

Package ‘msentropy’

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Type Package

Title Spectral Entropy for Mass Spectrometry Data

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Description Clean the MS/MS spectrum, calculate spectral entropy, unweighted entropy similarity, and entropy similarity for mass spectrometry data. The entropy similarity is a novel similarity measure for MS/MS spectra which outperform the widely used dot product similarity in compound identification. For more details, please refer to the paper: Yuanyue Li et al. (2021) ``Spectral entropy outperforms MS/MS dot product similarity for small-molecule compound identification" <[doi:10.1038/s41592-021-01331-z](https://doi.org/10.1038/s41592-021-01331-z)>.

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Suggests testthat

LinkingTo Rcpp

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URL <https://github.com/YuanyueLi/MSEntropy>

NeedsCompilation yes

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Contents

calculate_entropy_similarity	2
calculate_spectral_entropy	3
calculate_unweighted_entropy_similarity	4
clean_spectrum	5
msentropy_similarity	6

Index	9
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`calculate_entropy_similarity`*Entropy similarity between two spectra*

Description

Calculate the entropy similarity between two spectra

Usage

```
calculate_entropy_similarity(  
    peaks_a,  
    peaks_b,  
    ms2_tolerance_in_da,  
    ms2_tolerance_in_ppm,  
    clean_spectra,  
    min_mz,  
    max_mz,  
    noise_threshold,  
    max_peak_num  
)
```

Arguments

<code>peaks_a</code>	A matrix of spectral peaks, with two columns: m/z and intensity
<code>peaks_b</code>	A matrix of spectral peaks, with two columns: m/z and intensity
<code>ms2_tolerance_in_da</code>	The MS2 tolerance in Da, set to -1 to disable
<code>ms2_tolerance_in_ppm</code>	The MS2 tolerance in ppm, set to -1 to disable
<code>clean_spectra</code>	Whether to clean the spectra before calculating the entropy similarity, see clean_spectrum
<code>min_mz</code>	The minimum m/z value to keep, set to -1 to disable
<code>max_mz</code>	The maximum m/z value to keep, set to -1 to disable
<code>noise_threshold</code>	The noise threshold, set to -1 to disable, all peaks have intensity < noise_threshold * max_intensity will be removed
<code>max_peak_num</code>	The maximum number of peaks to keep, set to -1 to disable

Value

The entropy similarity

Examples

```
mz_a <- c(169.071, 186.066, 186.0769)
intensity_a <- c(7.917962, 1.021589, 100.0)
mz_b <- c(120.212, 169.071, 186.066)
intensity_b <- c(37.16, 66.83, 999.0)
peaks_a <- matrix(c(mz_a, intensity_a), ncol = 2, byrow = FALSE)
peaks_b <- matrix(c(mz_b, intensity_b), ncol = 2, byrow = FALSE)
calculate_entropy_similarity(peaks_a, peaks_b,
  ms2_tolerance_in_da = 0.02, ms2_tolerance_in_ppm = -1,
  clean_spectra = TRUE, min_mz = 0, max_mz = 1000,
  noise_threshold = 0.01,
  max_peak_num = 100)
```

calculate_spectral_entropy

Calculate spectral entropy of a spectrum

Description

Calculate spectral entropy of a spectrum

Usage

```
calculate_spectral_entropy(peaks)
```

Arguments

peaks A matrix of peaks, with two columns: m/z and intensity.

Value

A double value of spectral entropy.

