Rdisop

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`RcppVersion` *Rcpp Version and License Information*

Description

`RcppVersion` displays the version of Rcpp/RcppTemplate that was used to build this package.

Usage

```
RcppVersion()
```

Author(s)

Dominick Samperi

Examples

```
RcppVersion()
```

`addMolecules` *Add/subtract sum formulae*

Description

Simple arithmetic modifications of sum formulae.

Usage

```
addMolecules(formula1, formula2, elements = NULL)
subMolecules(formula1, formula2, elements = NULL)
```

Arguments

`formula1`	Sum formula
`formula2`	Sum formula
`elements`	list of allowed chemical elements, defaults to full periodic system of elements

Details

addMolecules() adds the second argument to the first. subMolecules() subtracts the second argument from the first.

This can be useful to revert e.g. adduct/fragment formation found in ESI mass spectrometry, or to mimick simple chemical reactions. No chemical checks are performed.

Value

	A list with the elements
formula	repeated sum formula
mass	exact mass of molecule
score	dummy value, always 1.0
isotopes	a list of isotopes

Author(s)

Steffen Neumann <sneumann@IPB-Halle.DE>

Examples

```
# For proton-Adduct of Ethanol:
subMolecules("C2H7O", "H")
```

---

decomposeIsotopes *Mass Decomposition of Isotope Patterns*

Description

Calculate the elementary compositions from an exact Mass or Isotope Pattern, obtained e.g. by FTICR or TOF mass spectrometers

Usage

```
decomposeMass(mass, ppm=2.0, mzabs=0.0001, elements=NULL, filter=NULL, z=0)
decomposeIsotopes(masses, intensities, ppm=2.0, mzabs=0.0001, elements=NULL, fil
isotopeScore(molecule, masses, intensities, elements = NULL, filter = NULL, z =
```

Arguments

mass	A single exact mass (or m/z value)
masses	A vector of masses (or m/z values) of an isotope cluster
intensities	Abolute or relative intensities of the masses peaks
ppm	allowed deviation of hypotheses from given mass
mzabs	absolute deviation in dalton (mzabs and ppm will be added)
z	charge z of m/z peaks for calculation of real mass. 0 is for auto-detection
elements	list of allowed chemical elements, defaults to CHNOPS
filter	NYI, will be a selection of DU, DBE and Nitrogen rules
molecule	a molecule as obtained from getMolecule() or decomposeMass / decomposeIsotopes

Details

Sum formulas are calculated which explain the given mass or isotope pattern.

Value

A list of molecules, which contain the sub-lists

`formula`	potential formulae
`mass`	exact mass of hypothesis
`score`	calculated score
`isotopes`	a list of isotopes

Author(s)

Steffen Neumann <sneumann@IPB-Halle.DE>

References

For a description of the underlying IMS see: WABI Paper

See Also

`decomposeMass`

Examples

```
# For Glutamate:
decomposeIsotopes(c(147.0529,148.0563), c(100.0,5.561173))
```

---

getMolecule *Calculate mass and isotope information for a molecule given as sum formula*

Description

Parse the sum formula and calculate the theoretical exact mass and the isotope distribution.

Usage

```
getMolecule(formula, elements = NULL, z = 0)
getMass(molecule)
getFormula(molecule)
getIsotope(molecule, index)
getScore(molecule)
getValid(molecule)
```

Arguments

formula	Sum formula
elements	list of allowed chemical elements, defaults to full periodic system of elements
z	charge z of molecule for exact mass calculation
molecule	an initialized molecule as returned by getMolecule() or the decomposeMass() and decomposeIsotope() functions
index	return the n-th isotope mass/abundance pair of the molecule

Details

getMolecule() Parse the sum formula and calculate the theoretical exact mass and the isotope distribution. For a given element, return the different mass values.

Value

getMolecule	A list with the elements
formula	repeated sum formula
mass	exact mass of molecule
score	probability, for given molecules a dummy value which is always 1.0
valid	result of neutrogen rule check
isotopes	a list of isotopes
getMass, getFormula and getScore	return the mass of the molecule as string or real value

Author(s)

Steffen Neumann <sneumann@IPB-Halle.DE>

References

For a description of the underlying IMS see: WABI Paper

Examples

```
# For Ethanol:
getMolecule("C2H6O")
```

---

initializeCHNOPS *Initialize (a subset of) elements of the periodic system of elements (PSE)*

Description

Initialize the information about name, mass and isotopes. To reduce the number of decomposition hypotheses, subsets of elements can be created.

Usage

```
initializeCHNOPS()
initializeCHNOPSMgKCaFe()
initializePSE()
initializeElements(names)
```

Arguments

names	vector of element names within PSE

Details

These functions return full, pre-defined or user-defined (sub-) lists of elements.

Value

A list with the elements

formula	repeated sum formula
mass	exact mass of molecule
isotopes	a list of isotopes

The initializeCharges() is special, since it allows to parse charges such as `getMolecule("H3O+", elements=c(initializeCHNOPS(),initializeCharges()))`

Author(s)

Steffen Neumann <sneumann@IPB-Halle.DE>

References

For a description of the underlying IMS see: WABI Paper

Isotope patterns obtained through wikipedia.org

See Also

`getMolecule`

Examples

```
# For Ethanol:
elements <- initializeCHNOPS()
```

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