Package 'gemma.R'

September 29, 2022

```
Description Low- and high-level wrappers for Gemma's RESTful API. They enable
     access to curated expression and differential expression data from over
     10,000 published studies. Gemma is a web site, database and a set of tools for the
     meta-analysis, re-use and sharing of genomics data, currently primarily
     targeted at the analysis of gene expression profiles.
URL https://pavlidislab.github.io/gemma.R/,
     https://github.com/PavlidisLab/gemma.R
License Apache License (>= 2)
Encoding UTF-8
Roxygen list(markdown = TRUE)
RoxygenNote 7.2.1
BugReports https://github.com/PavlidisLab/gemma.R/issues
Imports magrittr, glue, memoise, jsonlite, data.table, rlang,
     lubridate, utils, stringr, SummarizedExperiment, Biobase,
     tibble, tidyr, S4Vectors, httr, rappdirs, bit64
Suggests testthat (>= 2.0.0), rmarkdown, knitr, dplyr, covr, ggplot2,
     ggrepel, BiocStyle, microbenchmark, magick, purrr
Config/testthat/edition 2
VignetteBuilder knitr
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Maintainer Ogan Mancarci <ogan.mancarci@gmail.com>

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 ${\tt forget_gemma_memoised} \ \ {\it Clear gemma.R cache}$

Description

Forget past results from memoised calls to the Gemma API (ie. using functions with memoised = TRUE)

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Usage

```
forget_gemma_memoised()
```

Value

TRUE to indicate cache was cleared.

Examples

forget_gemma_memoised()

gemma.R

gemma.R package: Access curated gene expression data and differential expression analyses

Description

This package contains wrappers and convenience function for Gemma's RESTful API that enable access to curated expression and differential expression data from over 15,000 published studies (as of mid-2022). Gemma (https://gemma.msl.ubc.ca) is a web site, database and a set of tools for the meta-analysis, re-use and sharing of genomics data, currently primarily targeted at the analysis of gene expression profiles.

Details

The endpoints are classified as follows:

- Dataset endpoints: Access expression and differential expression data, and their associated properties and files.
- Platform endpoints: Access microarray or sequencing platforms and their associated properties and files.
- Gene endpoints: Access information about specific genes.

Most users will want to start with the high-level functions getDataset, getDatasetDE and getPlat-formAnnotations. Additional lower-level methods are available that directly map to the Gemma RESTful API methods.

For more information and detailed usage instructions check the README, the function reference and the vignette.

All software-related questions should be posted to the Bioconductor Support Site: https://support.bioconductor.org

Author(s)

Javier Castillo-Arnemann, Jordan Sicherman, Ogan Mancarci, Guillaume Poirier-Morency

References

Lim, N. et al., Curation of over 10 000 transcriptomic studies to enable data reuse, Database, 2021. https://doi.org/10.1093/database/baab006

gemma_call

Custom gemma call

Description

A minimal function to create custom calls. Can be used to acquire unimplemented endpoints and/or raw output without any processing. Refer to the API documentation.

Usage

```
gemma_call(call, ..., json = TRUE)
```

Arguments

call Gemma API endpoint.
... parameters included in the call
json If TRUE will parse the content as a list

Value

A list if json = TRUE and an httr response if FALSE

Examples

```
# get singular value decomposition for the dataset
gemma_call('datasets/{dataset}/svd',dataset = 1)
```

get_datasets_by_ids

Retrieve datasets by their identifiers

Description

Retrieve datasets by their identifiers

get_datasets_by_ids 5

Usage

```
get_datasets_by_ids(
  datasets = NA_character_,
  offset = 0L,
  limit = 20L,
  sort = "+id",
  raw = getOption("gemma.raw", FALSE),
  memoised = getOption("gemma.memoised", FALSE),
  file = getOption("gemma.file", NA_character_),
  overwrite = getOption("gemma.overwrite", FALSE),
  attributes = getOption("gemma.attributes", TRUE)
)
```

Arguments

| datasets | Numerical dataset identifiers or dataset short names |
|------------|---|
| offset | The offset of the first retrieved result. |
| limit | Optional, defaults to 20. Limits the result to specified amount of objects. Has a maximum value of 100. Use together with offset and the totalElements attribute in the output to compile all data if needed. |
| sort | Order results by the given property and direction. The '+' sign indicate ascending order whereas the '-' indicate descending. |
| raw | TRUE to receive results as-is from Gemma, or FALSE to enable parsing. Raw results usually contain additional fields and flags that are omitted in the parsed results. |
| memoised | Whether or not to save to cache for future calls with the same inputs and use the result saved in cache if a result is already saved. Doing options(gemma.memoised = TRUE) will ensure that the cache is always used. Use forget_gemma_memoised to clear the cache. |
| file | The name of a file to save the results to, or NULL to not write results to a file. If raw == TRUE, the output will be a JSON file. Otherwise, it will be a RDS file. |
| overwrite | Whether or not to overwrite if a file exists at the specified filename. |
| attributes | If TRUE additional information from the call will be added into the output object's attributes such as offset and available elements. |

Value

A data table with information about the queried dataset(s). A list if raw = TRUE. Returns an empty list if no datasets matched. A successful response may contain 'Geeq' information, which aims to provide a unified metric to measure experiments by the quality of their data, and their suitability for use in Gemma. You can read more about the geeq properties here.

The fields of the output data.table are:

- experiment. ShortName: Shortname given to the dataset within Gemma. Often corresponds to accession ID
- experiment.Name: Full title of the dataset

- experiment. ID: Internal ID of the dataset.
- experiment.Description: Description of the dataset
- experiment.Public: Is the dataset publicly available. Only useful for logged in users with access to non-public data
- experiment. Troubled: Did an automatic process within gemma or a curator mark the dataset as "troubled"
- experiment.Accession: Accession ID of the dataset in the external database it was taken from
- experiment. Database: The name of the database where the dataset was taken from
- experiment.URI: URI of the original database
- experiment. SampleCount: Number of samples in the dataset
- experiment.batchEffect: A text field describing whether the dataset has batch effects
- experiment.batchCorrected: Whether batch correction has been performed on the dataset. This is a text field with possible details about the batch correction. Use geeq.batchEffect if a more structured field is needed.
- geeq.batchConfound: 0 if batch info isn't available, -1 if batch counfoud is detected, 1 if batch information is available and no batch confound found
- geeq.batchEffect: -1 if batch p value < 0.0001, 1 if batch p value > 0.1, 0 if otherwise and when there is no batch information is available or when the data is confounded with batches.
- geeq.rawData: -1 if no raw data available, 1 if raw data was available. When available, Gemma reprocesses raw data to get expression values and batches
- geeq.qScore: Data quality score given to the dataset by Gemma.
- geeq.sScore: Suitability score given to the dataset by Gemma. Refers to factors like batches, platforms and other aspects of experimental design
- taxon.Name: The taxa of the study. In Gemma each study will include a single species. If
 the original source has samples from multiple species, they will be split into different studies
 within Gemma
- taxon. ID: Internal ID given to the taxon by Gemma

Examples

