

Package ‘RSnetwork’

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Title Ecological Networks Analysis with Reciprocal Scaling

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Description Analyze ecological networks and possibly species-environment tables using various correspondence analyses methods, and also quantify niches widths with reciprocal scaling.

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stringr,
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tidyr,
tidyselect

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| | |
|------------|-----------------------|
| add_params | <i>Add parameters</i> |
|------------|-----------------------|

Description

Adds parameters value to simulation results dataframe.

Usage

```
add_params(corsim, paramvec)
```

Arguments

| | |
|----------|--|
| corsim | a list with the correlation values (output of get_cor()). |
| paramvec | A vector of parameters (must be a 1-row dataframe) |

Value

The list corsim with added columns in each dataframe of the list. The added columns are named like the columns of paramvec and contain repeated values of paramvec

| | |
|-------------|-----------------------------|
| combine_cor | <i>Combine correlations</i> |
|-------------|-----------------------------|

Description

Combine lists formatted by [get_cor\(\)](#).

Usage

```
combine_cor(a, b)
```

Arguments

| | |
|---|--|
| a | First list formatted by get_cor() |
| b | Second list formatted by get_cor() |

Value

A list as formatted by [get_cor\(\)](#) (see this function's "Return" section).

compas2d

*Simulate species interactions***Description**

Simulate species interactions based on the matching between 2 traits. This is based on a code that originally simulated species abundances in different sites based on their environmental niches.

Usage

```
compas2d(
  nconsumer = 40,
  nresource = 100,
  le_grad = 100,
  ratio_grad = 0.8,
  consumer_ab = NULL,
  resource_ab = NULL,
  min_ab = 1,
  max_ab = 100,
  ninter = 100,
  delta = 1,
  mean_tol = 2,
  sd_tol = 10,
  buffer = 1,
  col_prefix = "c",
  row_prefix = "r",
  rowvar_prefix = "tr",
  remove_zeroes = FALSE,
  return_intermediate = FALSE
)
```

Arguments

| | |
|-------------|--|
| nconsumer | Number of consumer species |
| nresource | Number of resource species |
| le_grad | gradient length for traits. The values are comprised between 0 and le_grad. |
| ratio_grad | length of the second gradient as a ratio of the first one. |
| consumer_ab | Optional vector of abundances for consumers. If missing, abundances are taken from a uniform law between min_ab and max_ab. |
| resource_ab | Optional vector of abundances for resource species. If missing, abundances are taken from a uniform law between min_ab and max_ab. |
| min_ab | minimal abundance value for the uniform law (if consumer_ab and/or resource_ab are absent) |
| max_ab | maximal abundance value for the uniform law (if consumer_ab and/or resource_ab are absent) |
| ninter | Total number of observations in the matrix |
| delta | Exponent between 0 and 1 to give more (1) or less (0) weight to trait matching relatively to abundance. |

| | |
|---------------------|--|
| mean_tol | Mean value for the distribution of the normal law in which consumers niche standard deviation are drawn (for each trait). |
| sd_tol | Standard deviation for the distribution of the normal law in which consumers niche standard deviation are drawn (for each trait). |
| buffer | buffer to allow consumer species optima to fall outside the gradient by 0 - buffer ; le_grad + buffer. |
| col_prefix | Prefix for the column names of the matrix (which will be prefix + index). Defaults to "c" for "consumers". |
| row_prefix | Prefix for the row names of the matrix (which will be prefix + index). Defaults to "r" for "resources". |
| rowvar_prefix | Prefix for the column names of the row variable (which will be prefix + 1 or 2). Defaults to "tr" for "resource traits". |
| remove_zeroes | If there are unobserved species, should we keep them in the final matrix? |
| return_intermediate | Return intermediate results? If yes, will return the probability of interaction based on matching only, on abundances only and the predicted mix of matching and abundances (before sampling). |

Details

This function models species interactions as arising from the matching between their traits. Traits are drawn from a uniform law with parameters determined by `le_grad` and `ratio_grad` (for the first and second trait). The matching is based on a multivariate normal distribution with means corresponding to the difference between trait values, and variance determined by the variance of consumers' niches. It is then mixed to the abundances (using the parameter `delta`) and sampled to obtain `ninteractions`.

