# Package 'sRm'

January 16, 2017

Junuary 10, 2017
Type Package
<b>Title</b> Minimal parser for Selective Reaction Monitoring (SRM) mass spectrometry (MS) data
Version 0.1.2
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<b>Description</b> Functions for parsing SRM-MS data from .mzML files.
License GPL (>= 3)
<b>Depends</b> methods, xml2 (>= 1.0), ggplot2, grid, gridExtra
Imports ggrepel, base64enc, Rcpp, gtable, scales, munsell, colorspace, plyr
<pre>URL https://github.com/wilsontom/sRm</pre>
BugReports https://github.com/wilsontom/sRm/issues
LazyData TRUE
RoxygenNote 5.0.0
Suggests testthat
NeedsCompilation no
R topics documented:
combineTransitions meta openSRMfile plotAll plotMulti plotSRM show,SRM-method SRM-class transitions
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combineTransitions

Combine Transitions

#### **Description**

Combine Transitions

Each parent m/z will usually have multiple separate Q3 product ions. SRM transitions are stored separately in the mzML file. Each index is a transition from a single Q1 precursor ion to one of the specified Q3 ions. This method combines all Q3 product ion chromatograms into one, to produce a total ion chromatogram based on the parent m/z and all Q3 product ions assigned to that parent m/z

# Usage

```
combineTransitions(object)
## S4 method for signature 'SRM'
combineTransitions(object)
```

#### **Arguments**

object

a SRM object

#### Value

a transition class of SRM total ion chromatograms

#### Author(s)

meta

Show meta data

#### **Description**

Show meta data

#### Usage

```
meta(object)
## S4 method for signature 'SRM'
meta(object)
```

#### Arguments

object

a SRM object

openSRMfile 3

#### Author(s)

openSRMfile

Open a SRM-MS file

# Description

Open and parse mzML file into SRM object

# Usage

```
openSRMfile(filename)
```

#### **Arguments**

filename

a mzML file

#### Value

```
a SRM object (see SRM-class)
```

#### Author(s)

plotAll

Plot all

# Description

Plot the SRM-MS TIC and all individual Q3 traces for a selected parent scans'

# Usage

```
plotAll(x, n)
## S3 method for class 'transition'
plotAll(x, n)
```

# Arguments

x a transition class

n the number of the transition (index) to plot

#### Author(s)

4 plotSRM

plotMulti

plotMulti

#### **Description**

plotMulti

#### Usage

```
plotMulti(object, idn, labels = NULL)
## S4 method for signature 'SRM'
plotMulti(object, idn, labels = NULL)
```

#### **Arguments**

object a SRM object

idn the index number of a transition to plot

labels an optional character vector of label names. Default is NULL

#### Author(s)

plotSRM

Plot SRM

#### **Description**

Plot SRM

#### Usage

```
plotSRM(object, idn)
## S4 method for signature 'SRM'
plotSRM(object, idn)
```

# Arguments

object a SRM object

idn the index number of a transition to plot

#### Author(s)

show,SRM-method 5

show, SRM-method

show-SRM

#### **Description**

show-SRM

# Usage

```
## S4 method for signature 'SRM'
show(object)
```

# Arguments

object

a SRM object

#### Author(s)

SRM-class

**SRM** 

#### **Description**

peaks list of peak data

A S4 class to store single reaction monitoring (SRM) mass spectrometry (MS) data

# Slots

```
SHA1 character vector of the original SHA-1 checksum value
meta list containing file meta data
totalIonCount data.frame of sample total ion count (TIC)
index character vector of scan index's; ie SRM SIC 341.014,114.098
```

header data.frame of summary information for each scan index; parentMz, Q3mz, totalIonCount, basePeakInt

6 transitions

 ${\it transitions}$ 

Show transitions

# Description

Show transitions

# Usage

```
transitions(object)
## S4 method for signature 'SRM'
transitions(object)
```

# Arguments

object

a SRM object

# Author(s)

# **Index**