

Package ‘SEtools’

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Type Package

Title SEtools: tools for working with SummarizedExperiment

Version 1.10.0

Depends R (>= 4.0)

Description This includes a set of tools for working with the SummarizedExperiment class, including merging, melting, aggregation functions. Plotting functions historically in this package have been moved to the sechm package.

Imports BiocParallel, Matrix, SummarizedExperiment, DESeq2, S4Vectors, data.table, edgeR, openxlsx, stats, sva, sechm

Suggests BiocStyle, knitr, rmarkdown, ggplot2

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VignetteBuilder knitr

License GPL

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BugReports <https://github.com/plger/SEtools>

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aggSE	<i>aggSE</i>
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Description

Aggregates the rows of a ‘SummarizedExperiment’.

Usage

```
aggSE(x, by, assayFun = NULL, rowDatFuns = list())
```

Arguments

x	An object of class ‘SummarizedExperiment’
by	Vector by which to aggregate, or column of ‘rowData(x)’
assayFun	Function by which to aggregate, or a list of such functions (or vector of function names) of the same length as there are assays. If NULL will attempt to use an appropriate function (and notify the functions used), typically the mean.
rowDatFuns	A named list providing functions by which to aggregate each rowData columns. If a given column has no specified function, the default will be used, i.e. logical are transformed into a proportion, numerics are aggregated by median, and unique factors/characters are pasted together. Use ‘rowDataFuns=NULL’ to discard rowData.

Value

An object of class ‘SummarizedExperiment’

Examples

```
library(SummarizedExperiment)
data("SE", package="SEtools")
# arbitrary IDs for example aggregation:
rowData(SE)$otherID <- rep(LETTERS[1:10], each=10)
SE <- aggSE(SE, "otherID")
```

castSE	<i>castSE</i>
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Description

Casts a data.frame as a [SummarizedExperiment-class](#)

Usage

```
castSE(  
  x,  
  rowNames = NULL,  
  colNames = NULL,  
  assayNames = NULL,  
  colData = NULL,  
  rowData = NULL,  
  sparse = FALSE  
)
```

Arguments

x	A data.frame
rowNames	Column of 'x' containing the row.names (if omitted, will build from 'rowData')
colNames	Column of 'x' containing the column names (if omitted, will build from 'colData')
assayNames	Columns of 'x' to turn into assays
colData	Columns of 'x' to use as colData
rowData	Columns of 'x' to use as rowData
sparse	Local, whether to keep the assays sparse.

Value

A [SummarizedExperiment-class](#)

Examples

```
d <- data.frame(transcript=rep(LETTERS[1:10],each=2), gene=rep(LETTERS[1:5],each=4),  
               count=rpois(20, 10), sample=letters[1:2])  
head(d)  
castSE(d, rowData=c("transcript","gene"), colNames="sample")
```

data	<i>Example dataset</i>
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Description

A `SummarizedExperiment-class` containing (a subset of) whole-hippocampus RNAseq of mice after different stressors.

Value

a `SummarizedExperiment-class`.

References

Floriou-Servou et al. (2018). Distinct Proteomic, Transcriptomic, and Epigenetic Stress Responses in Dorsal and Ventral Hippocampus. *Biological Psychiatry*, **84**(7): 531-541. DOI: 10.1016/j.biopsych.2018.02.003.

flattenPB	<i>flattenPB</i>
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Description

Flattens a pseudo-bulk `SummarizedExperiment` as produced by ‘`muscat::aggregateData`’ so that all cell types are represented in a single assay. Optionally normalizes the data and calculates per-sample logFCs.

Usage

```
flattenPB(pb, norm = TRUE, lfc_group = NULL)
```

Arguments

pb	a pseudo-bulk <code>SummarizedExperiment</code> as produced by ‘ <code>muscat::aggregateData</code> ’, with different celltypes/clusters are assays.
norm	Logical; whether to calculate logcpm (TMM normalization).
lfc_group	the <code>colData</code> column to use to calculate foldchange. If <code>NULL</code> (default), no fold-change assay will be computed.

Value

A `SummarizedExperiment`

log2FC	<i>log2FC</i>
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Description

Generates log2(foldchange) matrix/assay, eventually on a per-batch fashion.

Usage

```
log2FC(
  x,
  fromAssay = NULL,
  controls,
  by = NULL,
  isLog = NULL,
  agFun = rowMeans,
  toAssay = "log2FC"
)
```

Arguments

<code>x</code>	A numeric matrix, or a ‘SummarizedExperiment’ object
<code>fromAssay</code>	The assay to use if ‘x’ is a ‘SummarizedExperiment’
<code>controls</code>	A vector of which samples should be used as controls for foldchange calculations.
<code>by</code>	An optional vector indicating groups/batches by which the controls will be averaged to calculate per-group foldchanges.
<code>isLog</code>	Logical; whether the data is log-transformed. If NULL, will attempt to figure it out from the data and/or assay name
<code>agFun</code>	Aggregation function for the baseline (default rowMeans)
<code>toAssay</code>	The name of the assay in which to save the output.

