Package 'fuzzySim'

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Type Package

Title Fuzzy similarity in species distributions

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Description Functions to calculate fuzzy versions of species' occurrence patterns based on presence-absence data (including inverse distance interpolation, trend surface analysis and prevalence-independent favourability GLM), and pair-wise fuzzy similarity (based on fuzzy versions of commonly used similarity indices) among those occurrence patterns. Includes also functions for data preparation, such as obtaining unique abbreviations of species names, converting species lists (long format) to presence-absence tables (wide format), transposing part of a data frame, assessing the false discovery rate, or analysing multicollinearity among variables. Includes also sample datasets for providing practical examples.

License GPL-3

R topics documented:

`fuzzySim-package`	*Fuzzy similarity in species distributions*

Description

Functions to calculate fuzzy versions of species' occurrence patterns based on presence-absence data (including inverse distance interpolation, trend surface analysis and prevalence-independent favourability GLM), and pair-wise fuzzy similarity (based on fuzzy versions of commonly used similarity indices) among those occurrence patterns. Includes also functions for data preparation, such as obtaining unique abbreviations of species names, converting species lists (long format) to presence-absence tables (wide format), transposing part of a data frame, assessing the false discovery rate, or analysing multicollinearity among variables. Includes also sample datasets for providing practical examples. A step-by-step illustrated tutorial is available from the package homepage (http://fuzzysim.r-forge.r-project.org).

Details

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Type:	Package
Version:	1.0
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License:	GPL-3

Author(s)

A. Marcia Barbosa

Maintainer: A. Marcia Barbosa <barbosa@uevora.pt>

References

Barbosa A.M. (2015) fuzzySim: applying fuzzy logic to binary similarity indices in ecology. Methods in Ecology and Evolution, in press.

See Also

RMACOQUI

Examples

```
data(rotifers)

head(rotifers)


# add column with species name abbreviations:

rotifers$spcode <- spCodes(rotifers$species, sep.species = "_", nchar.gen = 1,
nchar.sp = 5, nchar.ssp = 0)

head(rotifers)


# convert species list (long format) to presence-absence table (wide format):

rotifers.presabs <- splist2presabs(rotifers, sites.col = "TDWG4",
sp.col = "spcode", keep.n = FALSE)

head(rotifers.presabs)


# get 3rd-degree spatial trend surface for each species' distribution:

data(rotif.env)

names(rotif.env)

rotifers.tsa <- multTSA(rotif.env, sp.cols = 18:47,
coord.cols = c("Longitude", "Latitude"), id.col = 1)

head(rotifers.tsa)


# get inverse squared distance to presence for each species:

rotifers.isqd <- distPres(rotif.env, sp.cols = 18:47,
coord.cols = c("Longitude", "Latitude"), id.col = 1, p = 2, inv = TRUE)

head(rotifers.isqd)


# get prevalence-independent environmental favourability models for each species:

data(rotif.env)

names(rotif.env)

rotifers.fav <- multGLM(data = rotif.env, sp.cols = 18:47, var.cols = 5:17,
id.col = 1, step = FALSE, trim = TRUE, Favourability = TRUE)
```

get matrix of fuzzy similarity between rotifer species distributions:

# either based on inverse squared distance to presence:
rot.fuz.sim.mat <- simMat(rotifers.isqd[, -1], method = "Jaccard")

# or on environmental favourability for presence:
rot.fuz.sim.mat <- simMat(rotifers.fav$predictions[, 32:61], method = "Jaccard")

head(rot.fuz.sim.mat)

transpose fuzzy rotifer distribution data to compare
regional species composition rather than species' distributions:

names(rotifers.isqd)

rot.fuz.reg <- transpose(rotifers.fav$predictions, sp.cols = 32:61,
reg.names = 1)

head(rot.fuz.reg)

get matrix of fuzzy similarity between (some) regions' species compositions:

reg.fuz.sim.mat <- simMat(rot.fuz.reg[, 1:100], method = "Jaccard")

head(reg.fuz.sim.mat)

---

distPres   *(Inverse) distance to the nearest presence*

Description

This function takes a matrix or data frame containing species presence (1) and absence (0) data and
their spatial coordinates (optionally also a pre-calculated distance matrix between all localities), and
calculates the (inverse) distance from each locality to the nearest presence locality for each species.

