

Package ‘RBMRB’

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Type Package

Title BMRB Data Access and Visualization

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Description

The Biological Magnetic Resonance Data Bank (BMRB, <<http://www.bmrwisc.edu/>>) collects, annotates, archives, and disseminates (worldwide in the public domain) the important spectral and quantitative data derived from NMR (Nuclear Magnetic Resonance) spectroscopic investigations of biological macromolecules and metabolites. This package provides an interface to BMRB database for easy data access and includes a minimal set of data visualization functions. Users are encouraged to make their own data visualizations using BMRB data.

Depends R(>= 3.3.0)

License GPL-2

URL <https://github.com/uwbmrwisc/RBMRB>,
<https://github.com/kumar-physics/RBMRB>

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chemical_shift_corr *Chemical shift correlation between any two atoms from a single residue*

Description

Plots the correlated chemical shift distribution of any two atoms in a single residue for the 20 standard amino acids from BMRB database. By default it will generate interactive graphics using plotly library

Usage

```
chemical_shift_corr(atom1, atom2, res = NA, type = "c",
  interactive = TRUE)
```

Arguments

atom1	atom name in NMR-STAR nomenclature like CA,CB2
atom2	atom name in NMR_STAR nomenclature like HA,HB2
res	residue name like ALA,GLY (optional by default includes all possible amino acids)
type	'c' for contour plot and 's' for scatter plot default 'c'.scatter plot will be slow and heavy for large data set
interactive	TRUE/FALSE default=TRUE

Value

plot object

See Also

[fetch_atom_chemical_shifts](#)

Examples

```
plt<-chemical_shift_corr('HE21','HE22')
#plots the chemical shift distribution between HE21 and HE22
```

chemical_shift_hist *Plots chemical shift distribution*

Description

Plots the histogram (or) density of chemical shift distribution of a given amino acid (or) nucleic acid from BMRB database. Optionally particular atom can be specified in the parameter

Usage

```
chemical_shift_hist(res = "*", atm = "*", type = "count", bw = 0.1,
  cutoff = 8, interactive = TRUE)
```

Arguments

res	residue name in NMR-STAR atom nomenclature ; Example: ALA, GLY ; default '*' (includes everything)
atm	atom name in NMR-STAR nomenclature ; Example :CA, HB2 default '*' (includes all atoms)
type	count ; other than count will assume density plot
bw	binwidth for histogram; default value 0.1ppm
cutoff	values not within the cutoff time standard deviation from both sides of the mean will be excluded from the plot; default value 8
interactive	TRUE/FALSE default TRUE

Value

R plot object

See Also

[fetch_res_chemical_shifts](#), [filter_residue](#) and [chemical_shift_corr](#)

Examples

```
#plt<-chemical_shift_hist('ALA')
#plots the histogram of all atoms of ALA
#plt<-chemical_shift_hist('GLY', type='density')
#plots the density plot
```

chemical_shift_hists *Plots chemical shift distribution for a list of atoms*

Description

Plots the histogram (or) density of chemical shift distribution of a given list of atoms. Atoms from different residues can be specified as "residue-atom". Exammple "ALA-CA","GLN-HE21","GLN-HE*"

Usage

```
chemical_shift_hists(atm = NA, type = "count", bw = 0.1, cutoff = 8,  
  interactive = TRUE)
```

Arguments

atm	list Example: c("ALA-CA","GLY-CA")
type	count ; other than count will assume density plot
bw	binwidth for histogram; default value 0.1ppm
cutoff	values not with in the cutoff time standard deviation from both sides ofthe mean will be excluded from the plot;default value 8
interactive	TRUE/FALSE default TRUE

Value

R plot object

See Also

[fetch_res_chemical_shifts](#),[filter_residue](#) and [chemical_shift_corr](#)

Examples

```
#plt<-chemical_shift_hists(c('ALA-C*'))  
#plots the histogram of all atoms of ALA  
#plt<-chemical_shift_hists(c("GLY-H*","ALA-HA"),type='density')  
#plots the density plot
```

convert_cs_to_c13hsqc *Reformats chemical shift dataframe for easy plotting*

Description

Reformats the output dataframe from [fetch_entry_chemical_shifts](#) into a simple dataframe that contains proton and carbon chemical shifts in two columns. This will be helpful to plot 1H-13C HSQC(Hetronuclear Single Quantum Coherence) spectrum

Usage

```
convert_cs_to_c13hsqc(csdf)
```

Arguments

csdf chemical shift data frame from [fetch_entry_chemical_shifts](#)

Value

R data frame that contains proton and carbon chemical shifts in two columns for each residue

See Also

[convert_cs_to_n15hsqc](#) and [convert_cs_to_tocsy](#)

Examples

```
df<-fetch_entry_chemical_shifts(15060)
# Downloads data from BMRB
hsqc<-convert_cs_to_c13hsqc(df)
# Reformats for easy plotting
```

convert_cs_to_n15hsqc *Reformats chemical shift dataframe for easy plotting*

Description

Reformats the output dataframe from [fetch_entry_chemical_shifts](#) into a simple dataframe that contains algorithmically combined proton and nitrogen chemical shifts in two columns. This will be helpful to plot 1H-15N HSQC(Hetronuclear Single Quantum Coherence) spectrum.