Value

An object of class `compas`, which is a list with 3 components (4 if `return_intermediate`):

- `comm`: the interaction matrix of dimension `nresource` x `nconsumer`. Rows and columns are named like species (see `row_prefix` and `col_prefix`).
- `rowvar`: resource traits values. It is a `nresource` x 2 matrix. Rows are named like resource species (see `row_prefix`) and column are named like resource traits (see `rowvar_prefix`).
- `colvar`: consumer traits values. It is a `nconsumer` x 2 x 2 array. Rows are named like consumer species (see `col_prefix`). Column names in each array dimension are `mean` and `sd` (for the mean and standard deviation of the species niche, respectively).
- if `return_intermediate`, a fourth component named `intermediate` is returned. The first element is named `p_matching` and contains the matrix of probabilities of interactions based only on matching. The second `ab_neutral` contains the matrix of count of interactions as predicted by species abundances. The third `p_mix` contains the matrix of probabilities of interactions taking into account the mix of matching and abundances.

| | |
|-------------|--------------------|
| create_code | <i>Create code</i> |
|-------------|--------------------|

Description

Create a code for each (unique) value of a vector

Usage

```
create_code(col, code_pattern = NA, name = NA)
```

Arguments

| | |
|--------------|--|
| col | vector to code |
| code_pattern | first letter to add before code |
| name | optional column names for the output dataframe |

Value

dataframe with 2 columns: one with original vector (named col by default), the other with the corresponding codes (named col_code by default)

| | |
|--------------|--|
| df_to_matrix | <i>Interaction dataframe to matrix</i> |
|--------------|--|

Description

Transforms an interaction dataframe into an incidence matrix.

Usage

```
df_to_matrix(df, rows = "plant_species_code", columns = "animal_species_code")
```

Arguments

| | |
|---------|---|
| df | the dataframe to transform. Must have species names (columns which names are given by the parameters rows and columns), and a third column frequency with the interaction frequency/strength index. |
| rows | Name of the column with the species names that will be the rows of the matrix. |
| columns | Name of the column with the species names that will be the columns of the matrix. |

Value

Returns an incidence matrix with named rows and columns filled with the values in frequency.

| | |
|---------------|----------------------|
| filter_matrix | <i>Filter matrix</i> |
|---------------|----------------------|

Description

Filter incidence matrix to keep only species that interact with more partners that a given threshold.

Usage

```
filter_matrix(mat, thr)
```

Arguments

| | |
|-----|---|
| mat | The incidence matrix. |
| thr | The minimal number of different interacting partners a species must have (a species must have \geq thr different interacting partners). It is not weighted (for instance, if thr = 2 then a species interacting 100 times with a unique other partner will be removed.) |

Value

The interaction matrix but where the rows and/or columns with species that do not interact above the threshold removed.

| | |
|----------------|-----------------------|
| get_best_model | <i>Get best model</i> |
|----------------|-----------------------|

Description

Compare 2 linear models and returns the best

Usage

```
get_best_model(lmsimple, lm2, method = c("LRT", "AIC"), alpha = 0.05)
```

Arguments

| | |
|----------|---|
| lmsimple | First model (if using method = "LRT", must be nested in lm2) |
| lm2 | Second model |
| method | The method to use to compare models: likelihood ratio test (method = "LRT") for nested linear models comparison or AIC (method = "AIC") |
| alpha | Significance threshold to consider when using method = "LRT" |

Value

The best model

get_cor

*Get correlations***Description**

Measure correlations between niche obtained from CA or reciprocal scaling and the fundamental and realized niches.

Usage

```
get_cor(niches)
```

Arguments

niches An object containing the fundamental, realized and multivariate niches (output of `get_niches()`)

Value

A list of lists. The first element (named *fundamental*) contains the correlations with the fundamental niche values and the second element (named *realized*) contains the correlations with the realized niche values. Each sublist is a list of 2 dataframes:

- **mean_cor**: correlations of niches optima with the CA coordinates.
- **sd_cor**: correlations of niches breadths with reciprocal scaling standard deviations.

