

Package ‘alabaster.mae’

April 15, 2024

Title Load and Save MultiAssayExperiments

Version 1.2.0

Date 2023-08-17

License MIT + file LICENSE

Description Save MultiAssayExperiments 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 MultiAssayExperiment, alabaster.base

Imports methods, alabaster.se, S4Vectors

Suggests testthat, knitr, SummarizedExperiment, BiocParallel,
BiocStyle, rmarkdown

VignetteBuilder knitr

RoxygenNote 7.2.3

biocViews DataImport, DataRepresentation

git_url <https://git.bioconductor.org/packages/alabaster.mae>

git_branch RELEASE_3_18

git_last_commit 4636dc2

git_last_commit_date 2023-10-24

Repository Bioconductor 3.18

Date/Publication 2024-04-15

Author Aaron Lun [aut, cre]

Maintainer Aaron Lun <infinite.monkeys.with.keyboards@gmail.com>

R topics documented:

| | |
|---|---|
| loadMultiAssayExperiment | 2 |
| stageObject,MultiAssayExperiment-method | 3 |

| | |
|--------------|----------|
| Index | 5 |
|--------------|----------|

`loadMultiAssayExperiment`*Load a MultiAssayExperiment*

Description

Load a dataset as a [MultiAssayExperiment](#), based on the metadata generated by the corresponding [stageObject](#) method.

Usage

```
loadMultiAssayExperiment(  
  ds.info,  
  project,  
  experiments = NULL,  
  BPPARAM = NULL,  
  include.nested = TRUE  
)
```

Arguments

| | |
|-----------------------------|--|
| <code>ds.info</code> | Named list containing the metadata for this object. |
| <code>project</code> | Any argument accepted by the acquisition functions, see ?acquireFile . By default, this should be a string containing the path to a staging directory. |
| <code>experiments</code> | Character or integer vector specifying the subset of experiments to load. If NULL, all experiments are loaded. |
| <code>BPPARAM</code> | A BiocParallelParam object indicating how loading should be parallelized across multiple experiments. If NULL, loading is done serially. |
| <code>include.nested</code> | Logical scalar indicating whether to include nested DataFrames in the <code>colData</code> of the output. |

Value

A [MultiAssayExperiment](#) object.

Author(s)

Aaron Lun

Examples

```
library(SummarizedExperiment)  
  
# Mocking up an MAE  
mat <- matrix(rnorm(1000), ncol=10)  
colnames(mat) <- letters[1:10]  
rownames(mat) <- sprintf("GENE_%i", seq_len(nrow(mat)))
```

```
se <- SummarizedExperiment(list(counts=mat))

library(MultiAssayExperiment)
mae <- MultiAssayExperiment(list(gene=se))

# Staging it:
tmp <- tempfile()
dir.create(tmp)
info <- stageObject(mae, tmp, "dataset")

# Loading it back in:
loadMultiAssayExperiment(info, tmp)
```

stageObject,MultiAssayExperiment-method
Stage a dataset

Description

Save the metadata and annotations of a [MultiAssayExperiment](#) in a staging directory.

Usage

```
## S4 method for signature 'MultiAssayExperiment'
stageObject(
  x,
  dir,
  path,
  child = FALSE,
  sm.name = "sample_mapping",
  sd.name = "sample_data",
  meta.name = "dataset.json"
)
```

Arguments

| | |
|-----------|--|
| x | A MultiAssayExperiment 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. |
| sm.name | String containing the prefix of the sample mapping file. |
| sd.name | String containing the prefix of the sample data file. |
| meta.name | String containing the name of the metadata file. |

Details

`meta.name` is only needed to set up the output path, for consistency with the `stageObject` contract. Callers should make sure to write the metadata to the same document by using `.writeMetadata` with `meta.only=TRUE`.

Value

A named list containing the metadata for this dataset. The contents of `x` are saved into a path subdirectory inside `dir`.

Author(s)

Aaron Lun

Examples

```
# Mocking up an MAE
mat <- matrix(rnorm(1000), ncol=10)
colnames(mat) <- letters[1:10]
rownames(mat) <- sprintf("GENE_%i", seq_len(nrow(mat)))

library(SummarizedExperiment)
se <- SummarizedExperiment(list(counts=mat))

library(MultiAssayExperiment)
mae <- MultiAssayExperiment(list(gene=se))

# Staging it:
tmp <- tempfile()
dir.create(tmp)
stageObject(mae, tmp, "dataset")
```

Index

`.writeMetadata`, [4](#)
`acquireFile`, [2](#)
`loadMultiAssayExperiment`, [2](#)
`MultiAssayExperiment`, [2](#), [3](#)
`stageObject`, [2](#), [4](#)
`stageObject`, `MultiAssayExperiment`-method,
[3](#)
`stageObject`, `SampleMapFrame`-method
(`stageObject`, `MultiAssayExperiment`-method),
[3](#)