Package ‘MultiAssayExperiment’

May 4, 2024

Title Software for the integration of multi-omics experiments in Bioconductor

Version 1.30.1

Description Harmonize data management of multiple experimental assays performed on an overlapping set of specimens. It provides a familiar Bioconductor user experience by extending concepts from SummarizedExperiment, supporting an open-ended mix of standard data classes for individual assays, and allowing subsetting by genomic ranges or rownames. Facilities are provided for reshaping data into wide and long formats for adaptability to graphing and downstream analysis.

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URL http://waldronlab.io/MultiAssayExperiment/

BugReports https://github.com/waldronlab/MultiAssayExperiment/issues

Depends SummarizedExperiment (>= 1.3.81), R (>= 3.5.0)

Imports Biobase, BiocBaseUtils, BiocGenerics, DelayedArray, GenomicRanges (>= 1.25.93), IRanges, methods, S4Vectors (>= 0.23.19), tidyr, utils

Suggests BiocStyle, HDF5Array (>= 1.19.17), knitr, maftools (>= 2.7.10), R.rsp, RaggedExperiment, reshape2, rmarkdown, survival, survminer, testthat, UpSetR

VignetteBuilder knitr, R.rsp

biocViews Infrastructure, DataRepresentation

Encoding UTF-8

RoxygenNote 7.3.1

Roxygen list(markdown = TRUE)


Collate 'ExperimentList-class.R' 'MultiAssayExperiment-class.R'
'MatchedAssayExperiment-class.R' 'subsetBy-methods.R'
'MultiAssayExperiment-subset.R'
'MultiAssayExperiment-methods.R'
MultiAssayExperiment-package

MultiAssayExperiment: Build an integrative multi-assay container

Description

MultiAssayExperiment allows the manipulation of related multiassay datasets with partially overlapping samples, associated metadata at the level of an entire study, and at the level of the "biological unit". The biological unit may be a patient, plant, yeast strain, etc.

Details

The package hierarchy of information:

- study
- experiments
- samples

Author(s)

Maintainer: Marcel Ramos <marcel.ramos@roswellpark.org> (ORCID)

Authors:

- Martin Morgan [contributor]
- Vincent J Carey [contributor]
- Levi Waldron <lwaldron.research@gmail.com>

Other contributors:

- Lori Shepherd [contributor]
- Hervé Pagès [contributor]
- MultiAssay SIG <biocmultiassay@googlegroups.com> [contributor]

See Also

Useful links:

- http://waldronlab.io/MultiAssayExperiment/
- Report bugs at https://github.com/waldronlab/MultiAssayExperiment/issues
ExperimentList

Represent multiple experiments as a List-derivative ExperimentList

Description

The ExperimentList class can contain several different types of data. The only requirements for an ExperimentList class are that the objects contained have the following set of methods: dim, [ , dimnames

Usage

ExperimentList(...)

Arguments

... A named list class object

Value

A ExperimentList class object of experiment data

Examples

## Create an empty ExperimentList instance
ExperimentList()

## Create array matrix and AnnotatedDataFrame to create an ExpressionSet class
arraydat <- matrix(data = seq(101, length.out = 20), ncol = 4, 
dimnames = list( 
  c("ENST00000294241", "ENST00000355076", 
  "ENST00000383706", "ENST000000234812", "ENST00000383323"), 
  c("array1", "array2", "array3", "array4") 
))

colDat <- data.frame(slope53 = rnorm(4), 
  row.names = c("array1", "array2", "array3", "array4"))

## SummarizedExperiment constructor
exprdat <- SummarizedExperiment::SummarizedExperiment(arraydat, 
colData = colDat)

## Create a sample methylation dataset
methylDat <- matrix(data = seq(1, length.out = 25), ncol = 5, 
dimnames = list( 
  c("ENST00000355076", "ENST00000383706", 
  "ENST00000383323", "ENST000000234812", "ENST000000294241"), 
  c("methyl1", "methyl2", "methyl3", 
  "methyl4", "methyl5") 
))
## Create a sample RNASeqGene dataset

```r
rnadat <- matrix(
  data = sample(c(46851, 5, 19, 13, 2197, 507, 84318, 126, 17, 21, 23979, 614), size = 20, replace = TRUE),
  ncol = 4,
  dimnames = list(
    c("XIST", "RPS4Y1", "KDM5D", "ENST00000383323", "ENST00000234812"),
    c("samparray1", "samparray2", "samparray3", "samparray4")
  )
)
```

## Create a mock RangedSummarizedExperiment from a data.frame

```r
rangedat <- data.frame(chr="chr2", start = 11:15, end = 12:16,
    strand = c("+", "-", "+", "*", "."),
    samp0 = c(0,0,1,1,1), samp1 = c(1,0,1,0,1), samp2 = c(0,1,0,1,0),
    row.names = c(paste0("ENST", "00000", 135411:135414), "ENST00000383323"))
rangeSE <- SummarizedExperiment::makeSummarizedExperimentFromDataFrame(rangedat)
```

## Combine to a named list and call the ExperimentList constructor function

```r
assayList <- list(Affy = exprdat, Methyl450k = methyldat, RNASeqGene = rnadat,
              GISTIC = rangeSE)
ExpList <- ExperimentList(assayList)
```

### ExperimentList-class

**ExperimentList - A container for multi-experiment data**

#### Description

The `ExperimentList` class is a container that builds on the `SimpleList` with additional checks for consistency in experiment names and length. It contains a `SimpleList` of experiments with sample identifiers. One element present per experiment performed.

Convert from `SimpleList` or `list` to the multi-experiment data container. When using the `mergeReplicates` method, additional arguments are passed to the given `simplify` function argument (e.g., `na.rm = TRUE`)

#### Usage

```r
# S4 method for signature 'ExperimentList'
show(object)
```

```r
# S4 method for signature 'ExperimentList'
isEmpty(x)
```

```r
# S4 method for signature 'ExperimentList'
dimnames(x)
```

```r
# S4 method for signature 'ExperimentList'
```
ExperimentList-class

colnames(x, do.NULL = TRUE, prefix = "col")

## S4 method for signature 'ExperimentList'
rownames(x, do.NULL = TRUE, prefix = "row")

## S4 method for signature 'ExperimentList'
mergeReplicates(x, replicates = list(), simplify = BiocGenerics::mean, ...)

## S4 method for signature 'ANY,missing'
assay(x, i, withDimnames = TRUE, ...)

## S4 method for signature 'ExperimentList'
assays(x, withDimnames = TRUE, ...)

## S4 method for signature 'ExperimentList,missing'
assay(x, i, withDimnames = TRUE, ...)

## S4 method for signature 'ExperimentList,numeric'
assay(x, i, withDimnames = TRUE, ...)

## S4 method for signature 'ExperimentList,character'
assay(x, i, withDimnames = TRUE, ...)

Arguments

object, x An ExperimentList object
do.NULL, prefix See ?base::rownames for a description of these arguments.
replicates mergeReplicates: A list or LogicalList where each element represents a
sample and a vector of repeated measurements for the sample
simplify A function for merging columns where duplicates are indicated by replicates
... Additional arguments. See details for more information.
i A scalar character or integer index
withDimnames logical (default TRUE) whether to return dimension names

Value

An ExperimentList class object

Methods (by generic)

- show(ExperimentList): Show method for ExperimentList class
- isEmpty(ExperimentList): check for zero length across all experiments
- dimnames(ExperimentList): Get the dimension names for an ExperimentList using CharacterList
- colnames(ExperimentList): Get the column names for an ExperimentList as a CharacterList slightly more efficiently
- rownames(ExperimentList): Get the row names for an ExperimentList as a CharacterList slightly more efficiently
hasAssay

- `mergeReplicates(ExperimentList)`: Apply the mergeReplicates method on the ExperimentList elements
- `assay(x = ANY, i = missing)`: Obtain the specified assay with a numeric or character reference
- `assays(ExperimentList)`: Get the assay data from each element in the ExperimentList

**coercion**

Convert a list or S4 List to an ExperimentList using the `as()` function.

