Package ‘FAVA’

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Title Quantify Compositional Variability Across Relative Abundance Vectors

Version 1.0.0

Description Implements the statistic FAVA, an FST-based Assessment of Variability across vectors of relative Abundances, as well as a suite of helper functions which enable the visualization and statistical analysis of relative abundance data. The FAVA R package accompanies the paper, “Quantifying compositional variability in microbial communities with FAVA” by Morrison, Xue, and Rosenberg (2024) <doi:10.1101/2024.07.03.601929>.

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`bootstrap_fava` *Statistically compare FAVA values between pairs of relative abundance matrices.*

Description

`bootstrap_fava` uses bootstrapping to statistically compare FAVA values between pairs of relative abundance matrices. `bootstrap_fava` takes the same options as `fava`, so, as with `fava`, you can separately analyze multiple populations or groups of samples (specify `group`), and account for similarity among categories (specify `S`) or uneven weighting of rows (specify `w` or `time`). `bootstrap_fava` follows the bootstrapping procedure defined by Efron and Tibshirani (1993). Details on the bootstrapping procedure are available in the Methods section of the accompanying paper.

Usage

```
bootstrap_fava(
  relab_matrix,
  n_replicates = 1000,
  group,
  K = NULL,
  S = NULL,
  w = NULL,
  time = NULL,
  normalized = FALSE,
  seed = NULL,
  alternative = "two.sided"
)
```

Arguments

`relab_matrix` A matrix or data frame with rows containing non-negative entries that sum to 1. Each row represents a sample, each column represents a category, and each entry represents the abundance of that category in the sample. If `relab_matrix` contains any metadata, it must be on the left-hand side of the matrix, the right `K` entries of each row must sum to 1, and `K` must be specified. Otherwise, all entries of each row must sum to 1.

n_replicates	The number of bootstrap replicate matrices to generate. Default is n_replicates = 1000.
group	A string (or vector of strings) specifying the name(s) of the column(s) that describes which group(s) each row (sample) belongs to. Use if relab_matrix is a single matrix containing multiple groups of samples you wish to compare.
K	Optional; an integer specifying the number of categories in the data. Default is K=ncol(relab_matrix).
S	Optional; a K x K similarity matrix with diagonal elements equal to 1 and off-diagonal elements between 0 and 1. Entry S[i,k] for i!=k is the similarity between category and i and category k, equaling 1 if the categories are to be treated as identical and equaling 0 if they are to be treated as totally dissimilar. The default value is S = diag(ncol(relab_matrix)).
w	Optional; a vector of length I with non-negative entries that sum to 1. Entry w[i] represents the weight placed on row i in the computation of the mean abundance of each category across rows. The default value is w = rep(1/nrow(relab_matrix),nrow(rel
time	Optional; a string specifying the name of the column that describes the sampling time for each row. Include if you wish to weight FAVA by the distance between samples.
normalized	Optional; should normalized FAVA be used? Default is normalized = FALSE; use normalized = TRUE to compute normalized FAVA. FAVA can only be normalized if it is not weighted.
seed	Optional; an integer to be used as a random seed for the simulations.
alternative	Optional; do you want to do a one- or two.sided test? Default is alternative = "two.sided". If you wish to do a one-sided test, specify either alternative = "lesser" or alternative = "greater".

Value

A named list containing the following entries:

- p_values: The probability of observing the observed difference in variability between each pair of groups if there were no difference between groups. Computed as the fraction of bootstrap differences greater than or equal to the observed difference. Depends on what alternative is specified ("greater", "lesser", or "two.sided").
- bootstrap_distribution_plot: The distribution of bootstrap replicate differences in each variability value. The observed differences are shown in red. The further the red points are from 0, the more significant the statistical difference between groups.
- observed_stats: The observed diversity statistics for the groups.
- bootstrap_stats: The bootstrap replicate diversity statistics for the groups.
- bootstrap_replicates: The bootstrap replicate matrices, reported only if save_replicates = TRUE.

