

Package: fioRa (via r-universe)

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Title Mass-Spectra Prediction Using the FIORA Model

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Description Provides a wrapper for the python module 'FIORA' as well as a 'shiny'-App to facilitate data processing and visualization. 'FIORA' allows to predict Mass-Spectra based on the SMILES code of chemical compounds. It is described in the Nature Communications article by Nowatzky (2025) [doi:10.1038/s41467-025-57422-4](https://doi.org/10.1038/s41467-025-57422-4).

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Imports bslib, config, golem, InterpretMSSpectrum, rcdk, shiny, shinyjs, waiter

Encoding UTF-8

RoxygenNote 7.3.3

URL <https://github.com/janlisec/fioRa>

BugReports <https://github.com/janlisec/fioRa/issues>

Suggests reticulate, S4Vectors, Spectra, spelling, testthat (>= 3.0.0), vdiff, withr

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install_fiora	<i>Install the python module 'fiora' into a conda environment.</i>
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Description

This function will check and perform the installation of three components in the following order: reticulate, miniconda and fiora. It will ensure that a working conda environment 'fiora' is available. This is a prerequisite for both, [run_app](#) and [run_script](#).

Usage

```
install_fiora(conda_name = "fiora")
```

Arguments

conda_name The name of the conda environment where fiora shall be installed to.

Value

A list providing the current OS and path information on the current python executable and the fiora script.

Examples

```
## Not run:
# this will install packages and software on your machine
install_fiora()

## End(Not run)
```

read_fiora	<i>Read a fiora result file (mgf) into a R object.</i>
------------	--

Description

Allows to import a 'fioRa' output file (msp like format) to a list like format or simplified to a data frame or Spectra object.

Usage

```
read_fiora(fl, fmt = c("list", "df", "Spectra"))
```

Arguments

fl	file.
fmt	A named list object is returned if not specified otherwise. Set 'fmt=df' to simplify the return value to a data frame. Use 'fmt=Spectra' to convert to a [Spectra::Spectra()] object.

Value

A result list of length = n_compounds containing metadata fields as sub-lists and the predicted MS² spectrum as sub-list 'spec'. You can set parameter 'fmt' to 'df' in which case each compound list will be coerced to a data frame row (spectra will be encoded as 'mz1:int1 mz2:int2 ...' and information regarding SMILES, adduct or formula per peak will be lost).

See Also

[Spectra::Spectra()]

Examples

```
fl <- system.file("extdata/annotated_output.mgf", package = "fioRa")

# read as data.frame (return a simplified version)
str(fioRa::read_fiora(fl = fl, fmt = "df"))

# read as a list (standard case)
fioRa::read_fiora(fl = fl)

# read as Spectra object (requires Spectra-package)
if (requireNamespace("Spectra", quietly = TRUE)) {
  fioRa::read_fiora(fl = fl, fmt = "Spectra")
}
```

`run_app`*Run the Shiny Application.*

Description

Will open a Shiny App in the local browser.

Usage

```
run_app(  
  onStart = NULL,  
  options = list(),  
  enableBookmarking = NULL,  
  uiPattern = "/",  
  ...  
)
```

Arguments

<code>onStart</code>	A function that will be called before the app is actually run. This is only needed for <code>shinyAppObj</code> , since in the <code>shinyAppDir</code> case, a global <code>.R</code> file can be used for this purpose.
<code>options</code>	Named options that should be passed to the <code>runApp</code> call (these can be any of the following: "port", "launch.browser", "host", "quiet", "display.mode" and "test.mode"). You can also specify width and height parameters which provide a hint to the embedding environment about the ideal height/width for the app.
<code>enableBookmarking</code>	Can be one of "url", "server", or "disable". The default value, <code>NULL</code> , will respect the setting from any previous calls to <code>enableBookmarking()</code> . See <code>enableBookmarking()</code> for more information on bookmarking your app.
<code>uiPattern</code>	A regular expression that will be applied to each GET request to determine whether the <code>ui</code> should be used to handle the request. Note that the entire request path must match the regular expression in order for the match to be considered successful.
<code>...</code>	arguments to pass to <code>golem_opts</code> . See <code>'?golem::get_golem_options'</code> for more details.

Value

A `shinyApp` object. Will open a Shiny App in the local browser.

run_script	<i>Predict MS² fragment spectra from SMILES code.</i>
------------	--

Description

A wrapper around the python script 'fiora-predict' using the fiora open source model to generate a MS² spectra for a compound with known SMILES code.

Usage

```
run_script(
  x = data.frame(Name = "Example_0", SMILES = "CC1=CC(=O)OC2=CC(OS(O)(=O)=O)=CC=C12",
    Precursor_type = "[M-H]-", CE = 17, Instrument_type = "HCD"),
  min_prob = 0.001,
  annotation = FALSE,
  fiora_script = NULL,
  fmt = c("list", "df", "Spectra", "file_only"),
  file_out = NULL,
  verbose = TRUE
)
```

Arguments

x	A data frame containing columns Name (compound names), SMILES (SMILES code of the compounds, Precursor_type (currently "[M-H]-" or "[M+H]+"), CE (Collision energy) and Instrument_type (i.e. HCD).
min_prob	Minimum peak probability to be recorded in the spectrum.
annotation	Return SMILES for fragments if TRUE.
fiora_script	Path to python script fiora-predict.
fmt	Set fmt to 'df' to simplify the return value to a data frame (named list otherwise). You may also use 'file_only' to omit output to the console if you specify a valid path in 'file_out'.
file_out	Specify a path to a file to store the FIORA result permanently.
verbose	Set to FALSE to omit messages.

Details

This wrapper will generate a fiora ready input file (csv-format) based on the user parameters which is stored as a temp file. It will ensure that the current version of the fiora package is installed in a respective python environment. It will use 'system2()' to run the python script 'fiora-predict' and import its result back into R using function 'read_fiora()'. You can try different installed version of 'fiora' by providing the path the the script explicitly.

Value

A list with the fiora results for the specified compound(s).

Examples

```
## Not run:
# !!! running this example will install the python module `fiora`
td <- fioRa::test_data
x <- setNames(data.frame(
  t(sapply(td[2:11], function(x) { strsplit(x, ",")[[1]] })),
  strsplit(td[1], ",")[[1]]
)
foo <- run_script(x = x)
foo[[1]][["spec"]]
# modify parameters
run_script(x = x[1,,drop=FALSE], min_prob = 0.05)

# you may also return a Spectra object
run_script(x = x[1,,drop=FALSE], min_prob = 0.05, annotation = TRUE, fmt = "Spectra")

# use a different fiora environment/model
s_pth <- "c:/Users/jlisec/AppData/Local/r-miniconda/envs/fiora-0.1.0/Scripts/fiora-predict"

# this setup will be used internally
fioRa::find_fiora_predict_paths(default_path = dirname(s_pth), script_name = basename(s_pth))

# run the script
foo2 <- run_script(x = x, fiora_script = s_pth)

foo2[[1]][["spec"]]

for (i in 1:length(foo)) {
  cat("\n")
  print(names(foo)[i])#'
  print(foo[[i]][["spec"]])
  print(foo2[[i]][["spec"]])
}

## End(Not run)
```

test_data

The example set of test compounds provided with FIORA.

Description

The example set of test compounds provided with FIORA.

Usage

```
data(test_data)
```

Format

A character vector of length = 12 containing the readLines equivalent of a fiora input file in csv format, with header, 10 example compounds, defined via their SMILES, and an empty final line (or vector element respectively).

Source

https://github.com/BAMeScience/fiora/blob/main/examples/example_input.csv

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