In the following example, `x` is either a list or `List`:

```r
as(x, "ExperimentList")
```

**Examples**

```r
ExperimentList()
```

---

**Description**

The `hasAssay` function is intended for developers who would like to include new classes into a MultiAssayExperiment instance. It checks the methods tables of the assay function for the specified class of the argument.

**Usage**

```r
hasAssay(object)
```

**Arguments**

- `object` A MultiAssayExperiment or named list object instance

**Value**

A logical value indicating method availability

**Examples**

```r
lst <- structure(list(), .Names=character())
hasAssay(lst)
```
listToMap

Convert map from data.frame or DataFrame to list and vice versa

Description

The mapToList function provides a convenient way of reordering a data.frame to a list. The listToMap function does the opposite by taking a list and converting it to DataFrame.

Usage

listToMap(listmap, fill = TRUE)

mapToList(dfmap, assayCol = "assay")

Arguments

- **listmap**: A named list object containing DataFrames with "primary" and "colname" columns
- **fill**: logical(1) Whether to fill the map with an empty DataFrame when empty elements are present in the input list
- **dfmap**: A data.frame or DataFrame object with identifiers in the first column
- **assayCol**: A character vector of length one indicating the assay names column

Value

- A DataFrame class object of names
- A list object of DataFrames for each assay

Functions

- `listToMap()`: The inverse of the listToMap operation

Examples

```r
example("MultiAssayExperiment")

## Create a sampleMap from a list using the listToMap function
sampMap <- listToMap(maplist)

## The inverse operation is also available
maplist <- mapToList(sampMap)
```
MatchedAssayExperiment-class

**MatchedAssayExperiment** - A matched-samples MultiAssayExperiment class

---

**Description**

This class supports the use of matched samples where an equal number of observations per biological unit are present in all assays.

**Usage**

```r
MatchedAssayExperiment(...)```

**Arguments**

```r
... Either a single MultiAssayExperiment or the components to create a valid MultiAssayExperiment```

**Value**

A MatchedAssayExperiment object

**Functions**

- `MatchedAssayExperiment()`: Construct a MatchedAssayExperiment class from `MultiAssayExperiment`

**See Also**

`MultiAssayExperiment`

**Examples**

```r
data("miniACC")
acc <- as(miniACC, "MatchedAssayExperiment")
acc```
Description

A MultiAssayExperiment object providing a reduced version of the TCGA ACC dataset for all 92 patients. RNA-seq, copy number, and somatic mutations are included only for genes whose proteins are included in the reverse-phase protein array. The MicroRNA-seq dataset is also included, with infrequently expressed microRNA removed. Clinical, pathological, and subtype information are provided by colData(miniACC), and some additional details are provided by metadata(miniACC).

Usage

data("miniACC")

Format

A MultiAssayExperiment with 5 experiments, providing:

- **RNASeq2GeneNorm** RNA-seq count data: an ExpressionSet with 198 rows and 79 columns
- **gistict** Recurrent copy number lesions identified by GISTIC2: a SummarizedExperiment with 198 rows and 90 columns
- **RPPAArray** Reverse Phase Protein Array: an ExpressionSet with 33 rows and 46 columns. Rows are indexed by genes, but protein annotations are available from featureData(miniACC["RPPAArray"]). The source of these annotations is noted in abstract(miniACC["RPPAArray"]).
- **Mutations** Somatic mutations: a matrix with 223 rows and 90 columns. 1 for any kind of non-silent mutation, zero for silent (synonymous) or no mutation.
- **miRNASeqGene** microRNA sequencing: an ExpressionSet with 471 rows and 80 columns. Rows not having at least 5 counts in at least 5 samples were removed.

Author(s)

Levi Waldron <lwaldron.research@gmail.com>

Source

https://github.com/waldronlab/multiassayexperiment-tcga

References

Examples

```r
data("miniACC")
metadata(miniACC)
colnames(colData(miniACC))
table(miniACC$vital_status)
longFormat(
  miniACC["MAPK3", , ],
  colDataCols = c("vital_status", "days_to_death")
)
wideFormat(
  miniACC["MAPK3", , ],
  colDataCols = c("vital_status", "days_to_death")
)
```

MultiAssayExperiment  Construct an integrative representation of multi-omic data with MultiAssayExperiment

Description

The constructor function for the `MultiAssayExperiment` combines multiple data elements from the different hierarchies of data (study, experiments, and samples). It can create instances where neither a `sampleMap` or a `colData` set is provided. Please see the MultiAssayExperiment API documentation for more information.

Usage

```r
MultiAssayExperiment(
  experiments = ExperimentList(),
  colData = S4Vectors::DataFrame(),
  sampleMap = S4Vectors::DataFrame(assay = factor(), primary = character(), colname = character()),
  metadata = list(),
  drops = list()
)
```

Arguments

- `experiments` A list or `ExperimentList` of all combined experiments
- `colData` A `DataFrame` or `data.frame` of characteristics for all biological units
- `sampleMap` A `DataFrame` or `data.frame` of assay names, sample identifiers, and `colname` samples
- `metadata` An optional argument of "ANY" class (usually list) for content describing the experiments
- `drops` A list of unmatched information (included after subsetting)
MultiAssayExperiment

Value

A MultiAssayExperiment object that can store experiment and phenotype data

colData

The colData input can be either DataFrame or data.frame with subsequent coercion to DataFrame. The rownames in the colData must match the colnames in the experiments if no sampleMap is provided.

experiments

The experiments input can be of class SimpleList or list. This input becomes the ExperimentList. Each element of the input list or list must be named, rectangular with two dimensions, and have dimnames.

sampleMap

The sampleMap can either be input as DataFrame or data.frame with eventual coercion to DataFrame. The sampleMap relates biological units and biological measurements within each assay. Each row in the sampleMap is a single such link. The standard column names of the sampleMap are "assay", "primary", and "colname". Note that the "assay" column is a factor corresponding to the names of each experiment in the ExperimentList. In the case where these names do not match between the sampleMap and the experiments, the documented experiments in the sampleMap take precedence and experiments are dropped by the harmonization procedure. The constructor function will generate a sampleMap in the case where it is not provided and this method may be a 'safer' alternative for creating the MultiAssayExperiment (so long as the rownames are identical in the colData, if provided). An empty sampleMap may produce empty experiments if the levels of the "assay" factor in the sampleMap do not match the names in the ExperimentList.

See Also

MultiAssayExperiment

Examples

