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`cwt`

*Continuous Wavelet Transform (CWT)*

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## Description

CWT(Continuous Wavelet Transform) with Mexican Hat wavelet (by default) to match the peaks in Mass Spectrometry spectrum

## Usage

```
cwt(ms, scales = 1, wavelet = "mexh")
```

## Arguments

`ms` Mass Spectrometry spectrum (a vector of MS intensities)  
`scales` a vector represents the scales at which to perform CWT.  
`wavelet` The wavelet base, Mexican Hat by default. User can provide wavelet  $\Psi(x)$  as a form of two row matrix. The first row is the  $x$  value, and the second row is  $\Psi(x)$  corresponding to  $x$ .

## Value

The return is the 2-D CWT coefficient matrix, with column names as the scale. Each column is the CWT coefficients at that scale.

## Author(s)

Pan Du, Simon Lin

## Examples

```
data(exampleMS)
scales <- seq(1, 64, 3)
wCoefs <- cwt(exampleMS[5000:11000], scales=scales, wavelet='mexh')

## Plot the 2-D CWT coefficients as image (It may take a while!)
xTickInterval <- 1000
image(5000:11000, scales, wCoefs, col=terrain.colors(256), axes=FALSE, xlab='m/z index', ylab='CWT')
axis(1, at=seq(5000, 11000, by=xTickInterval))
axis(2, at=c(1, seq(10, 64, by=10)))
box()
```

---

`exampleMS`

*An example mass spectrum*

---

### Description

An example mass spectrum from CAMDA 2006. All-in-1 Protein Standard II (CIPHERGEN Cat. C100-0007) were measured on CIPHERGEN NP20 chips. There are 7 polypeptides in the sample with  $m/z$  values of 7034, 12230, 16951, 29023, 46671, 66433, 147300.

### Usage

```
data(exampleMS)
```

### Format

A numeric vector represents the mass spectrum with equal sample intervals.

### Source

CAMDA, CAMDA 2006 Competition Data Set. 2006, <http://camda.duke.edu>.

---

`extend.nBase`

*Extend the row number of a matrix as the exponential of base N*

---

### Description

Extend the data as the exponential of base N by increasing row number.

### Usage

```
extend.nBase(x, nLevel=1, base=2, ...)
```

### Arguments

<code>x</code>	data matrix
<code>nLevel</code>	the level of DWT decomposition. Basically, it is equivalent to changing the 'base' as $\text{base}^{\text{nLevel}}$
<code>base</code>	the base, 2 by default
<code>...</code>	other parameters of used by <code>extendLength</code>

### Details

The method 'open' is padding the the matrix with the last row.

**Value**

Return a extended matrix

**Author(s)**

Pan Du

**See Also**

[extendLength](#)

**Examples**

```
a = matrix(rnorm(9), 3)
extend.nBase(a)
```

---

<code>extendLength</code>	<i>Extend the length of a signal or matrix</i>
---------------------------	--

---

**Description**

Extend the length of a signal or matrix by row

**Usage**

```
extendLength(x, addLength = NULL, method = c("reflection", "open", "circular"), direction = c('
```

**Arguments**

<code>x</code>	a vector or matrix with column with each column as a signal
<code>addLength</code>	the length to be extended
<code>method</code>	three methods available, <code>c("reflection", "open", "circular")</code> . By default, it is "reflection".
<code>direction</code>	three options available: <code>c("right", "left", "both")</code>

**Value**

return the extended vector or matrix.

**Author(s)**

Pan Du

**See Also**

[extend.nBase](#)

## Examples

```
a = matrix(rnorm(9), 3)
extendLength(a, 3, direction='right')
```

---

```
get.localMaximum.cwt
```

*Identify the local maximum of each column in 2-D CWT coefficients matrix*

---

## Description

Identify the local maximum of each column in 2-D CWT coefficients matrix by using a slide window. The size of slide window linearly changes from the coarse scale (bigger window size) to detail scale. The scale of CWT increases with the column index.