Usage

```
convert_cs_to_n15hsqc(csdf)
```

Arguments

csdf Chemical shift data frame from [fetch_entry_chemical_shifts](#)

Value

R data frame that contains proton and nitrogen chemical shifts in two columns for each residue

See Also

[convert_cs_to_c13hsqc](#) and [convert_cs_to_tocsy](#)

Examples

```
df<-fetch_entry_chemical_shifts(15060)
#Downloads the chemical shift data from BMRB
hsqc<-convert_cs_to_n15hsqc(df)
#Reformats for easy plotting
```

`convert_cs_to_tocsy` *Reformats chemical shift dataframe for easy plotting*

Description

Reformats the output dataframe from [fetch_entry_chemical_shifts](#) into a simple dataframe that contains algorithmically combined proton shifts in two columns. This will be helpful to plot TOCSY(TOtal Correlation Spectroscopy) spectrum

Usage

```
convert_cs_to_tocsy(csdf)
```

Arguments

csdf chemical shift data frame from [fetch_entry_chemical_shifts](#)

Value

R data frame that contains all possible combinations of proton chemical shifts in two columns

See Also

[convert_cs_to_c13hsqc](#) and [convert_cs_to_n15hsqc](#)

Examples

```
df<-fetch_entry_chemical_shifts(15060)
# Downloads data from BMRB
tocsy<-convert_cs_to_tocsy(df)
# Reformats for easy plotting
```

export_star_data	<i>Exports NMR-STAR file to BMRB API server</i>
------------------	---

Description

Exports NMR-STAR file to BMRB API server, so that local data can be visualized using RBMRB library. This function will return a token, which can be used to access the data through RBMRB package. The token will expire after 7 days

Usage

```
export_star_data(filename)
```

Arguments

filename	filename with correct path
----------	----------------------------

Value

Temporary token to access the file

See Also

[fetch_atom_chemical_shifts](#), [fetch_entry_chemical_shifts](#) [fetch_res_chemical_shifts](#)

Examples

```
# ent_id <- export_star_data('/nmrdata/hpr.str')  
# Exports hpr.str file to BMRB API server and gets a temporary token
```

fetch_atom_chemical_shifts	<i>Fetchs atom specific NMR chemical shift data from BMRB database</i>
----------------------------	--

Description

Downloads the full list of chemical shifts from BMRB macromolecular/metabolomics database for a given atom

Usage

```
fetch_atom_chemical_shifts(atom, db = "macromolecules")
```

Arguments

atom atom name in NMR-STAR atom nomenclature ; Example: CA,CB2
 db macromolecules, metabolomics (optional, by default will fetch from macromolecules database)

Value

R data frame that contains full chemical shift list for a given atom

See Also

[fetch_entry_chemical_shifts](#), [fetch_res_chemical_shifts](#), [filter_residue](#) and [chemical_shift_corr](#)

Examples

```
df<-fetch_atom_chemical_shifts('CG2','macromolecules')
# Downloads CB2 chemical shifts from macromolecules database at BMRB
df<-fetch_atom_chemical_shifts('C1','metabolomics')
# Downloads C1 chemical shifts from metabolomics database at BMRB
```

fetch_entry_chemical_shifts

Fetchs entry specific NMR chemical shift data from BMRB database

Description

Downloads NMR chemical shift data from BMRB database for a given Entry ID or list of Entry IDs

Usage

```
fetch_entry_chemical_shifts(BMRBidlist)
```

Arguments

BMRBidlist sinlge BMRB ID (or) list of BMRB IDs in csv format For metabolomics entries
 entry id should have 'bmse' prefix example: c('bmse000034','bmse000035','bmse000036')

Value

R data frame that contains Atom_chem_shift data for a given list of entries

See Also

[fetch_atom_chemical_shifts](#) and [fetch_res_chemical_shifts](#)

Examples

```
df<-fetch_entry_chemical_shifts(15060)
# Downloads NMR chemical shifts of a single entry from BMRB
df<-fetch_entry_chemical_shifts(c(17074,17076,17077))
# Downloads NMR chemical shifts of multiple entries from BMRB
df<-fetch_entry_chemical_shifts(c('bmse000034','bmse000035','bmse000036'))
# Downloads data from BMRB metabolomics database
```

fetch_res_chemical_shifts

Fetchs residue specific NMR chemical shift data from BMRB database

Description

Downloads the full list of chemical shifts from BMRB macromolecular database for a given amino acid (or) nucleic acid. Optionally particular atom can be specified in the parameter