```
get_datasets_by_ids("GSE2018")
get_datasets_by_ids(c("GSE2018", "GSE2872"))
```

get_dataset_annotations

Retrieve the annotations analysis of a dataset

Description

Retrieve the annotations analysis of a dataset

get_dataset_annotations

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Usage

```
get_dataset_annotations(
  dataset,
  raw = getOption("gemma.raw", FALSE),
  memoised = getOption("gemma.memoised", FALSE),
  file = getOption("gemma.file", NA_character_),
  overwrite = getOption("gemma.overwrite", FALSE),
  attributes = getOption("gemma.attributes", TRUE)
)
```

Arguments

| dataset | A numerical dataset identifier or a dataset short name |
|------------|---|
| raw | TRUE to receive results as-is from Gemma, or FALSE to enable parsing. Raw results usually contain additional fields and flags that are omitted in the parsed results. |
| memoised | Whether or not to save to cache for future calls with the same inputs and use the result saved in cache if a result is already saved. Doing options(gemma.memoised = TRUE) will ensure that the cache is always used. Use forget_gemma_memoised to clear the cache. |
| file | The name of a file to save the results to, or NULL to not write results to a file. If raw == TRUE, the output will be a JSON file. Otherwise, it will be a RDS file. |
| overwrite | Whether or not to overwrite if a file exists at the specified filename. |
| attributes | If TRUE additional information from the call will be added into the output object's attributes such as offset and available elements. |

Value

A data table with information about the annotations of the queried dataset. A list if raw = TRUE.A 404 error if the given identifier does not map to any object.

The fields of the output data.table are:

- class. Type: Type of the annotation class
- class. Name: Name of the annotation class (e.g. organism part)
- class. URI: URI for the annotation class
- term.Name: Name of the annotation term (e.g. lung)
- term. URI: URI for the annotation term

Examples

```
get_dataset_annotations("GSE2018")
```

8 get_dataset_design

Description

Retrieve the design of a dataset

Usage

```
get_dataset_design(
  dataset,
  raw = getOption("gemma.raw", FALSE),
  memoised = getOption("gemma.memoised", FALSE),
  file = getOption("gemma.file", NA_character_),
  overwrite = getOption("gemma.overwrite", FALSE),
  attributes = getOption("gemma.attributes", TRUE)
)
```

Arguments

| dataset | A numerical dataset identifier or a dataset short name |
|------------|---|
| raw | TRUE to receive results as-is from Gemma, or FALSE to enable parsing. Raw results usually contain additional fields and flags that are omitted in the parsed results. |
| memoised | Whether or not to save to cache for future calls with the same inputs and use the result saved in cache if a result is already saved. Doing options(gemma.memoised = TRUE) will ensure that the cache is always used. Use forget_gemma_memoised to clear the cache. |
| file | The name of a file to save the results to, or NULL to not write results to a file. If raw == TRUE, the output will be a JSON file. Otherwise, it will be a RDS file. |
| overwrite | Whether or not to overwrite if a file exists at the specified filename. |
| attributes | If TRUE additional information from the call will be added into the output object's attributes such as offset and available elements. |

Value

A data table of the design matrix for the queried dataset. A 404 error if the given identifier does not map to any object

Examples

```
head(get_dataset_design("GSE2018"))
```

Description

Retrieve the differential analyses of a dataset

Usage

```
get_dataset_differential_expression_analyses(
  dataset,
  raw = getOption("gemma.raw", FALSE),
  memoised = getOption("gemma.memoised", FALSE),
  file = getOption("gemma.file", NA_character_),
  overwrite = getOption("gemma.overwrite", FALSE),
  attributes = getOption("gemma.attributes", TRUE)
)
```

Arguments

| dataset | A numerical dataset identifier or a dataset short name |
|------------|---|
| raw | TRUE to receive results as-is from Gemma, or FALSE to enable parsing. Raw results usually contain additional fields and flags that are omitted in the parsed results. |
| memoised | Whether or not to save to cache for future calls with the same inputs and use the result saved in cache if a result is already saved. Doing options(gemma.memoised = TRUE) will ensure that the cache is always used. Use forget_gemma_memoised to clear the cache. |
| file | The name of a file to save the results to, or NULL to not write results to a file. If raw == TRUE, the output will be a JSON file. Otherwise, it will be a RDS file. |
| overwrite | Whether or not to overwrite if a file exists at the specified filename. |
| attributes | If TRUE additional information from the call will be added into the output object's attributes such as offset and available elements. |

Value

A data table with information about the differential expression analysis of the queried dataset. Note that this funciton does not return differential expression values themselves. Use get_differential_expression_values to get differential expression values (see examples).

The fields of the output data.table are:

• result.ID: Result set ID of the differential expression analysis. May represent multiple factors in a single model.

- contrast.ID: Id of the specific contrast factor. Together with the result.ID they uniquely represent a given contrast.
- experiment. ID: Id of the source experiment
- baseline.category: Category for the contrast
- baseline.categoryURI: URI for the baseline category
- baseline.factorValue: Factor value assigned as the baseline in the contrast. Typically represent control samples
- baseline.factorValueURI: URI for the baseline.factorValue
- experimental.factorValue: Factor value assigned to the experimental group.
- experimental.factorValueURI: URI for the experimental.factorValue
- subsetFactor.subset: TRUE if the result set belong to a subset, FALSE if not. Subsets are created when performing differential expression to avoid unhelpful comparisons.
- subsetFactor.category: Category of the subset
- subsetFactor.categoryURI: URI of the subset category
- subsetFactor.factorValue: Factor Value of the subset
- subsetFactor.factorValueURI: URI of the subset factor value
- probes . Analyzed: Number of probesets represented in the contrast
- genes. Analyzed: Number of genes represented in the contrast
- platform. ID: Platform id for the contrast

Examples

