

Package ‘alabaster.sce’

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Title Load and Save SingleCellExperiment from File

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Description Save SingleCellExperiment into file artifacts, and load them back into memory.

This is a more portable alternative to serialization of such objects into RDS files.

Each artifact is associated with metadata for further interpretation;

downstream applications can enrich this metadata with context-specific properties.

Depends SingleCellExperiment, alabaster.base

Imports methods, alabaster.se

Suggests knitr, testthat, BiocStyle, rmarkdown

VignetteBuilder knitr

RoxygenNote 7.2.3

biocViews DataImport, DataRepresentation

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loadSingleCellExperiment

Load a SingleCellExperiment

Description

Load a [SingleCellExperiment](#) object from file, using the metadata generated by the corresponding [stageObject](#) method.

Usage

```
loadSingleCellExperiment(exp.info, project, ...)
```

Arguments

exp.info	Named list containing the metadata for this experiment.
project	Any argument accepted by the acquisition functions, see ?acquireFile . By default, this should be a string containing the path to a staging directory.
...	Further arguments to pass to loadSummarizedExperiment .

Value

A [SingleCellExperiment](#) object.

Author(s)

Aaron Lun

Examples

```
# Mocking up an SCE:
mat <- matrix(rpois(10000, 10), ncol=10)
colnames(mat) <- letters[1:10]
rownames(mat) <- sprintf("GENE_%i", seq_len(nrow(mat)))

se <- SingleCellExperiment(list(counts=mat))
se$stuff <- LETTERS[1:10]
se$blah <- runif(10)
reducedDims(se) <- list(
  PCA=matrix(rnorm(ncol(se)*10), ncol=10),
  TSNE=matrix(rnorm(ncol(se)*2), ncol=2)
)
altExps(se) <- list(spikes=SummarizedExperiment(list(counts=mat[1:2,])))

# Staging it:
tmp <- tempfile()
dir.create(tmp)
info <- stageObject(se, dir=tmp, "rna-seq")
```

```
# Loading it back into memory:
loadSingleCellExperiment(info, tmp)
```

```
stageObject, SingleCellExperiment-method
Stage an experiment
```

Description

Save a [SingleCellExperiment](#) to file inside the staging directory.

Usage

```
## S4 method for signature 'SingleCellExperiment'
stageObject(x, dir, path, child = FALSE, rd.name = "dimreds", ...)
```

Arguments

x	A SingleCellExperiment object or one of its subclasses.
dir	String containing the path to the staging directory.
path	String containing a prefix of the relative path inside dir where x is to be saved. The actual path used to save x may include additional components, see Details.
child	Logical scalar indicating whether x is a child of a larger object.
rd.name	String containing the prefix of the file to save the reduced dimensions.
...	Further arguments to pass to the RangedSummarizedExperiment method.

Value

A named list of metadata that follows the single_cell_experiment schema. The contents of x are saved into a path subdirectory inside dir.

Author(s)

Aaron Lun

Examples

```
# Mocking up an SCE:
mat <- matrix(rpois(10000, 10), ncol=10)
colnames(mat) <- letters[1:10]
rownames(mat) <- sprintf("GENE_%i", seq_len(nrow(mat)))

se <- SingleCellExperiment(list(counts=mat))
se$stuff <- LETTERS[1:10]
se$blah <- runif(10)
```

```
reducedDims(se) <- list(
  PCA=matrix(rnorm(ncol(se)*10), ncol=10),
  TSNE=matrix(rnorm(ncol(se)*2), ncol=2)
)
altExps(se) <- list(spikes=SummarizedExperiment(list(counts=mat[1:2,])))

# Staging it:
tmp <- tempfile()
dir.create(tmp)
stageObject(se, dir=tmp, "rna-seq")
list.files(file.path(tmp, "rna-seq"))
```

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