## Usage

```
get.localMaximum.cwt(wCoefs, winSize0 = 5, amp.Th = 0)
```

## Arguments

<code>wCoefs</code>	2-D CWT coefficients, each column corresponding to CWT coefficient at one scale. The column name is the scale.
<code>winSize0</code>	The slide window size at the 0 scale, i.e., original MS spectrum.
<code>amp.Th</code>	The minimum peak amplitude.

## Value

return a matrix with same dimension as CWT coefficient matrix, `wCoefs`. The local maxima are marked as 1, others are 0.

## Author(s)

Pan Du

## See Also

[localMaximum](#)

## Examples

```
data(exampleMS)
scales <- seq(1, 64, 3)
wCoefs <- cwt(exampleMS[5000:11000], scales=scales, wavelet='mexh')

localMax <- get.localMaximum.cwt(wCoefs)
plot.localMax(localMax)
```

---

`get.ridge`

*Identify ridges based on the local maximum matrix*

---

## Description

Identify ridges by connecting the local maximum of 2-D CWT coefficients from the coarse scale to detail scale. The local maximum matrix is returned from [get.localMaximum](#)

## Usage

```
get.ridge(localMax, iInit = ncol(localMax), step = -1, iFinal = 1, endWinSize = 5, gapTh = 3, skip = 0)
```

## Arguments

<code>localMax</code>	The local maximum matrix is returned from <a href="#">get.localMaximum</a> with 1 represents maximum, others are 0.
<code>iInit</code>	The start column to search ridge. By default, it starts from the coarsest scale level.
<code>step</code>	Search step. -1 by default, which means searching from coarse scale to detail scale column by column.
<code>iFinal</code>	The final column index of search ridge.
<code>endWinSize</code>	The smallest slide window size allowed.
<code>gapTh</code>	The gap allowed during searching for ridge. 3 by default.
<code>skip</code>	The column to be skipped during search.

## Value

Return a list of ridge, with m/z index of peak at the raw spectrum as the names of the ridges.

## Author(s)

Pan Du, Simon Lin

## References

Pan Du, Warren A. Kibbe, Simon M. Lin, Improved peak detection in mass spectrum by incorporating continuous wavelet transform-based pattern matching, *Bioinformatics Advance* Access published online on July 4, 2006 (<http://bioinformatics.oxfordjournals.org/cgi/content/abstract/bt1355?ijke>)

## See Also

[get.localMaximum](#), [identify.majorPeaks](#)

## Examples

```
data(exampleMS)
scales <- seq(1, 64, 3)
wCoefs <- cwt(exampleMS[5000:11000], scales=scales, wavelet='mexh')

localMax <- get.localMaximum.cwt(wCoefs)
ridgeList <- get.ridge(localMax)
plot(ridgeList)
```

---

`get.ridgeLength`      *Estimate the length of the ridge*

---

## Description

Estimate the length of the ridge line, which is composed of local maxima at adjacent CWT scales. The ridge line is cut off at the end point, whose amplitude divided by the maximum ridge amplitude is larger than the cutoff amplitude ratio threshold (0.5 by default).

## Usage

```
get.ridgeLength(ridgeList, Th = 0.5)
```

## Arguments

`ridgeList`      a list of identified ridges  
`Th`              the cutoff amplitude ratio (the amplitude divided by the maximum amplitude of the ridge) threshold of the ridge line end.

## Value

a vector of estimated ridge length

## Author(s)

Pan Du

---

`get.ridgeValue`            *Get the CWT coefficient values corresponding to the peak ridge*

---

### Usage

```
get.ridgeValue(ridgeList, wCoefs, skip = 0)
```

### Arguments

<code>ridgeList</code>	a list of ridge lines
<code>wCoefs</code>	2-D CWT coefficients
<code>skip</code>	the CWT scale level to be skipped, by default the 0 scale level (raw spectrum) is skipped.

### Value

A list of ridge values corresponding to the input `ridgeList`.