```
result <- get_dataset_differential_expression_analyses("GSE2018")
get_differential_expression_values(resultSet = result$result.ID)</pre>
```

```
get_dataset_expression
```

Retrieve the expression data of a dataset

Description

Retrieve the expression data of a dataset

Usage

```
get_dataset_expression(
  dataset,
  filter = FALSE,
  raw = getOption("gemma.raw", FALSE),
  memoised = getOption("gemma.memoised", FALSE),
  file = getOption("gemma.file", NA_character_),
  overwrite = getOption("gemma.overwrite", FALSE),
  attributes = getOption("gemma.attributes", TRUE)
)
```

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Arguments

| dataset | A numerical dataset identifier or a dataset short name |
|------------|---|
| filter | The filtered version (filter = TRUE) corresponds to what is used in most Gemma analyses, removing some probes/elements. Unfiltered includes all elements. |
| raw | TRUE to receive results as-is from Gemma, or FALSE to enable parsing. Raw results usually contain additional fields and flags that are omitted in the parsed results. |
| memoised | Whether or not to save to cache for future calls with the same inputs and use the result saved in cache if a result is already saved. Doing options(gemma.memoised = TRUE) will ensure that the cache is always used. Use forget_gemma_memoised to clear the cache. |
| file | The name of a file to save the results to, or NULL to not write results to a file. If raw == TRUE, the output will be a JSON file. Otherwise, it will be a RDS file. |
| overwrite | Whether or not to overwrite if a file exists at the specified filename. |
| attributes | If TRUE additional information from the call will be added into the output object's attributes such as offset and available elements. |

Value

If raw is FALSE (default), a data table of the expression matrix for the queried dataset. If raw is TRUE, returns the binary file in raw form.

Examples

Description

Return an annotated Bioconductor-compatible data structure or a long form tibble of the queried dataset, including expression data and the experimental design.

Usage

```
get_dataset_object(
  dataset,
  filter = FALSE,
  type = "se",
  memoised = getOption("gemma.memoised", FALSE)
)
```

Arguments

dataset A dataset identifier.

filter The filtered version corresponds to what is used in most Gemma analyses, re-

moving some probes/elements. Unfiltered includes all elements.

type "se"for a SummarizedExperiment or "eset" for Expression Set. We recommend

using SummarizedExperiments which are more recent. See the Summarized

experiment vignette or the ExpressionSet vignette for more details.

memoised Whether or not to save to cache for future calls with the same inputs and use the

result saved in cache if a result is already saved. Doing options(gemma.memoised = TRUE) will ensure that the cache is always used. Use forget_gemma_memoised

to clear the cache.

Value

A SummarizedExperiment, ExpressionSet or tibble containing metadata and expression data for the queried dataset.

Examples

```
get_dataset_object("GSE2018")
```

get_dataset_platforms Retrieve the platform of a dataset

Description

Retrieve the platform of a dataset

Usage

```
get_dataset_platforms(
  dataset,
  raw = getOption("gemma.raw", FALSE),
  memoised = getOption("gemma.memoised", FALSE),
  file = getOption("gemma.file", NA_character_),
  overwrite = getOption("gemma.overwrite", FALSE),
  attributes = getOption("gemma.attributes", TRUE)
)
```

Arguments

dataset A numerical dataset identifier or a dataset short name

raw TRUE to receive results as-is from Gemma, or FALSE to enable parsing. Raw

results usually contain additional fields and flags that are omitted in the parsed

results.

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| memoised | Whether or not to save to cache for future calls with the same inputs and use the result saved in cache if a result is already saved. Doing options(gemma.memoised = TRUE) will ensure that the cache is always used. Use forget_gemma_memoised to clear the cache. |
|------------|---|
| file | The name of a file to save the results to, or NULL to not write results to a file. If raw == TRUE, the output will be a JSON file. Otherwise, it will be a RDS file. |
| overwrite | Whether or not to overwrite if a file exists at the specified filename. |
| attributes | If TRUE additional information from the call will be added into the output object's attributes such as offset and available elements. |

Value

A data table with information about the platform(s). A list if raw = TRUE. A 404 error if the given identifier does not map to any object

The fields of the output data.table are:

- platform. ID: Internal identifier of the platform
- platform. ShortName: Shortname of the platform.
- platform. Name: Full name of the platform.
- platform. Description: Free text description of the platform
- platform. Troubled: Whether or not the platform was marked "troubled" by a Gemma process or a curator
- platform. ExperimentCount: Number of experiments using the platform within Gemma
- taxon. Name: Name of the species platform was made for
- taxon. ID: Internal identifier given to the species by Gemma
- platform. Type: Technology type for the platform.

Examples

```
get_dataset_platforms("GSE2018")
```

get_dataset_samples

Retrieve the samples of a dataset

Description

Retrieve the samples of a dataset

Usage

```
get_dataset_samples(
  dataset,
  raw = getOption("gemma.raw", FALSE),
  memoised = getOption("gemma.memoised", FALSE),
  file = getOption("gemma.file", NA_character_),
  overwrite = getOption("gemma.overwrite", FALSE),
  attributes = getOption("gemma.attributes", TRUE)
)
```

Arguments

| dataset | A numerical dataset identifier or a dataset short name |
|------------|---|
| raw | TRUE to receive results as-is from Gemma, or FALSE to enable parsing. Raw results usually contain additional fields and flags that are omitted in the parsed results. |
| memoised | Whether or not to save to cache for future calls with the same inputs and use the result saved in cache if a result is already saved. Doing options(gemma.memoised = TRUE) will ensure that the cache is always used. Use forget_gemma_memoised to clear the cache. |
| file | The name of a file to save the results to, or NULL to not write results to a file. If raw == TRUE, the output will be a JSON file. Otherwise, it will be a RDS file. |
| overwrite | Whether or not to overwrite if a file exists at the specified filename. |
| attributes | If TRUE additional information from the call will be added into the output object's attributes such as offset and available elements. |

Value

A data table with information about the samples of the queried dataset. A list if raw = TRUE. A 404 error if the given identifier does not map to any object.

The fields of the output data.table are:

- sample. Name: Internal name given to the sample.
- sample. ID: Internal ID of the sample
- sample.Description: Free text description of the sample
- sample.Outlier: Whether or not the sample is marked as an outlier
- sample. Accession: Accession ID of the sample in it's original database
- sample.Database: Database of origin for the sample
- sample. Characteristics: Characteristics of the sample. This field is a data table
- sample. Factor Values: Experimental factor values of the sample. This field is a data table

Examples