Examples

```
# Statistically compare values of FAVA between
# subjects in the xue_microbiome_sample data:

 boot_out = bootstrap_fava(relab_matrix = xue_microbiome_sample,
               n_replicates = 100,
               seed = 1,
```

```
                group = "subject",
                K = 524,
                S = xue_species_similarity)

# Table of P-values comparing values of FAVA between group 1 and group 2:
 boot_out$P_values

 # Plots of the bootstrap distributions of differences in FAVA between each pair of matrices,
 # and how the true observed differences (red dots) compare to the distribution.
 boot_out$bootstrap_distribution_plot
```

`fava` *Compute the Fst of a matrix of compositional vectors*

Description

This function computes the population genetic statistic Fst on any matrix with rows that sum to 1. Values of 0 are achieved when each row is a permutation of (1,0,..., 0) and at least two categories have non-zero abundance across all rows. The value equals 1 when each row is identical.

Usage

```
fava(
  relab_matrix,
  K = NULL,
  S = NULL,
  w = NULL,
  time = NULL,
  group = NULL,
  normalized = FALSE
)
```

Arguments

`relab_matrix`	A matrix or data frame with rows containing non-negative entries that sum to 1. Each row represents a sample, each column represents a category, and each entry represents the abundance of that category in the sample. If `relab_matrix` contains any metadata, it must be on the left-hand side of the matrix, the right `K` entries of each row must sum to 1, and `K` must be specified. Otherwise, all entries of each row must sum to 1.
`K`	Optional; an integer specifying the number of categories in the data. Default is `K=ncol(relab_matrix)`.
`S`	Optional; a K x K similarity matrix with diagonal elements equal to 1 and off-diagonal elements between 0 and 1. Entry `S[i,k]` for `i!=k` is the similarity between category and `i` and category `k`, equaling 1 if the categories are to be treated as identical and equaling 0 if they are to be treated as totally dissimilar. The default value is `S = diag(ncol(relab_matrix))`.
`w`	Optional; a vector of length `I` with non-negative entries that sum to 1. Entry `w[i]` represents the weight placed on row `i` in the computation of the mean abundance of each category across rows. The default value is `w = rep(1/nrow(relab_matrix),nrow(rel`

`time`	Optional; a string specifying the name of the column that describes the sampling time for each row. Include if you wish to weight FAVA by the distance between samples.
`group`	Optional; a string (or vector of strings) specifying the name(s) of the column(s) that describes which group(s) each row (sample) belongs to. Use if `relab_matrix` is a single matrix containing multiple groups of samples you wish to compare.
`normalized`	Optional; should normalized FAVA be used? Default is `normalized = FALSE`; use `normalized = TRUE` to compute normalized FAVA. FAVA can only be normalized if it is not weighted.

Value

A numeric value between 0 and 1.

Examples

```
# Compute the Fst of
# the following compositional vectors:
q1 = c(1,   0,   0,   0)
q2 = c(0.5, 0.5, 0,   0)
q3 = c(1/4, 1/4, 1/4, 1/4)
q4 = c(0,   0,   1,   0)
relative_abundances = matrix(c(q1, q2, q3, q4),
                  byrow = TRUE, nrow = 4)

fava(relative_abundances)

# Incoporating weights:

# Compute fava ignoring
# rows 2 and 3
row_weights = c(0.5, 0, 0, 0.5)
fava(relative_abundances, w = row_weights)

# Compute fava assuming that
# categories 1 and 2 are identical:
similarity_matrix = diag(4)
similarity_matrix[1,2] = 1
similarity_matrix[2,1] = 1
fava(relative_abundances, S = similarity_matrix)

# Assume categories 1 and 2 are identical AND
# ignore rows 2 and 4:
row_weights = c(0.5, 0, 0.5, 0)
fava(relative_abundances, w = row_weights, S = similarity_matrix)
```

`fava_norm` *Compute the normalized Fst of a matrix of compositional vectors*

Description

This function computes the normalized Fst given the number of rows and the mean abundance of the most abundant category. We employ the normalization employed in the FSTruct package by Morrison, Alcala, and Rosenberg (2020) doi: 10.1111/17550998.13647.

Usage

```
fava_norm(relab_matrix, K = ncol(relab_matrix))
```

Arguments

`relab_matrix`	A matrix with `I=nrow(relab_matrix)` rows, each containing `K=ncol(relab_matrix)` non-negative entries that sum to 1. If `relab_matrix` contains any metadata, it must be on the left-hand side of the matrix and the number of entries that sum to 1 (`K`) must be specified.
`K`	Optional; an integer specifying the number of categories in the data. Default is `K=ncol(relab_matrix)`.

Value

A numeric value between 0 and 1.

