

MSnbase development

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Abstract

This vignette describes the classes implemented in *MSnbase* package. It is intended as a starting point for developers or users who would like to learn more or further develop/extend *pSet*.

Keywords: Mass Spectrometry (MS), proteomics, infrastructure.

Contents

1	Introduction	2
2	MSnbase classes	2
2.1	<i>pSet</i> : a virtual class for raw mass spectrometry data and meta data	3
2.2	<i>MSnExp</i> : a class for MS experiments	3
2.3	<i>MSnSet</i> : a class for quantitative proteomics data	4
2.4	<i>MSnProcess</i> : a class for logging processing meta data	5
2.5	<i>MIAPE</i> : Minimum Information About a Proteomics Experiment	6
2.6	<i>Spectrum et al.</i> : classes for MS spectra	8
2.7	<i>ReporterIons</i> : a class for isobaric tags	9
2.8	<i>NAnnotatedDataFrame</i> : multiplexed <i>AnnotatedDataFrames</i>	10
2.9	Other classes	10
3	Miscellaneous	10
4	Session information	11

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Foreword

MSnbase is under active development; current functionality is evolving and new features will be added. This software is free and open-source software. If you use it, please support the project by citing it in publications:

Laurent Gatto and Kathryn S. Lilley. *MSnbase - an R/Bioconductor package for isobaric tagged mass spectrometry data visualization, processing and quantitation*. Bioinformatics 28, 288-289 (2011).

Questions and bugs

You are welcome to contact me directly about *MSnbase*. For bugs, typos, suggestions or other questions, please file an issue in our tracking system¹ providing as much information as possible, a reproducible example and the output of `sessionInfo()`.

If you wish to reach a broader audience for general questions about proteomics analysis using R, you may want to use the Bioconductor support site: <https://support.bioconductor.org/>.

1 Introduction

This document is not a replacement for the individual manual pages, that document the slots of the *MSnbase* classes. It is a centralised high-level description of the package design.

MSnbase aims at being compatible with the *Biobase* infrastructure [1]. Many meta data structures that are used in *eSet* and associated classes are also used here. As such, knowledge of the *Biobase development and the new eSet vignette*² would be beneficial.

The initial goal is to use the *MSnbase* infrastructure for labelled quantitation using reporter ions (iTRAQ [2] and TMT [3]). Spectral counting should be trivial to apply with current features, as long as identification data is at hand. Currently, no effort is invested to streamline label-free quantitative proteomics, although some effort has been done to keep the infrastructure flexible enough to accommodate more designs.

2 MSnbase classes

All classes have a `...classVersion...` slot, of class *Versioned* from the *Biobase* package. This slot documents the class version for any instance to be used for debugging and object update purposes. Any

¹<https://github.com/lgatto/MSnbase/issues>

²The vignette can directly be accessed with `vignette("BiobaseDevelopment", package="Biobase")` once *Biobase* is loaded.

change in a class implementation should trigger a version change.

2.1 pSet: a virtual class for raw mass spectrometry data and meta data

This virtual class is the main container for mass spectrometry data, i.e spectra, and meta data. It is based on the *eSet* implementation for genomic data. The main difference with *eSet* is that the *assayData* slot is an environment containing any number of *Spectrum* instances (see section 2.6).

One new slot is introduced, namely *processingData*, that contains one *MSnProcess* instance (see section 2.4). and the *experimentData* slot is now expected to contain *MIAPE* data (see section 2.5). The annotation slot has not been implemented, as no prior feature annotation is known in shotgun proteomics.

```
getClass("pSet")
Virtual Class "pSet" [package "MSnbase"]

Slots:

Name:          assayData          phenoData
Class:          environment NAnnotatedDataFrame

Name:          featureData         experimentData
Class: AnnotatedDataFrame          MIAxE

Name:          protocolData        processingData
Class: AnnotatedDataFrame          MSnProcess

Name:          .cache      .__classVersion__
Class:          environment          Versions

Extends: "Versioned"

Known Subclasses: "MSnExp"
```

Future work Currently, few setters have been implemented.