```
head(get_dataset_samples("GSE2018"))
```

Description

Retrieves the differential expression result set(s) associated with the dataset. To get more information about the contrasts in individual resultSets and annotation terms associated them, use get_dataset_differential_expression_analyses()

Usage

```
get_differential_expression_values(
  dataset = NA_character_,
  resultSet = NA_integer_,
  readableContrasts = FALSE,
  memoised = getOption("gemma.memoised", FALSE)
```

Arguments

dataset A dataset identifier.

resultSet A resultSet identifier.

readableContrasts

If FALSE (default), the returned columns will use internal constrasts IDs as names. Details about the contrasts can be accessed using get_dataset_differential_expression_ana

If TRUE IDs will be replaced with human readable contrast information.

memoised

Whether or not to save to cache for future calls with the same inputs and use the result saved in cache if a result is already saved. Doing options(gemma.memoised = TRUE) will ensure that the cache is always used. Use forget_gemma_memoised

to clear the cache.

Details

In Gemma each result set corresponds to the estimated effects associated with a single factor in the design, and each can have multiple contrasts (for each level compared to baseline). Thus a dataset with a 2x3 factorial design will have two result sets, one of which will have one contrast, and one having two contrasts.

The methodology for differential expression is explained in Curation of over 10000 transcriptomic studies to enable data reuse. Briefly, differential expression analysis is performed on the dataset based on the annotated experimental design with up two three potentially nested factors. Gemma attempts to automatically assign baseline conditions for each factor. In the absence of a clear control condition, a baseline is arbitrarily selected. A generalized linear model with empirical Bayes shrinkage of t-statistics is fit to the data for each platform element (probe/gene) using an implementation of the limma algorithm. For RNA-seq data, we use weighted regression, applying the voom algorithm to compute weights from the mean–variance relationship of the data. Contrasts of each

get_genes

condition are then computed compared to the selected baseline. In some situations, Gemma will split the data into subsets for analysis. A typical such situation is when a 'batch' factor is present and confounded with another factor, the subsets being determined by the levels of the confounding factor.

Value

A list of data tables with differential expression values per result set.

Examples

```
get_differential_expression_values("GSE2018")
```

get_genes

Retrieve genes matching a gene identifier

Description

Retrieve genes matching a gene identifier

Usage

```
get_genes(
   genes,
   raw = getOption("gemma.raw", FALSE),
   memoised = getOption("gemma.memoised", FALSE),
   file = getOption("gemma.file", NA_character_),
   overwrite = getOption("gemma.overwrite", FALSE),
   attributes = getOption("gemma.attributes", TRUE)
)
```

Arguments

| genes | An ensembl gene identifier which typically starts with ensg or an ncbi gene identifier or an official gene symbol approved by hgnc |
|------------|---|
| raw | TRUE to receive results as-is from Gemma, or FALSE to enable parsing. Raw results usually contain additional fields and flags that are omitted in the parsed results. |
| memoised | Whether or not to save to cache for future calls with the same inputs and use the result saved in cache if a result is already saved. Doing options(gemma.memoised = TRUE) will ensure that the cache is always used. Use forget_gemma_memoised to clear the cache. |
| file | The name of a file to save the results to, or NULL to not write results to a file. If raw == TRUE, the output will be a JSON file. Otherwise, it will be a RDS file. |
| overwrite | Whether or not to overwrite if a file exists at the specified filename. |
| attributes | If TRUE additional information from the call will be added into the output object's attributes such as offset and available elements. |

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Value

A data table with information about the querried gene(s) A list if raw = TRUE.

The fields of the output data.table are:

- gene. Symbol: Symbol for the gene
- gene. Ensembl: Ensembl ID for the gene
- gene. NCBI: NCBI id for the gene
- gene. Name: Name of the gene
- taxon.name: Name of the taxon of origin
- taxon. ID: Gemma ID for the taxon
- taxon. Scientific: Scientific name for the taxon

Examples

```
get_genes("DYRK1A")
get_genes(c("DYRK1A", "PTEN"))
```

get_gene_go_terms

Retrieve the GO terms associated to a gene

Description

Retrieve the GO terms associated to a gene

Usage

```
get_gene_go_terms(
  gene,
  raw = getOption("gemma.raw", FALSE),
 memoised = getOption("gemma.memoised", FALSE),
  file = getOption("gemma.file", NA_character_),
  overwrite = getOption("gemma.overwrite", FALSE),
  attributes = getOption("gemma.attributes", TRUE)
)
```

Arguments

gene

An ensembl gene identifier which typically starts with ensg or an ncbi gene identifier or an official gene symbol approved by hgnc

raw

TRUE to receive results as-is from Gemma, or FALSE to enable parsing. Raw results usually contain additional fields and flags that are omitted in the parsed

results.

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| memoised | Whether or not to save to cache for future calls with the same inputs and use the result saved in cache if a result is already saved. Doing options(gemma.memoised = TRUE) will ensure that the cache is always used. Use forget_gemma_memoised to clear the cache. |
|------------|---|
| file | The name of a file to save the results to, or NULL to not write results to a file. If raw == TRUE, the output will be a JSON file. Otherwise, it will be a RDS file. |
| overwrite | Whether or not to overwrite if a file exists at the specified filename. |
| attributes | If TRUE additional information from the call will be added into the output object's attributes such as offset and available elements. |

Value

A data table with information about the GO terms assigned to the queried gene. A list if raw = TRUE. A 404 error if the given identifier does not map to any object. Go terms were updated on June 10 2022.

The fields of the output data.table are:

```
term.Name: Name of the term
term.ID: ID of the term
term.URI: URI of the term
```

Examples