### Author(s)

Pan Du

### Examples

---

`i2u`                            *Transfer m/z index to m/z value*

---

### Description

Transfer m/z index to m/z value

### Usage

```
i2u(i, t0 = 9e-08, a = 336301655.051424, b = 0, U = 20000, delta = 4e-09)
```

### Arguments

<code>i</code>	a vector of m/z index
<code>t0</code>	initial time
<code>a</code>	accelerate rate
<code>b</code>	constant shift
<code>U</code>	Unit
<code>delta</code>	sample interval in second

## Details

Here is the conversion equation:  $t = i * \text{delta}$ , (unit in seconds),  $u = (a*(t-t_0)^2 + b) * U$ , (here is m/z at 150KDa)

## Value

a vector of m/z values

## Author(s)

Simon Lin

## See Also

[u2i](#)

---

`identify.majorPeaks` *Identify peaks based on the ridges in 2-D CWT coefficient matrix*

---

## Description

Identify the peaks based on the ridge list (returned by [get.ridge](#)) in 2-D CWT coefficient matrix and estimated Signal to Noise Ratio (SNR)

## Usage

```
identify.majorPeaks(ridgeList, wCoefs, scales = as.numeric(colnames(wCoefs)), SNR.Th = 3, peakScaleRange)
```

## Arguments

<code>ridgeList</code>	returned by <a href="#">get.ridge</a>
<code>wCoefs</code>	2-D CWT coefficients
<code>scales</code>	scales of CWT, by default it is the colnames of wCoefs
<code>SNR.Th</code>	threshold of SNR
<code>peakScaleRange</code>	the CWT scale range of the peak.
<code>ridgeLength</code>	the maximum ridge scale of the major peaks.
<code>nearbyPeak</code>	determine whether to include the small peaks close to large major peaks
<code>nearbyWinSize</code>	the window size to determine the nearby peaks. Only effective when nearbyPeak is true.
<code>winSize.noise</code>	the local window size to estimate the noise level.
<code>SNR.method</code>	method to estimate noise level. Currently, only 95 percentage quantile is supported.
<code>minNoiseLevel</code>	the minimum noise level used in calculating SNR, i.e., if the estimated noise level is less than "minNoiseLevel", it will use "minNoiseLevel" instead. If the noise level is less than 0.5, it will be treated as the ratio to the maximum amplitude of the spectrum.

## Details

The determination of the peaks is based on three rules: Rule 1: The maximum ridge scale of the peak should larger than a certain threshold Rule 2: Based on the scale of the peak (corresponding to the maximum value of the peak ridge) should be within certain range Rule 3: Based on the peak SNR

## Value

Return a list with following elements:

**peakIndex** the m/z indexes of the identified peaks  
**peakCenterIndex** the m/z indexes of peak centers, which correspond to the maximum on the ridge. **peakCenterIndex** includes all the peaks, not just the identified major peaks.  
**peakCenterValue** the CWT coefficients (the maximum on the ridge) corresponding to **peakCenterIndex**  
**peakSNR** the SNR of the peak, which is the ratio of **peakCenterValue** and noise level  
**peakScale** the estimated scale of the peak, which corresponds to the **peakCenterIndex**  
**potentialPeakIndex** the m/z indexes of all potential peaks, which satisfy all requirements of a peak without considering its SNR. Useful, if you want to change to a lower SNR threshold later.

## Author(s)

Pan Du, Simon Lin

## References

Pan Du, Warren A. Kibbe, Simon M. Lin, Improved peak detection in mass spectrum by incorporating continuous wavelet transform-based pattern matching, *Bioinformatics Advance* Access published online on July 4, 2006 (<http://bioinformatics.oxfordjournals.org/cgi/content/abstract/bt1355?ijke>)

## See Also

[main.peakDetection](#), [tuneIn.peakInfo](#)

