

# Package: gemma.R (via r-universe)

June 23, 2024

**Title** A wrapper for Gemma's Restful API to access curated gene expression data and differential expression analyses

**Version** 3.1.6

**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

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**BugReports** <https://github.com/PavlidisLab/gemma.R/issues>

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**Repository** <https://bioc.r-universe.dev>

**RemoteUrl** <https://github.com/bioc/gemma.R>

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filter_properties	<i>Return all supported filter properties</i>
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**Description**

Some functions such as [get\\_datasets](#) and [get\\_platforms\\_by\\_ids](#) include a filter argument that allows creation of more complex queries. This function returns a list of supported properties to be used in those filters

**Usage**

```
filter_properties()
```

**Value**

A list of data.tables that contain supported properties and their data types

**Examples**

```
filter_properties()
```

---

forget_gemma_memoised	<i>Clear gemma.R cache</i>
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---

**Description**

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

**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 functions for Gemma's RESTful API that enables 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 transcriptomics data, currently primarily targeted at the analysis of gene expression profiles.

## Details

Most users will want to start with the high-level functions like [get\\_dataset\\_object](#), [get\\_differential\\_expression\\_val](#) and [get\\_platform\\_annotations](#). 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>

## See Also

Useful links:

- <https://pavlidislab.github.io/gemma.R/>
- <https://github.com/PavlidisLab/gemma.R>
- Report bugs at <https://github.com/PavlidisLab/gemma.R/issues>

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gemma_call	<i>Custom gemma call</i>
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---

**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)
```

---

gemma_kable	<i>Create printable tables out of gemma.R outputs</i>
-------------	---

---

**Description**

Creates a [kable](#) where certain columns are automatically shortened to better fit a document.

**Usage**

```
gemma_kable(table)
```

**Arguments**

table	A data.table or data.frame outputted by a gemma.R function
-------	--

---

gemma_memoise	<i>Enable and disable memoisation of gemma.R functions</i>
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---

**Description**

Enable and disable memoisation of gemma.R functions

**Usage**

```
gemma_memoise(  
  memoised = FALSE,  
  cache = rappdirs::user_cache_dir(appname = "gemmaR")  
)
```

**Arguments**

memoised	boolean. If TRUE memoisation will be enabled
cache	File path or "cache_in_memory". File path will chose a location to save the cache files for memoisation. "cache_in_memory" will store the cache in the current R session

---

get_all_pages	<i>Get all pages of a paginated call</i>
---------------	--

---

**Description**

Given a Gemma.R output from a function with offset and limit arguments, returns the output from all pages. All arguments other than offset, limit

**Usage**

```
get_all_pages(  
  query,  
  step_size = 100,  
  binder = rbind,  
  directory = NULL,  
  file = getOption("gemma.file", NA_character_),  
  overwrite = getOption("gemma.overwrite", FALSE)  
)
```

**Arguments**

query	Output from a gemma.R function with offset and limit argument
step_size	Size of individual calls to the server. 100 is the maximum value
binder	Binding function for the calls. If raw = FALSE use rbind to combine the data.tables. If not, use c to combine lists
directory	Directory to save the output from the individual calls to. If provided, each page is saved to separate files.
file	The name of a file to save the results to, or NULL to not write results to a file. This function always saves the output as an RDS file. Otherwise, it will be a RDS file.
overwrite	Whether or not to overwrite if a file exists at the specified filename.

**Value**

A data.table or a list containing data from all pages.

---

get\_child\_terms      *Return child terms of a term*

---

**Description**

When querying for ontology terms, Gemma propagates these terms to include any datasets with their child terms in the results. This function returns these children for any number of terms, including all children and the terms itself in the output vector

**Usage**

```
get_child_terms(terms)
```

**Arguments**

terms	An array of terms
-------	-------------------

**Value**

An array containing descendends of the annotation terms, including the terms themselves

**Examples**

```
get_child_terms("http://purl.obolibrary.org/obo/MONDO_0000408")
```

---

get_datasets	<i>Retrieve all datasets</i>
--------------	------------------------------

---

## Description

Retrieve all datasets

## Usage