```
get_gene_go_terms("DYRK1A")
```

Description

Retrieve the physical locations of a given gene

Usage

```
get_gene_locations(
   gene,
   raw = getOption("gemma.raw", FALSE),
   memoised = getOption("gemma.memoised", FALSE),
   file = getOption("gemma.file", NA_character_),
   overwrite = getOption("gemma.overwrite", FALSE),
   attributes = getOption("gemma.attributes", TRUE)
)
```

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Arguments

| gene | An ensembl gene identifier which typically starts with ensg or an ncbi gene identifier or an official gene symbol approved by hgnc |
|------------|---|
| raw | TRUE to receive results as-is from Gemma, or FALSE to enable parsing. Raw results usually contain additional fields and flags that are omitted in the parsed results. |
| memoised | Whether or not to save to cache for future calls with the same inputs and use the result saved in cache if a result is already saved. Doing options(gemma.memoised = TRUE) will ensure that the cache is always used. Use forget_gemma_memoised to clear the cache. |
| file | The name of a file to save the results to, or NULL to not write results to a file. If raw == TRUE, the output will be a JSON file. Otherwise, it will be a RDS file. |
| overwrite | Whether or not to overwrite if a file exists at the specified filename. |
| attributes | If TRUE additional information from the call will be added into the output object's attributes such as offset and available elements. |

Value

A data table with information about the physical location of the queried gene. A list if raw = TRUE. A 404 error if the given identifier does not map to any object.

The fields of the output data.table are:

- chromosome: Name of the chromosome the gene is located
- strand: Which strand the gene is located
- nucleotide: Nucleotide number for the gene
- length: Gene length
- taxon.name: Name of the taxon
- taxon. Scientific: Scientific name for the taxon
- taxon. ID: Internal ID for the taxon given by Gemma
- taxon.NCBI: NCBI ID for the taxon
- taxon.Database.Name: Name of the database used in Gemma for the taxon

Examples

```
get_gene_locations("DYRK1A")
```

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get_gene_probes Retrieve the probes associated to a genes

Description

Retrieve the probes associated to a genes

Usage

```
get_gene_probes(
   gene,
   offset = 0L,
   limit = 20L,
   raw = getOption("gemma.raw", FALSE),
   memoised = getOption("gemma.memoised", FALSE),
   file = getOption("gemma.file", NA_character_),
   overwrite = getOption("gemma.overwrite", FALSE),
   attributes = getOption("gemma.attributes", TRUE)
)
```

Arguments

| gene | An ensembl gene identifier which typically starts with ensg or an ncbi gene identifier or an official gene symbol approved by hgnc |
|------------|---|
| offset | The offset of the first retrieved result. |
| limit | Optional, defaults to 20. Limits the result to specified amount of objects. Has a maximum value of 100. Use together with offset and the totalElements attribute in the output to compile all data if needed. |
| raw | TRUE to receive results as-is from Gemma, or FALSE to enable parsing. Raw results usually contain additional fields and flags that are omitted in the parsed results. |
| memoised | Whether or not to save to cache for future calls with the same inputs and use the result saved in cache if a result is already saved. Doing options(gemma.memoised = TRUE) will ensure that the cache is always used. Use forget_gemma_memoised to clear the cache. |
| file | The name of a file to save the results to, or NULL to not write results to a file. If raw == TRUE, the output will be a JSON file. Otherwise, it will be a RDS file. |
| overwrite | Whether or not to overwrite if a file exists at the specified filename. |
| attributes | If TRUE additional information from the call will be added into the output object's attributes such as offset and available elements. |

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Value

A data table with information about the probes representing a gene across all platrofms. A list if raw = TRUE. A 404 error if the given identifier does not map to any genes.

The fields of the output data.table are:

- mapping.name: Name of the mapping. Typically the probeset name
- mappping.description: A free text field providing optional information about the mapping
- platform. ShortName: Shortname of the platform given by Gemma. Typically the GPL identifier.
- platform. Name: Full name of the platform
- platform. ID: Id number of the platform given by Gemma
- platform. Taxon: Species the platform was designed for
- platform. TaxonID: Id number of the species given by Gemma
- platform. Type: Type of the platform.
- platform. Description: Free text field describing the platform.
- platform. Troubled: Whether the platform is marked as troubled by a Gemma curator.

Examples

```
get_gene_probes("DYRK1A")
```

get_platforms_by_ids Retrieve all platforms matching a set of platform identifiers

Description

Retrieve all platforms matching a set of platform identifiers

Usage

```
get_platforms_by_ids(
  platforms = NA_character_,
  offset = 0L,
  limit = 20L,
  sort = "+id",
  raw = getOption("gemma.raw", FALSE),
  memoised = getOption("gemma.memoised", FALSE),
  file = getOption("gemma.file", NA_character_),
  overwrite = getOption("gemma.overwrite", FALSE),
  attributes = getOption("gemma.attributes", TRUE)
)
```

Arguments

| platforms | Platform numerical identifiers or platform short names |
|------------|---|
| offset | The offset of the first retrieved result. |
| limit | Optional, defaults to 20. Limits the result to specified amount of objects. Has a maximum value of 100. Use together with offset and the totalElements attribute in the output to compile all data if needed. |
| sort | Order results by the given property and direction. The '+' sign indicate ascending order whereas the '-' indicate descending. |
| raw | TRUE to receive results as-is from Gemma, or FALSE to enable parsing. Raw results usually contain additional fields and flags that are omitted in the parsed results. |
| memoised | Whether or not to save to cache for future calls with the same inputs and use the result saved in cache if a result is already saved. Doing options(gemma.memoised = TRUE) will ensure that the cache is always used. Use forget_gemma_memoised to clear the cache. |
| file | The name of a file to save the results to, or NULL to not write results to a file. If raw == TRUE, the output will be a JSON file. Otherwise, it will be a RDS file. |
| overwrite | Whether or not to overwrite if a file exists at the specified filename. |
| attributes | If TRUE additional information from the call will be added into the output object's attributes such as offset and available elements. |

Value

A data table with information about the platform(s). A list if raw = TRUE. A 404 error if the given identifier does not map to any object

The fields of the output data.table are:

- $\bullet\,$ platform. ID: Internal identifier of the platform
- platform. ShortName: Shortname of the platform.
- platform. Name: Full name of the platform.
- platform. Description: Free text description of the platform
- platform. Troubled: Whether or not the platform was marked "troubled" by a Gemma process or a curator
- platform. ExperimentCount: Number of experiments using the platform within Gemma
- taxon.Name: Name of the species platform was made for
- taxon. ID: Internal identifier given to the species by Gemma
- platform. Type: Technology type for the platform.

Examples