```r
## Run the example ExperimentList
example("ExperimentList")

## Create sample maps for each experiment
exprmap <- data.frame(
  primary = c("Jack", "Jill", "Barbara", "Bob"),
  colname = c("array1", "array2", "array3", "array4"),
  stringsAsFactors = FALSE)
methylmap <- data.frame(
  primary = c("Jack", "Jack", "Jill", "Barbara", "Bob"),
  colname = c("methyl1", "methyl2", "methyl3", "methyl4", "methyl5"),
  stringsAsFactors = FALSE)
rnamap <- data.frame(
)```
MultiAssayExperiment-class

primary = c("Jack", "Jill", "Bob", "Barbara"),
colname = c("samparray1", "samparray2", "samparray3", "samparray4"),
stringsAsFactors = FALSE)
gistmap <- data.frame(
  primary = c("Jack", "Bob", "Jill"),
colname = c("samp0", "samp1", "samp2"),
stringsAsFactors = FALSE)

## Combine as a named list and convert to a DataFrame
maplist <- list(Affy = exprmap, Methyl450k = methylmap,
                 RNASeqGene = rnamap, GISTIC = gistmap)

## Create a sampleMap
sampMap <- listToMap(maplist)

## Create an example phenotype data
colDat <- data.frame(sex = c("M", "F", "M", "F"), age = 38:41,
                      row.names = c("Jack", "Jill", "Bob", "Barbara"))

## Create a MultiAssayExperiment instance
mae <- MultiAssayExperiment(experiments = ExpList, colData = colDat,
                             sampleMap = sampMap)

MultiAssayExperiment-class

MultiAssayExperiment - An integrative multi-assay class for experiment data

Description

The MultiAssayExperiment class can be used to manage results of diverse assays on a collection of specimen. Currently, the class can handle assays that are organized instances of SummarizedExperiment, ExpressionSet, matrix, RaggedExperiment (inherits from GRangesList), and RangedVcfStack. Create new MultiAssayExperiment instances with the homonymous constructor, minimally with the argument ExperimentList, potentially also with the arguments colData (see section below) and sampleMap.

Usage

## S4 method for signature 'MultiAssayExperiment'
show(object)

## S4 method for signature 'MultiAssayExperiment'
length(x)

## S4 method for signature 'MultiAssayExperiment'
names(x)

## S4 method for signature 'MultiAssayExperiment'
MultiAssayExperiment-class

updateObject(object, ..., verbose = FALSE)

## S4 method for signature 'MultiAssayExperiment'

dimnames(x)

## S4 method for signature 'MultiAssayExperiment'
c(x, ..., sampleMap = NULL, mapFrom = NULL)

## S4 method for signature 'MultiAssayExperiment'
exportClass(
  object,
  dir = tempdir(),
  fmt,
  ext,
  match = FALSE,
  verbose = TRUE,
  ...
)

## S4 method for signature 'MultiAssayExperiment'
assays(x, withDimnames = TRUE, ...)

## S4 method for signature 'MultiAssayExperiment,missing'
assay(x, i, withDimnames = TRUE, ...)

## S4 method for signature 'MultiAssayExperiment,numeric'
assay(x, i, withDimnames = TRUE, ...)

## S4 method for signature 'MultiAssayExperiment(character'
assay(x, i, withDimnames = TRUE, ...)

Arguments

object, x A MultiAssayExperiment object

... Additional arguments for supporting functions. See details.

verbose logical(1) Whether to print additional information (default TRUE)

sampleMap c method: a sampleMap list or DataFrame to guide merge

mapFrom Either a logical, character, or integer vector indicating the experiment(s)
that have an identical colname order as the experiment input(s). If using a char-
acter input, the name must match exactly.

dir character(1) A directory for saving exported data (default: tempdir())

fmt character(1) or function() Either a format character atomic as supported by
write.table either ('csv', or 'tsv') or a function whose first two arguments are
'object to save' and 'file location'

ext character(1) A file extension supported by the format argument

match logical(1) Whether to coerce the current object to a 'MatchedAssayExperi-
ment' object (default: FALSE)
MultiAssayExperiment-class

withDimnames logical (default TRUE) whether to return dimension names included in the object

An integer or character scalar indicating the assay to return

Details

The dots (...) argument allows the user to specify additional arguments in several instances.

- subsetting []: additional arguments sent to findOverlaps.
- mergeReplicates: used to specify arguments for the simplify functional argument
- assay: may contain withDimnames, which is forwarded to assays
- combining c: compatible MultiAssayExperiment classes passed on to the ExperimentList constructor, can be a list, List, or a series of named arguments. See the examples below.

Value

A MultiAssayExperiment object

Methods (by generic)

- show(MultiAssayExperiment): Show method for a MultiAssayExperiment
- length(MultiAssayExperiment): Get the length of ExperimentList
- names(MultiAssayExperiment): Get the names of the ExperimentList
- updateObject(MultiAssayExperiment): Update old serialized MultiAssayExperiment objects to new API
- dimnames(MultiAssayExperiment): Get the dimension names for a MultiAssayExperiment object
- c(MultiAssayExperiment): Add a supported data class to the ExperimentList
- exportClass(MultiAssayExperiment): Export data from class to a series of text files
- assays(MultiAssayExperiment): Obtain a SimpleList of assay data for all available experiments in the object
- assay(x = MultiAssayExperiment, i = missing): Convenience function for extracting the assay of the first element (default) in the ExperimentList. A numeric or character index can also be provided

Slots

ExperimentList A ExperimentList class object for each assay dataset
colData A DataFrame of all clinical/specimen data available across experiments
sampleMap A DataFrame of translatable identifiers of samples and participants
metadata Additional data describing the MultiAssayExperiment object
drops A metadata list of dropped information
colData

The colData slot is a collection of primary specimen data valid across all experiments. This slot is strictly of class DataFrame but arguments for the constructor function allow arguments to be of class data.frame and subsequently coerced.

ExperimentList

The ExperimentList slot is designed to contain results from each experiment/assay. It contains a SimpleList.

sampleMap

The sampleMap contains a DataFrame of translatable identifiers of samples and participants or biological units. The standard column names of the sampleMap are "assay", "primary", and "col-name". Note that the "assay" column is a factor corresponding to the names of each experiment in the ExperimentList. In the case where these names do not match between the sampleMap and the experiments, the documented experiments in the sampleMap take precedence and experiments are dropped by the harmonization procedure. The constructor function will generate a sampleMap in the case where it is not provided and this method may be a 'safer' alternative for creating the MultiAssayExperiment (so long as the rownames are identical in the colData, if provided). An empty sampleMap may produce empty experiments if the levels of the "assay" factor in the sampleMap do not match the names in the ExperimentList.

coection

Convert a list or S4 List to a MultiAssayExperiment object using the methods::as function.
In the following example, x is either a list or List:
as(x, "MultiAssayExperiment")
Convert a MultiAssayExperiment to MAF class object using the methods::as function.
In the following example, x is a MultiAssayExperiment:
MultiAssayExperimentToMAF(x)

See Also

MultiAssayExperiment-methods for slot modifying methods, MultiAssayExperiment API

Examples

example("MultiAssayExperiment")

## Subsetting
# Rows (i) Rows/Features in each experiment
mae[, , ]
mae[c(TRUE, FALSE), , ]

# Columns (j) Rows in colData
mae[, rownames(colData(mae))[3:2], ]
## MultiAssayExperiment-helpers

A group of helper functions for manipulating and cleaning a MultiAssayExperiment

### Description

A set of helper functions were created to help clean and manipulate a MultiAssayExperiment object. intersectRows also works for ExperimentList objects.

