2
intensity(refPeaks) # c(1, 1)
```

---

removeBaseline-methods

*Removes the baseline of a MassSpectrum object.*

---

## Description

This method removes the baseline of mass spectrometry data (represented by a [MassSpectrum](#) object).

The intensity of the mass spectrometry data would be reduced by baseline.

## Usage

```
## S4 method for signature 'MassSpectrum'
removeBaseline(object,
  method=c("SNIP", "TopHat", "ConvexHull", "median"),
  ...)
## S4 method for signature 'list'
removeBaseline(object, ...)
```

## Arguments

|        |  |
|--------|--|
| object | <a href="#">MassSpectrum</a> object or a list of <a href="#">MassSpectrum</a> objects.   |
| method | used baseline estimation method, one of "SNIP", "TopHat", "ConvexHull" or "median". See <a href="#">estimateBaseline, MassSpectrum-method</a> for details. |
| ...    | arguments to be passed to <a href="#">estimateBaseline, MassSpectrum-method</a> . If object is a list <code>mc.cores</code> is also supported.             |



**Value**

Returns a modified [MassSpectrum](#) object with reduced intensities.

**Author(s)**

Sebastian Gibb <mail@sebastiangibb.de>

**See Also**

[MassSpectrum](#), [estimateBaseline](#), [MassSpectrum-method](#)

`demo("baseline")`

Website: <https://strimmerlab.github.io/software/maldiquant/>

**Examples**

```
## load package
library("MALDIquant")

## load example data
data("fiedler2009subset", package="MALDIquant")

## choose only the first mass spectrum
s <- fiedler2009subset[[1]]

## plot spectrum
plot(s)

## subtract baseline
b <- removeBaseline(s, method="SNIP")

## draw modified spectrum on the plot
lines(b, col="blue")
```

---

smoothIntensity-methods

*Smooths intensities of a MassSpectrum object.*

---

**Description**

This method smoothes the intensity values of a [MassSpectrum](#) object.

**Usage**

```
## S4 method for signature 'MassSpectrum'
smoothIntensity(object,
  method=c("SavitzkyGolay", "MovingAverage"), halfWindowSize,
  ...)
```

## Arguments

|                |   |
|----------------|---|
| object         | <a href="#">AbstractMassObject</a> object or a list of <a href="#">AbstractMassObject</a> objects.  |
| method         | used smoothing method, one of "SavitzkyGolay" or "MovingAverage".   |
| halfWindowSize | half window size. The resulting window reaches from <code>mass[currentIndex-halfWindowSize]</code> to <code>mass[currentIndex+halfWindowSize]</code> (window size is $2*halfWindowSize+1$ ). The best size differs depending on the selected method.  |
| ...            | arguments to be passed to method. SavitzkyGolay has an additional <code>polynomialOrder</code> argument (default: 3) to control the order of the filter. MovingAverage has an additional <code>weighted</code> argument (default: FALSE) to indicate if the average should be equal weight (default) or if it should have weights depending on the distance from the center as calculated as $1/2^{abs(-halfWindowSize:halfWindowSize)}$ with the sum of all weights normalized to 1. |

## Details

`halfWindowSize`: Depends on the selected method. For the SavitzkyGolay the `halfWindowSize` should be smaller than *FWHM* of the peaks (full width at half maximum; please find details in Bromba and Ziegler 1981). In general the `halfWindowSize` for the MovingAverage has to be much smaller than for SavitzkyGolay to conserve the peak shape.

## Author(s)

Sebastian Gibb <[mail@sebastiangibb.de](mailto:mail@sebastiangibb.de)>

Weighted moving average: Sigurdur Smarason

## References

A. Savitzky and M. J. Golay. 1964. Smoothing and differentiation of data by simplified least squares procedures. *Analytical chemistry*, 36(8), 1627-1639.

M. U. Bromba and H. Ziegler. 1981. Application hints for Savitzky-Golay digital smoothing filters. *Analytical Chemistry*, 53(11), 1583-1586.

## See Also

[MassSpectrum](#)

Website: <https://strimmerlab.github.io/software/maldiquant/>

## Examples

```
## load package
library("MALDIquant")

## load example data
data("fiedler2009subset", package="MALDIquant")

## smooth spectra
s <- smoothIntensity(fiedler2009subset, method="MovingAverage",
                    halfWindowSize=2)
```

```
## or
s <- smoothIntensity(fiedler2009subset, method="MovingAverage",
                     halfWindowSize=2, weighted=TRUE)
## or
s <- smoothIntensity(fiedler2009subset, method="SavitzkyGolay",
                     halfWindowSize=10)
```

---

transformIntensity-methods

*Transforms intensities of an AbstractMassObject object.*

---

## Description

This method performs a transformation (e.g. sqrt-transformation) on the intensities of an [AbstractMassObject](#) object.

## Usage