Usage

distPres(data, sp.cols, coord.cols = NULL, id.col = NULL, dist.mat = NULL,
method = "euclidian", suffix = "_D", p = 1, inv = TRUE)

Arguments

data   a matrix or data frame containing, at least, two columns with spatial coordinates,
                and one column per species containing their presence (1) and absence (0) data,
                with localities in rows.

sp.cols	names or index numbers of the columns containing the species presences and absences in data. It must contain only zeros (0) for absences and ones (1) for presences.
coord.cols	names or index numbers of the columns containing the spatial coordinates in data (in this order, x and y, or longitude and latitude).
id.col	optionally, the name or index number of a column (to be included in the output) containing locality identifiers in data.
dist.mat	optionally, if you do not want distances calculated with any of the methods available in dist, you may provide a distance matrix calculated elsewhere for the localities in data.
method	the method with which to calculate distances between localities. Available options are those of dist. The default is "euclidian".
suffix	character indicating the suffix to add to the distance columns in the resulting data frame. The default is "_D".
p	the power to which distance should be raised. The default is 1; use 2 or higher if you want more conservative distances.
inv	logical value indicating whether distance should be inverted, so that it varies between 0 and 1 and higher values mean closer to presence. The default is TRUE, which is adequate as a fuzzy version of presence-absence (for using e.g. with fuzSim and simMat). In this case, presences maintain the value 1, and inverse distance to presence is calculated only for absence localities.

Value

distPres returns a matrix or data frame containing the identifier column (if provided in id.col) and one column per species containing the distance (inverse squared by default) from each locality to the nearest presence of that species.

Author(s)

A. Marcia Barbosa

See Also

dist

Examples

```
data(rotif.env)

head(rotif.env)

names(rotif.env)


# calculate plain distance to presence:

rotifers.dist <- distPres(rotif.env, sp.cols = 18:47, coord.cols = c("Longitude",
```

```
"Latitude"), id.col = 1, p = 1, inv = FALSE, suffix = "_D")

head(rotifers.dist)


# calculate inverse squared distance to presence:

rotifers.invd2 <- distPres(rotif.env, sp.cols = 18:47, coord.cols = c("Longitude",
"Latitude"), id.col = 1, p = 2, inv = TRUE, suffix = "_iDsq")

head(rotifers.invd2)
```

Fav   *Favourability*

Description

Environmental (prevalence-independent) favourability for a species' presence

Usage

```
Fav(obs = NULL, pred = NULL, n1n0 = NULL, model = NULL, sample.preval = NULL,
method = "RBV", true.preval = NULL)
```

Arguments

obs	a vector of 1-0 values of a modelled binary variable.
pred	a vector of predicted probability values for obs, given e.g. by logistic regression.
n1n0	alternatively to obs, an integer vector of length 2 providing the total number of ones and zeros, in this order. This argument is ignored if the obs vector is provided.
model	alternatively to all of the above, a model object of class "glm". If provided, will override any values provided in the arguments described above.
sample.preval	alternatively to obs or n1n0, the prevalence (proportion of positive cases) of the modelled binary variable in the modelled data.
method	either "RBV" for the original Real, Barbosa & Vargas (2006) procedure, or "AT"" for the modification proposed by Albert & Thuiller (2008) (but see Acevedo & Real 2012)
true.preval	the true prevalence (as opposed to sample prevalence), necessary if you want to use the AT method.

Details

Logistic regression (Generalised Linear Model with binomial error distribution and a logit link) is widely used for modelling species' potential distributions using presence/absence data and a set of categorical or continuous predictor variables. However, this GLM incorporates the prevalence (proportion of presences) of the species in the training sample, which affects the probability values produced. Barbosa (2006) and Real, Barbosa & Vargas (2006) proposed an environmental favourability function which is based on logistic regression but cancels out uneven proportions of presences and absences in the modelled data. Favourability thus assesses the extent to which the environmental conditions change the probability of occurrence of a species with respect to its overall prevalence in the study area. Model predictions become, therefore, directly comparable among species with different prevalences. The favourability function is implemented in the **fuzzySim** package and is also in the SAM (Spatial Analysis in Macroecology) software (Rangel et al. 2010).

Using simulated data, Albert & Thuiller (2008) proposed a modification to the favourability function, but it requires knowing the true prevalence of the species (not just the prevalence in the studied sample), which is rarely possible in real-world modelling. Besides, this suggestion was based on the misunderstanding that the favourability function was a way to obtain the probability of occurrence when prevalence differs from 50%, which is incorrect (see Acevedo & Real 2012).