Examples

```
# Compute the weighted fava of
# the following compositional vectors:
q1 = c(1,   0,   0,   0)
q2 = c(0.5, 0.5, 0,   0)
q3 = c(1/4, 1/4, 1/4, 1/4)
q4 = c(0,   0,   1,   0)
relative_abundances = matrix(c(q1, q2, q3, q4),
                  byrow = TRUE, nrow = 4)

fava_norm(relative_abundances)
```

`gini_simpson` *Compute the Gini-Simpson index of a compositional vector*

Description

This function computes the Gini-Simpson index, a statistical measure of variability also known as the Gini-Simpson index, of vector of non-negative entries which sum to 1. The function returns a number between 0 and 1 which quantifies the variability of the vector. Values of 0 are achieved when the vector is a permutation of (1,0,..., 0). The value approaches 1 as the number of categories K increases when the vector is equal to (1/K, 1/K, ..., 1/K).

Usage

```
gini_simpson(q, K = length(q), S = diag(K))
```

Arguments

`q`	A vector with `K=length(q)` non-negative entries that sum to 1.
`K`	Optional; an integer specifying the number of categories in the data. Default is `K=length(q)`.

`S` Optional; a K x K similarity matrix with diagonal elements equal to 1 and off-diagonal elements between 0 and 1. Entry `S[i,k]` for `i!=k` is the similarity between category and `i` and category `k`, equalling 1 if the categories are to be treated as identical and equaling 0 if they are to be treated as totally dissimilar. The default value is `S = diag(ncol(q))`.

Value

A numeric value between 0 and 1.

Examples

```
# Compute unweighted Gini-Simpson index:
gini_simpson(q = c(0.4, 0.3, 0.3))

# Compute Gini-Simpson index assuming that
# categories 1 and 2 are identical:
similarity_matrix = diag(3)
similarity_matrix[1,2] = 1
similarity_matrix[2,1] = 1
gini_simpson(q = c(0.4, 0.3, 0.3), S = similarity_matrix)
```

`gini_simpson_mean` *Compute the mean Gini-Simpson index of the rows in a matrix of compositional vectors*

Description

This function computes the mean Gini-Simpson index, a statistical measure of variability also known as the Gini-Simpson index, of a set of vectors of non-negative entries which sum to 1. The function returns a number between 0 and 1 which quantifies the mean variability of the vectors. Values of 0 are achieved when each vector is a permutation of (1,0,..., 0). The value approaches 1 as the number of categories K increases when the vectors are equal to (1/K, 1/K, ..., 1/K).

Usage

```
gini_simpson_mean(
  relab_matrix,
  K = NULL,
  S = NULL,
  w = NULL,
  time = NULL,
  group = NULL
)
```

Arguments

`relab_matrix` A matrix or data frame with rows containing non-negative entries that sum to 1. Each row represents a sample, each column represents a category, and each entry represents the abundance of that category in the sample. If `relab_matrix` contains any metadata, it must be on the left-hand side of the matrix, the right `K` entries of each row must sum to 1, and `K` must be specified. Otherwise, all entries of each row must sum to 1.

K	Optional; an integer specifying the number of categories in the data. Default is `K=ncol(relab_matrix)`.
S	Optional; a K x K similarity matrix with diagonal elements equal to 1 and off-diagonal elements between 0 and 1. Entry `S[i,k]` for `i!=k` is the similarity between category and `i` and category `k`, equalling 1 if the categories are to be treated as identical and equaling 0 if they are to be treated as totally dissimilar. The default value is `S = diag(ncol(relab_matrix))`.
w	Optional; a vector of length `I` with non-negative entries that sum to 1. Entry `w[i]` represents the weight placed on row `i` in the computation of the mean abundance of each category across rows. The default value is `w = rep(1/nrow(relab_matrix),nrow(rel`
time	Optional; a string specifying the name of the column that describes the sampling time for each row. Include if you wish to weight FAVA by the distance between samples.
group	Optional; a string (or vector of strings) specifying the name(s) of the column(s) that describes which group(s) each row (sample) belongs to. Use if `relab_matrix` is a single matrix containing multiple groups of samples you wish to compare.

Value

A numeric value between 0 and 1.