- `complete.cases`: Returns a logical vector corresponding to `colData` rows that have data across all experiments
- `isEmpty`: Returns a logical TRUE value for zero length MultiAssayExperiment objects
- `intersectRows`: Takes all common rows across experiments, excludes experiments with empty rownames
- `intersectColumns`: A wrapper for `complete.cases` to return a MultiAssayExperiment with only those biological units that have measurements across all experiments
- `replicated`: Identifies, via logical vectors, colnames that originate from a single biological unit within each assay
- `replicates`: Provides the replicate colnames found with the replicated function by their name, empty list if none
• anyReplicated: Whether the assay has replicate measurements
• showReplicated: Displays the actual columns that are replicated per assay and biological unit, i.e., primary value (colData rowname) in the sampleMap
• mergeReplicates: A function that combines replicated / repeated measurements across all experiments and is guided by the replicated return value
• longFormat: A MultiAssayExperiment method that returns a small and skinny DataFrame. The colDataCols arguments allows the user to append colData columns to the data.
• wideFormat: A function to reshape the data in a MultiAssayExperiment to a "wide" format DataFrame. Each row in the DataFrame represents an observation (corresponding to an entry in the colData). If replicates are present, their data will be appended at the end of the corresponding row and will generate additional NA data. It is recommended to remove or consolidate technical replicates with mergeReplicates. Optional colDataCols can be added when the original object is a MultiAssayExperiment.
• hasRowRanges: A function that identifies ExperimentList elements that have a rowRanges method
• getWithColData: A convenience function for extracting an assay and associated colData
• renamePrimary: A convenience function to rename the primary biological units as represented in the rownames(colData)
• renameColname: A convenience function to rename the colnames of a particular assay

Usage

```
## S4 method for signature 'MultiAssayExperiment'
complete.cases(...)  # complete cases

## S4 method for signature 'MultiAssayExperiment'
isEmpty(x)  # is empty

intersectRows(x)

intersectColumns(x)

replicated(x)

## S4 method for signature 'MultiAssayExperiment'
replicated(x)

anyReplicated(x)

## S4 method for signature 'MultiAssayExperiment'
anyReplicated(x)

showReplicated(x)

## S4 method for signature 'MultiAssayExperiment'
showReplicated(x)
```
replicates(x, ...)

## S4 method for signature 'MultiAssayExperiment'
replicates(x, ...)

mergeReplicates(x, replicates = list(), simplify = BiocGenerics::mean, ...)

## S4 method for signature 'MultiAssayExperiment'
mergeReplicates(
  x,
  replicates = replicated(x),
  simplify = BiocGenerics::mean,
  ...
)

## S4 method for signature 'ANY'
mergeReplicates(x, replicates = list(), simplify = BiocGenerics::mean, ...)

longFormat(object, colDataCols = NULL, i = 1L)

wideFormat(
  object,
  colDataCols = NULL,
  check.names = TRUE,
  collapse = "_",
  i = 1L
)

hasRowRanges(x)

## S4 method for signature 'MultiAssayExperiment'
hasRowRanges(x)

## S4 method for signature 'ExperimentList'
hasRowRanges(x)

getchWithColData(x, i, mode = c("append", "replace"), verbose = FALSE)

renamePrimary(x, value)

tenameColname(x, i, value)

splitAssays(x, hitList)

## S4 method for signature 'MultiAssayExperiment'
splitAssays(x, hitList)
makeHitList(x, patternList)

Arguments

... Additional arguments. See details for more information.

x A MultiAssayExperiment or ExperimentList

replicates A list of LogicalLists indicating multiple / duplicate entries for each biological unit per assay, see replicated (default replicated(x)).

simplify A function for merging repeat measurements in experiments as indicated by the replicated method for MultiAssayExperiment

object Any supported class object

colDataCols A character, logical, or numeric index for colData columns to be included

i longFormat: The i-th assay in SummarizedExperiment-like objects. A vector input is supported in the case that the SummarizedExperiment object(s) has more than one assay (default 1L), renameColname: Either a numeric or character index indicating the assay whose colnames are to be renamed

check.names (logical default TRUE) Column names of the output DataFrame will be checked for syntactic validity and made unique, if necessary

collapse (character default ".") A single string delimiter for output column names. In wideFormat, experiments and rownames (and when replicate samples are present, colnames) are separated by this delimiter

mode String indicating how MultiAssayExperiment column-level metadata should be added to the SummarizedExperiment colData.

verbose logical(1) Whether to suppressMessages on subsetting operations in getWithColData (default FALSE)

value renamePrimary: A character vector of the same length as the existing rownames(colData) to use for replacement, renameColname: A CharacterList or list with matching lengths to replace colnames(x)

hitList a named list or List of logical vectors that indicate groupings in the assays

patternList a named list or List of atomic character vectors that are the input to grepl for identifying groupings in the assays

Details

The replicated function finds replicate measurements in each assay and returns a list of LogicalLists. Each element in a single LogicalList corresponds to a biological or primary unit as in the sampleMap. Below is a small graphic for one particular biological unit in one assay, where the logical vector corresponds to the number of measurements/samples in the assay:

```r
> replicated(MultiAssayExperiment)
(list str) '--- $ AssayName
(LogicalList str) '--- [[ "Biological Unit" ]] Replicated if sum(...) > 1 '--- TRUE TRUE FALSE FALSE
```
anyReplicated determines if any of the assays have at least one replicate. Note. These methods are not available for the ExperimentList class due to a missing sampleMap structure (by design). showReplicated returns a list of CharacterLists where each element corresponds to the the biological or primary units that are replicated in that assay element. The values in the inner list are the colnames in the assay that are technical replicates.

The replicates function (noun) returns the colnames from the sampleMap that were identified as replicates. It returns a list of CharacterLists for each assay present in the MultiAssayExperiment and an inner entry for each biological unit that has replicate observations in that assay.

The mergeReplicates function is a house-keeping method for a MultiAssayExperiment where only complete cases are returned. This by-assay operation averages replicate measurements (by default) and columns are aligned by the row order in colData. Users can provide their own function for merging replicates with the simplify functional argument. Additional inputs ... are sent to the 'simplify' function.

The mergeReplicates "ANY" method consolidates duplicate measurements for rectangular data structures, returns object of the same class (endomorphic). The ellipsis or ... argument allows the user to provide additional arguments to the simplify functional argument.