To get environmental favourability with either the Real, Barbosa & Vargas ("RBV") or the Albert & Thuiller ("AT") method, you just need to get a probabilistic model (e.g. logistic regression) from your data and then use the Fav function. Input data for this function are either a model object resulting from the glm function, or the presences-absences (1-0) of your species and the corresponding presence probability values, obtained e.g. with predict(mymodel, mydata, type = "response"). Alternatively to the presences-absences, you can provide either the sample prevalence or the numbers of presences and absences. In case you want to use the "AT" method, you also need to provide the true (absolute) prevalence of your species.

Value

A numeric vector of the favourability values corresponding to the input probability values.

Author(s)

A. Marcia Barbosa

References

Acevedo P. & Real R. (2012) Favourability: concept, distinctive characteristics and potential usefulness. Naturwissenschaften 99: 515-522

Albert C.H. & Thuiller W. (2008) Favourability functions versus probability of presence: advantages and misuses. Ecography 31: 417-422.

Barbosa A.M.E. (2006) Modelacion de relaciones biogeograficas entre predadores, presas y parasitos: implicaciones para la conservacion de mamiferos en la Peninsula Iberica. PhD Thesis, University of Malaga (Spain).

Rangel T.F.L.V.B, Diniz-Filho J.A.F & Bini L.M. (2010) SAM: a comprehensive application for Spatial Analysis in Macroecology. Ecography 33: 46-50.

Real R., Barbosa A.M. & Vargas J.M. (2006) Obtaining environmental favourability functions from logistic regression. Environmental and Ecological Statistics 13: 237-245.

See Also

glm, multGLM

Examples

```
# obtain a probability model and its predictions:

data(rotif.env)

names(rotif.env)

mod <- with(rotif.env, glm(Abrigh ~ Area + Altitude + AltitudeRange +
HabitatDiversity + HumanPopulation, family = binomial))

prob <- predict(mod, data = rotif.env, type = "response")


# obtain predicted favourability in different ways:

Fav(model = mod)

Fav(obs = rotif.env$Abrigh, pred = prob)

Fav(pred = mod$fitted.values, n1n0 = c(112, 179))

Fav(pred = mod$fitted.values, sample.preval = 0.38)
```

FDR   *False Discovery Rate*

Description

Calculate the false discovery rate (type I error) under repeated testing and determine which variables to select and to exclude from multivariate analysis.

Usage

```
FDR(data = NULL, sp.cols = NULL, var.cols = NULL, pvalues = NULL, model.type,
family = "binomial", correction = "fdr", q = 0.05, verbose = TRUE)
```

Arguments

data	a data frame containing the response and predictor variables (one in each column).
sp.cols	index number of the column containing the response variable (currently implemented for only one response variable at a time).
var.cols	index numbers of the columns containing the predictor variables.

pvalues	optionally, instead of data, sp.cols and var.cols, a data frame with the names of the predictor variables in the first column and their bivariate p-values (obtained elsewhere) in the second column. Example: pvalues <- data.frame(var = letters[1:5], pval = c(0.02, 0.004, 0.07, 0.03, 0.05)).
model.type	either "LM" (linear model, for continuous response variables) or "GLM" (generalized linear models, for binary or other variables for which such models are more appropriate)
family	if model.type = "GLM", the error distribution and link function (see glm or family for details); defaults to "binomial" (for binary logistic regression).
correction	the correction procedure to apply to the p-values; see p.adjust.methods for available options and p.adjust for more information. The default is "fdr".
q	the threshold value of FDR-corrected significance above which to reject variables. Defaults to 0.05.
verbose	logical, whether to report a short description of the results.

Details

It is common in ecology to search for statistical relationships between species' occurrence and a set of predictor variables. However, when a large number of variables is analysed (compared to the number of observations), false findings may arise due to repeated testing. Garcia (2003) recommended controlling the false discovery rate (FDR; Benjamini & Hochberg 1995) in ecological studies. The p.adjust R function performs this and other corrections to the significance (p) values of variables under repeated testing. The FDR function performs repeated regressions (either linear or binary logistic) or uses already-obtained p values for a set of variables; calculates the FDR with p.adjust; and shows which variables should be retained for or excluded from further multivariate analysis according to their corrected p values (see, for example, Barbosa, Real & Vargas 2009).