## Examples

```
data(exampleMS)
scales <- seq(1, 64, 3)
wCoefs <- cwt(exampleMS, scales=scales, wavelet='mexh')

localMax <- get.localMaximum.cwt(wCoefs)
ridgeList <- get.ridge(localMax)
```

```
SNR.Th <- 3
majorPeakInfo <- identify.majorPeaks(ridgeList, wCoefs, SNR.Th=SNR.Th)
## Plot the identified peaks
peakIndex <- majorPeakInfo$peakIndex
plot.peak(exampleMS, peakIndex, main=paste('Identified peaks with SNR >', SNR.Th))
```

---

localMaximum	<i>Identify local maximum within a slide window.</i>
--------------	--

---

## Description

Find local maximum by transform the vector as matrix, then get the the maximum of each column. This operation is performed twice with vector shifted half of the winSize.

## Usage

```
localMaximum(x, winSize = 5)
```

## Arguments

x	a vector represents a signal profile
winSize	the slide window size, 5 by default.

## Details

Instead of find the local maximum by a slide window, which slide all possible positions, we find local maximum by transform the vector as matrix, then get the the maximum of each column. This operation is performed twice with vector shifted half of the winSize. The main purpose of this is to increase the efficiency of the algorithm.

## Value

Return a vector with the same length of the input x. The position of local maximum is set as 1, 0 else where.

## Author(s)

Pan Du

## See Also

[get.localMaximum](#)

## Examples

```
x <- rnorm(200)
lmax <- localMaximum(x, 5)
maxInd <- which(lmax > 0)
plot(x, type='l')
points(maxInd, x[maxInd], col='red')
```

---

MassSpecWavelet-package

*Peak detection of mass spectrum by Wavelet transform based methods*

---

## Description

Process Mass Spectrum (MS) by Wavelet Transforms-based algorithms

## Details

Package: MassSpecWavelet  
Type: Package  
Version: 1.1  
Date: 2006-07-05  
License: GPL 2 or newer

MassSpecWavelet R package is aimed to process Mass Spectrometry (MS) data mainly based on Wavelet Transforms. The current version only supports the peak detection based on Continuous Wavelet Transform (CWT). Future versions will include more functions covering entire MS data processes.

## Author(s)

Pan Du, Simon Lin

Maintainer: Pan Du <dupan@northwestern.edu>

## References

Pan Du, Warren A. Kibbe, Simon M. Lin, Improved peak detection in mass spectrum by incorporating continuous wavelet transform-based pattern matching, *Bioinformatics Advance* Access published online on July 4, 2006 (<http://bioinformatics.oxfordjournals.org/cgi/content/abstract/bt1355?ijke>)

## Examples

```
data(exampleMS)
SNR.Th <- 3
peakInfo <- peakDetection.cwt(exampleMS, SNR.Th=SNR.Th)
majorPeakInfo = peakInfo$majorPeakInfo
peakIndex <- majorPeakInfo$peakIndex
plot.peak(exampleMS, peakIndex, main=paste('Identified peaks with SNR >', SNR.Th))
```

---

mzInd2vRange

*Match m/z index to m/z value with a certain error range*

---

### Description

Match m/z index to m/z value with a certain error range

### Usage

```
mzInd2vRange(mzInd, error = 0.003)
```

### Arguments

mzInd	a vector of m/z index
error	error range

### Value

return a vector of sorted m/z values

### Author(s)

Pan Du

### See Also

[mzV2indRange](#)

---

mzV2indRange

*Match m/z value to m/z index with a certain error range*

---

### Description

Match m/z value to m/z index with a certain error range

### Usage

```
mzV2indRange(mzV, error = 0.003)
```

### Arguments

mzV	a vector of m/z value
error	error range

### Value

return a vector of sorted m/z indexes

## Author(s)

Pan Du

## See Also

[mzInd2vRange](#)

---

<code>peakDetection.cwt</code>	<i>The main function of peak detection by CWT based pattern matching</i>
--------------------------------	--

---

## Description

This function is a wrapper of [cwt](#), [get.localMaximum](#), [get.ridge](#), [identify.majorPeaks](#)