```
get_datasets(
  query = NA_character_,
  filter = NA_character_,
  taxa = NA_character_,
  uris = 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)
)
```

## Arguments

query	The search query. Queries can include plain text or ontology terms. They also support conjunctions ("alpha AND beta"), disjunctions ("alpha OR beta") grouping ("(alpha OR beta) AND gamma"), prefixing ("alpha*"), wildcard characters ("BRCA?") and fuzzy matches ("alpha~").
filter	Filter results by matching expression. Use <a href="#">filter_properties</a> function to get a list of all available parameters. These properties can be combined using "and" "or" clauses and may contain common operators such as "=", "<" or "in". (e.g. "taxon.commonName = human", "taxon.commonName in (human,mouse)", "id < 1000")
taxa	A vector of taxon common names (e.g. human, mouse, rat). Providing multiple species will return results for all species. These are appended to the filter and equivalent to filtering for taxon.commonName property
uris	A vector of ontology term URIs. Providing multiple terms will return results containing any of the terms and their children. These are appended to the filter and equivalent to filtering for allCharacteristics.valueUri
offset	The offset of the first retrieved result.
limit	Defaults to 20. Limits the result to specified amount of objects. Has a maximum value of 100. Use together with offset and the totalElements <a href="#">attribute</a> 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 <code>options(gemma.memoised = TRUE)</code> will ensure that the cache is always used. Use <a href="#">forget_gemma_memoised</a> 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 <code>raw == TRUE</code> , the output will be the raw endpoint from the API, likely a JSON or a gzip file. Otherwise, it will be a RDS file.
overwrite	Whether or not to overwrite if a file exists at the specified filename.

### Value

A data table with information about the queried dataset(s). A list if `raw = TRUE`. Returns an empty list if no datasets matched.

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.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.batchEffectText`: A text field describing whether the dataset has batch effects
- `experiment.batchCorrected`: Whether batch correction has been performed on the dataset.
- `experiment.batchConfound`: 0 if batch info isn't available, -1 if batch confound is detected, 1 if batch information is available and no batch confound found
- `experiment.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.
- `experiment.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`: Name of the species

- `taxon.scientific`: Scientific name for the taxon
- `taxon.ID`: Internal identifier given to the species by Gemma
- `taxon.NCBI`: NCBI ID of the taxon
- `taxon.database.name`: Underlying database used in Gemma for the taxon
- `taxon.database.ID`: ID of the underlying database used in Gemma for the taxon

## Examples

```
get_datasets()
get_datasets(taxa = c("mouse", "human"), uris = "http://purl.obolibrary.org/obo/UBERON_0002048")
# filter below is equivalent to the call above
get_datasets(filter = "taxon.commonName in (mouse,human) and allCharacteristics.valueUri = http://purl.obolibrary.org/obo/UBERON_0002048")
get_datasets(query = "lung")
```

---

`get_datasets_by_ids`    *Retrieve datasets by their identifiers*

---

## Description

Retrieve datasets by their identifiers

## Usage

```
get_datasets_by_ids(
  datasets = NA_character_,
  filter = NA_character_,
  taxa = NA_character_,
  uris = 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)
)
```

## Arguments

<code>datasets</code>	Numerical dataset identifiers or dataset short names. If not specified, all datasets will be returned instead
<code>filter</code>	Filter results by matching expression. Use <a href="#">filter_properties</a> function to get a list of all available parameters. These properties can be combined using "and" "or" clauses and may contain common operators such as "=", "<" or "in". (e.g. "taxon.commonName = human", "taxon.commonName in (human,mouse)", "id < 1000")

taxa	A vector of taxon common names (e.g. human, mouse, rat). Providing multiple species will return results for all species. These are appended to the filter and equivalent to filtering for <code>taxon.commonName</code> property
uris	A vector of ontology term URIs. Providing multiple terms will return results containing any of the terms and their children. These are appended to the filter and equivalent to filtering for <code>allCharacteristics.valueUri</code>
offset	The offset of the first retrieved result.
limit	Defaults to 20. Limits the result to specified amount of objects. Has a maximum value of 100. Use together with <code>offset</code> and the <code>totalElements</code> 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 <code>options(gemma.memoised = TRUE)</code> will ensure that the cache is always used. Use <code>forget_gemma_memoised</code> 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 <code>raw == TRUE</code> , the output will be the raw endpoint from the API, likely a JSON or a gzip file. Otherwise, it will be a RDS file.
overwrite	Whether or not to overwrite if a file exists at the specified filename.

### Value

A data table with information about the queried dataset(s). A list if `raw = TRUE`. Returns an empty list if no datasets matched.

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.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.batchEffectText`: A text field describing whether the dataset has batch effects

- `experiment.batchCorrected`: Whether batch correction has been performed on the dataset.
- `experiment.batchConfound`: 0 if batch info isn't available, -1 if batch confound is detected, 1 if batch information is available and no batch confound found
- `experiment.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.
- `experiment.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`: Name of the species
- `taxon.scientific`: Scientific name for the taxon
- `taxon.ID`: Internal identifier given to the species by Gemma
- `taxon.NCBI`: NCBI ID of the taxon
- `taxon.database.name`: Underlying database used in Gemma for the taxon
- `taxon.database.ID`: ID of the underlying database used in Gemma for the taxon

## Examples

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

---