The FDR function uses the Benjamini & Hochberg ("BH") correction by default, but check the p.adjust documentation for other available methods. Input data may be the response variable (for example, the presence-absence or abundance of a species) and the predictors (a table with one independent variable in each column, with the same number of rows and in the same order as the response); there should be no missing values in the data. Model type can be either "LM" (linear model) for continuous response variables such as abundance, or "GLM" (generalized linear model) for binary responses such as presence-absence. Alternatively, you may already have performed the univariate regressions and have a set of variables and corresponding p values which you want to correct with FDR; in this case, get a table with your variables' names in the first column and their p values in the second column, and supply it as the pvalues argument to the FDR function (no need to provide response or predictors in this case).

Value

A list with the following components:

exclude	a data frame of the variables to exclude under the chosen criteria, including the variables' names, their bivariate coefficients against the response, their p-value and adjusted p-value.
select	a data frame similar to the above for the variables to select.

Author(s)

A. Marcia Barbosa

References

Barbosa A.M., Real R. & Vargas J.M (2009) Transferability of environmental favourability models in geographic space: The case of the Iberian desman (Galemys pyrenaicus) in Portugal and Spain. Ecological Modelling 220: 747-754

Benjamini Y. & Hochberg Y. (1995) Controlling the false discovery rate: a practical and powerful approach to multiple testing. Journal of the Royal Statistical Society, Series B 57: 289-300

Garcia L.V. (2003) Controlling the false discovery rate in ecological research. Trends in Ecology and Evolution 18: 553-554

See Also

p.adjust

Examples

```
data(rotif.env)

names(rotif.env)

FDR(data = rotif.env, sp.cols = 18, var.cols = 5:17, model.type = "GLM")
```

---

fuzSim   *Fuzzy similarity*

Description

This function calculates fuzzy similarity, based on a fuzzy version of the binary similarity index specified in method, between two binary (0 or 1) or fuzzy (between 0 and 1) variables.

Usage

```
fuzSim(x, y, method)
```

Arguments

x	a vector of (optionally fuzzy) presence-absence data, with 1 meaning presence, 0 meaning absence, and values in between meaning fuzzy presence (or the degree to which each locality belongs to the set of species presences, or to which each species belongs to the locality; Zadeh, 1965). Fuzzy presence-absence can be obtained, for example, with functions multGLM, distPres or multTSA in this package.
y	a vector similar to x, of the same length and in the same order.
method	the similarity index to use. Currently available options are 'Jaccard', 'Sorensen', 'Simpson' and 'Baroni' (see Details).

Details

Similarity between ecological communities, beta diversity patterns, biotic regions, and distributional relationships among species are commonly determined based on pair-wise (dis)similarities in species' occurrence patterns. Some of the most commonly employed similarity indices are those of Jaccard (1901), Sorensen (1948), Simpson (1960) and Baroni-Urbani & Buser (1976), which are here implemented in their fuzzy versions (Barbosa, 2015), able to deal with both binary and fuzzy data.

Value

The function returns a value between 0 and 1 representing the fuzzy similarity between x and y. Note, for example, that Jaccard similarity can be converted to dissimilarity (or Jaccard distance) if subtracted from 1, while 1-Sorensen is not a proper distance metric as it lacks the property of triangle inequality (see http://en.wikipedia.org/wiki/S%C3%B8rensen%E2%80%93Dice_coefficient).

Note

The formulas used in this function may look slighty different from some of their published versions (e.g. Baroni-Urbani & Buser 1976), not only because the letters are switched, but because here the A and B are the numbers of attributes present in each element, whether or not they are also present in the other one. Thus, our 'A+B' is equivalent to 'A+B+C' in formulas where A and B are the numbers of attributes present in one but not the other element, and our A+B-C is equivalent to their A+B+C. The formulas used here (adapted from Olivero et al. 1998) are faster to calculate, visibly for large datasets.

Author(s)

A. Marcia Barbosa

References

Barbosa A.M. (2015) fuzzySim: applying fuzzy logic to binary similarity indices in ecology. Methods in Ecology and Evolution, in press.

Baroni-Urbani C. & Buser M.W. (1976) Similarity of Binary Data. Systematic Zoology, 25: 251-259

Jaccard P. (1901) Etude comparative de la distribution florale dans une portion des Alpes et des Jura. Memoires de la Societe Vaudoise des Sciences Naturelles, 37: 547-579

Olivero J., Real R. & Vargas J.M. (1998) Distribution of breeding, wintering and resident waterbirds in Europe: biotic regions and the macroclimate. Ornis Fennica, 75: 153-175

Simpson, G.G. (1960) Notes on the measurement of faunal resemblance. Amer. J. Sci. 258A, 300-311