2.2 MSnExp: a class for MS experiments

MSnExp extends *pSet* to store MS experiments. It does not add any new slots to *pSet*. Accessors and setters are all inherited from *pSet* and new ones should be implemented for *pSet*. Methods that manipulate actual data in experiments are implemented for *MSnExp* objects.

```
getClass("MSnExp")
Class "MSnExp" [package "MSnbase"]

Slots:

Name:          assayData          phenoData
Class:          environment NAnnotatedDataFrame

Name:          featureData          experimentData
Class:  AnnotatedDataFrame          MIAxE

Name:          protocolData          processingData
Class:  AnnotatedDataFrame          MSnProcess

Name:          .cache          .__classVersion__
Class:          environment          Versions

Extends:
Class "pSet", directly
Class "Versioned", by class "pSet", distance 2
```

2.3 MSnSet: a class for quantitative proteomics data

This class stores quantitation data and meta data after running `quantify` on an *MSnExp* object or by creating an *MSnSet* instance from an external file, as described in the *MSnbase-io* vignette and in `?readMSnSet`, `readMzTabData`, etc. The quantitative data is in form of a $m \times n$ matrix, where m is the number of features/spectra originally in the *MSnExp* used as parameter in `quantify` and n is the number of reporter ions (see section 2.7). If read from an external file, n corresponds to the number of features (protein groups, proteins, peptides, spectra) in the file and m is the number of columns with quantitative data (samples) in the file.

This prompted to keep a similar implementation as the *ExpressionSet* class, while adding the proteomics-specific annotation slot introduced in the *pSet* class, namely `processingData` for objects of class *MSnProcess* (see section 2.4).

```
getClass("MSnSet")
Class "MSnSet" [package "MSnbase"]

Slots:

Name:          experimentData          processingData          qual
Class:          MIAPE          MSnProcess          data.frame
```

```

Name:      assayData      phenoData      featureData
Class:      AssayData AnnotatedDataFrame AnnotatedDataFrame

Name:      annotation      protocolData  .__classVersion__
Class:      character AnnotatedDataFrame      Versions

Extends:
Class "eSet", directly
Class "VersionedBiobase", by class "eSet", distance 2
Class "Versioned", by class "eSet", distance 3

```

The *MSnSet* class extends the virtual *eSet* class to provide compatibility for *ExpressionSet*-like behaviour. The experiment meta-data in *experimentData* is also of class *MIAPE* (see section 2.5). The annotation slot, inherited from *eSet* is not used. As a result, it is easy to convert *ExpressionSet* data from/to *MSnSet* objects with the coercion method `as`.

```

data(itraqdata)
class(msnset)

[1] "MSnSet"
attr("package")
[1] "MSnbase"

class(as(msnset, "ExpressionSet"))

[1] "ExpressionSet"
attr("package")
[1] "Biobase"

data(sample.ExpressionSet)
class(sample.ExpressionSet)

[1] "ExpressionSet"
attr("package")
[1] "Biobase"

class(as(sample.ExpressionSet, "MSnSet"))

[1] "MSnSet"
attr("package")
[1] "MSnbase"

```

2.4 MSnProcess: a class for logging processing meta data

This class aims at recording specific manipulations applied to *MSnExp* or *MSnSet* instances. The processing slot is a character vector that describes major processing. Most other slots are of class `logical` that indicate whether the data has been centroided, smoothed, ...although many of the

functionality is not implemented yet. Any new processing that is implemented should be documented and logged here.

It also documents the raw data file from which the data originates (files slot) and the *MSnbase* version that was in use when the *MSnProcess* instance, and hence the *MSnExp*/*MSnSet* objects, were originally created.

```
getClass("MSnProcess")
```

```
Class "MSnProcess" [package "MSnbase"]
```

```
Slots:
```

Name:	files	processing	merged
Class:	character	character	logical

Name:	cleaned	removedPeaks	smoothed
Class:	logical	character	logical

Name:	trimmed	normalised	MSnbaseVersion
Class:	numeric	logical	character

Name:	.__classVersion__
Class:	Versions

```
Extends: "Versioned"
```

2.5 MIAPE: Minimum Information About a Proteomics Experiment

The Minimum Information About a Proteomics Experiment [4, 5] *MIAPE* class describes the experiment, including contact details, information about the mass spectrometer and control and analysis software.

```
getClass("MIAPE")
```

```
Class "MIAPE" [package "MSnbase"]
```

```
Slots:
```