Examples

```
# To compute the mean Gini-Simpson index of
# the following compositional vectors...
q1 = c(1,   0,   0,   0)
q2 = c(0.5, 0.5, 0,   0)
q3 = c(1/4, 1/4, 1/4, 1/4)
q4 = c(0,   0,   1,   0)

# we could compute the mean manually:
mean(sapply(list(q1, q2, q3, q4), gini_simpson))

# Or we could use gini_simpson_mean:
relative_abundances = matrix(c(q1, q2, q3, q4),
                  byrow = TRUE, nrow = 4)

gini_simpson_mean(relative_abundances)

# Incoporating weights:

# Compute mean Gini-Simpson index ignoring
# rows 2 and 3
row_weights = c(0.5, 0, 0, 0.5)
gini_simpson_mean(relative_abundances, w = row_weights)

# Compute mean Gini-Simpson index assuming that
# categories 1 and 2 are identical:
similarity_matrix = diag(4)
similarity_matrix[1,2] = 1
similarity_matrix[2,1] = 1
gini_simpson_mean(relative_abundances, S = similarity_matrix)

# Assume categories 1 and 2 are identical AND
```

```
# ignore rows 2 and 4:
row_weights = c(0.5, 0, 0.5, 0)
gini_simpson_mean(relative_abundances, w = row_weights, S = similarity_matrix)
```

`gini_simpson_pooled` *Compute the pooled Gini-Simpson index of the rows in a matrix of compositional vectors*

Description

This function computes the Gini-Simpson index of a "pooled" vector equal to `colMeans(relab_matrix)`. Values of 0 are achieved when this pooled vector is a permutation of (1,0,..., 0). The value approaches 1 as the number of categories K increases when this pooled vector is equal to (1/K, 1/K, ..., 1/K).

Usage

```
gini_simpson_pooled(
  relab_matrix,
  K = NULL,
  S = NULL,
  w = NULL,
  time = NULL,
  group = NULL
)
```

Arguments

`relab_matrix`	A matrix or data frame with rows containing non-negative entries that sum to 1. Each row represents a sample, each column represents a category, and each entry represents the abundance of that category in the sample. If `relab_matrix` contains any metadata, it must be on the left-hand side of the matrix, the right `K` entries of each row must sum to 1, and `K` must be specified. Otherwise, all entries of each row must sum to 1.
`K`	Optional; an integer specifying the number of categories in the data. Default is `K=ncol(relab_matrix)`.
`S`	Optional; a K x K similarity matrix with diagonal elements equal to 1 and off-diagonal elements between 0 and 1. Entry `S[i,k]` for `i!=k` is the similarity between category and `i` and category `k`, equalling 1 if the categories are to be treated as identical and equaling 0 if they are to be treated as totally dissimilar. The default value is `S = diag(ncol(relab_matrix))`.
`w`	Optional; a vector of length `I` with non-negative entries that sum to 1. Entry `w[i]` represents the weight placed on row `i` in the computation of the mean abundance of each category across rows. The default value is `w = rep(1/nrow(relab_matrix),nrow(rel`
`time`	Optional; a string specifying the name of the column that describes the sampling time for each row. Include if you wish to weight FAVA by the distance between samples.
`group`	Optional; a string (or vector of strings) specifying the name(s) of the column(s) that describes which group(s) each row (sample) belongs to. Use if `relab_matrix` is a single matrix containing multiple groups of samples you wish to compare.

Value

A numeric value between 0 and 1.

Examples

```
# To compute the pooled Gini-Simpson index of
# the following compositional vectors...
q1 = c(1,   0,   0,   0)
q2 = c(0.5, 0.5, 0,   0)
q3 = c(1/4, 1/4, 1/4, 1/4)
q4 = c(0,   0,   1,   0)

# we could compute the mean manually:
qPooled = (q1 + q2 + q3 + q4)/4
gini_simpson(qPooled)

# Or we could use gini_simpson_pooled:
relative_abundances = matrix(c(q1, q2, q3, q4),
                  byrow = TRUE, nrow = 4)

gini_simpson_pooled(relative_abundances)

# Incoporating weights:

# Compute pooled Gini-Simpson index ignoring
# rows 2 and 3
row_weights = c(0.5, 0, 0, 0.5)
gini_simpson_pooled(relative_abundances, w = row_weights)

# Compute pooled Gini-Simpson index assuming that
# categories 1 and 2 are identical:
similarity_matrix = diag(4)
similarity_matrix[1,2] = 1
similarity_matrix[2,1] = 1
gini_simpson_pooled(relative_abundances, S = similarity_matrix)

# Assume categories 1 and 2 are identical AND
# ignore rows 2 and 4:
row_weights = c(0.5, 0, 0.5, 0)
gini_simpson_pooled(relative_abundances, w = row_weights, S = similarity_matrix)
```

---

plot_relabund *Visualize a relative abundance matrix as a stacked bar plot.*

Description

This function enables graphical visualization of a matrix of compostional data. In the output plot, each vertical bar represents a single vector; the height of each color in the bar corresponds to the abundance of each category in that vector. Because this function produces a ggplot object, its output can be modified using standard ggplot2 syntax.