The longFormat "ANY" class method, works with classes such as ExpressionSet and SummarizedExperiment as well as matrix to provide a consistent long and skinny DataFrame.

The hasRowRanges method identifies assays that support a rowRanges method and return a GRanges object.

Value
See the itemized list in the description section for details.

mergeReplicates
The mergeReplicates function makes use of the output from replicated which will point out the duplicate measurements by biological unit in the MultiAssayExperiment. This function will return a MultiAssayExperiment with merged replicates. Additional arguments can be provided to the simplify argument via the ellipsis (...). For example, when replicates "TCGA-B" and "TCGA-A" are found in the assay, the name of the first appearing replicate is taken (i.e., "B"). Note that a typical use case of merging replicates occurs when there are multiple measurements on the same sample (within the same assay) and can therefore be averaged.

longFormat
The 'longFormat' method takes data from the ExperimentList in a MultiAssayExperiment and returns a uniform DataFrame. The resulting DataFrame has columns indicating primary, rowname, colname and value. This method can optionally include columns of the MultiAssayExperiment colData named by colDataCols character vector argument. (MultiAssayExperiment method only). The i argument allows the user to specify the assay value for the SummarizedExperiment assay function’s i argument.

wideFormat
The wideFormat function returns standardized wide DataFrame where each row represents a biological unit as in the colData. Depending on the data and setup, biological units can be patients,
tumors, specimens, etc. Metadata columns are generated based on the names produced in the wide format DataFrame. These can be accessed via the `mcols()` function. See the wideFormat section for description of the `colDataCols` and `i` arguments.

**hasRowRanges**

The `hasRowRanges` method identifies assays with associated ranged row data by directly testing the method on the object. The result from the test must be a `GRanges` class object to satisfy the test.

**getWithColData**

The `getWithColData` function allows the user to conveniently extract a particular assay as indicated by the `i` index argument. It will also attempt to provide the `colData` along with the extracted object using the `colData<-` replacement method when possible. Typically, this method is available for `SummarizedExperiment` and `RaggedExperiment` classes.

The setting of `mode` determines how the `colData` is added. If `mode="append"`, the `MultiAssayExperiment` metadata is appended onto that of the `SummarizedExperiment`. If any fields are duplicated by name, the values in the `SummarizedExperiment` are retained, with a warning emitted if the values are different. For `mode="replace"`, the `MultiAssayExperiment` metadata replaces that of the `SummarizedExperiment`, while for `mode="none"`, no replacement or appending is performed.

**rename**

The `renamePrimary` function allows the user to conveniently change the actual names of the primary biological units as seen in `rownames(colData)`. `renameColname` allows the user to change the names of a particular assay based on index `i`. `i` can either be a single numeric or character value. See `colnames<-` method for renaming multiple `colnames` in a `MultiAssayExperiment`.

**splitAssays**

The `splitAssays` method separates columns in each of the assays based on the `hitList` input. The `hitList` can be generated using the `makeHitList` helper function. To use the `makeHitList` helper, the user should input a list of patterns that will match on the column names of each assay. These matches should be mutually exclusive as to avoid repetition of columns across assays. See the examples section.

**Examples**

```r
example(MultiAssayExperiment)
complete.cases(mae)
isEmpty(MultiAssayExperiment())
```

```r
## renaming biological units (primary)
mae2 <- renamePrimary(mae, paste0("pt", 1:4))
colData(mae2)
sampleMap(mae2)
```
```r
## renaming observational units (colname)
mae2 <- renameColname(mae, i = "Affy", paste0("ARRAY", 1:4))
colnames(mae2)
sampleMap(mae2)

patts <- list(
    normals = "TCGA-[A-Z0-9]{2}-[A-Z0-9]{4}-11",
    tumors = "TCGA-[A-Z0-9]{2}-[A-Z0-9]{4}-01"
)
data("miniACC")

hits <- makeHitList(miniACC, patts)

## only tumors present
splitAssays(miniACC, hits)
```

---

**MultiAssayExperiment-methods**

*Accessing and modifying information in MultiAssayExperiment*

---

**Description**

A set of accessor and setter generic functions to extract either the `sampleMap`, the `ExperimentList`, `colData`, or `metadata` slots of a MultiAssayExperiment object.

**Usage**

```r
## S4 method for signature 'MultiAssayExperiment'
sampleMap(x)

## S4 method for signature 'MultiAssayExperiment'
experiments(x)

## S4 method for signature 'MultiAssayExperiment'
colData(x, ...)

## S4 method for signature 'MultiAssayExperiment'
drops(x)

## S4 replacement method for signature 'MultiAssayExperiment,DataFrame'
sampleMap(object) <- value
```
## S4 replacement method for signature 'MultiAssayExperiment,ANY'
sampleMap(object) <- value
drops(x, ...) <- value

## S4 replacement method for signature 'MultiAssayExperiment,ExperimentList'
experiments(object) <- value

## S4 replacement method for signature 'MultiAssayExperiment,List'
experiments(object) <- value

## S4 replacement method for signature 'MultiAssayExperiment,DataFrame'
colData(x) <- value

colData(x) <- value

colData(x) <- value
drops(x, ...) <- value

## S4 replacement method for signature 'MultiAssayExperiment'
x$name <- value

names(x) <- value

## S4 replacement method for signature 'MultiAssayExperiment,List'
colnames(x) <- value
colnames(x) <- value
colnames(x) <- value

colnames(x) <- value

colnames(x) <- value

colnames(x) <- value

## S4 method for signature 'MultiAssayExperiment'
x$name

colnames(x)

## S4 method for signature 'MultiAssayExperiment'
metadata(x, ...)

## S4 replacement method for signature 'MultiAssayExperiment'
metadata(x, ...) <- value

### Arguments

- **...**  
  Argument not in use

- **object, x**  
  A MultiAssayExperiment object

- **value**  
  See details.

- **name**  
  A column in colData
Value

Accessors: Either a sampleMap, ExperimentList, or DataFrame object
Setters: A MultiAssayExperiment object

Accessors

Eponymous names for accessing MultiAssayExperiment slots with the exception of the ExperimentList accessor named experiments.

- colData: Access the colData slot
- sampleMap: Access the sampleMap slot
- experiments: Access the ExperimentList slot
- [[: Access the ExperimentList slot
- $: Access a column in colData
- drops: Get a vector of dropped ExperimentList names

Setters

Setter method values (i.e., 'function(x) <- value'):

- experiments<-: An ExperimentList object containing experiment data of supported classes
- sampleMap<-: A DataFrame object relating samples to biological units and assays
- colData<-: A DataFrame object describing the biological units
- metadata<-: A list object of metadata
- [[<-: Equivalent to the experiments<- setter method for convenience
- $<-: A vector to replace the indicated column in colData
- drops<-: Trace ExperimentList names that have been removed

Examples

```r
## Load example MultiAssayExperiment
defer2(MultiAssayExperiment)

## Access the sampleMap
defer2(sampleMap(mae))

## Replacement method for a MultiAssayExperiment sampleMap
defer2(sampleMap(mae) <- S4Vectors::DataFrame())

## Access the ExperimentList
defer2(experiments(mae))

## Replace with an empty ExperimentList
defer2(experiments(mae) <- ExperimentList())

## Access the metadata
defer2(metadata(mae))
```
## Replace metadata with a list

```r
metadata(mae) <- list(runDate = format(Sys.time(), "%B %d, %Y"))
```

## Access the colData

```r
colData(mae)
```

## Access a column in colData

```r
mae$age
```

## Replace a column in colData

```r
mae$age <- mae$age + 1
```

---

### MultiAssayExperimentToMAF

*Convert MultiAssayExperiment to MAF class*

#### Description

Take a `MultiAssayExperiment` object with specific mutation assays and convert these into a `maftools` representation. The names provided via `synAssay` and `nonSynAssay` must match exactly those assays in the `MultiAssayExperiment`.