```
get_dataset_annotations
```

*Retrieve the annotations of a dataset*

---

## Description

Retrieve the annotations of a dataset

## 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)  
)
```

**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 <code>options(gemma.memoised = TRUE)</code> will ensure that the cache is always used. Use <code>forget_gemma_memoised</code> 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 <code>raw == TRUE</code> , the output will be the raw endpoint from the API, likely a JSON or a gzip file. Otherwise, it will be a RDS file.
overwrite	Whether or not to overwrite if a file exists at the specified filename.

**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.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
- `object.class`: Class of object that the term originated from.

**Examples**

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

---

```
get_dataset_design      Retrieve the design of a dataset
```

---

**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)
)
```

### 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 <code>options(gemma.memoised = TRUE)</code> will ensure that the cache is always used. Use <code>forget_gemma_memoised</code> 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 <code>raw == TRUE</code> , the output will be the raw endpoint from the API, likely a JSON or a gzip file. Otherwise, it will be a RDS file.
overwrite	Whether or not to overwrite if a file exists at the specified filename.

### 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"))
```

---

```
get_dataset_differential_expression_analyses
```

*Retrieve annotations and surface level stats for a dataset's differential analyses*

---

### Description

Retrieve annotations and surface level stats for a dataset's differential analyses

### 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)  
)
```

**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 <code>options(gemma.memoised = TRUE)</code> will ensure that the cache is always used. Use <a href="#">forget_gemma_memoised</a> 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 <code>raw == TRUE</code> , the output will be the raw endpoint from the API, likely a JSON or a gzip file. Otherwise, it will be a RDS file.
overwrite	Whether or not to overwrite if a file exists at the specified filename.

**Value**

A data table with information about the differential expression analysis of the queried dataset. Note that this function 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
- `factor.category`: Category for the contrast
- `factor.category.URI`: URI for the contrast category
- `factor.ID`: ID of the factor
- `baseline.factors`: Characteristics of the baseline. This field is a `data.table`
- `experimental.factors`: Characteristics of the experimental group. This field is a `data.table`
- `isSubset`: TRUE if the result set belong to a subset, FALSE if not. Subsets are created when performing differential expression to avoid unhelpful comparisons.
- `subsetFactor`: Characteristics of the subset. This field is a `data.table`
- `probes.analyzed`: Number of probesets represented in the contrast
- `genes.analyzed`: Number of genes represented in the contrast

**Examples**

```
result <- get_dataset_differential_expression_analyses("GSE2872")
get_differential_expression_values(resultSet = result$result.ID[1])
```

---

```
get_dataset_expression_for_genes
```

*Retrieve the expression data matrix of a set of datasets and genes*

---

### Description

Retrieve the expression data matrix of a set of datasets and genes

### Usage

```
get_dataset_expression_for_genes(
  datasets,
  genes,
  keepNonSpecific = FALSE,
  consolidate = NA_character_,
  raw = getOption("gemma.raw", FALSE),
  memoised = getOption("gemma.memoised", FALSE),
  file = getOption("gemma.file", NA_character_),
  overwrite = getOption("gemma.overwrite", FALSE)
)
```

### Arguments

datasets	A numerical dataset identifier or a dataset short name
genes	An ensembl gene identifier which typically starts with ensg or an ncbi gene identifier or an official gene symbol approved by hgnc
keepNonSpecific	logical. FALSE by default. If TRUE, results from probesets that are not specific to the gene will also be returned.
consolidate	An option for gene expression level consolidation. If empty, will return every probe for the genes. "pickmax" to pick the probe with the highest expression, "pickvar" to pick the prove with the highest variance and "average" for returning the average expression
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 <code>options(gemma.memoised = TRUE)</code> will ensure that the cache is always used. Use <code>forget_gemma_memoised</code> 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 <code>raw == TRUE</code> , the output will be the raw endpoint from the API, likely a JSON or a gzip file. Otherwise, it will be a RDS file.
overwrite	Whether or not to overwrite if a file exists at the specified filename.



**Value**

A list of data frames

**Examples**

