Package ‘SEtools’

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Type Package

Title SEtools: tools for working with SummarizedExperiment

Version 1.16.0

Depends R (>= 4.0), SummarizedExperiment, sechm

Description This includes a set of convenience functions for working with the SummarizedExperiment class. Note that plotting functions historically in this package have been moved to the sechm package (see vignette for details).

Imports BiocParallel, Matrix, DESeq2, S4Vectors, data.table, edgeR, openxlsx, pheatmap, stats, circlize, methods, sva

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\textbf{R topics documented:}

\begin{itemize}
  \item \texttt{aggSE} \hspace{1cm} 2
  \item \texttt{castSE} \hspace{1cm} 3
  \item \texttt{data} \hspace{1cm} 4
  \item \texttt{flattenPB} \hspace{1cm} 4
  \item \texttt{log2FC} \hspace{1cm} 5
  \item \texttt{mergeSEs} \hspace{1cm} 6
  \item \texttt{resetAllSEtoolsOptions} \hspace{1cm} 7
  \item \texttt{se2xls} \hspace{1cm} 7
  \item \texttt{sehm} \hspace{1cm} 8
  \item \texttt{svacor} \hspace{1cm} 9
\end{itemize}

\section*{Index}

\begin{tabular}{ll}
  \texttt{aggSE} & \texttt{aggSE} \\
\end{tabular}

\section*{Description}

Aggregates the rows of a `SummarizedExperiment`.

\section*{Usage}

\texttt{aggSE(x, by, assayFun = NULL, rowDatFuns = list())}

\section*{Arguments}

\begin{itemize}
  \item \texttt{x} \hspace{1cm} An object of class `SummarizedExperiment`
  \item \texttt{by} \hspace{1cm} Vector by which to aggregate, or column of `rowData(x)`
  \item \texttt{assayFun} \hspace{1cm} 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.
  \item \texttt{rowDatFuns} \hspace{1cm} 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.
\end{itemize}

\section*{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")

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")
**flattenPB**

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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


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**flattenPB**

---

### 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

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

### Arguments

- **pb**
  - A pseudo-bulk SummarizedExperiment as produced by `muscat::aggregateData`, with different celltypes/clusters as assays.

- **norm**
  - Logical; whether to calculate logcpm (TMM normalization).

- **lfc_group**
  - The colData column to use to calculate foldchange. If NULL (default), no fold-change assay will be computed.

### Value

A SummarizedExperiment
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

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

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.
resetAllSEtoolsOptions

Description
Resets all global options relative to SEtools.

Examples

resetAllSEtoolsOptions()

Value
None

Description

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

Usage
se2xls(se, filename, addSheets = NULL)

Examples

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

Description

Deprecated pheatmap wrapper for SummarizedExperiment-class. **This function has been replaced by the sehm function from the 'sechm' package and is retained here solely for backward compatibility.**

Usage

sehm(
    se,
    genes,
    do.scale = FALSE,
    assayName = .getDef("assayName"),
    sortRowsOn = seq_len(ncol(se)),
    cluster_cols = FALSE,
    cluster_rows = is.null(sortRowsOn),
    toporder = NULL,
    hmcols = NULL,
    breaks = .getDef("breaks"),
    gaps_at = .getDef("gaps_at"),
    gaps_row = NULL,
    anno_rows = .getDef("anno_rows"),
    anno_columns = .getDef("anno_columns"),
    anno_colors = NULL,
    show_rownames = NULL,
    show_colnames = FALSE,
    ...
)
### Arguments

- **se**: A `SummarizedExperiment-class`.
- **genes**: An optional vector of genes (i.e. row names of `se`).
- **do.scale**: Logical; whether to scale rows (default FALSE).
- **assayName**: An optional vector of assayNames to use. The first available will be used, or the first assay if NULL.
- **sortRowsOn**: Sort rows by MDS polar order using the specified columns (default all)
- **cluster_cols**: Whether to cluster columns (default F)
- **cluster_rows**: Whether to cluster rows; default FALSE if `do.sortRows=TRUE`.
- **toporder**: Optional vector of categories on which to supra-order when sorting rows, or name of a `rowData` column to use for this purpose.
- **hmcols**: Colors for the heatmap.
- **breaks**: Breaks for the heatmap colors. Alternatively, symmetrical breaks can be generated automatically by setting ‘breaks’ to a numerical value between 0 and 1. The value is passed as the ‘split.prop’ argument to the `getBreaks` function, and indicates the proportion of the points to map to a linear scale, while the more extreme values will be plotted on a quantile scale. ‘breaks=FALSE‘ will disable symmetrical scale and quantile capping, while retaining automatic breaks. ‘breaks=1’ will produce a symmetrical scale without quantile capping.
- **gaps_at**: Columns of `colData` to use to establish gaps between columns.
- **gaps_row**: Passed to the heatmap function; if missing, will be set automatically according to toporder.
- **anno_rows**: Columns of `rowData` to use for left annotation.
- **anno_columns**: Columns of `colData` to use for top annotation.
- **anno_colors**: List of colors to use for annotation.
- **show_rownames**: Whether to show row names (default TRUE if less than 50 rows to plot).
- **show_colnames**: Whether to show column names (default FALSE).
- **...**: Further arguments passed to `pheatmap`

### Value

A heatmap.

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### svacor

#### 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,
  form$ = ~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
form$       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 ‘form$’.
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)
Index

aggSE, 2, 6
castSE, 3
data, 4
flattenPB, 4
getBreaks, 9
log2FC, 5
mergeSEs, 6
resetAllSEtoolsOptions, 7
SE (data), 4
se2xls, 7
sechm, 8
sehm, 8
sva, 10
svacor, 9