These dataframes have the following columns:

- **axis**: correlations measured on which multivariate axis?
- **type**: the type of individuals for which the measure is (rows or columns?). Column or row codes are chosen with *rowname* and *colname* from the parameter *niches*.
- **cor**: the correlation value

get_mean_sd

*Get mean and standard deviation***Description**

Return the mean and standard deviation from a reciprocal scaling analysis

Usage

```
get_mean_sd(recscal, ax = 1:2)
```

Arguments

recscal A dataframe (expected to be the output of `reciprocal.coa()`). The first columns must contain the reciprocal scaling scores and the last 3 columns are *Row*, *Col* and *Weight*.

ax the axes for which to compute mean and variance (the function will use the n-th first columns from the *recscal* dataframe).

Value

A list of 4 matrices, each containing the mean or standard deviation per axes. Each matrix has one column per axis given in ax and as many rows as there are grouping levels.

- rowstd is the standard deviation per row
- rowmean is the mean per row
- colstd is the standard deviation per column
- colmean is the mean per column

Matrices rownames are the row or column groups (species names). Column names are the same as in recscal.

| | |
|------------|-------------------|
| get_niches | <i>Get niches</i> |
|------------|-------------------|

Description

Returns the fundamental and realized niches and the niches returned by reciprocal scaling.

Usage

```
get_niches(
  rec,
  comm,
  consumer_niche,
  resource_traits,
  rowname = "row",
  colname = "col"
)
```

Arguments

| | |
|-----------------|--|
| rec | A dataframe (expected to be the output of <code>reciprocal.coa()</code>). The first 2 columns must contain the reciprocal scaling scores (there can be more axes, but only the first 2 will be used.) The last 3 columns are Row, Col and Weight. |
| comm | The observed community matrix. It is used to get the resource species niche width. |
| consumer_niche | A nx2x2 array with n rows (n species), 2 columns corresponding to species niche optima and standard deviation (corresponding to their multivariate normal niche) and 2 traits stored in the dimension. There is 1 row per species. |
| resource_traits | A mx2 matrix. Each row corresponds to a species and each column to a trait value. |
| rowname | How to name the first element of the list (resources/rows) (e.g. could be "plants") |
| colname | How to name the second element of the list (resources/rows) (e.g. could be "birds") |

Details

The fundamental niche optimum of a consumer is the mean of its niche (from the `consumer_niche` table). For a resource, it corresponds to the traits from the `resource_traits` table. The fundamental niche breadths are computed for consumers only, and corresponds to the breadths of their niche (from the `consumer_niche` table).

The realized niche optima are computed as the weighted means of the traits of species' interacting partners: therefore, the realized niche of a consumer is the weighted mean of the traits of the resource species it interacts with, and reciprocally. The realized niche breadths are computed as the weighted standard deviations of the traits of a species' interacting partners. Therefore, the realized niche breadth of a consumer is the weighted standard deviation of the traits of the resource species it interacts with, and reciprocally.

Niche optima computed with multivariate methods correspond to the reciprocal scaling mean and niche breadths correspond to the reciprocal scaling standard deviation.

Value

A list of lists of lists. The first element (named `rowname`) contains the niches for the resources and the second element (named `colname`) contains the niches for the consumers. Each element contains 2 lists:

- `mean`: niche optima
- `sd`: niche breadths

For each of these sublists, there are 3 matrices elements named:

- `fundamental`: the fundamental niche based on species traits
- `realized`: the realized niche based on interacting species traits
- `mvar`: niche parameters corresponding to the mean and standard deviations given by reciprocal scaling.

These matrices describe the value for each species (in rows) along each trait/axis (in column).

get_pred

Get the predictions of a linear model

Description

Returns the model prediction on given data range.