```
get_dataset_expression_for_genes("GSE2018", genes = c(10225, 2841))
```

---

get\_dataset\_object      *Compile gene expression data and metadata*

---

**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(  
  datasets,  
  genes = NULL,  
  keepNonSpecific = FALSE,  
  consolidate = NA_character_,  
  resultSets = NULL,  
  contrasts = NULL,  
  metaType = "text",  
  type = "se",  
  memoised = getOption("gemma.memoised", FALSE)  
)
```

**Arguments**

datasets	A numerical dataset identifier or a dataset short name
genes	An ensembl gene identifier which typically starts with ensg or an ncbi gene identifier or an official gene symbol approved by hgnc
keepNonSpecific	logical. FALSE by default. If TRUE, results from probesets that are not specific to the gene will also be returned.
consolidate	An option for gene expression level consolidation. If empty, will return every probe for the genes. "pickmax" to pick the probe with the highest expression, "pickvar" to pick the probe with the highest variance and "average" for returning the average expression
resultSets	Result set IDs of the a differential expression analysis. Optional. If provided, the output will only include the samples from the subset used in the result set ID. Must be the same length as datasets.'

contrasts	Contrast IDs of a differential expression contrast. Optional. Need resultSets to be defined to work. If provided, the output will only include samples relevant to the specific contrasts.
metaType	How should the metadata information should be included. Can be "text", "uri" or "both". "text" and "uri" options
type	"se" for a SummarizedExperiment or "eset" for Expression Set. We recommend using SummarizedExperiments which are more recent. See the Summarized experiment <a href="#">vignette</a> or the ExpressionSet <a href="#">vignette</a> for more details. "tidy" for a long form data frame compatible with tidyverse functions. 'list' to return a list containing individual data frames containing expression values, design and the experiment.
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 <code>options(gemma.memoised = TRUE)</code> will ensure that the cache is always used. Use <a href="#">forget_gemma_memoised</a> to clear the cache.

### Value

A list of [SummarizedExperiments](#), [ExpressionSets](#) or a tibble containing metadata and expression data for the queried datasets and genes. Metadata will be expanded to include a variable number of factors that annotates samples from a dataset but will always include single "factorValues" column that houses data.tables that include all annotations for a given sample.

### Examples

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

---

`get_dataset_platforms` *Retrieve the platforms of a dataset*

---

### Description

Retrieve the platforms 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)
)
```

### 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 <code>options(gemma.memoised = TRUE)</code> will ensure that the cache is always used. Use <a href="#">forget_gemma_memoised</a> 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 <code>raw == TRUE</code> , the output will be the raw endpoint from the API, likely a JSON or a gzip file. Otherwise, it will be a RDS file.
overwrite	Whether or not to overwrite if a file exists at the specified filename.

### 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
- `platform.type`: Technology type for the platform.
- `taxon.name`: Name of the species platform was made for
- `taxon.scientific`: Scientific name for the taxon
- `taxon.ID`: Internal identifier given to the species by Gemma
- `taxon.NCBI`: NCBI ID of the taxon
- `taxon.database.name`: Underlying database used in Gemma for the taxon
- `taxon.database.ID`: ID of the underlying database used in Gemma for the taxon

### Examples

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

---

`get_dataset_processed_expression`*Retrieve processed expression data of a dataset*

---

## Description

Retrieve processed expression data of a dataset

## Usage

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

## Arguments

<code>dataset</code>	A numerical dataset identifier or a dataset short name
<code>raw</code>	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.
<code>memoised</code>	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 <code>options(gemma.memoised = TRUE)</code> will ensure that the cache is always used. Use <a href="#">forget_gemma_memoised</a> to clear the cache.
<code>file</code>	The name of a file to save the results to, or NULL to not write results to a file. If <code>raw == TRUE</code> , the output will be the raw endpoint from the API, likely a JSON or a gzip file. Otherwise, it will be a RDS file.
<code>overwrite</code>	Whether or not to overwrite if a file exists at the specified filename.

## 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