Usage

```
plot_relabund(
  relab_matrix,
  group = NULL,
  time = NULL,
  w = NULL,
  K = NULL,
  arrange = FALSE
)
```

Arguments

`relab_matrix`	A matrix with `I=nrow(relab_matrix)` rows, each containing `K=ncol(relab_matrix)` non-negative entries that sum to 1. If `relab_matrix` contains any metadata, it must be on the left-hand side of the matrix and the number of entries that sum to 1 (`K`) must be specified.
`group`	Optional; a string specifying the name of the column that describes which group each row (sample) belongs to. Use if `matrices` is a single matrix containing multiple groups of samples you wish to compare.
`time`	Optional; a string specifying the name of the column that describes the sampling time for each row. Include if you wish to weight FAVA by the distance between samples.
`w`	Optional; a vector of length `I` with non-negative entries that sum to 1. Entry `w[i]` represents the weight placed on row `i` in the computation of the mean abundance of each category across rows. The default value is `w = rep(1/nrow(relab_matrix),nrow(rel`
`K`	Optional; an integer specifying the number of categories in the data. Default is `K=ncol(relab_matrix)`.
`arrange`	Optional; controls horizontal ordering of samples and vertical ordering of categories. If `arrange = TRUE` or `arrange = "both"`, samples are ordered by the categories of greatest abundance and categories are ordered in decreasing abundance. If `arrange = "vertical"`, sample order is unchanged but categories are ordered in decreasing abundance. If `arrange = "horizontal"`, samples are ordered by the most abundant categories, but category order is unchanged. If `arrange` is missing or `arrange = FALSE`, neither order is changed.

Value

A ggplot object containing a bar plot visualization of the relative abundance matrix.

Examples

```
# Make an example matrix of compositional data
# Each row is an individual. Rows sum to 1.
population_A = matrix(c(
    .5, .3, .2,
    .4, .2, .4,
    .5, .4, .1,
    .6, .1, .3,
    .2, 0, .8
  ),
  nrow = 5,
  byrow = TRUE
```

```
  )

  plot_relabund(relab_matrix = population_A,
              K = 3, # How many categories per vector?
              arrange = FALSE
              )
  plot_relabund(relab_matrix = population_A,
              K = 3, # How many categories per vector?
              arrange = "horizontal"
              )
  plot_relabund(relab_matrix = population_A,
              K = 3, # How many categories per vector?
              arrange = "vertical"
              )
   plot_relabund(relab_matrix = population_A,
              K = 3, # How many categories per vector?
              arrange = TRUE  # could also be "both"
              )


# You can modify the plot as you would any ggplot2 object
plot_relabund(relab_matrix = population_A,
              K = 3, # How many categories per vector?
              arrange = TRUE
              ) +
  # Below are example, optional modifications to the default plot
  ggplot2::ggtitle("Population A") +
  ggplot2::scale_fill_brewer("Blues") +
  ggplot2::scale_color_brewer("Blues") +
  ggplot2::xlab("Individuals")
  # Note that both scale_fill and scale_color are needed to change the color of the bars.


  # Plot a dataset which has 2 populations

  population_B = matrix(c(
    .9, 0, .1,
    .6, .4, 0,
    .7, 0, .3,
    .3, .4, .3,
    .5, .3, .2
  ),
  nrow = 5,
  byrow = TRUE
  )


  populations_AB = cbind(data.frame(c("A", "A", "A", "A", "A",
                                      "B", "B", "B", "B", "B")),
                         rbind(population_A, population_B))
  colnames(populations_AB) = c("population", "category_1", "category_2", "category_3")


 plot_relabund(relab_matrix = populations_AB, group = "population")
 plot_relabund(relab_matrix = populations_AB, group = "population", arrange = "vertical")
 plot_relabund(relab_matrix = populations_AB, group = "population", arrange = "horizontal")
 plot_relabund(relab_matrix = populations_AB, group = "population", arrange = "both")
```

`relab_phyloseq` *Generate a relative abundance matrix with sample meta data and OTU abundances from a phyloseq object.*

Description

The R package phyloseq streamlines the storage and analysis of microbiome sequence data. This function takes a phyloseq object and extracts the OTU table and the sample meta data and combines them into one relative abundance matrix with rows corresponding to samples, meta data on the left-hand side, and OTU relative abundances on the right-hand side.