## Usage

```
peakDetection.cwt(ms, scales = c(1, seq(2, 30, 2), seq(32, 64, 4)), SNR.Th = 3, nearbyPeak = TRUE)
```

## Arguments

<code>ms</code>	the mass spectrometry spectrum
<code>scales</code>	scales of CWT
<code>SNR.Th</code>	SNR (Signal to Noise Ratio) threshold
<code>nearbyPeak</code>	Determine whether to include the nearby small peaks of major peaks. TRUE by default
<code>peakScaleRange</code>	the scale range of the peak. larger than 5 by default.
<code>amp.Th</code>	the minimum required amplitude of the peak
<code>minNoiseLevel</code>	the minimum noise level used in computing the SNR
<code>ridgeLength</code>	the minimum highest scale of the peak in 2-D CWT coefficient matrix
<code>tuneIn</code>	determine whether to tune in the parameter estimation of the detected peaks
<code>...</code>	other parameters of used by <a href="#">get.localMaximum</a> , <a href="#">get.ridge</a> , <a href="#">identify.majorPeaks</a>

## Value

<code>majorPeakInfo</code>	return of <a href="#">identify.majorPeaks</a>
<code>ridgeList</code>	return of <a href="#">get.ridge</a>
<code>localMax</code>	return of <a href="#">get.localMaximum</a>
<code>wCoefs</code>	2-D CWT coefficient matrix, see <a href="#">cwt</a> for details.

## Author(s)

Pan Du, Simon Lin

## References

Pan Du, Warren A. Kibbe, Simon M. Lin, Improved peak detection in mass spectrum by incorporating continuous wavelet transform-based pattern matching, *Bioinformatics Advance*  
Access published online on July 4, 2006 (<http://bioinformatics.oxfordjournals.org/cgi/content/abstract/bt1355?ijke>)

## See Also

[cwt](#), [get.localMaximum](#), [get.ridge](#), [identify.majorPeaks](#)

## Examples

```
data(exampleMS)
SNR.Th <- 3
peakInfo <- peakDetection.cwt(exampleMS, SNR.Th=SNR.Th)
majorPeakInfo = peakInfo$majorPeakInfo
peakIndex <- majorPeakInfo$peakIndex
plot.peak(exampleMS, peakIndex, main=paste('Identified peaks with SNR >', SNR.Th))
```

---

`peakDetection.dwt`      *Peak detection based on DWT denoising and baseline removal*

---

## Description

### Usage

```
peakDetection.dwt(ms, nLevel = 6, SNR.Th = 3, amp.Th = 0.001, minNoiseLevel = amp.Th/SNR.Th, winSize.localMax)
```

### Arguments

<code>ms</code>	a vector representing the mass spectrum
<code>nLevel</code>	the level of DWT decomposition
<code>SNR.Th</code>	SNR (Signal to Noise Ratio) threshold
<code>amp.Th</code>	the minimum required amplitude of the peak
<code>minNoiseLevel</code>	the minimum noise level used in computing the SNR
<code>winSize.localMax</code>	the slide window size for detecting local maximum
<code>localWinSize</code>	local window size for estimate local noise threshold
<code>smLocalNoiseTh</code>	the local noise threshold for smoothing (denoising)
<code>twoSectionMode</code>	determine whether treating the spectrum as two sections and adopting different parameters for them.
<code>...</code>	parameters used by <a href="#">smooth.dwt</a>

## Details

### Value

<code>peakInd</code>	the index of the peak position
<code>smoothMS</code>	the smoothed mass spectrum
<code>localMaxInd</code>	all local maximum indexes, including SNR less than the SNR threshold
<code>SNR</code>	the SNR corresponding to <code>localMaxInd</code>
<code>globalSNR</code>	the global SNR

### Author(s)

Pan Du

### References

### See Also

[smooth.dwt](#), [peakDetection.cwt](#)