```
get_dataset_processed_expression("GSE2018")
```

---

get\_dataset\_quantitation\_types  
*Retrieve quantitation types of a dataset*

---

## Description

Retrieve quantitation types of a dataset

## Usage

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

## 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 <code>options(gemma.memoised = TRUE)</code> will ensure that the cache is always used. Use <a href="#">forget_gemma_memoised</a> 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 <code>raw == TRUE</code> , the output will be the raw endpoint from the API, likely a JSON or a gzip file. Otherwise, it will be a RDS file.
overwrite	Whether or not to overwrite if a file exists at the specified filename.

## Value

A `data.table` containing the quantitation types

The fields of the output `data.table` are:

- `id`: Id of the quantitation type. Any raw quantitation type can be accessed by [get\\_dataset\\_raw\\_expression](#) function using this id.
- `name`: Name of the quantitation type
- `description`: Description of the quantitation type
- `type`: Type of the quantitation type. Either raw or processed. Each dataset will have one processed quantitation type which is the data returned using [get\\_dataset\\_processed\\_expression](#)

- **ratio**: Whether or not the quantitation type is a ratio of multiple quantitation types. Typically TRUE for processed TWOCOLOR quantitation type.
- **preferred**: The preferred raw quantitation type. This version is used in generation of the processed data within gemma.
- **recomputed**: If TRUE this quantitation type is generated by recomputing raw data files Gemma had access to.

## Examples

```
get_dataset_quantitation_types("GSE59918")
```

---

```
get_dataset_raw_expression
```

*Retrieve raw expression data of a dataset*

---

## Description

Retrieve raw expression data of a dataset

## Usage

```
get_dataset_raw_expression(
  dataset,
  quantitationType,
  raw = getOption("gemma.raw", FALSE),
  memoised = getOption("gemma.memoised", FALSE),
  file = getOption("gemma.file", NA_character_),
  overwrite = getOption("gemma.overwrite", FALSE)
)
```

## Arguments

dataset	A numerical dataset identifier or a dataset short name
quantitationType	Quantitation type id. These can be acquired using <a href="#">get_dataset_quantitation_types</a> function. This endpoint can only return non-processed quantitation types.
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 <code>options(gemma.memoised = TRUE)</code> will ensure that the cache is always used. Use <a href="#">forget_gemma_memoised</a> 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 <code>raw == TRUE</code> , the output will be the raw endpoint from the API, likely a JSON or a gzip file. Otherwise, it will be a RDS file.
overwrite	Whether or not to overwrite if a file exists at the specified filename.

**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**

```
q_types <- get_dataset_quantitation_types("GSE59918")
get_dataset_raw_expression("GSE59918", q_types$id[q_types$name == "Counts"])
```

---

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)
)
```

**Arguments**

dataset	A numerical dataset identifier or a dataset short name
raw	<code>TRUE</code> to receive results as-is from Gemma, or <code>FALSE</code> 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 <code>options(gemma.memoised = TRUE)</code> will ensure that the cache is always used. Use <code>forget_gemma_memoised</code> to clear the cache.
file	The name of a file to save the results to, or <code>NULL</code> to not write results to a file. If <code>raw == TRUE</code> , the output will be the raw endpoint from the API, likely a JSON or a gzip file. Otherwise, it will be a RDS file.
overwrite	Whether or not to overwrite if a file exists at the specified filename.

**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.factorValues: Experimental factor values of the sample. This field is a data table

**Examples**

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

---

```
get_differential_expression_values
```

*Retrieve differential expression results*

---

**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_,
  resultSets = NA_integer_,
  readableContrasts = FALSE,
  memoised = getOption("gemma.memoised", FALSE)
)
```

**Arguments**

dataset	A dataset identifier.
resultSets	resultSet identifiers. If a dataset is not provided, all result sets will be downloaded. If it is provided it will only be used to ensure all result sets belong to the dataset.



**readableContrasts**

If FALSE (default), the returned columns will use internal contrasts IDs as names. Details about the contrasts can be accessed using [get\\_dataset\\_differential\\_expression\\_analysis](#). 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 to two or 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 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 gene identifiers*

---

**Description**

Retrieve genes matching gene identifiers

**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)
)

```

**Arguments**

genes	An ensembl gene identifier which typically starts with <code>ensg</code> 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 <code>options(gemma.memoised = TRUE)</code> will ensure that the cache is always used. Use <a href="#">forget_gemma_memoised</a> 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 <code>raw == TRUE</code> , the output will be the raw endpoint from the API, likely a JSON or a gzip file. Otherwise, it will be a RDS file.
overwrite	Whether or not to overwrite if a file exists at the specified filename.