Usage

```
get_pred(
  dat_predict,
  lmpred,
  by = 0.001,
  level = 0.95,
  interval = c("confidence", "prediction")
)
```

Arguments

| | |
|--------------------------|---|
| <code>dat_predict</code> | The data to predict |
| <code>lmpred</code> | The model. It must have only one explanatory variable named mean. |
| <code>by</code> | The step to use for the range of <code>dat_predict</code> values |
| <code>level</code> | Confidence level used for the prediction. Defaults to 0.95. |
| <code>interval</code> | The type of interval to use: the argument is used in <code>stats::predict.lm()</code> and has the same interpretation. <code>confidence</code> gives the confidence interval around the mean of the observations whereas <code>prediction</code> gives the confidence interval of predicted values. |

Value

The model's prediction as a dataframe with columns:

- `fit`: the predicted value
- `lwr`: lower bound of the confidence interval (see arguments `level` and `interval`)
- `upr`: upper bound of the confidence interval (see arguments `level` and `interval`)
- `x`: the explanatory variable

`lm_labels`
Labels of a linear model

Description

Get the labels of a linear model to display on a plot.

Usage

```
lm_labels(mod, a)
```

Arguments

| | |
|------------------|--|
| <code>mod</code> | The linear model. It is expected to have 2 or 3 coefficients of the form $y = ax + b$ or $y = ax + cx^2 + b$. |
| <code>a</code> | The axis to consider. Used for the subscript of the variables. |

Details

Inspired from <https://r-graphics.org/recipe-annotate-facet>

Value

A dataframe with columns `formula` and `r2` Containing respectively the model equation and coefficient of determination. The formulas are written to be parsed later in the plot. The variables' names are s_a for the predicted value and m_a for the predictor.

| | |
|--------------|--|
| matrix_to_df | <i>Interaction matrix to dataframe</i> |
|--------------|--|

Description

Transform interaction matrix (incidence matrix) to a dataframe

Usage

```
matrix_to_df(mat, colnames = "animals", rownames = "plants", tofactor = TRUE)
```

Arguments

| | |
|----------|---|
| mat | Interaction matrix. The function assumes this matrix has row and column names, that are then used to fill the output dataframe columns. |
| colnames | The name to give to the column containing the matrix column names (defaults to "animals"). |
| rownames | The name to give to the column containing the matrix row names (defaults to "plants"). |
| tofactor | Should the final columns colnames and rownames be of type factor? |

Value

A dataframe with 3 columns:

- colnames containing the initial matrix column names
- rownames containing the initial matrix row names
- value containing the matrix values

| | |
|-----------|------------------------------------|
| multiplot | <i>Plot bi, tri or quadriplots</i> |
|-----------|------------------------------------|

Description

Plot bi, tri or quadriplots from a multivariate analysis

Usage

```
multiplot(
  indiv_row = NULL,
  indiv_col = NULL,
  indiv_row_lab = rownames(indiv_row),
  indiv_col_lab = rownames(indiv_col),
  var_row = NULL,
  var_col = NULL,
  var_row_lab = rownames(var_row),
  var_col_lab = rownames(var_col),
  row_color = "black",
```

```

    col_color = "black",
    eig = NULL,
    x = 1,
    y = 2,
    xlim = NULL,
    ylim = NULL,
    grad = 4,
    mult = 1,
    title = NULL,
    max.overlaps = 20,
    labsize = 3,
    alphapoints = 0.8
  )

```

Arguments

| | |
|----------------------------|--|
| <code>indiv_row</code> | Matrix of individuals coordinates for rows (type <code>dudi\$li</code>) |
| <code>indiv_col</code> | Matrix of individuals coordinates for columns (type <code>dudi\$co</code>) |
| <code>indiv_row_lab</code> | Labels for row individuals |
| <code>indiv_col_lab</code> | Labels for columns individuals |
| <code>var_row</code> | Matrix of variables coordinates for rows (type <code>dudi\$corR</code>) |
| <code>var_col</code> | Matrix of variables coordinates for columns (type <code>dudi\$corQ</code>) |
| <code>var_row_lab</code> | Labels for row variables |
| <code>var_col_lab</code> | Labels for columns variables |
| <code>row_color</code> | Color for the row individuals or variables |
| <code>col_color</code> | Color for the columns individuals or variables |
| <code>eig</code> | Eigenvalues vector |
| <code>x</code> | Axis to use for the x-axis |
| <code>y</code> | Axis to use for the y-axis |
| <code>xlim</code> | x-axis limits. Defaults to data range if not specified |
| <code>ylim</code> | y-axis limits. Defaults to data range if not specified |
| <code>grad</code> | graduations for the major.grid |
| <code>mult</code> | factor to multiply the vectors (arrows on the plot) |
| <code>title</code> | plot title |
| <code>max.overlaps</code> | <code>max.overlaps</code> argument for <code>ggrepel::geom_text_repel</code> |
| <code>labsize</code> | text size for point labels |
| <code>alphapoints</code> | Transparency value for points |