Examples

```
mz <- c(100.212, 300.321, 535.325)
intensity <- c(37.16, 66.83, 999.0)
peaks <- matrix(c(mz, intensity), ncol = 2, byrow = FALSE)
calculate_spectral_entropy(peaks)
```

`calculate_unweighted_entropy_similarity`*Unweighted entropy similarity between two spectra*

Description

Calculate the unweighted entropy similarity between two spectra

Usage

```
calculate_unweighted_entropy_similarity(  
    peaks_a,  
    peaks_b,  
    ms2_tolerance_in_da,  
    ms2_tolerance_in_ppm,  
    clean_spectra,  
    min_mz,  
    max_mz,  
    noise_threshold,  
    max_peak_num  
)
```

Arguments

<code>peaks_a</code>	A matrix of spectral peaks, with two columns: m/z and intensity
<code>peaks_b</code>	A matrix of spectral peaks, with two columns: m/z and intensity
<code>ms2_tolerance_in_da</code>	The MS2 tolerance in Da, set to -1 to disable
<code>ms2_tolerance_in_ppm</code>	The MS2 tolerance in ppm, set to -1 to disable
<code>clean_spectra</code>	Whether to clean the spectra before calculating the entropy similarity, see clean_spectrum
<code>min_mz</code>	The minimum m/z value to keep, set to -1 to disable
<code>max_mz</code>	The maximum m/z value to keep, set to -1 to disable
<code>noise_threshold</code>	The noise threshold, set to -1 to disable, all peaks have intensity < noise_threshold * max_intensity will be removed
<code>max_peak_num</code>	The maximum number of peaks to keep, set to -1 to disable

Value

The unweighted entropy similarity

Examples

```

mz_a <- c(169.071, 186.066, 186.0769)
intensity_a <- c(7.917962, 1.021589, 100.0)
mz_b <- c(120.212, 169.071, 186.066)
intensity_b <- c(37.16, 66.83, 999.0)
peaks_a <- matrix(c(mz_a, intensity_a), ncol = 2, byrow = FALSE)
peaks_b <- matrix(c(mz_b, intensity_b), ncol = 2, byrow = FALSE)
calculate_unweighted_entropy_similarity(peaks_a, peaks_b,
                                       ms2_tolerance_in_da = 0.02, ms2_tolerance_in_ppm = -1,
                                       clean_spectra = TRUE, min_mz = 0, max_mz = 1000,
                                       noise_threshold = 0.01,
                                       max_peak_num = 100)

```

clean_spectrum	<i>Clean a spectrum</i>
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Description

Clean a spectrum

This function will clean the peaks by the following steps: 1. Remove empty peaks (mz <= 0 or intensity <= 0). 2. Remove peaks with mz >= max_mz or mz < min_mz. 3. Centroid the spectrum by merging peaks within min_ms2_difference_in_da or min_ms2_difference_in_ppm. 4. Remove peaks with intensity < noise_threshold * max_intensity. 5. Keep only the top max_peak_num peaks. 6. Normalize the intensity to sum to 1.

Note: The only one of min_ms2_difference_in_da and min_ms2_difference_in_ppm should be positive.

Usage

```

clean_spectrum(
  peaks,
  min_mz,
  max_mz,
  noise_threshold,
  min_ms2_difference_in_da,
  min_ms2_difference_in_ppm,
  max_peak_num,
  normalize_intensity
)

```

Arguments

peaks	A matrix of spectral peaks, with two columns: mz and intensity
min_mz	The minimum mz value to keep, set to -1 to disable
max_mz	The maximum mz value to keep, set to -1 to disable

noise_threshold The noise threshold, set to -1 to disable, all peaks have intensity < noise_threshold * max_intensity will be removed

min_ms2_difference_in_da The minimum mz difference in Da to merge peaks, set to -1 to disable, any two peaks with mz difference < min_ms2_difference_in_da will be merged

min_ms2_difference_in_ppm The minimum mz difference in ppm to merge peaks, set to -1 to disable, any two peaks with mz difference < min_ms2_difference_in_ppm will be merged