**Value**

A data table with information about the queried 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
- `gene.aliases`: Gene aliases. Each row includes a vector
- `gene.MFX.rank`: Multifunctionality rank for the gene
- `taxon.name`: Name of the species
- `taxon.scientific`: Scientific name for the taxon
- `taxon.ID`: Internal identifier given to the species by Gemma
- `taxon.NCBI`: NCBI ID of the taxon
- `taxon.database.name`: Underlying database used in Gemma for the taxon
- `taxon.database.ID`: ID of the underlying database used in Gemma 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)
)
```

**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 <code>options(gemma.memoised = TRUE)</code> will ensure that the cache is always used. Use <a href="#">forget_gemma_memoised</a> 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 <code>raw == TRUE</code> , the output will be the raw endpoint from the API, likely a JSON or a gzip file. Otherwise, it will be a RDS file.
overwrite	Whether or not to overwrite if a file exists at the specified filename.

**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.

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(3091)
```

---

```
get_gene_locations      Retrieve the physical locations of a given gene
```

---

**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)
)
```

**Arguments**

gene	An ensembl gene identifier which typically starts with <code>ensg</code> 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 <code>options(gemma.memoised = TRUE)</code> will ensure that the cache is always used. Use <a href="#">forget_gemma_memoised</a> 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 <code>raw == TRUE</code> , the output will be the raw endpoint from the API, likely a JSON or a gzip file. Otherwise, it will be a RDS file.
overwrite	Whether or not to overwrite if a file exists at the specified filename.

**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")
get_gene_locations(1859)
```

---

get_gene_probes	<i>Retrieve the probes associated to a genes across all platforms</i>
-----------------	---

---

### Description

Retrieve the probes associated to a genes across all platforms

### 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)
)
```

### Arguments

gene	An ensembl gene identifier which typically starts with <code>ensg</code> or an ncbi gene identifier or an official gene symbol approved by hgnc
offset	The offset of the first retrieved result.
limit	Defaults to 20. Limits the result to specified amount of objects. Has a maximum value of 100. Use together with <code>offset</code> and the <code>totalElements</code> 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 <code>options(gemma.memoised = TRUE)</code> will ensure that the cache is always used. Use <code>forget_gemma_memoised</code> 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 the raw endpoint from the API, likely a JSON or a gzip file. Otherwise, it will be a RDS file.
overwrite	Whether or not to overwrite if a file exists at the specified filename.

**Value**

A data table with information about the probes representing a gene across all platforms. 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:

- element.name: Name of the element. Typically the probeset name
- element.description: A free text field providing optional information about the element
- 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.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.
- taxon.name: Name of the species platform was made for
- taxon.scientific: Scientific name for the taxon
- taxon.ID: Internal identifier given to the species by Gemma
- taxon.NCBI: NCBI ID of the taxon
- taxon.database.name: Underlying database used in Gemma for the taxon
- taxon.database.ID: ID of the underlying database used in Gemma for the taxon

**Examples**

```
get_gene_probes(1859)
```

---

```
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_,
  filter = NA_character_,
  taxa = 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)
)

```

**Arguments**

platforms	Platform numerical identifiers or platform short names. If not specified, all platforms will be returned instead
filter	Filter results by matching expression. Use <a href="#">filter_properties</a> function to get a list of all available parameters. These properties can be combined using "and" "or" clauses and may contain common operators such as "=", "<" or "in". (e.g. "taxon.commonName = human", "taxon.commonName in (human,mouse)", "id < 1000")
taxa	A vector of taxon common names (e.g. human, mouse, rat). Providing multiple species will return results for all species. These are appended to the filter and equivalent to filtering for taxon.commonName property
offset	The offset of the first retrieved result.
limit	Defaults to 20. Limits the result to specified amount of objects. Has a maximum value of 100. Use together with offset and the totalElements <a href="#">attribute</a> 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 <a href="#">forget_gemma_memoised</a> 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 the raw endpoint from the API, likely a JSON or a gzip file. Otherwise, it will be a RDS file.
overwrite	Whether or not to overwrite if a file exists at the specified filename.

**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
- platform.type: Technology type for the platform.
- taxon.name: Name of the species platform was made for
- taxon.scientific: Scientific name for the taxon
- taxon.ID: Internal identifier given to the species by Gemma
- taxon.NCBI: NCBI ID of the taxon
- taxon.database.name: Underlying database used in Gemma for the taxon
- taxon.database.ID: ID of the underlying database used in Gemma for the taxon

**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 numerical identifiers or platform short name.
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 <code>options(gemma.memoised = TRUE)</code> will ensure that the cache is always used. Use <code>forget_gemma_memoised</code> 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 "|"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.

**Examples**

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

---

`get_platform_datasets` *Retrieve all experiments using a given platform*

---

**Description**

Retrieve all experiments using 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)
)
```

**Arguments**

platform	A platform numerical identifier or a platform short name
offset	The offset of the first retrieved result.
limit	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 the raw endpoint from the API, likely a JSON or a gzip file. Otherwise, it will be a RDS file.
overwrite	Whether or not to overwrite if a file exists at the specified filename.