Name:	title	url
Class:	character	character

Name:	abstract	pubMedIds
Class:	character	character

Name:	samples	preprocessing
Class:	list	list

```

Name:          other          dateStamp
Class:         list          character

Name:          name          lab
Class:         character     character

Name:          contact       email
Class:         character     character

Name:          instrumentModel instrumentManufacturer
Class:         character     character

Name: instrumentCustomisations softwareName
Class:         character     character

Name:          softwareVersion switchingCriteria
Class:         character     character

Name:          isolationWidth parameterFile
Class:         numeric      character

Name:          ionSource      ionSourceDetails
Class:         character     character

Name:          analyser       analyserDetails
Class:         character     character

Name:          collisionGas    collisionPressure
Class:         character     numeric

Name:          collisionEnergy detectorType
Class:         character     character

Name:          detectorSensitivity .__classVersion__
Class:         character     Versions

Extends:
Class "MIAxE", directly
Class "Versioned", by class "MIAxE", distance 2

```

2.6 Spectrum et al.: classes for MS spectra

Spectrum is a virtual class that defines common attributes to all types of spectra. MS1 and MS2 specific attributes are defined in the *Spectrum1* and *Spectrum2* classes, that directly extend *Spectrum*.

```
getClass("Spectrum")
```

```
Virtual Class "Spectrum" [package "MSnbase"]
```

```
Slots:
```

Name:	msLevel	peaksCount	rt
Class:	integer	integer	numeric

Name:	acquisitionNum	scanIndex	tic
Class:	integer	integer	numeric

Name:	mz	intensity	fromFile
Class:	numeric	numeric	integer

Name:	centroided	__classVersion__
Class:	logical	Versions

```
Extends: "Versioned"
```

```
Known Subclasses: "Spectrum2", "Spectrum1"
```

```
getClass("Spectrum1")
```

```
Class "Spectrum1" [package "MSnbase"]
```

```
Slots:
```

Name:	polarity	msLevel	peaksCount
Class:	integer	integer	integer

Name:	rt	acquisitionNum	scanIndex
Class:	numeric	integer	integer

Name:	tic	mz	intensity
Class:	numeric	numeric	numeric

Name:	fromFile	centroided	__classVersion__
Class:	integer	logical	Versions

```
Extends:
```



```
Class "Spectrum", directly
Class "Versioned", by class "Spectrum", distance 2
```

```
getClass("Spectrum2")
```

```
Class "Spectrum2" [package "MSnbase"]
```

```
Slots:
```

Name:	merged	precScanNum	precursorMz
Class:	numeric	integer	numeric

Name:	precursorIntensity	precursorCharge	collisionEnergy
Class:	numeric	integer	numeric

Name:	msLevel	peaksCount	rt
Class:	integer	integer	numeric

Name:	acquisitionNum	scanIndex	tic
Class:	integer	integer	numeric

Name:	mz	intensity	fromFile
Class:	numeric	numeric	integer

Name:	centroided	.__classVersion__
Class:	logical	Versions

```
Extends:
```

```
Class "Spectrum", directly
```

```
Class "Versioned", by class "Spectrum", distance 2
```

2.7 ReporterIons: a class for isobaric tags

The iTRAQ and TMT (or any other peak of interest) are implemented *ReporterIons* instances, that essentially defines an expected MZ position for the peak and a width around this value as well a names for the reporters.

```
getClass("ReporterIons")
```

```
Class "ReporterIons" [package "MSnbase"]
```

```
Slots:
```

Name:	name	reporterNames	description
Class:	character	character	character

```

Name:          mz          col          width
Class:         numeric     character    numeric

Name:  __classVersion__
Class:  Versions

Extends: "Versioned"

```

2.8 NAnnotatedDataFrame: multiplexed AnnotatedDataFrames

The simple expansion of the *AnnotatedDataFrame* classes adds the `multiplex` and `multiLabel` slots to document the number and names of multiplexed samples.

```

getClass("NAnnotatedDataFrame")

Class "NAnnotatedDataFrame" [package "MSnbase"]

Slots:

Name:          multiplex      multiLabels      varMetadata
Class:         numeric       character        data.frame

Name:          data          dimLabels  __classVersion__
Class:         data.frame    character    Versions

Extends:
Class "AnnotatedDataFrame", directly
Class "Versioned", by class "AnnotatedDataFrame", distance 2

```

2.9 Other classes

Lists of MSnSet instances

3 Miscellaneous

Unit tests *MSnbase* implements unit tests with the *testthat* package.