max_peak_num The maximum number of peaks to keep, set to -1 to disable

normalize_intensity Whether to normalize the intensity to sum to 1

Value

A matrix of spectral peaks, with two columns: mz and intensity

Examples

```
mz <- c(100.212, 169.071, 169.078, 300.321)
intensity <- c(0.3716, 7.917962, 100., 66.83)
peaks <- matrix(c(mz, intensity), ncol = 2, byrow = FALSE)
clean_spectrum(peaks, min_mz = 0, max_mz = 1000, noise_threshold = 0.01,
               min_ms2_difference_in_da = 0.02, min_ms2_difference_in_ppm = -1,
               max_peak_num = 100, normalize_intensity = TRUE)
```

msentropy_similarity *Calculate spectral entropy similarity between two spectra*

Description

msentropy_similarity calculates the spectral entropy between two spectra (Li et al. 2021). It is a wrapper function defining defaults for parameters and calling the [calculate_entropy_similarity\(\)](#) or [calculate_unweighted_entropy_similarity\(\)](#) functions to perform the calculation.

Usage

```
msentropy_similarity(
  peaks_a,
  peaks_b,
  ms2_tolerance_in_da = 0.02,
  ms2_tolerance_in_ppm = -1,
  clean_spectra = TRUE,
  min_mz = 0,
  max_mz = 1000,
  noise_threshold = 0.01,
```

```

    max_peak_num = 100,
    weighted = TRUE,
    ...
)

```

Arguments

peaks_a	A two-column numeric matrix with the m/z and intensity values for peaks of one spectrum.
peaks_b	A two-column numeric matrix with the m/z and intensity values for peaks of one spectrum.
ms2_tolerance_in_da	The MS2 tolerance in Da, set to -1 to disable. Defaults to <code>ms2_tolerance_in_da = 0.02</code> .
ms2_tolerance_in_ppm	The MS2 tolerance in ppm, set to -1 to disable. Defaults to <code>ms2_tolerance_in_ppm = -1</code> .
clean_spectra	Whether to clean the spectra before calculating the entropy similarity, see <code>clean_spectrum()</code> .
min_mz	The minimum mz value to keep, set to -1 to disable. Defaults to <code>min_mz = 0</code> .
max_mz	The maximum mz value to keep, set to -1 to disable. Defaults to <code>max_mz = 1000</code> .
noise_threshold	The noise threshold, set to -1 to disable, all peaks have intensity $< \text{noise_threshold} * \text{max_intensity}$ will be removed. Defaults to <code>noise_threshold = 0.01</code> , thus, by default, all peaks with an intensity less than 1% of the maximum intensity of a spectrum will be removed.
max_peak_num	The maximum number of peaks to keep, set to -1 to disable. Defaults to <code>max_peak_num = 1000</code> .
weighted	<code>logical(1)</code> whether the weighted or unweighted entropy similarity should be calculated. Defaults to <code>weighted = TRUE</code> , thus <code>calculate_entropy_similarity()</code> is used for the calculation. For <code>weighted = FALSE</code> <code>calculate_unweighted_entropy_similarity()</code> is used instead.
...	Optional additional parameters (currently ignored)

Value

The entropy similarity

References

Li, Y., Kind, T., Folz, J. et al. (2021) Spectral entropy outperforms MS/MS dot product similarity for small-molecule compound identification. *Nat Methods* 18, 1524-1531. doi: [10.1038/s41592-02101331z](https://doi.org/10.1038/s41592-02101331z).

Examples

```
peaks_a <- cbind(mz = c(169.071, 186.066, 186.0769),
  intensity = c(7.917962, 1.021589, 100.0))
peaks_b <- cbind(mz = c(120.212, 169.071, 186.066),
  intensity <- c(37.16, 66.83, 999.0))
msentropy_similarity(peaks_a, peaks_b, ms2_tolerance_in_da = 0.02)
```

Index

`calculate_entropy_similarity`, 2
`calculate_entropy_similarity()`, 6, 7
`calculate_spectral_entropy`, 3
`calculate_unweighted_entropy_similarity`,
4
`calculate_unweighted_entropy_similarity()`,
6, 7
`clean_spectrum`, 2, 4, 5
`clean_spectrum()`, 7
`msentropy_similarity`, 6