**Value**

A data table with information about the queried dataset(s). A list if raw = TRUE. Returns an empty list if no datasets matched.

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.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.batchEffectText: A text field describing whether the dataset has batch effects
- experiment.batchCorrected: Whether batch correction has been performed on the dataset.
- experiment.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
- experiment.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.

- `experiment.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`: Name of the species
- `taxon.scientific`: Scientific name for the taxon
- `taxon.ID`: Internal identifier given to the species by Gemma
- `taxon.NCBI`: NCBI ID of the taxon
- `taxon.database.name`: Underlying database used in Gemma for the taxon
- `taxon.database.ID`: ID of the underlying database used in Gemma for the taxon

### 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)  
)
```

### Arguments

<code>platform</code>	A platform numerical identifier or a platform short name
<code>probe</code>	A probe name or its numerical identifier
<code>offset</code>	The offset of the first retrieved result.
<code>limit</code>	Defaults to 20. Limits the result to specified amount of objects. Has a maximum value of 100. Use together with <code>offset</code> and the <code>totalElements</code> 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 <code>options(gemma.memoised = TRUE)</code> will ensure that the cache is always used. Use <code>forget_gemma_memoised</code> 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 <code>raw == TRUE</code> , the output will be the raw endpoint from the API, likely a JSON or a gzip file. Otherwise, it will be a RDS file.
overwrite	Whether or not to overwrite if a file exists at the specified filename.

### Value

A data table with information about the queried 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
- `gene.aliases`: Gene aliases. Each row includes a vector
- `gene.MFX.rank`: Multifunctionality rank for the gene
- `taxon.name`: Name of the species
- `taxon.scientific`: Scientific name for the taxon
- `taxon.ID`: Internal identifier given to the species by Gemma
- `taxon.NCBI`: NCBI ID of the taxon
- `taxon.database.name`: Underlying database used in Gemma for the taxon
- `taxon.database.ID`: ID of the underlying database used in Gemma for the taxon

### Examples

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

---

```
get_result_sets
```

*Retrieve all result sets matching the provided criteria*

---

### Description

Returns queried result set

**Usage**

```

get_result_sets(
  datasets = NA_character_,
  resultSets = NA_character_,
  filter = NA_character_,
  offset = 0,
  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)
)

```

**Arguments**

datasets	A numerical dataset identifier or a dataset short name
resultSets	A resultSet identifier. Note that result set identifiers are not static and can change when Gemma re-runs analyses internally. When using these as inputs, try to make sure you access a currently existing result set ID by basing them on result sets returned for a particular dataset or filter used in <a href="#">get_result_sets</a>
filter	Filter results by matching expression. Use <a href="#">filter_properties</a> function to get a list of all available parameters. These properties can be combined using "and" "or" clauses and may contain common operators such as "=", "<" or "in". (e.g. "taxon.commonName = human", "taxon.commonName in (human,mouse)", "id < 1000")
offset	The offset of the first retrieved result.
limit	Defaults to 20. Limits the result to specified amount of objects. Has a maximum value of 100. Use together with offset and the totalElements <a href="#">attribute</a> 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 <code>options(gemma.memoised = TRUE)</code> will ensure that the cache is always used. Use <a href="#">forget_gemma_memoised</a> 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 <code>raw == TRUE</code> , the output will be the raw endpoint from the API, likely a JSON or a gzip file. Otherwise, it will be a RDS file.
overwrite	Whether or not to overwrite if a file exists at the specified filename.

**Details**

Output and usage of this function is mostly identical to [get\\_dataset\\_differential\\_expression\\_analyses](#). The principal difference being the ability to restrict your result sets, being able to query across multiple datasets and being able to use the filter argument to search based on result set properties.

**Value**

A data table with information about the queried result sets. Note that this function does not return differential expression values themselves. Use [get\\_differential\\_expression\\_values](#) to get differential expression values

- `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
- `factor.category`: Category for the contrast
- `factor.category.URI`: URI for the contrast category
- `factor.ID`: ID of the factor
- `baseline.factors`: Characteristics of the baseline. This field is a `data.table`
- `experimental.factors`: Characteristics of the experimental group. This field is a `data.table`
- `isSubset`: TRUE if the result set belong to a subset, FALSE if not. Subsets are created when performing differential expression to avoid unhelpful comparisons.
- `subsetFactor`: Characteristics of the subset. This field is a `data.table`

**Examples**