Value

a `ggplot` representing row and/or column individuals as points, possibly including row and/or column variables as vectors

| | |
|----------------|--------------------------------|
| plot_corcircle | <i>Plot correlation circle</i> |
|----------------|--------------------------------|

Description

Plot a correlation circle

Usage

```
plot_corcircle(
  cor,
  label = rownames(cor),
  col = "black",
  lty = "solid",
  cor2 = NULL,
  label2 = rownames(cor2),
  col2 = "black",
  lty2 = "solid",
  xax = 1,
  yax = 2,
  xlab = NULL,
  ylab = NULL,
  eig = NULL,
  col_bg = "grey30",
  mar = 0.01,
  labsize = 3.88,
  lim = 1,
  clip = "off"
)
```

Arguments

| | |
|--------|--|
| cor | dataframe containing the coordinates of the variables for each axis in columns |
| label | label to display on the arrows for the first set of variables |
| col | color for the first set of arrows |
| lty | linetype for the first set of variables |
| cor2 | optional second set of variables (for double constrained analyses) |
| label2 | label to display on the arrows for the second set of variables |
| col2 | color for the second set of arrows |
| lty2 | linetype for the second set of variables |
| xax | index of the column of the matrix to use for the x-axis |
| yax | index of the column of the matrix to use for the y-axis |
| xlab | Custom x-label |
| ylab | Custom y-label |
| eig | Eigenvalues vector (to display percentage of variation explained by each axis) |
| col_bg | color to use for background elements (circle and $y = 0$ and $x = 0$ lines) |
| mar | margin between arrow tips and label |

| | |
|---------|---|
| labsize | Size for arrow labels. |
| lim | limits for the x and y axes: a unique number that gives the distance between the plot limit and zero (so that the circle is centered) |
| clip | constrain the labels to be inside the plot? If yes, change to "on". |

Value

A ggplot with the correlation circle where one or two sets of variables are represented as arrows inside the circle of radius 1.

| | |
|----------|-------------------------|
| plot_eig | <i>Plot eigenvalues</i> |
|----------|-------------------------|

Description

Barplot of eigenvalues

Usage

```
plot_eig(eigenvalues, showrank = FALSE)
```

Arguments

| | |
|-------------|--|
| eigenvalues | The eigenvalues |
| showrank | Should the x-axis display the rank of the eigenvalues? |

Value

A ggplot with the eigenvalues displayed as a barplot.

| | |
|-----------------|--------------------------|
| plot_lm_mean_sd | <i>Plot linear model</i> |
|-----------------|--------------------------|

Description

Plot the linear models that predicts standard deviation with mean of the reciprocal scaling.

Usage

```
plot_lm_mean_sd(
  df,
  dat_pred,
  col = "black",
  ylab = "Standard deviation",
  xlab = "Mean",
  lab = NULL,
  text_size = 3,
  points_size = 1,
  facet = NULL,
```

```

    facet_dir = c("h", "v"),
    strip.position = "top",
    nudge_x = NULL,
    max.overlaps = 5
  )

