Package 'biodbLipidmaps'

May 22, 2022

Title biodbLipidmaps, a library for connecting to the Lipidmaps Structure database

Version 1.2.0

Description The biodbLipidmaps library provides access to the Lipidmaps Structure Database, using biodb package framework. It allows to retrieve entries by their accession number, and run web the services lmsdSearch and lmsdRecord.

URL https://github.com/pkrog/biodbLipidmaps

BugReports https://github.com/pkrog/biodbLipidmaps/issues

biocViews Software, Infrastructure, DataImport

License AGPL-3

Encoding UTF-8

VignetteBuilder knitr

Depends R (>= 4.1)

Imports biodb (>= 1.3.2), lifecycle, R6

Suggests BiocStyle, lgr, roxygen2, devtools, testthat (>= 2.0.0), knitr, rmarkdown, covr

RoxygenNote 7.1.2

Collate 'LipidmapsStructureConn.R' 'LipidmapsStructureEntry.R' 'package.R'

git_url https://git.bioconductor.org/packages/biodbLipidmaps

git_branch RELEASE_3_15

git_last_commit 484c60e

git_last_commit_date 2022-04-26

Date/Publication 2022-05-22

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R topics documented:

`biodbLipidmaps-package`

biodbLipidmaps: biodbLipidmaps, a library for connecting to the Lipidmaps Structure database

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Description

The biodbLipidmaps library provides access to the Lipidmaps Structure Database, using biodb package framework. It allows to retrieve entries by their accession number, and run web the services lmsdSearch and lmsdRecord.

Details

See vignette intro: "‘ vignette('intro', package='biodbLipidmaps') "‘

Author(s)

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See Also

LipidmapsStructureConn.

`LipidmapsStructureConn`

Lipidmaps Structure connector class.

Description

Lipidmaps Structure connector class.

Lipidmaps Structure connector class.

Details

Connector class for Lipidmaps Structure.

Super classes

`biodb::BiodbConnBase` -> `biodb::BiodbConn` -> `LipidmapsStructureConn`

Methods

Public methods:

- LipidmapsStructureConn$wsLmsdSearch()
- LipidmapsStructureConn$wsLmsd()
- LipidmapsStructureConn$wsLmsdRecord()
- LipidmapsStructureConn$clone()

Method `wsLmsdSearch()`**:** Calls LMSDSearch web service. See https://www.lipidmaps.org/data/structure/programmatica for details.

Usage:

```
LipidmapsStructureConn$wsLmsdSearch(
  mode = NULL,
  output.mode = NULL,
  output.type = NULL,
  output.delimiter = NULL,
  output.quote = NULL,
  output.column.header = NULL,
  lmid = NULL,
  name = NULL,
  formula = NULL,
  search.type = NULL,
  smiles.string = NULL,
  exact.mass = NULL,
  exact.mass.offset = NULL,
  core.class = NULL,
  main.class = NULL,
  sub.class = NULL,
  retfmt = c("plain", "request", "parsed", "ids")
)
```

Arguments:

`mode` The search mode: 'ProcessStrSearch', 'ProcessTextSearch' or 'ProcessTextOntologySearch'. Compulsory.

`output.mode` If set to 'File', will output a in format 'output.type', otherwise will output HTML.

`output.type` The output format: 'TSV', 'CSV' or 'SDF'.

`output.delimiter` The delimiter for TSV or CSV formats: 'Tab', 'Comma', 'Semicolon'.

`output.quote` If quotes are to be used: 'Yes' or 'No'.

`output.column.header` If header must be output: 'Yes' or 'No'.

`lmid` a Lipidmaps ID.

`name` The name to search for.

`formula` The chemical formula to search for.

`search.type` The search type: 'SubStructure' or 'ExactMatch'.

`smiles.string` A SMILES to search for.

`exact.mass` The mass to search for.

`exact.mass.offset` The tolerance on the mass search.

`core.class` An integer number from 1 to 8.

