## Package 'IsoCorrectoR'

November 13, 2024

**Title** Correction for natural isotope abundance and tracer purity in MS and MS/MS data from stable isotope labeling experiments

Version 1.24.0

- **Imports** dplyr, magrittr, methods, quadprog, readr, readxl, stringr, tibble, tools, utils, pracma, WriteXLS
- **Description** IsoCorrectoR performs the correction of mass spectrometry data from stable isotope labeling/tracing metabolomics experiments with regard to natural isotope abundance and tracer impurity. Data from both MS and MS/MS measurements can be corrected (with any tracer isotope: 13C, 15N, 18O...), as well as ultra-high resolution MS data from multiple-tracer experiments (e.g. 13C and 15N used simultaneously). See the Bioconductor package IsoCorrectoRGUI for a graphical user interface to IsoCorrectoR. NOTE: With R version 4.0.0, writing correction results to Excel files may currently not work on Windows. However, writing results to csv works as before.

**Depends** R (>= 3.5)

URL https://genomics.ur.de/files/IsoCorrectoR/

License GPL-3

LazyData TRUE

NeedsCompilation no

**biocViews** Software, Metabolomics, MassSpectrometry, Preprocessing, ImmunoOncology

RoxygenNote 6.1.1

Suggests IsoCorrectoRGUI, knitr, rmarkdown, testthat, BiocStyle

VignetteBuilder knitr

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IsoCorrection Algorithm For Natural Isotope Abundance And Tracer Purity Correction Of Data From Stable Isotope Labeling Experiments

#### Description

IsoCorrection is the main function of the IsoCorrectoR package. It performs the correction of mass spectrometry data from stable isotope labeling experiments with regard to natural abundance and tracer purity. Data from both MS and MS/MS experiments can be corrected (with any tracer isotope: 13C, 15N, 18O...), as well as high resolution data from multiple-tracer experiments (e.g. 13C and 15N used simultaneously).

#### Usage

```
IsoCorrection(MeasurementFile = NA, ElementFile = NA, MoleculeFile = NA,
CorrectTracerImpurity = FALSE, CorrectTracerElementCore = TRUE,
CalculateMeanEnrichment = TRUE, UltraHighRes = FALSE, DirOut = ".",
FileOut = "result", FileOutFormat = "csv", ReturnResultsObject = TRUE,
CorrectAlsoMonoisotopic = FALSE, CalculationThreshold = 10^-8,
CalculationThreshold_UHR = 8, verbose = FALSE, Testmode = FALSE)
```

#### Arguments

	MeasurementFile	2							
		Required. The file that contains the measured data to be corrected. Only ".xls", ".xlsx" and ".csv" file formats are supported.							
	ElementFile	Required. The file that contains the element information required for correction. Only ".xls", ".xlsx" and ".csv" file formats are supported.							
	MoleculeFile	Required. The file that contains the information on the molecules for which data is to be corrected. Only ".xls", ".xlsx" and ".csv" file formats are supported.							
CorrectTracerImpurity									
		Logical. If TRUE, correction for isotopic impurity of the tracer substrate is performed.							
	CorrectTracerElementCore								
		Logical. If TRUE (recommended!), the tracer element atoms in the core module (usually the part of the molecule that does not come from derivatization) are considered when correcting.							
	CalculateMeanEr	richment							
		Logical. If TRUE, the mean isotopic enrichment is calculated for each molecule.							
	UltraHighRes	Logical. If TRUE, high resolution correction is performed on the data. Should only be set to TRUE, if you know that you have high resolution data.							
	DirOut	Character String. Defines the directory the corrected data and log-file should be written to. Default directory is set to current working directoy ('.').							