```

Arguments

| | |
|----------------|---|
| df | The dataframe containing the true data. It must have columns mean (used for the x-axis) and sd (y-axis). |
| dat_pred | The dataframe containing the predicted data (assumed to be over the true data range). It is assumed to be obtained with get_pred() . It must have columns x, fit, lwr, upr. |
| col | Color of the data points and prediction line. If there are facets, it should have length 2 (one color per facet). |
| ylab | Label of the y-axis |
| xlab | Label of the x-axis |
| lab | A label to add to the graph. If provided, will be added in the top-left corner. It must be a dataframe with at least one column R2 (to create this dataframe, see lm_labels()). If facet is not NULL, it must have a column type too. If it has a column formula, the formula will be displayed along with the coefficient of determination. |
| text_size | Size of the text (points labels and R squared). Other text (axes titles, panels) is scaled to be bigger. |
| points_size | Size of the points |
| facet | A column to use for facetting. Defaults to NULL (no facets). |
| facet_dir | Direction of the facet grid: "h" for "horizontal and "v" for "vertical" (useful only if facet is not NULL) |
| strip.position | Strip position (useful only if facet is not NULL). See the documentation of ggplot2::facet_wrap() (argument strip.position). |
| nudge_x | Nudge to use for top left corner labels. If NULL, defaults to 1 % of the range. |
| max.overlaps | max.overlaps argument for ggrepel::geom_text_repel() |

Value

A ggplot object showing standard deviation vs mean, along with the predicted values of the linear model

| | |
|-------------|--------------------|
| plot_matrix | <i>Plot matrix</i> |
|-------------|--------------------|

Description

Plot an interaction matrix

Usage

```
plot_matrix(
  mat,
  tile_fill = "white",
  col = "black",
  alpha = FALSE,
  max_size = max(mat) * 0.7,
  base_size = 12,
  xlab = "Animals",
  ylab = "Plants",
  trans = "identity",
  breaks = waiver(),
  legend_title = "Frequency"
)
```

Arguments

| | |
|--------------|---|
| mat | The matrix. It represents the interaction frequency of species interactions (it is assumed to be plants x animals). |
| tile_fill | Color fill of tiles. Use NA for no fill. |
| col | Color of the points |
| alpha | If true, will make less abundant interactions more transparent |
| max_size | The maximum circle area to be used in scale_size_area. Defaults to 0.7 times the actual maximum. |
| base_size | Base text size |
| xlab | Title of the x axis |
| ylab | Title of the y axis |
| trans | Transformation to apply to the scaling of points. |
| breaks | Breaks to use for the legend of the points size. |
| legend_title | Title of the legend |

Value

A ggplot representing the interaction matrix as a rectangle with points whose sizes correspond to the interaction frequencies

| | |
|-----------------|--------------------------------|
| plot_reciprocal | <i>Plot reciprocal scaling</i> |
|-----------------|--------------------------------|

Description

Plot reciprocal scaling

Usage

```
plot_reciprocal(
  dudi,
  recscal,
  xax = 1,
  yax = 2,
  labsize = 1,
  psize = 1,
  xlab = NULL,
  ylab = NULL,
  group = c("li", "co"),
  col = "cornflowerblue",
  alpha = 0.2,
  plot_arrows = TRUE,
  plot_points = TRUE,
  plot_labels = TRUE,
  ...
)
```

Arguments

| | |
|-------------|---|
| dudi | An object of class <code>ade4::dudi()</code> to pair with the analysis. Used for the eigenvalues (always) and for the arrows of explanatory variables if required. It must be of class <code>coa</code> , <code>pcaiv</code> or <code>dpcaiv</code> (obtained with the new function <code>dpcaiv2</code>). |
| recscal | A dataframe (expected to be the output of <code>reciprocal.coa</code> function). The first columns must contain the reciprocal scaling scores and the last 3 columns are <code>Row</code> , <code>Col</code> and <code>Weight</code> . |
| xax | The index of the multivariate axis to plot on the x-axis (the column <code>xax</code> from the <code>recscal</code> dataframe). |
| yax | The index of the multivariate axis to plot on the y-axis (the column <code>yax</code> from the <code>recscal</code> dataframe). |
| labsize | Size of the ellipses labels |
| psize | Size of the points |
| xlab | Custom x-label |
| ylab | Custom y-label |
| group | The group to use for the ellipses (<code>co</code> or <code>li</code>) |
| col | The ellipses colors |
| alpha | The ellipses transparency |
| plot_arrows | Whether to plot the arrows for the variables |
| plot_points | Plot the data points? |
| plot_labels | Plot ellipses labels? |
| ... | Additional parameters passed to <code>s.class</code> |

Value

A plot with the points grouped by ellipses.

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