```
## S4 method for signature 'AbstractMassObject'
transformIntensity(object,
                  method=c("sqrt", "log", "log2", "log10"))
## S4 method for signature 'list'
transformIntensity(object, ...)
```

## Arguments

|        |  |
|--------|--|
| object | <a href="#">AbstractMassObject</a> object or a list of <a href="#">AbstractMassObject</a> objects.           |
| method | used transformation method.  |
| ...    | arguments to be passed to underlying functions. If object is a list <code>mc.cores</code> is also supported. |

## Author(s)

Sebastian Gibb <[mail@sebastiangibb.de](mailto:mail@sebastiangibb.de)>

## See Also

[AbstractMassObject](#), [MassSpectrum](#)

Website: <https://strimmerlab.github.io/software/maldiquant/>

## Examples

```
## load package
library("MALDIquant")

## load example data
data("fiedler2009subset", package="MALDIquant")
```

```
## choose only the first mass spectrum
s <- fiedler2009subset[[1]]

## transform spectrum
t <- transformIntensity(s, method="sqrt")

## plot spectrum
par(mfrow=c(2, 1))
plot(s, main="raw spectrum")
plot(t, main="transformed spectrum")
par(mfrow=c(1, 1))
```

---

trim-methods

*Trim an AbstractMassObject object.*

---

### Description

This method trims an [AbstractMassObject](#) object. That is useful if some mass ranges should be excluded from further analysis.

### Usage

```
## S4 method for signature 'AbstractMassObject,numeric'
trim(object, range)
## S4 method for signature 'list,numeric'
trim(object, range, ...)
## S4 method for signature 'list,missing'
trim(object, range, ...)
```

### Arguments

|        |  |
|--------|--|
| object | <a href="#">AbstractMassObject</a> object or a list of <a href="#">AbstractMassObject</a> objects.   |
| range  | numeric, limits of trimming (left/minimal mass, right/maximal mass). If missing it is automatically determined (largest overlapping mass range) for a list of <a href="#">AbstractMassObject</a> . |
| ...    | arguments to be passed to underlying functions (currently only <code>mc.cores</code> is supported).  |

### Author(s)

Sebastian Gibb <[mail@sebastiangibb.de](mailto:mail@sebastiangibb.de)>

### See Also

[AbstractMassObject](#), [MassPeaks](#), [MassSpectrum](#)

Website: <https://strimmerlab.github.io/software/maldiquant/>

## Examples

```
## load package
library("MALDIquant")

## load example data
data("fiedler2009subset", package="MALDIquant")

## select only one spectrum
s <- fiedler2009subset[[1]]

## remove all mass lower 3000
trim(s, range=c(3000, Inf))

## remove all mass higher 8000
trim(s, range=c(0, 8000))

## remove all mass lower 3000 and higher 8000
trim(s, range=c(3000, 8000))

## choose largest overlapping mass range for all spectra
trim(fiedler2009subset)
```

---

warpMassSpectra

*Run warping functions on AbstractMassObject objects.*

---

## Description

These functions run warping functions on [AbstractMassObject](#) objects (warping is also known as *phase correction*).

## Usage

```
warpMassPeaks(l, w, emptyNoMatches=FALSE)
```

```
warpMassSpectra(l, w, emptyNoMatches=FALSE)
```

## Arguments

**l** list, list of [MassPeaks](#) or [MassSpectrum](#) objects.

**w** a list of warping functions determined by [determineWarpingFunctions](#). Has to be of the same length as **l**.

**emptyNoMatches** logical, if TRUE (default: FALSE) the intensity values of [MassSpectrum](#) or [MassPeaks](#) objects with missing (NA) warping functions are set to zero.

## Details

The warping function *w* is called in the following way:

$$newMass = oldMass + w(oldMass)$$

**Value**

Returns a list of warped `MassPeaks` or `MassSpectrum` objects.

**Author(s)**

Sebastian Gibb <mail@sebastiangibb.de>

**See Also**

[determineWarpingFunctions](#), [MassPeaks](#), [MassSpectrum](#)

Website: <https://strimmerlab.github.io/software/maldiquant/>

**Examples**

```
## load package
library("MALDIquant")

## create a MassPeaks object
p <- createMassPeaks(mass=1:5, intensity=1:5)

## stupid warping function for demonstration
## (please use determineWarpingFunctions in real life applications)
simpleWarp <- function(x) { return(1) }

## run warping function
w <- warpMassPeaks(list(p), list(simpleWarp))[[1]]

## compare results
all(mass(w) == mass(p)+1) # TRUE

## no warping (MassPeaks object is not changed)
warpMassPeaks(list(p), list(NA))

## no warping (intensity values of MassPeaks object are set to zero)
warpMassPeaks(list(p), list(NA), emptyNoMatches=TRUE)
```

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