```
get_platforms_by_ids("GPL1355")
get_platforms_by_ids(c("GPL1355", "GPL96"))
```

```
get_platform_annotations
```

Retrieve Platform Annotations by Gemma

Description

Gets Gemma's platform annotations including mappings of microarray probes to genes.

Usage

```
get_platform_annotations(
  platform,
  annotType = c("noParents", "allParents", "bioProcess"),
  file = getOption("gemma.file", NA_character_),
  overwrite = getOption("gemma.overwrite", FALSE),
  memoised = getOption("gemma.memoise", FALSE),
  unzip = FALSE
)
```

Arguments

| platform | A platform identifier @seealso getPlatforms |
|-----------|---|
| annotType | Which GO terms should the output include |
| file | Where to save the annotation file to, or empty to just load into memory |
| overwrite | Whether or not to overwrite an existing file |
| memoised | Whether or not to save to cache for future calls with the same inputs and use the result saved in cache if a result is already saved. Doing options(gemma.memoised = TRUE) will ensure that the cache is always used. Use forget_gemma_memoised to clear the cache. |
| unzip | Whether or not to unzip the file (if @param file is not empty) |

Value

A table of annotations

- ProbeName: Probeset names provided by the platform. Gene symbols for generic annotations
- GeneSymbols: Genes that were found to be aligned to the probe sequence. Note that it is possible for probes to be non-specific. Alignment to multiple genes are indicated with gene symbols separated by "I"s
- GeneNames: Name of the gene
- GOTerms: GO Terms associated with the genes. annotType argument can be used to choose which terms should be included.
- GemmaIDs and NCBIids: respective IDs for the genes.

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Examples

```
head(get_platform_annotations("GPL96"))
head(get_platform_annotations('Generic_human'))
```

get_platform_datasets Retrieve all experiments within a given platform

Description

Retrieve all experiments within a given platform

Usage

```
get_platform_datasets(
  platform,
  offset = 0L,
  limit = 20L,
  raw = getOption("gemma.raw", FALSE),
  memoised = getOption("gemma.memoised", FALSE),
  file = getOption("gemma.file", NA_character_),
  overwrite = getOption("gemma.overwrite", FALSE),
  attributes = getOption("gemma.attributes", TRUE)
)
```

Arguments

| platform | A platform numerical identifier or a platform short name |
|------------|---|
| offset | The offset of the first retrieved result. |
| limit | Optional, defaults to 20. Limits the result to specified amount of objects. Has a maximum value of 100. Use together with offset and the totalElements attribute in the output to compile all data if needed. |
| raw | TRUE to receive results as-is from Gemma, or FALSE to enable parsing. Raw results usually contain additional fields and flags that are omitted in the parsed results. |
| memoised | Whether or not to save to cache for future calls with the same inputs and use the result saved in cache if a result is already saved. Doing options(gemma.memoised = TRUE) will ensure that the cache is always used. Use forget_gemma_memoised to clear the cache. |
| file | The name of a file to save the results to, or NULL to not write results to a file. If raw == TRUE, the output will be a JSON file. Otherwise, it will be a RDS file. |
| overwrite | Whether or not to overwrite if a file exists at the specified filename. |
| attributes | If TRUE additional information from the call will be added into the output object's attributes such as offset and available elements. |

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Value

A data table with information about the queried dataset(s). A list if raw = TRUE. Returns an empty list if no datasets matched. A successful response may contain 'Geeq' information, which aims to provide a unified metric to measure experiments by the quality of their data, and their suitability for use in Gemma. You can read more about the geeq properties here.

The fields of the output data.table are:

- experiment. ShortName: Shortname given to the dataset within Gemma. Often corresponds to accession ID
- experiment.Name: Full title of the dataset
- experiment. ID: Internal ID of the dataset.
- experiment.Description: Description of the dataset
- experiment.Public: Is the dataset publicly available. Only useful for logged in users with access to non-public data
- experiment. Troubled: Did an automatic process within gemma or a curator mark the dataset as "troubled"
- experiment.Accession: Accession ID of the dataset in the external database it was taken from
- experiment.Database: The name of the database where the dataset was taken from
- experiment. URI: URI of the original database
- experiment.SampleCount: Number of samples in the dataset
- experiment.batchEffect: A text field describing whether the dataset has batch effects
- experiment.batchCorrected: Whether batch correction has been performed on the dataset. This is a text field with possible details about the batch correction. Use geeq.batchEffect if a more structured field is needed.
- geeq.batchConfound: 0 if batch info isn't available, -1 if batch counfoud is detected, 1 if batch information is available and no batch confound found
- geeq.batchEffect: -1 if batch p value < 0.0001, 1 if batch p value > 0.1, 0 if otherwise and when there is no batch information is available or when the data is confounded with batches.
- geeq.rawData: -1 if no raw data available, 1 if raw data was available. When available, Gemma reprocesses raw data to get expression values and batches
- geeq.qScore: Data quality score given to the dataset by Gemma.
- geeq.sScore: Suitability score given to the dataset by Gemma. Refers to factors like batches, platforms and other aspects of experimental design
- taxon.Name: The taxa of the study. In Gemma each study will include a single species. If the original source has samples from multiple species, they will be split into different studies within Gemma
- taxon. ID: Internal ID given to the taxon by Gemma

Examples

head(get_platform_datasets("GPL1355"))

```
get_platform_element_genes
```

Retrieve the genes associated to a probe in a given platform

Description

Retrieve the genes associated to a probe in a given platform

Usage