### Examples

```
library(waveslim)
data(exampleMS)
SNR.Th <- 3
peakInfo <- peakDetection.dwt(exampleMS, SNR.Th=SNR.Th)
peakIndex <- peakInfo$peakIndex
plot.peak(exampleMS, peakIndex, main=paste('Identified peaks with SNR >', SNR.Th))
```

---

`plot.localMax` *Plot the local maximum matrix*

---

### Description

Plot the local maximum matrix of 2-D CWT coefficients returned by [get.localMaximum.cwt](#)

### Usage

```
plot.localMax(localMax, wCoefs = NULL, range = c(1, nrow(localMax)), colorMap = "RYB", main = M
```

## Arguments

<code>localMax</code>	local maximum matrix of 2-D CWT coefficients returned by <code>get.localMaximum.cwt</code>
<code>wCoefs</code>	2-D CWT coefficients
<code>range</code>	plot range of m/z index
<code>colorMap</code>	the colormap used in plotting tht points
<code>main</code>	parameter of <code>plot</code>
<code>cex</code>	parameter of <code>plot</code>
<code>pch</code>	parameter of <code>plot</code>
<code>...</code>	other parameters of <code>points</code>

## Author(s)

Pan Du

## See Also

[get.localMaximum.cwt](#)

## Examples

```
data(exampleMS)
scales <- seq(1, 64, 3)
wCoefs <- cwt(exampleMS[5000:11000], scales=scales, wavelet='mexh')

localMax <- get.localMaximum.cwt(wCoefs)
plot.localMax(localMax)
```

---

`plot.peak`

*Plot the identified peaks over the spectrum*

---

## Description

Plot the identified peaks over the spectrum. The identified peaks are returned by `main.peakDetection` or `identify.majorPeaks`

## Usage

```
plot.peak(ms, peakIndex = NULL, mz = 1:length(ms), range = c(min(mz), max(mz)), method = c("p",
```

## Arguments

<code>ms</code>	the MS spectrum
<code>peakIndex</code>	m/z indexes of the identified peaks
<code>mz</code>	m/z value correspond to m/z index
<code>range</code>	the plot range of m/z value
<code>method</code>	plot method of the identified peaks. method 'p' plot circles on the peaks; method 'l' add vertical lines over the peaks.
<code>main</code>	parameter of <code>plot</code>
<code>log</code>	parameter of <code>plot</code>
<code>...</code>	other parameters of <code>points</code>

## Author(s)

Pan Du

## See Also

`main.peakDetection`, `identify.majorPeaks`

## Examples

```
data(exampleMS)
SNR.Th <- 3
peakInfo <- peakDetection.cwt(exampleMS, SNR.Th=SNR.Th)
majorPeakInfo = peakInfo$majorPeakInfo
peakIndex <- majorPeakInfo$peakIndex
plot.peak(exampleMS, peakIndex, main=paste('Identified peaks with SNR >', SNR.Th))
```

---

`plot.ridgeList`      *Plot the ridge list*

---

## Description

Plot the ridge list returned by `get.ridge`

## Usage

```
plot.ridgeList(ridgeList, wCoefs = NULL, range = NULL, colorMap = "RYB", main = NULL, pch = ".')
```

## Arguments

<code>ridgeList</code>	returned by <code>get.ridge</code>
<code>wCoefs</code>	2-D CWT coefficients
<code>range</code>	plot range of m/z index
<code>colorMap</code>	colorMap to plot the points of local maximum
<code>main</code>	parameter of <code>plot</code>
<code>pch</code>	parameter of <code>plot</code>
<code>cex</code>	parameter of <code>plot</code>
<code>...</code>	other parameters of <code>points</code>

## Author(s)

Pan Du

## See Also

[get.ridge](#)

## Examples

```
data(exampleMS)
scales <- seq(1, 64, 3)
wCoefs <- cwt(exampleMS[5000:11000], scales=scales, wavelet='mexh')

localMax <- get.localMaximum.cwt(wCoefs)
ridgeList <- get.ridge(localMax)
plot(ridgeList)
```

---

<code>smooth.dwt</code>	<i>smooth (denoise) the spectrum by DWT (Discrete Wavelet Transform)</i>
-------------------------	--