Usage

```
relab_phyloseq(phyloseq_object)
```

Arguments

`phyloseq_object`
: A phyloseq object containing both an OTU table (`otu_table`) and sample metadata (`sample_data`).

Value

A data frame with rows representing samples and columns representing sample data categories or OTU relative abundances.

Examples

```
otu_table = relab_phyloseq(FAVA::xue_phyloseq)
otu_table[1:10, 1:6]
```

`time_weights` *Compute a normalized weighting vector based on a vector of sampling times.*

Description

This function takes a vector of sampling times, $t = (t_1, t_2, \ldots, t_I)$ and computes a normalized vector which can be used to weight each sample based on the time between the subsequent and the preceding samples. The weighting vector w is defined such that each entry, $w_i = d_i/2T$, where $T = t_I - t_1$ and $d_i = t_{i+1} - t_{i-1}$ for i not equal to 1 or I. $d_1 = t_2 - t_1$ and $d_I = t_I - t_{I-1}$.

Usage

```
time_weights(times, group = NULL)
```

Arguments

`times`	A numeric vector of sampling times. Each entry must be greater than the previous entry.
`group`	Optional; a character vector specifying the group identity of each sampling time. Use if there are samples from multiple replicates or subjects in one data set.

Value

A numeric vector. Each entry provides a weight for each entry in the provided `times` vector. If `group` is not specified, the vector sums to 1. If `group` is specified, the vector sums to the number of distinct groups.

Examples

```
time_vector = c(1, 8, 22, 23, 24, 25, 26, 27, 28, 29, 30, 31,
                32, 33, 34, 35, 36, 37, 38, 39, 44, 50, 57, 64)

time_weights(times = time_vector)
```

`window_fava` *Compute FAVA in sliding windows.*

Description

This function computes FAVA in sliding window slices of a data set.

Usage

```
window_fava(
  relab_matrix,
  window_size,
  window_step = 1,
  group = NULL,
  index = NULL,
  time = NULL,
  w = NULL,
  S = NULL,
  K = NULL,
  normalized = FALSE,
  alpha = 0.5
)
```

Arguments

`relab_matrix`	matrix or data frame with rows containing non-negative entries that sum to 1. Each row represents a sample, each column represents a category, and each entry represents the abundance of that category in the sample. If `relab_matrix` contains any metadata, it must be on the left-hand side of the matrix, the right `K` entries of each row must sum to 1, and `K` must be specified. Otherwise, all entries of each row must sum to 1.
`window_size`	An integer number specifying the number of samples per window.

`window_step`	Optional; an integer specifying the distance between the first entry of adjacent windows. Default is `window_step=1`.
`group`	Optional; a string specifying the name of the column that describes which group each row (sample) belongs to. Use if `relab_matrix` is a single matrix containing multiple groups of samples you wish to compare.
`index`	Optional; a string specifying the name of the column in `relab_matrix` containing an index for each sample. For example, if `relab_matrix` contains time series data, `index` would be the column containing the time of each sample. If `index` is not specified but `time` is, `time` is by default used as the index.
`time`	Optional; a string specifying the name of the column that describes the sampling time for each row. Include if you wish to weight FAVA by the distance between samples.
`w`	Optional; a vector of length `I` with non-negative entries that sum to 1. Entry `w[i]` represents the weight placed on row `i` in the computation of the mean abundance of each category across rows. The default value is `w = rep(1/nrow(relab_matrix),nrow(rel`
`S`	Optional; a K x K similarity matrix with diagonal elements equal to 1 and off-diagonal elements between 0 and 1. Entry `S[i,k]` for `i!=k` is the similarity between category and `i` and category `k`, equaling 1 if the categories are to be treated as identical and equaling 0 if they are to be treated as totally dissimilar. The default value is `S = diag(ncol(relab_matrix))`.
`K`	Optional; an integer specifying the number of categories in the data. Default is `K=ncol(relab_matrix)`.
`normalized`	Optional; should normalized FAVA be used? Default is `normalized = FALSE`; use `normalized = TRUE` to compute normalized FAVA. FAVA can only be normalized if it is not weighted.
`alpha`	Optional; number between 0 and 1 specifying the opacity of the horizontal lines plotted. Default is `alpha = 0.5`.