Value

An object of same class as ‘x’; if a ‘SummarizedExperiment’, will have the additional assay named from ‘toAssay’.

Examples

```
log2FC( matrix(rnorm(40), ncol=4), controls=1:2 )
```

meltSE	<i>meltSE</i>
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Description

Melts a SE object into a [ggplot](#)-ready long data.frame.

Usage

```
meltSE(x, genes, assayName = NULL, colDat.columns = NULL, rowDat.columns = NA)
```

Arguments

x	An object of class SummarizedExperiment-class
genes	A vector of genes to include. Use 'genes=NULL' to include all.
assayName	The name(s) of the assay(s) to use. If NULL and the assays are named, all of them will be included (if they are not named, the first one will be used).
colDat.columns	The colData columns to include (defaults includes all). Use 'colDat.columns=NA' in order not to include any.
rowDat.columns	The rowData columns to include (none included by default). Use 'rowData=NULL' to include all.

Value

A data.frame.

Examples

```
data("SE", package="SEtools")
head(meltSE(SE, "Fos"))
```

mergeSEs	<i>mergeSEs</i>
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Description

Merges a list of [SummarizedExperiment-class](#), either by row.names or through specified rowData fields. In cases of many-to-many (or one-to-many) mappings, 'aggFun' determines whether the records are aggregated by linking ID (if an aggregation method is given) or all combinations are returned (if 'aggFun=NULL' - default).

Usage

```
mergeSEs(
  ll,
  use.assays = NULL,
  do.scale = TRUE,
  commonOnly = TRUE,
  colColumns = NULL,
  mergeBy = NULL,
  aggFun = NULL,
  addDatasetPrefix = TRUE,
  defValues = list(),
  keepRowData = TRUE,
  BPPARAM = SerialParam()
)
```

Arguments

ll	A (named) list of SummarizedExperiment-class
use.assays	Names (or indexes) of the assays to use. By default, all common assays are used.
do.scale	A logical vector indicating (globally or for each assay) whether to perform row unit-variance scaling on each dataset before merging (default TRUE).
commonOnly	Logical; whether to restrict to rows present in all datasets (default TRUE).
colColumns	A character vector specifying ‘colData’ columns to include (if available in at least one of the datasets). If NULL, everything is kept.
mergeBy	The ‘rowData’ column to merge with. If NULL, row.names are used.
aggFun	The aggregation function to use when multiple rows have the same ‘mergeBy’ value. If merging multiple assays, a different function per assay can be passed as a named list (see aggSE). If NULL (default), entries will be reused to have each combination.
addDatasetPrefix	Logical; whether the name of the dataset should be appended to the sample names (default TRUE).
defValues	An optional named list of default ‘colData’ values when some columns are missing from some SEs.
keepRowData	Logical, whether to keep the rowData (default TRUE).
BPPARAM	For multithreading the aggregation step.

Value

An object of class [SummarizedExperiment-class](#)

Examples

```
data("SE", package="SEtools")
mergeSEs( list( se1=SE[,1:10], se2=SE[,11:20] ) )
```

```
resetAllSEtoolsOptions
      resetAllSEtoolsOptions
```

Description

Resets all global options relative to SEtools.

Usage

```
resetAllSEtoolsOptions()
```

Value

None

Examples

```
resetAllSEtoolsOptions()
```

```
se2xls          se2xlsx
```

Description

Writes a SummarizedExperiment to an excel/xlsx file. Requires the 'openxlsx' package.

Usage

```
se2xls(se, filename, addSheets = NULL)
```

Arguments

se	The 'SummarizedExperiment'
filename	xlsx file name
addSheets	An optional list of additional tables to save as sheets.

Value

Saves to file.

Examples

```
data("SE", package="SEtools")
# not run
# se2xls(SE, filename="SE.xlsx")
```

svacor	<i>svacor</i>
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Description

A wrapper around SVA-based correction, providing a corrected assay. If this is RNAseq data or similar, use a count assay assay with ‘useVST=TRUE’; otherwise (e.g. proteomics) a log-normalized assay is recommended.

Usage

```
svacor(
  SE,
  form,
  form0 = ~1,
  assayName = NULL,
  regressOutNull = TRUE,
  useVST = TRUE,
  n.sv = NULL,
  ...
)
```

Arguments

SE	An object of class ‘SummarizedExperiment’.
form	The formula of the differential expression model
form0	An optional formula for the null model
assayName	The name (or index) of the assay to use.
regressOutNull	Logical; whether to regress out the variables of ‘form0’.
useVST	Logical; whether to use DESeq2’s variance-stabilizing transformation; (for count data!)
n.sv	The number of surrogate variables (if omitted, sva will attempt to estimate it)
...	Any other argument passed to the sva command.

Value

Returns the ‘SummarizedExperiment’ with a ‘corrected’ assay and the surrogate variables in ‘col-Data’.

Examples

```
data("SE", package="SEtools")
SE <- svacor(SE, ~Condition)
```

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