```
get_platform_element_genes(
  platform,
  probe,
  offset = 0L,
  limit = 20L,
  raw = getOption("gemma.raw", FALSE),
  memoised = getOption("gemma.memoised", FALSE),
  file = getOption("gemma.file", NA_character_),
  overwrite = getOption("gemma.overwrite", FALSE),
  attributes = getOption("gemma.attributes", TRUE)
)
```

Arguments

| platform | A platform numerical identifier or a platform short name |
|------------|---|
| probe | A probe name or it's numerical identifier |
| offset | The offset of the first retrieved result. |
| limit | Optional, defaults to 20. Limits the result to specified amount of objects. Has a maximum value of 100. Use together with offset and the totalElements attribute in the output to compile all data if needed. |
| raw | TRUE to receive results as-is from Gemma, or FALSE to enable parsing. Raw results usually contain additional fields and flags that are omitted in the parsed results. |
| memoised | Whether or not to save to cache for future calls with the same inputs and use the result saved in cache if a result is already saved. Doing options(gemma.memoised = TRUE) will ensure that the cache is always used. Use forget_gemma_memoised to clear the cache. |
| file | The name of a file to save the results to, or NULL to not write results to a file. If raw == TRUE, the output will be a JSON file. Otherwise, it will be a RDS file. |
| overwrite | Whether or not to overwrite if a file exists at the specified filename. |
| attributes | If TRUE additional information from the call will be added into the output object's attributes such as offset and available elements. |

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Value

A data table with information about the querried gene(s) A list if raw = TRUE.

The fields of the output data.table are:

• gene.Symbol: Symbol for the gene

• gene.Ensembl: Ensembl ID for the gene

• gene. NCBI: NCBI id for the gene

• gene.Name: Name of the gene

• taxon.name: Name of the taxon of origin

• taxon. ID: Gemma ID for the taxon

• taxon. Scientific: Scientific name for the taxon

Examples

```
get_platform_element_genes("GPL1355", "AFFX_Rat_beta-actin_M_at")
```

get_taxa

Get taxa

Description

Returns taxa and their versions used in Gemma

Usage

```
get_taxa(memoised = getOption("gemma.memoised", FALSE))
```

Arguments

memoised

Whether or not to save to cache for future calls with the same inputs and use the result saved in cache if a result is already saved. Doing options(gemma.memoised = TRUE) will ensure that the cache is always used. Use forget_gemma_memoised to clear the cache.

Value

A data frame including the names, IDs and database information about the taxons

Examples

```
get_taxa()
```

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get_taxon_datasets

Retrieve the datasets for a given taxon

Description

Retrieve the datasets for a given taxon

Usage

```
get_taxon_datasets(
  taxon,
  offset = 0L,
  limit = 20,
  sort = "+id",
  raw = getOption("gemma.raw", FALSE),
  memoised = getOption("gemma.memoised", FALSE),
  file = getOption("gemma.file", NA_character_),
  overwrite = getOption("gemma.overwrite", FALSE),
  attributes = getOption("gemma.attributes", TRUE)
)
```

Arguments

taxon

can either be Taxon ID, Taxon NCBI ID, or one of its string identifiers: scientific name, common name. It is recommended to use Taxon ID for efficiency. Please note, that not all taxa have all the possible identifiers available. Use the get_taxa_by_ids function to retrieve the necessary information. For convenience, below is a list of officially supported taxa:

| ID | Comm.name | Scient.name | NcbiID |
|----|-----------|--------------------------|--------|
| 1 | human | Homo sapiens | 9606 |
| 2 | mouse | Mus musculus | 10090 |
| 3 | rat | Rattus norvegicus | 10116 |
| 11 | yeast | Saccharomyces cerevisiae | 4932 |
| 12 | zebrafish | Danio rerio | 7955 |
| 13 | fly | Drosophila melanogaster | 7227 |
| 14 | worm | Caenorhabditis elegans | 6239 |

offset

The offset of the first retrieved result.

limit

Optional, defaults to 20. Limits the result to specified amount of objects. Has a maximum value of 100. Use together with offset and the totalElements attribute in the output to compile all data if needed.

sort

Order results by the given property and direction. The '+' sign indicate ascending order whereas the '-' indicate descending.

raw

TRUE to receive results as-is from Gemma, or FALSE to enable parsing. Raw results usually contain additional fields and flags that are omitted in the parsed

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results.

memoised Whether or not to save to cache for future calls with the same inputs and use the

result saved in cache if a result is already saved. Doing options(gemma.memoised = TRUE) will ensure that the cache is always used. Use forget_gemma_memoised

to clear the cache.

file The name of a file to save the results to, or NULL to not write results to a file. If

raw == TRUE, the output will be a JSON file. Otherwise, it will be a RDS file.

overwrite Whether or not to overwrite if a file exists at the specified filename.

attributes If TRUE additional information from the call will be added into the output ob-

ject's attributes such as offset and available elements.

Value

A data table with information about the queried dataset(s). A list if raw = TRUE. Returns an empty list if no datasets matched. A successful response may contain 'Geeq' information, which aims to provide a unified metric to measure experiments by the quality of their data, and their suitability for use in Gemma. You can read more about the geeq properties here.

The fields of the output data.table are:

- experiment.ShortName: Shortname given to the dataset within Gemma. Often corresponds to accession ID
- experiment.Name: Full title of the dataset
- experiment. ID: Internal ID of the dataset.
- experiment.Description: Description of the dataset
- experiment.Public: Is the dataset publicly available. Only useful for logged in users with access to non-public data
- experiment.Troubled: Did an automatic process within gemma or a curator mark the dataset as "troubled"
- experiment.Accession: Accession ID of the dataset in the external database it was taken from
- experiment.Database: The name of the database where the dataset was taken from
- experiment.URI: URI of the original database
- experiment.SampleCount: Number of samples in the dataset
- experiment.batchEffect: A text field describing whether the dataset has batch effects
- experiment.batchCorrected: Whether batch correction has been performed on the dataset.
 This is a text field with possible details about the batch correction. Use geeq.batchEffect if a more structured field is needed.
- geeq.batchConfound: 0 if batch info isn't available, -1 if batch counfoud is detected, 1 if batch information is available and no batch confound found
- geeq.batchEffect: -1 if batch p value < 0.0001, 1 if batch p value > 0.1, 0 if otherwise and when there is no batch information is available or when the data is confounded with batches.
- geeq.rawData: -1 if no raw data available, 1 if raw data was available. When available, Gemma reprocesses raw data to get expression values and batches

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- geeq. qScore: Data quality score given to the dataset by Gemma.
- geeq.sScore: Suitability score given to the dataset by Gemma. Refers to factors like batches, platforms and other aspects of experimental design
- taxon.Name: The taxa of the study. In Gemma each study will include a single species. If the original source has samples from multiple species, they will be split into different studies within Gemma
- taxon. ID: Internal ID given to the taxon by Gemma

Examples