#### Usage

```r
MultiAssayExperimentToMAF(x, synAssay = "maf_syn", nonSynAssay = "maf_nonSyn")
```

#### Arguments

- `x` A `MultiAssayExperiment` object
- `synAssay` character(1) The name of the `ExperimentList` element in the `MultiAssayExperiment` that identifies synonymous variant classifications.
- `nonSynAssay` character(1) The name of the `ExperimentList` element in the `MultiAssayExperiment` that identifies non-synonymous variant classifications.

---

### prepMultiAssay

*Prepare a MultiAssayExperiment instance*

#### Description

The purpose of this helper function is to facilitate the creation of a `MultiAssayExperiment` object by detecting any inconsistencies with all types of names in either the `ExperimentList`, the `colData`, or `sampleMap`. 
**Usage**

```
prepMultiAssay(ExperimentList, colData, sampleMap, ...)
```

**Arguments**

- **ExperimentList**  
  A list of all combined experiments
- **colData**  
  A `DataFrame` of the phenotype data for all participants
- **sampleMap**  
  A `DataFrame` of sample identifiers, assay samples, and assay names
- **...**  
  Optional arguments for the `MultiAssayExperiment` constructor function such as metadata and `drops`.

**Value**

A list containing all the essential components of a `MultiAssayExperiment` as well as a "drops" metadata element that indicates non-matched names. The names of the resulting list correspond to the arguments of the `MultiAssayExperiment` constructor function.

**Checks**

The `prepMultiAssay` function checks that all columns in the `sampleMap` are character.

It checks that all names and lengths match in both the `ExperimentList` and in the unique assay names of the `sampleMap`.

If `ExperimentList` names and assay names only differ by case and are not duplicated, the function will standardize all names to lowercase.

If names cannot be matched between the colname column of the `sampleMap` and the colnames of the `ExperimentList`, those unmatched will be dropped and found in the "drops" element of the resulting list.

Names in the "primary" column of the `sampleMap`, will be matched to those in the `colData`. Unmatched "primary" column rows will be dropped from the `sampleMap`. Suggestions for name fixes in either the `ExperimentList` or colnames will be made when necessary.

**Examples**

```r
## Run example
example("MultiAssayExperiment")

## Check if there are any inconsistencies within the different names
preparedMAE <- prepMultiAssay(ExpList, colDat, sampMap)

## Results in a list of components for the MultiAssayExperiment constructor function
MultiAssayExperiment(preparedMAE$experiments, preparedMAE$colData, preparedMAE$sampleMap)

## Alternatively, use the do.call function
do.call(MultiAssayExperiment, preparedMAE)
```
Description

These objects are imported from other packages. Click on the function name to see their documentation.

- S4Vectors: DataFrame

Examples

DataFrame()

---

saveHDF5MultiAssayExperiment

Save a MultiAssayExperiment class object to HDF5 and Rds files

Description

This function takes a MultiAssayExperiment object and uses the assays functionality to obtain data matrices out of the experiments. These are then saved into the .h5 file format. This function relies heavily on the HDF5Array package whose installation is required before use. saveHDF5MultiAssayExperiment preserves the classes contained in the ExperimentList with the exception of matrix which is converted to HDF5Matrix. Internal SummarizedExperiment assays are converted to HDF5-backed assays as in HDF5Array::saveHDF5SummarizedExperiment. SummarizedExperiment objects with multiple i-th assays will have the first assay take precedence and others assays will be dropped with a warning. If the first assay in a SummarizedExperiment contains an array, the array is preserved in the process of saving and loading the HDF5-backed MultiAssayExperiment.

Usage

```r
saveHDF5MultiAssayExperiment(x, 
  dir = "h5_mae", 
  prefix = NULL, 
  replace = FALSE, 
  chunkdim = NULL, 
  level = NULL, 
  as.sparse = NA, 
  verbose = NA)
```

loadHDF5MultiAssayExperiment(dir = "h5_mae", prefix = NULL)
Arguments

- **x**: A `MultiAssayExperiment` object or derivative
- **dir**: The path (as a single string) to the directory where to save the HDF5-based `MultiAssayExperiment` object or to load it from.
  When saving, the directory will be created if it doesn’t already exist. If the directory already exists and no prefix is specified and replace is set to `TRUE`, then it’s replaced with an empty directory.
- **prefix**: An optional prefix to add to the names of the files created inside `dir`. This allows saving more than one object in the same directory. When the prefix is `NULL`, the name of the `x` input `MultiAssayExperiment` is used. To avoid the default setting use an empty character string i.e., "". An underscore (_) is appended to the prefix when provided; therefore, typical inputs should be words, e.g., "test".
- **replace**: When no prefix is specified, should a pre-existing directory be replaced with a new empty one? The content of the pre-existing directory will be lost!
- **chunkdim, level**: The dimensions of the chunks and the compression level to use for writing the assay data to disk.
  Passed to the internal calls to `writeHDF5Array`. See `?writeHDF5Array` for more information.
- **as.sparse**: Whether the assay data should be flagged as sparse or not. If set to `NA` (the default), then the specific `as.sparse` value to use for each assay is determined by calling `is_sparse()` on them.
  Passed to the internal calls to `writeHDF5Array`. See `?writeHDF5Array` for more information and an IMPORTANT NOTE.
- **verbose**: Set to `TRUE` to make the function display progress.
  In the case of `saveHDF5MultiAssayExperiment()`, `verbose` is set to `NA` by default, in which case verbosity is controlled by `DelayedArray.verbose.block.processing` option. Setting `verbose` to `TRUE` or `FALSE` overrides the option.

Examples

data("miniACC")

testDir <- file.path(tempdir(), "test_mae")

saveHDF5MultiAssayExperiment(
  miniACC, dir = testDir, verbose = TRUE, replace = TRUE
)

# inspect the files in the dir
list.files(testDir)

loadHDF5MultiAssayExperiment(
  dir = testDir
)

# remove example files
unlink(testDir, recursive = TRUE)
subsetBy

Subsetting a MultiAssayExperiment object

Description
A set of functions for extracting and dividing a MultiAssayExperiment

Usage
subsetByRow(x, y, ...)
subsetByColData(x, y)
subsetByColumn(x, y)
subsetByAssay(x, y)

## S4 method for signature 'ExperimentList,ANY'
subsetByRow(x, y, ...)

## S4 method for signature 'ExperimentList,list'
subsetByRow(x, y)

## S4 method for signature 'ExperimentList,List'
subsetByRow(x, y)

## S4 method for signature 'ExperimentList,logical'
subsetByRow(x, y)

## S4 method for signature 'ExperimentList,list'
subsetByColumn(x, y)

## S4 method for signature 'ExperimentList,List'
subsetByColumn(x, y)

## S4 method for signature 'ExperimentList,logical'
subsetByColumn(x, y)

## S4 method for signature 'ExperimentList'
subsetByAssay(x, y)

## S4 method for signature 'MultiAssayExperiment,ANY'
subsetByColData(x, y)

## S4 method for signature 'MultiAssayExperiment,character'
subsetByColData(x, y)
## S4 method for signature 'MultiAssayExperiment,ANY'
subsetByRow(x, y, ...)