---

## Description

## Usage

```
smooth.dwt(ms, nLevel = 6, wf = "la8", localNoiseTh = seq(1, 0, by = -0.2), localWinSize = 500,
```

## Arguments

<code>ms</code>	a vector representing the mass spectrum
<code>nLevel</code>	the level of DWT decomposition
<code>wf</code>	the name of wavelet for DWT
<code>localNoiseTh</code>	local noise level threshold
<code>localWinSize</code>	local window size for estimate local noise threshold
<code>globalNoiseTh</code>	global noise level threshold
<code>smoothMethod</code>	the method used for denoising. 'hard' means keeping the dwt coefficients higher than the threshold unchanged; "soft" means the dwt coefficients higher than the threshold were subtracted by the threshold.
<code>method</code>	'dwt' or 'modwt' used for decomposition

## Details

## Value

return the smoothed mass spectrum with the 'detail' component of DWT as an attribute 'detail'.

## Author(s)

Pan Du

## References

## See Also

[peakDetect.dwt](#)

## Examples

---

<code>tuneIn.peakInfo</code>	<i>Tune in the peak information: peak position and peak scale</i>
------------------------------	---

---

## Description

Based on the identified peak position, more precise estimation of the peak information, i.e., peak position and peak scale, can be got by this function. The basic idea is to cut the segment of spectrum near the identified peaks, and then do similar procedures as [main.peakDetection](#), but with more detailed scales around the estimated peak scale.

## Usage

```
tuneIn.peakInfo(ms, majorPeakInfo = NULL, peakIndex = NULL, peakScale = NULL, maxScale = 128, .
```

## Arguments

<code>ms</code>	the mass spectrometry spectrum
<code>majorPeakInfo</code>	return of <a href="#">identify.majorPeaks</a>
<code>peakIndex</code>	the m/z index of the identified peaks
<code>peakScale</code>	the scales of the identified peaks
<code>maxScale</code>	the maximum scale allowed for the peak
<code>...</code>	other parameters of used by <a href="#">get.localMaximum</a> , <a href="#">get.ridge</a> , <a href="#">identify.majorPeaks</a>

## Details

The `majorPeakInfo` or `peakIndex` and `peakScale` must be provided.

## Value

<code>peakCenterIndex</code>	the updated peak center m/z index
<code>peakScale</code>	the updated peak scale
<code>peakValue</code>	the corresponding peak value

## Author(s)

Pan Du

## References

Pan Du, Warren A. Kibbe, Simon M. Lin, Improved peak detection in mass spectrum by incorporating continuous wavelet transform-based pattern matching, *Bioinformatics Advance* Access published online on July 4, 2006 (<http://bioinformatics.oxfordjournals.org/cgi/content/abstract/bt1355?ijke>)

## See Also

[main.peakDetection](#)

## Examples

```
data(exampleMS)
SNR.Th <- 3
peakInfo <- peakDetection.cwt(exampleMS, SNR.Th=SNR.Th)
majorPeakInfo <- peakInfo$majorPeakInfo
betterPeakInfo <- tuneIn.peakInfo(exampleMS, majorPeakInfo)
plot(500:length(exampleMS), exampleMS[500:length(exampleMS)], type='l', log='x')
abline(v=betterPeakInfo$peakCenterIndex, col='red')
```

---

u2i	<i>Transfer m/z value to m/z index</i>
-----	--

---

## Description

Transfer m/z value to m/z index

## Usage

```
u2i(u, t0 = 9e-08, a = 336301655.051424, b = 0, U = 20000, delta = 4e-09)
```

## Arguments

u	a vector of m/z values
t0	initial time
a	accelerate rate
b	constant shift
U	Unit
delta	sample interval in second

## Details

Here is the conversion equation:  $i = (((u/U-b)/a)^{1/2} + t0)/\text{delta}$

## Value

a vector of m/z indexes

## Author(s)

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## See Also

[i2u](#)