Processing methods Methods that process raw data, i.e. spectra should be implemented for *Spectrum* objects first and then eapply'ed (or similar) to the `assayData` slot of an *MSnExp* instance in the specific method.

4 Session information

- R version 3.2.2 Patched (2015-10-08 r69496), x86_64-apple-darwin10.8.0
- Locale: en_US.UTF-8/en_US.UTF-8/en_US.UTF-8/C/en_US.UTF-8/en_US.UTF-8
- Base packages: base, datasets, graphics, grDevices, grid, methods, parallel, stats, stats4, utils
- Other packages: annotate 1.48.0, AnnotationDbi 1.32.0, Biobase 2.30.0, BiocGenerics 0.16.0, BiocParallel 1.4.0, cluster 2.0.3, ggplot2 1.0.1, gplots 2.17.0, IRanges 2.4.0, knitr 1.11, MLInterfaces 1.50.0, MSnbase 1.18.0, mzR 2.4.0, pRoloc 1.10.0, pRolocdata 1.7.4, ProtGenerics 1.2.0, Rcpp 0.12.1, RcppClassic 0.9.6, Rdisop 1.30.0, reshape2 1.4.1, S4Vectors 0.8.0, XML 3.98-1.3, zoo 1.7-12
- Loaded via a namespace (and not attached): affy 1.48.0, affyio 1.40.0, assertthat 0.1, base64enc 0.1-3, BiocInstaller 1.20.0, BiocStyle 1.8.0, biomaRt 2.26.0, bitops 1.0-6, car 2.1-0, caret 6.0-57, caTools 1.17.1, class 7.3-14, codetools 0.2-14, colorspace 1.2-6, DBI 0.3.1, DEoptimR 1.0-3, digest 0.6.8, diptest 0.75-7, doParallel 1.0.10, dplyr 0.4.3, e1071 1.6-7, evaluate 0.8, flexmix 2.3-13, FNN 1.1, foreach 1.4.3, formatR 1.2.1, fpc 2.1-10, futile.logger 1.4.1, futile.options 1.0.0, gbm 2.1.1, gdata 2.17.0, genefilter 1.52.0, ggvis 0.4.2, gtable 0.1.2, gtools 3.5.0, highr 0.5.1, htmltools 0.2.6, htmlwidgets 0.5, httpuv 1.3.3, hwriter 1.3.2, impute 1.44.0, iterators 1.0.8, kernlab 0.9-22, KernSmooth 2.23-15, labeling 0.3, lambda.r 1.1.7, lattice 0.20-33, limma 3.26.0, lme4 1.1-10, lpSolve 5.6.13, magrittr 1.5, MALDIquant 1.13, MASS 7.3-44, Matrix 1.2-2, MatrixModels 0.4-1, mclust 5.0.2, mgcv 1.8-7, mime 0.4, minqa 1.2.4, mlbench 2.1-1, modeltools 0.2-21, munsell 0.4.2, mvtnorm 1.0-3, mzID 1.8.0, nlme 3.1-122, nloptr 1.0.4, nnet 7.3-11, pbkrtest 0.4-2, pcaMethods 1.60.0, pls 2.5-0, plyr 1.8.3, prabclus 2.2-6, preprocessCore 1.32.0, proto 0.3-10, proxy 0.4-15, quantreg 5.19, R6 2.1.1, randomForest 4.6-12, RColorBrewer 1.1-2, RCurl 1.95-4.7, rda 1.0.2-2, rgl 0.95.1367, robustbase 0.92-5, rpart 4.1-10, RSQLite 1.0.0, sampling 2.7, scales 0.3.0, sfsmisc 1.0-28, shiny 0.12.2, SparseM 1.7, splines 3.2.2, stringi 0.5-5, stringr 1.0.0, survival 2.38-3, threejs 0.2.1, tools 3.2.2, trimcluster 0.1-2, vsn 3.38.0, xtable 1.7-4, zlibbioc 1.16.0

References

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- reagents. *Mol Cell Proteomics*, 3(12):1154–1169, Dec 2004. URL: <http://dx.doi.org/10.1074/mcp.M400129-MCP200>, doi:10.1074/mcp.M400129-MCP200.
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