## S4 method for signature 'MultiAssayExperiment,ANY'
subsetByColumn(x, y)

## S4 method for signature 'MultiAssayExperiment'
subsetByAssay(x, y)

## S4 method for signature 'MultiAssayExperiment,ANY,ANY'
x[i, j, k, ..., drop = FALSE]

## S4 method for signature 'MultiAssayExperiment,ANY'
x[[i, j, ...]]

## S4 replacement method for signature 'MultiAssayExperiment,ANY,ANY'
x[[i, j, ...]] <- value

## S4 replacement method for signature 'MultiAssayExperiment,ANY,ANY'
x[i, j, ...] <- value

### Arguments

**x**  
A MultiAssayExperiment or ExperimentList

**y**  
Any argument used for subsetting, can be a character, logical, integer, list or List vector

**...**  
Additional arguments passed on to lower level functions.

**i**  
Either a character, integer, logical or GRanges object for subsetting by rows

**j**  
Either a character, logical, or numeric vector for subsetting by colData rows. See details for more information.

**k**  
Either a character, logical, or numeric vector for subsetting by assays

**drop**  
logical (default FALSE) whether to drop all empty assay elements in the ExperimentList

**value**  
An assay compatible with the MultiAssayExperiment API

### Details

Subsetting a MultiAssayExperiment by the j index can yield a call to either subsetByColData or subsetByColumn. For vector inputs, the subset will be applied to the colData rows. For List-type inputs, the List will be applied to each of the elements in the ExperimentList. The order of the subsetting elements in the List must match that of the ExperimentList in the MultiAssayExperiment.

- subsetBycolData: Select biological units by vector input types
- subsetByColumn: Select observations by assay or for each assay
- subsetByRow: Select rows by assay or for each assay
- subsetByAssay: Select experiments
Value

subsetBy* operations are endomorphic and return either MultiAssayExperiment or ExperimentList depending on the input.

Examples

```r
## Load the example MultiAssayExperiment
dput(mae)

## Using experiment names
dputByAssay(mae, "Affy")

## Using numeric indices
dputByAssay(mae, 1:2)

## Using a logical vector
dputByAssay(mae, c(TRUE, FALSE, TRUE))

## Subset by character vector (Jack)
dputByColData(mae, "Jack")

## Subset by numeric index of colData rows (Jack and Bob)
dputByColData(mae, c(1, 3))

## Subset by logical indicator of colData rows (Jack and Jill)
dputByColData(mae, c(TRUE, TRUE, FALSE, FALSE))

dputByColumn(mae, list(Affy = 1:2,  
                   Methyl450k = c(3,5,2), RNASeqGene = 2:4, GISTIC = 1))

# Use a GRanges object to subset rows where ranged data present  
egr <- GenomicRanges::GRanges(seqnames = "chr2",  
                         IRanges::IRanges(start = 11, end = 13), strand = 
system("-")

# Use a logical vector (recycling used)
dputByRow(mae, c(TRUE, FALSE))

# Use a character vector  
dputByRow(mae, "ENST00000355076")
```

| upsetSamples | Create a generalized Venn Diagram analog for sample membership in multiple assays, using the upset algorithm in UpSetR |
Description

Create a generalized Venn Diagram analog for sample membership in multiple assays, using the upset algorithm in UpSetR

Usage

upsetSamples(
  MultiAssayExperiment,
  nsets = NULL,
  sets = names(MultiAssayExperiment),
  nintersects = NA_integer_,
  order.by = "freq",
  check.names = FALSE,
  ...
)

Arguments

MultiAssayExperiment

A MultiAssayExperiment object

nsets numeric(1) The number of sets to analyze. If specified, sets will be ignored.

sets character() A character vector of names in MultiAssayExperiment to use. If specified, nsets will be ignored.

nintersects numeric(1) The number of intersections to plot. By default, all intersections will be plotted.

order.by How the intersections in the matrix should be ordered by. Options include frequency (entered as "freq"), degree, or both in any order.

check.names logical(1) Whether to munge names as in the data.frame() constructor (default FALSE).

... parameters passed to UpSetR::upset

Value

Produces a visualization of set intersections using the UpSet matrix design

Note

This function is intended to provide convenient visualization of assay availability configurations in MultiAssayExperiment instances. The UpSetR::upset function requires data.frame input and has many parameters to tune appearance of the result. Assay name handling is important for interpretability.

Author(s)

Vincent J Carey
Examples

```r
data(miniACC)
upsetSamples(miniACC)
upsetSamples(miniACC, nsets = 3, nintersects = 3)
```
Index

* data

  miniACC, 10
  [,MultiAssayExperiment,ANY,ANY,ANY-method (subsetBy), 30
  [,MultiAssayExperiment,ANY-method (subsetBy), 30
  [<-,MultiAssayExperiment,ANY,ANY,ANY-method (subsetBy), 30
  [[,MultiAssayExperiment,ANY,ANY-method (subsetBy), 30
  [[<-,MultiAssayExperiment,ANY,ANY-method (subsetBy), 30
  $,MultiAssayExperiment-method (MultiAssayExperiment-methods), 23
  $<-,MultiAssayExperiment-method (MultiAssayExperiment-methods), 23

  anyReplicated (MultiAssayExperiment-methods), 17
  anyReplicated,MultiAssayExperiment-method (MultiAssayExperiment-methods), 17
  assay,ANY,missing-method (ExperimentList-class), 5
  assay,ExperimentList,character-method (ExperimentList-class), 5
  assay,ExperimentList,missing-method (ExperimentList-class), 5
  assay,ExperimentList,numeric-method (ExperimentList-class), 5
  assay,MultiAssayExperiment,character-method (MultiAssayExperiment-class), 13
  assay,MultiAssayExperiment,missing-method (MultiAssayExperiment-class), 13
  assay,MultiAssayExperiment,numeric-method (MultiAssayExperiment-class), 13
  assays,ExperimentList-method (ExperimentList-class), 5
  assays,MultiAssayExperiment-method (MultiAssayExperiment-class), 13
  c,MultiAssayExperiment-method (MultiAssayExperiment-class), 13
  CharacterList, 6, 21
  coerce,ExperimentList-method (ExperimentList-class), 5
  coerce,MultiAssayExperiment-method (MultiAssayExperiment-class), 13
  colData, 22
  colData,MultiAssayExperiment-method (MultiAssayExperiment-methods), 23
  colData<-,MultiAssayExperiment,ANY-method (MultiAssayExperiment-methods), 23
  colData<-,MultiAssayExperiment,DataFrame-method (MultiAssayExperiment-methods), 23