Value

A list of values of fava for each window.

Examples

```
A = matrix(c(.3,.7,0,.1,0,.9,.2,.5,.3,.1,.8,.1,.3,.4,.3,.6,.4,0,0,.5,.5),
           ncol = 3, byrow = TRUE)
window_out = window_fava(relab_matrix = A, window_size = 4, normalized = TRUE)
```

`window_list` *Generate sliding windows of specified length given the maximum number of samples*

Description

This function generates a list of of sliding windows conditional on two parameters: the length of each window (number of samples) and the total number of samples present in the data.

Usage

```
window_list(window_size, length, window_step = 1)
```

Arguments

`window_size`	An integer number specifying the number of samples per window.
`length`	An integer number specifying the total number of samples.
`window_step`	Optional; an integer number specifying the distance between the first entry of adjacent windows. Default is `window_step=1`.

Value

A list of samples of sample indices. Each list entry represents one window.

Examples

```
window_list(window_size = 6, length = 40)
window_list(window_size = 6, length = 40, window_step = 2)
```

`window_plot` *Generate a plot of FAVA in sliding windows.*

Description

This function generates a plot of normalized or unnormalized, weighted or unweighted FAVA computed in sliding windows across samples for one or many groups of samples.

Usage

```
window_plot(window_fava, alpha = 0.5)
```

Arguments

`window_fava`	The output of `window_fava`.
`alpha`	Optional; number between 0 and 1 specifying the opacity of the horizontal lines plotted. Default is `alpha = 0.5`.

Value

A ggplot2 object.

Examples

```
A = matrix(c(.3,.7,0,.1,0,.9,.2,.5,.3,.1,.8,.1,.3,.4,.3,.6,.4,0,0,.5,.5),
           ncol = 3, byrow = TRUE)
window_out = window_fava(relab_matrix = A, window_size = 4, normalized = TRUE)
window_out$window_data
window_out$window_plot
```

xue_microbiome_sample *Temporal microbiome composition data*

Description

A subset of the data generated by Katherine Xue et al. detailing longitudinal composition of the human gut microbiome for three subjects who experience an antiobitic perturbation between days 29 and 34. We include only the subjects XAA, XBA, and XCA.

Usage

```
xue_microbiome_sample
```

Format

`xue_microbiome_sample`**:**

A data frame with 75 rows and 1,348 columns:

subject Subject ID: XBA, XDA, or XMA

timepoint Time (days) of sample collection

... Species names ...

xue_phyloseq *Phyloseq object containing the information in xue_microbiome_sample*

Description

A Phyloseq object containing an OTU table and the corresponding sample information (subject and timepoint).

Usage

```
xue_phyloseq
```

Format

`xue_phyloseq`**:**

#' A phyloseq-class experiment-level object with two elements with an OTU table and sample data.

xue_species_info *Table of species information*

Description

A data frame providing taxonomic information for the species included in xue_species_tree.

Usage

```
xue_species_info
```

Format

`xue_species_info`**:**

#' A data frame with 1346 rows and 9 columns:

species_id The species_id given in xue_microbiome_sample

kingdom, phylum, class, order, family, genus, species The corresponding taxonomic category for each species

species_id_number The numeric code associated with each species, as used in xue_species_tree

xue_species_similarity *Species similarity matrix for the species included in xue_microbiome_sample*

Description

A similarity matrix, with entry (i,j) corresponding to the pairwise similarity between species i and species j. This similarity matrix was derived from a phylogenetic distance matrix, inferred from the tree xue_species_tree, using the expression s(i,j) = exp(-d(i,j)), where d(i,j) is the phylogenetic distance between species i and j.

Usage

```
xue_species_similarity
```

Format

`xue_species_similarity`**:**

A dataframe with 524 rows and 524 columns, each corresponding to one species.

`xue_species_tree` *Phylogenetic tree for the species included in xue_microbiome_sample*

Description

A phylogenetic tree in the Newick format.

Usage

```
xue_species_tree
```

Format

`xue_species_tree`:
A Newick tree.

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