Usage

```
fetch_res_chemical_shifts(res = "*", atm = "*")
```

Arguments

res	residue name in NMR-STAR atom nomenclature ; Example: ALA, GLY ; default '*' (includes everything)
atm	atom name in NMR-STAR nomenclature ; Example :CA, HB2; default * (includes all atoms)

Value

R data frame that contains full chemical shift list for a given atom

See Also

[fetch_atom_chemical_shifts](#), [filter_residue](#) and [chemical_shift_hist](#)

Examples

```
#df<-fetch_res_chemical_shifts('GLY')
# Downloads chemical shift data of all atoms of GLY
#df<-fetch_res_chemical_shifts('ALA','CA')
# Downloads C alpha chemical shifts of ALA from macromolecules database at BMRB
```

filter_residue	<i>Filter for standard 20 amino acids</i>
----------------	---

Description

Filters out non standard amino acids using Comp_ID. The data frame should contain three letter amino acid code in COMP_ID column.

Usage

```
filter_residue(df)
```

Arguments

df	data frame with amino acid information in Comp_ID column
----	--

Value

R data frame that contains information from only standard 20 amino acids.

See Also

[fetch_atom_chemical_shifts](#)

Examples

```
df<-filter_residue(fetch_atom_chemical_shifts("CG2"))
#Downloads all CG2 chemical shifts and removes non standard amino acids
```

HSQC_13C	<i>Simulates H1-C13 HSQC spectra for a given entry or list of entries from BMRB</i>
----------	---

Description

Simulates H1-C13 HSQC(Hetronuclear Single Quantum Coherence) spectra directly from BMRB database. By default it will generate interactive graphics using plotly library

Usage

```
HSQC_13C(idlist, type = "scatter", interactive = TRUE)
```

Arguments

idlist	list of bmr ids in csv
type	scatter/line default=scatter
interactive	TRUE/FALSE default=TRUE

Value

R plot object

See Also

[HSQC_15N](#) and [TOCSY](#)

Examples

```
plot_hsqc<-HSQC_13C(c(17074,17076,17077))
#Simulates C13-HSQC spectra form the given list of entries
plot_hsqc<-HSQC_13C(c(17074,17076,17077),'line')
#Simulates C13-HSQC and connects the peaks with same sequence number
plot_hsqc<-HSQC_13C(c(17074,17076,17077),interactive=FALSE)
#Example for non interactive plot
```

HSQC_15N	<i>Simulates H1-N15 HSQC spectra for a given entry or list of entries from BMRB</i>
----------	---

Description

Simulates H1-N15 HSQC(Hetronuclear Single Quantum Coherence) spectra directly from BMRB database. Default plot type will be 'scatter'.Peaks from different spectra(entries) can be connected based on residue numbers by specifying plot type as 'line'. By default it will generate interactive graphics using plotly library

Usage

```
HSQC_15N(idlist, type = "scatter", interactive = TRUE)
```

Arguments

idlist	list of bmr ids in csv
type	scatter/line default=scatter
interactive	TRUE/FALSE default=TRUE

Value

R plot object

See Also

[HSQC_13C](#) and [TOCSY](#)

Examples

```

plot_hsqc<-HSQC_15N(c(17074,17076,17077))
#simulates N15-HSQC spectra for the given 3 entreis
plot_hsqc<-HSQC_15N(18857,'line')
#simulates the N15-HSQC spectra from many chemical shift lists from a single entry
plot_hsqc<-HSQC_15N(c(17074,17076,17077),interactive=FALSE)
#example for non interactive plots

```

makeRandomString	<i>Generates random string of fixed length(for internal use in RBMRB)</i>
------------------	---

Description

Local files may not have Entry_ID, in that case random Entry_ID is assigned using this function. It is an internal function used only by RBMRB package

Usage

```
makeRandomString()
```

TOCSY	<i>Simulates TOCSY spectra for a given entry or a list of entries from BMRB</i>
-------	---

Description

Simulates TOCSY(TOtal Correlation SpectroscopY) spectra directly from BMRB database. By default it will generate interactive graphics using plotly library

Usage

```
TOCSY(idlist, interactive = TRUE)
```

Arguments

idlist	list of bmr ids c(17074,17076,17077)
interactive	TRUE/FALSE default=TRUE

Value

plot object

See Also

[HSQC_15N](#) and [HSQC_13C](#)

Examples

```
plot_tocsy<-TOCSY(c(17074,17076,17077))  
#Simulates TOCSY spectra for the given 3 entries  
plot_tocsy<-TOCSY(c(17074,17076,17077),interactive=FALSE)  
# Example to disable interactive plot feature
```

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