  35
colnames, ExperimentList-method (ExperimentList-class), 5
colnames<-, MultiAssayExperiment, List-method (MultiAssayExperiment-methods), 23
colnames<-, MultiAssayExperiment, list-method (MultiAssayExperiment-methods), 23
complete.cases, MultiAssayExperiment-method (MultiAssayExperiment-methods), 17
DataFrame, 8, 11, 16, 18, 21, 25, 27, 28
DataFrame (reexports), 28
dimnames, ExperimentList-method (ExperimentList-class), 5
dimnames, MultiAssayExperiment-method (MultiAssayExperiment-class), 13
drops (MultiAssayExperiment-methods), 23
drops, MultiAssayExperiment-method (MultiAssayExperiment-methods), 23
drops<-, MultiAssayExperiment-method (MultiAssayExperiment-methods), 23
drops<-, MultiAssayExperiment, List-method (MultiAssayExperiment-methods), 23
ExperimentList, 4, 6, 7, 11–13, 15, 16, 21, 23, 25–28
ExperimentList-class, 5
experiments (MultiAssayExperiment-methods), 23
experiments, MultiAssayExperiment-method (MultiAssayExperiment-methods), 23
experiments<-, MultiAssayExperiment-method (MultiAssayExperiment-methods), 23
exportClass, MultiAssayExperiment-method (MultiAssayExperiment-class), 13
exportClass, MultiAssayExperiment-method (MultiAssayExperiment-class), 13
ExpressionSet, 13, 21
findOverlaps, 15
getWithColData (MultiAssayExperiment-methods), 17
GRanges, 21, 22
GRangesList, 13
hasAssay, 7
hasRowRanges (MultiAssayExperiment-methods), 17
hasRowRanges, ExperimentList-method (MultiAssayExperiment-methods), 17
hasRowRanges, MultiAssayExperiment-method (MultiAssayExperiment-methods), 17
HDF5MultiAssayExperiment (saveHDF5MultiAssayExperiment), 28
intersectColumns (MultiAssayExperiment-methods), 17
intersectRows (MultiAssayExperiment-methods), 17
isEmpty, ExperimentList-method (ExperimentList-class), 5
isEmpty, MultiAssayExperiment-method (MultiAssayExperiment-methods), 17
length, MultiAssayExperiment-method (MultiAssayExperiment-class), 13
List, 7, 15, 16
ListToMap, 8
loadHDF5MultiAssayExperiment (saveHDF5MultiAssayExperiment), 28
LogicalList, 6, 20
longFormat
  (MultiAssayExperiment-helpers), 17
makeHitList
  (MultiAssayExperiment-helpers), 17
makeMatchList
  (MultiAssayExperiment-helpers), 17
mapToList(listToMap), 8
MatchedAssayExperiment
  (MatchedAssayExperiment-class), 9
MatchedAssayExperiment-class, 9
mcols(), 22
mergeReplicates
  (MultiAssayExperiment-helpers), 17
mergeReplicates, ANY-method
  (MultiAssayExperiment-helpers), 17
mergeReplicates, ExperimentList-method
  (ExperimentList-class), 5
mergeReplicates, MultiAssayExperiment-method
  (MultiAssayExperiment-helpers), 17
metadata, MultiAssayExperiment-method
  (MultiAssayExperiment-methods), 23
metadata<-, MultiAssayExperiment-method
  (MultiAssayExperiment-methods), 23
methods::as, 16
miniACC, 10
MultiAssayExperiment, 9–11, 12, 23, 26, 27, 29
MultiAssayExperiment-class, 13
MultiAssayExperiment-helpers, 17
MultiAssayExperiment-methods, 16, 23
MultiAssayExperiment-package, 3
MultiAssayExperimentToMAF, 26
names, MultiAssayExperiment-method
  (MultiAssayExperiment-class), 13
names<-, MultiAssayExperiment-method
  (MultiAssayExperiment-methods), 23
prepMultiAssay, 26
reexports, 28
renameColname
  (MultiAssayExperiment-helpers), 17
renamePrimary
  (MultiAssayExperiment-helpers), 17
replicated
  (MultiAssayExperiment-helpers), 17
replicated, MultiAssayExperiment-method
  (MultiAssayExperiment-helpers), 17
replicates
  (MultiAssayExperiment-helpers), 17
replicates, MultiAssayExperiment-method
  (MultiAssayExperiment-helpers), 17
rownames, 6
rownames, ExperimentList-method
  (ExperimentList-class), 5
rowRanges, 18, 21
sampleMap. 12, 13, 16, 26, 27
sampleMap
  (MultiAssayExperiment-methods), 23
sampleMap, MultiAssayExperiment-method
  (MultiAssayExperiment-methods), 23
sampleMap<-
  (MultiAssayExperiment-methods), 23
sampleMap<-, MultiAssayExperiment, ANY-method
  (MultiAssayExperiment-methods), 23
sampleMap<-, MultiAssayExperiment, DataFrame-method
  (MultiAssayExperiment-methods), 23
saveHDF5MultiAssayExperiment, 28
show, ExperimentList-method
  (ExperimentList-class), 5
show, MultiAssayExperiment-method
  (MultiAssayExperiment-class), 13
showReplicated
  (MultiAssayExperiment-helpers), 17

showReplicated,MultiAssayExperiment-method
  (MultiAssayExperiment-helpers), 17

SimpleList, 12, 15, 16

splitAssays
  (MultiAssayExperiment-helpers), 17

splitAssays,MultiAssayExperiment-method
  (MultiAssayExperiment-helpers), 17

subset (subsetBy), 30
subsetBy, 30
subsetByAssay (subsetBy), 30
subsetByAssay,ExperimentList-method
  (subsetBy), 30
subsetByAssay,MultiAssayExperiment-method
  (subsetBy), 30
subsetByColData (subsetBy), 30
subsetByColData,MultiAssayExperiment,ANY-method
  (subsetBy), 30
subsetByColData,MultiAssayExperiment,character-method
  (subsetBy), 30
subsetByColumn (subsetBy), 30
subsetByColumn,ExperimentList,List-method
  (subsetBy), 30
subsetByColumn,ExperimentList,list-method
  (subsetBy), 30
subsetByColumn,ExperimentList,logical-method
  (subsetBy), 30
subsetByColumn,MultiAssayExperiment,ANY-method
  (subsetBy), 30
subsetByRow (subsetBy), 30
subsetByRow,ExperimentList,ANY-method
  (subsetBy), 30
subsetByRow,ExperimentList,List-method
  (subsetBy), 30
subsetByRow,ExperimentList,list-method
  (subsetBy), 30
subsetByRow,ExperimentList,logical-method
  (subsetBy), 30
subsetByRow,MultiAssayExperiment,ANY-method
  (subsetBy), 30
SummarizedExperiment, 13, 21, 22

updateObject,MultiAssayExperiment-method
  (MultiAssayExperiment-class),