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FileOut	Character String. Defines the name of the file that contains the corrected data. The name of the fill will be IsoCorrectoR_ <fileout>.<fileformat>. If the format is set to "csv", the name will also contain the type of the corrected data in the respective file.</fileformat></fileout>									
FileOutFormat	Character String. Defines the format of the files that contain the corrected data. Can either be "csv" or "xls". If set to "csv", multiple files will be generated, one for each type of corrected data (eg. corrected data, fractions, mean enrich- ment). If set to "xls", all correction results are provided in one excel file in different sheets.									
ReturnResults0	ReturnResultsObject									
	Logical. If TRUE, the correction results are returned as a list in the current									
	R_session in addition to writing the results to a file. This is useful if the corrected data has to be further processed directly in R.									
CorrectAlsoMon	CorrectAlsoMonoisotopic									
	Logical. If TRUE, monoisotopic correction results are also provided.									
CalculationThr	CalculationThreshold									
	(Advanced Option) Numeric. Defines a threshold to stop probability calcula- tions at for making correction faster (normal resolution mode). Should be left at the default value.									
CalculationThr	eshold_UHR									
	(Advanced Option) Numeric. Defines a threshold to stop probability calcula- tions at for making correction faster (high resolution mode). Should be left at the default value.									
verbose	Logical. If TRUE, status messages are sent to standard output.									
Testmode	Logical. If TRUE, starts a testmode for development purposes. Not required for users of IsoCorrectoR.									

#### Value

The function returns a list with 4 elements

- success: string that is "TRUE" if the correction was successful, "FALSE" if an error occured and "WARNINGS" if warnings occured
- **results:** a list containing a dataframe for each type of corrected data (normal, fractions, mean enrichment ...). Will be NA if ReturnResultsObject is set to FALSE
- **log:** list containing log information on the correction run (parameters, file names and paths, warnings and errors)
- error: contains a string with the associated error message if an error occurred, empty otherwise

#### References

See Reference 1 LinktoIsoCorrectoR-Paper

#### Examples

```
# Normal resolution data
```

```
# 1) get path of example files
path.molecule <- system.file("extdata","normal_resolution","MoleculeFile.csv",
    package = "IsoCorrectoR", mustWork = TRUE);
path.element <- system.file("extdata","normal_resolution","ElementFile.csv",
    package = "IsoCorrectoR", mustWork = TRUE);</pre>
```

```
path.measurement <- system.file("extdata", "normal_resolution", "MeasurementFile.csv",</pre>
      package = "IsoCorrectoR", mustWork = TRUE);
   # 2) run correction algorithm and save results in variable
   correctionResults <- IsoCorrection(MeasurementFile=path.measurement,</pre>
      ElementFile=path.element,
     MoleculeFile=path.molecule)
# High resolution data
   # 1) get path of example files
   path.molecule <- system.file("extdata", "high_resolution", "MoleculeFile.csv",</pre>
      package = "IsoCorrectoR", mustWork = TRUE);
   path.element <- system.file("extdata", "high_resolution", "ElementFile.csv",</pre>
      package = "IsoCorrectoR", mustWork = TRUE);
   path.measurement <- system.file("extdata", "high_resolution", "MeasurementFile.csv",</pre>
      package = "IsoCorrectoR", mustWork = TRUE);
   # 2) run correction algorithm and save results in variable
   correctionResults <- IsoCorrection(MeasurementFile=path.measurement,</pre>
      ElementFile=path.element,
```

```
MoleculeFile=path.molecule,UltraHighRes=TRUE)
```

IsoCorrectoR

IsoCorrectoR example data

#### Description

Example data for the IsoCorrectoR package

#### Usage

data(IsoCorrectoR)

#### Format

A list with four elements, namely

- **tool\_features:** overview of correction features as provided by various tools IsoCorrectoR is compared with.
- **element\_file:** example data for an Element file. This information is independent of the resolution type (normal or high)
- **normal\_resolution:** list with two elements, containing examples for normal resolution data: molecule information and measurement data
- **high\_resolution:** list with two elements, containing examples for high resolution data: molecule information and measurement data

#### Author(s)

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