```
get_result_sets(dataset = 1)
# get all contrasts comparing disease states. use filter_properties to see available options
get_result_sets(filter = "baselineGroup.characteristics.value = disease")
```

---

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()
```

---

make_design	<i>Make simplified design frames</i>
-------------	--------------------------------------

---

**Description**

Using on the output of `get_dataset_samples`, this function creates a simplified design table, granting one column to each experimental variable

**Usage**

```
make_design(samples, metaType = "text")
```

**Arguments**

`samples` An output from `get_dataset_samples`. The output should not be raw

`metaType` Type of metadata to include in the output. "text", "uri" or "both"

**Value**

A data.frame including the design table for the dataset

**Examples**

```
samples <- get_dataset_samples('GSE46416')
make_design(samples)
```

---

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)
)
```

## Arguments

query	The search query. Queries can include plain text or ontology terms They also support conjunctions ("alpha AND beta"), disjunctions ("alpha OR beta") grouping ("(alpha OR beta) AND gamma"), prefixing ("alpha*"), wildcard characters ("BRCA?") and fuzzy matches ("alpha~").
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 <code>options(gemma.memoised = TRUE)</code> will ensure that the cache is always used. Use <a href="#">forget_gemma_memoised</a> 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 <code>raw == TRUE</code> , the output will be the raw endpoint from the API, likely a JSON or a gzip file. Otherwise, it will be a RDS file.
overwrite	Whether or not to overwrite if a file exists at the specified filename.

## 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_annotatons("traumatic")
```

---

```
search_gemma
```

---

*Search everything in Gemma*

---

**Description**

Search everything in Gemma

**Usage**

```
search_gemma(
  query,
  taxon = NA_character_,
  platform = NA_character_,
  limit = 100,
  resultType = "experiment",
  raw = getOption("gemma.raw", FALSE),
  memoised = getOption("gemma.memoised", FALSE),
  file = getOption("gemma.file", NA_character_),
  overwrite = getOption("gemma.overwrite", FALSE)
)
```

**Arguments**

query	The search query. Queries can include plain text or ontology terms They also support conjunctions ("alpha AND beta"), disjunctions ("alpha OR beta") grouping ("(alpha OR beta) AND gamma"), prefixing ("alpha*"), wildcard characters ("BRCA?") and fuzzy matches ("alpha~").
taxon	A numerical taxon identifier or an ncbi taxon identifier or a taxon identifier that matches either its scientific or common name
platform	A platform numerical identifier or a platform short name
limit	Defaults to 100 with a maximum value of 2000. Limits the number of returned results. Note that this function does not support pagination.
resultType	The kind of results that should be included in the output. Can be experiment, gene, platform or a long object type name, documented in the API documentation.
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 <a href="#">forget_gemma_memoised</a> 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 the raw endpoint from the API, likely a JSON or a gzip file. Otherwise, it will be a RDS file.
overwrite	Whether or not to overwrite if a file exists at the specified filename.

**Value**

If raw = FALSE and resultType is experiment, gene or platform, a data.table containing the search results. If it is any other type, a list of results. A list with additional details about the search if raw = TRUE

**Examples**

```
search_gemma("bipolar")
```

---

set_gemma_user	<i>Authentication by user name</i>
----------------	------------------------------------

---

**Description**

Allows the user to access information that requires logging in to Gemma. To log out, run set\_gemma\_user 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

---

update_result	<i>Update result</i>
---------------	----------------------

---

**Description**

Re-runs the function used to create a gemma.R output to update the data at hand. Useful if you have a reason to believe parts of the data has changed since your last accession and you wish to update while decoupling the update process from your original code used to generate the data.

**Usage**

```
update_result(query)
```

**Arguments**

query	Output from a gemma.R function
-------	--------------------------------

**Details**

Note that if you have used the file and overwrite arguments with the original call, this will also repeat to regenerate the file based on your initial preference

**Examples**

```
annots <- get_dataset_annotations(1)
# wait for a couple of years..
# wonder if the results are the same
updated_annots <- update_result(annots)

# also works with outputs of get_all_pages
platforms <- get_all_pages(get_platforms_by_ids())
updated_platforms <- update_result(platforms)
```

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