```
get_taxon_datasets("human")
```

search_annotations

Search for annotation tags

Description

Search for annotation tags

Usage

```
search_annotations(
  query,
  raw = getOption("gemma.raw", FALSE),
  memoised = getOption("gemma.memoised", FALSE),
  file = getOption("gemma.file", NA_character_),
  overwrite = getOption("gemma.overwrite", FALSE),
  attributes = getOption("gemma.attributes", TRUE)
)
```

Arguments

| query | The search query |
|------------|---|
| raw | TRUE to receive results as-is from Gemma, or FALSE to enable parsing. Raw results usually contain additional fields and flags that are omitted in the parsed results. |
| memoised | Whether or not to save to cache for future calls with the same inputs and use the result saved in cache if a result is already saved. Doing options(gemma.memoised = TRUE) will ensure that the cache is always used. Use forget_gemma_memoised to clear the cache. |
| file | The name of a file to save the results to, or NULL to not write results to a file. If raw == TRUE, the output will be a JSON file. Otherwise, it will be a RDS file. |
| overwrite | Whether or not to overwrite if a file exists at the specified filename. |
| attributes | If TRUE additional information from the call will be added into the output object's attributes such as offset and available elements. |

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Value

A data table with annotations (annotation search result value objects) matching the given identifiers. A list if raw = TRUE. A 400 error if required parameters are missing.

The fields of the output data.table are:

• category. Name: Category that the annotation belongs to

• category.URI: URI for the category.Name

• value.Name: Annotation term

• value.URI: URI for the value.Name

Examples

```
search_annotations("traumatic")
```

search_datasets

Retrieve datasets within a given taxa associated to an annotation tags search

Description

Retrieve datasets within a given taxa associated to an annotation tags search

Usage

```
search_datasets(
  query,
  taxon = NA_character_,
  offset = 0L,
  limit = 20L,
  sort = "+id",
  raw = getOption("gemma.raw", FALSE),
  memoised = getOption("gemma.memoised", FALSE),
  file = getOption("gemma.file", NA_character_),
  overwrite = getOption("gemma.overwrite", FALSE),
  attributes = getOption("gemma.attributes", TRUE)
)
```

Arguments

| query | The search query. Either plain text ('traumatic'), or an ontology term (UBERON_0002048). |
|--------|--|
| | Datasets that contain the given string in their short or full name will also be |
| | matched. Can be multiple identifiers separated by commas. |
| taxon | A numerical taxon identifier or an ncbi taxon identifier or a taxon identifier that matches either its scientific or common name |
| offset | The offset of the first retrieved result. |

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| limit | Optional, defaults to 20. Limits the result to specified amount of objects. Has a maximum value of 100. Use together with offset and the totalElements attribute in the output to compile all data if needed. |
|------------|---|
| sort | Order results by the given property and direction. The '+' sign indicate ascending order whereas the '-' indicate descending. |
| raw | TRUE to receive results as-is from Gemma, or FALSE to enable parsing. Raw results usually contain additional fields and flags that are omitted in the parsed results. |
| memoised | Whether or not to save to cache for future calls with the same inputs and use the result saved in cache if a result is already saved. Doing options(gemma.memoised = TRUE) will ensure that the cache is always used. Use forget_gemma_memoised to clear the cache. |
| file | The name of a file to save the results to, or NULL to not write results to a file. If raw == TRUE, the output will be a JSON file. Otherwise, it will be a RDS file. |
| overwrite | Whether or not to overwrite if a file exists at the specified filename. |
| attributes | If TRUE additional information from the call will be added into the output object's attributes such as offset and available elements. |

Value

A data table with information about the queried dataset(s). A list if raw = TRUE. Returns an empty list if no datasets matched. A successful response may contain 'Geeq' information, which aims to provide a unified metric to measure experiments by the quality of their data, and their suitability for use in Gemma. You can read more about the geeq properties here.

The fields of the output data.table are:

- experiment. ShortName: Shortname given to the dataset within Gemma. Often corresponds to accession ID
- experiment.Name: Full title of the dataset
- experiment. ID: Internal ID of the dataset.
- experiment.Description: Description of the dataset
- experiment.Public: Is the dataset publicly available. Only useful for logged in users with access to non-public data
- experiment. Troubled: Did an automatic process within gemma or a curator mark the dataset as "troubled"
- experiment.Accession: Accession ID of the dataset in the external database it was taken from
- experiment.Database: The name of the database where the dataset was taken from
- experiment.URI: URI of the original database
- experiment.SampleCount: Number of samples in the dataset
- experiment.batchEffect: A text field describing whether the dataset has batch effects
- experiment.batchCorrected: Whether batch correction has been performed on the dataset. This is a text field with possible details about the batch correction. Use geeq.batchEffect if a more structured field is needed.

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• geeq.batchConfound: 0 if batch info isn't available, -1 if batch counfoud is detected, 1 if batch information is available and no batch confound found

- geeq.batchEffect: -1 if batch p value < 0.0001, 1 if batch p value > 0.1, 0 if otherwise and when there is no batch information is available or when the data is confounded with batches.
- geeq.rawData: -1 if no raw data available, 1 if raw data was available. When available, Gemma reprocesses raw data to get expression values and batches
- geeq.qScore: Data quality score given to the dataset by Gemma.
- geeq.sScore: Suitability score given to the dataset by Gemma. Refers to factors like batches, platforms and other aspects of experimental design
- taxon.Name: The taxa of the study. In Gemma each study will include a single species. If the original source has samples from multiple species, they will be split into different studies within Gemma
- taxon. ID: Internal ID given to the taxon by Gemma

Examples

```
search_datasets("bipolar")
```

set_gemma_user

Authentication by user name

Description

Allows the user to access information that requires logging in to Gemma. To log out, run setGemmaUser without specifying the username or password.

Usage

```
set_gemma_user(username = NULL, password = NULL)
```

Arguments

username Your username (or empty, if logging out)
password Your password (or empty, if logging out)

Value

TRUE if authentication is successful, FALSE if not

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