`main.class` An integer number. See Lipidmaps documentation.

`sub.class` An integer number. See Lipidmaps documentation.

`retfmt` Use to set the format of the returned value. 'plain' will return the raw results from the server, as a character value. 'request' will return the request that would have been sent, as a BiodbRequest object. 'parsed' will return data frame. 'ids' will return a character vector containing the IDs of the matching entries.

Returns: Depending on 'retfmt'.

Method `wsLmsd()`**:** Calls LMSD web service for downloading one entry.

Usage:

```
LipidmapsStructureConn$wsLmsd(
  lmid,
  format = c("tsv", "csv"),
  retfmt = c("plain", "request", "parsed")
)
```

Arguments:

`lmid` The accession number of the entry to retrieve.

`format` The output format (either 'tsv' or 'csv').

`retfmt` Use to set the format of the returned value. 'plain' will return the raw results from the server, as a character value. 'request' will return the request that would have been sent, as a BiodbRequest object. 'parsed' will return data frame.

Returns: Depending on 'retfmt'.

Method `wsLmsdRecord()`**:** Calls LMSDRecord web service. See http://www.lipidmaps.org/data/structure/programmatica

Usage:

```
LipidmapsStructureConn$wsLmsdRecord(
  lmid,
  mode = NULL,
  output.type = NULL,
  output.delimiter = NULL,
  output.quote = NULL,
  output.column.header = NULL,
  retfmt = c("plain", "request", "parsed")
)
```

Arguments:

`lmid` A character vector containing the IDs of the wanted entries.

`mode` If set to 'File', will output a in format 'output.type', otherwise will output HTML.

`output.type` The output format: 'TSV', 'CSV' or 'SDF'.

`output.delimiter` The delimiter for TSV or CSV formats: 'Tab', 'Comma', 'Semicolon'.

`output.quote` If quotes are to be used: 'Yes' or 'No'.

`output.column.header` If header must be output: 'Yes' or 'No'.

`retfmt` Use to set the format of the returned value. 'plain' will return the raw results from the server, as a character value. 'request' will return the request that would have been sent, as a BiodbRequest object. 'parsed' will return data frame.

Returns: Depending on 'retfmt'.

Method `clone()`**:** The objects of this class are cloneable with this method.

Usage:

```
LipidmapsStructureConn$clone(deep = FALSE)
```

Arguments:

`deep` Whether to make a deep clone.

Examples

```
# Create an instance with default settings:
mybiodb <- biodb::newInst()

# Create a connector
conn <- mybiodb$getFactory()$createConn('lipidmaps.structure')

# Get an entry
e <- conn$getEntry('LMFA00000001')

# Terminate instance.
mybiodb$terminate()
```

`LipidmapsStructureEntry`

Lipidmaps Structure entry class.

Description

Lipidmaps Structure entry class.

Lipidmaps Structure entry class.

Details

Entry class for Lipidmaps Structure.

Super classes

`biodb::BiodbEntry` -> `biodb::BiodbCsvEntry` -> `LipidmapsStructureEntry`

Methods

Public methods:

- `LipidmapsStructureEntry$new()`
- `LipidmapsStructureEntry$clone()`

Method `new()`**:** New instance initializer. Entry classes must not be instantiated directly. Instead, you must use the getEntry() method of the connector class.

Usage:

```
LipidmapsStructureEntry$new(...)
```

Arguments:

`...` All parameters are passed to super class' initializer.

Returns: Nothing.

Method `clone()`**:** The objects of this class are cloneable with this method.

Usage:

```
LipidmapsStructureEntry$clone(deep = FALSE)
```

Arguments:

`deep` Whether to make a deep clone.

Examples

```
# Create an instance with default settings:
mybiodb <- biodb::newInst()

# Create a connector
conn <- mybiodb$getFactory()$createConn('lipidmaps.structure')

# Get an entry
e <- conn$getEntry('LMFA00000001')

# Terminate instance.
mybiodb$terminate()
```

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