

# Package ‘q3ML’

September 16, 2021

**Type** Package

**Title** Open and Parse SRM mzML files from Thermo Triple Quadrupole Mass Spectrometers

**Version** 0.1.0

**Description** More about what it does (maybe more than one line)

Use four spaces when indenting paragraphs within the Description.

**License** GPL (>= 3)

**Encoding** UTF-8

**LazyData** true

**Depends** magrittr

**Imports** xml2,

dplyr,

tidyr,

purrr,

tibble,

stringr,

base64enc,

crayon,

cli

**RoxygenNote** 7.1.1

**Suggests** covr,

testthat (>= 3.0.0),

mzR

**Config/testthat/edition** 3

**Roxygen** list(markdown = TRUE)

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`decodePeaks`*Decode peaks*

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**Description**

Decode a `binaryDataArray` into a vector of either time or intensity

**Usage**

```
decodePeaks(x, compression = "none", size)
```

**Arguments**

<code>x</code>	a base64 encoded vector
<code>compression</code>	compression type (default = "none")
<code>size</code>	a numeric value for the number of bytes per element in the byte stream

**Value**

a numeric vector

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`idRefs`*Extract mzML idRefs*

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**Description**

Extract all the `idRefs` in the specified `mzML` file. The `idRefs` will correspond to the individual scan events.

**Usage**

```
idRefs(xmlDoc)
```

**Arguments**

<code>xmlDoc</code>	a xml document
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**Value**

a character vector of all available `idRefs`

**Examples**

```
## Not run:
library(xml2)
xmlDoc <- read_xml("example_file.mzML")
xmlRefs <- idRefs(xmlDoc)
head(xmlRefs)
[1] "TIC" "SRM SIC 153.01,65.271"
[3] "SRM SIC 153.01,67.232" "SRM SIC 153.01,109.094"
[5] "SRM SIC 179.022,107.007" "SRM SIC 179.022,134.006"

## End(Not run)
```

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openFile*Open .mzML File*

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**Description**

Open and parse a .mzML file containing Selective Reaction Monitoring (SRM) Mass Spectrometry (MS) data using xml2

**Usage**

```
openFile(mzml_file)
```

**Arguments**

mzml\_file      the absolute file path of a valid .mzML file

**Value**

a list of two elements containing peaks and header

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parseChromNode*Parse the XML nodeset of a SRM chromatogram*

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**Description**

Parse the XML nodeset of a SRM chromatogram

**Usage**

```
parseChromNode(x, mode)
```

**Arguments**

x                      a xml2 nodeset for a Chromatogram block  
mode                    a character of either SRM or TIC

**Value**

a list of two elements

- **time**
  - attributes
  - raw
- **intensity**
  - attributes
  - raw

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