Sorensen T. (1948) A method of establishing groups of equal amplitude in plant sociology based on similarity of species and its application to analyses of the vegetation on Danish commons. Kongelige Danske Videnskabernes Selskab, 5(4): 1-34

Zadeh L.A. (1965) Fuzzy sets. Information and Control, 8: 338-353

See Also

simMat

Examples

data(rotif.env)

head(rotif.env)

names(rotif.env)

fuzSim(rotif.env[,18], rotif.env[,19], method = "Jaccard")

fuzSim(rotif.env[,18], rotif.env[,19], method = "Sorensen")

fuzSim(rotif.env[,18], rotif.env[,19], method = "Simpson")

fuzSim(rotif.env[,18], rotif.env[,19], method = "Baroni")

---

getPreds   *Get model predictions*

Description

This function allows getting the predictions of multiple models when applied to a given dataset. It can be useful if you have a list of model objects (e.g. resulting from multGLM) and want to apply them to a new data set containing the same variables for another region or time period. There are options to include the logit link (Y) and/or Favourability (see Fav).

Usage

getPreds(data, models, id.col = NULL, Y = FALSE, P = TRUE, Favourability = TRUE,
incl.input = TRUE)

Arguments

data	the data frame to which to apply the models to get their predictions; must contain all variables (with the same names, case-sensitive) included in any of the models.
models	a list of model objects obtained e.g. with function glm or multGLM.
id.col	optionally, the index number of a column of data containing row identifiers, to be included in the result. Ignored if incl.input = TRUE.
Y	logical, whether to include the logit link (y) value in the predictions.
P	logical, whether to include the probability value in the predictions.
Favourability	logical, whether to include Favourability in the predictions (see Fav).
incl.input	logical, whether to include input columns in the output. The default is TRUE.

Value

This function returns a data frame containing the model predictions, next to the id.col if provided, and to the input data if incl.input = TRUE.

Author(s)

A. Marcia Barbosa

See Also

multGLM, predict

Examples

```
data(rotif.env)

names(rotif.env)


# identify rotifer data corresponding to the Eastern and Western hemispheres:

unique(rotif.env$CONTINENT)

rotif.env$HEMISPHERE <- "Eastern"

rotif.env$HEMISPHERE[rotif.env$CONTINENT %in%
c("NORTHERN_AMERICA", "SOUTHERN_AMERICA")] <- "Western"

head(rotif.env)


# separate the rotifer data into hemispheres

east.hem <- rotif.env[rotif.env$HEMISPHERE == "Eastern", ]
west.hem <- rotif.env[rotif.env$HEMISPHERE == "Western", ]


# make models for 3 of the species in rotif.env based on their distribution
# in the Eastern hemisphere:

mods <- multGLM(east.hem, sp.cols = 18:20, var.cols = 5:17, id.col = 1,
step = FALSE, FDR = FALSE, trim = FALSE)


# get the models' predictions for the Western hemisphere dataset:

preds <- getPreds(west.hem, models = mods$models, P = TRUE, Favourability = TRUE)

head(preds)
```

integerCols   *Classify integer columns*

Description

This function detects which numeric columns in a data frame contain only whole numbers, and converts those columns to integer class, so that they take up less space.

Usage

integerCols(data)

Arguments

data   a data frame containing possibly integer columns classified as numeric.

Value

The function returns a data frame with the same columns as data, but with those that are numeric and contain only whole numbers (possibly including NA) now classified as integer.

Author(s)

A. Marcia Barbosa

See Also

is.integer, as.integer, multConvert

Examples

```
dat <- data.frame(
  var1 = 1:10,
  var2 = as.numeric(1:10),
  var3 = as.numeric(c(1:4, NA, 6:10)),
  var4 = as.numeric(c(1:3, NaN, 5, Inf, 7, -Inf, 9:10)),
  var5 = as.character(1:10),
  var6 = seq(0.1, 1, by = 0.1),
  var7 = letters[1:10]
)  # creates a sample data frame

dat

str(dat)
# var2 classified as 'numeric' but contains only whole numbers
# var3 same as var2 but containing also NA values
# var4 same as var2 but containing also NaN and infinite values
# var5 contains only whole numbers but initially classified as factor
```

```
dat <- integerCols(dat)

str(dat)
# var2 and var3 now classified as 'integer'
# var4 remains as numeric because contains infinite and NaN (not integer) values
# var5 remains as factor
```

| modelTrim | *Trim off non-significant variables from a model* |

Description

This function performs a stepwise removal of non-significant variables from a model.

Usage

```
modelTrim(model, method = "summary", alpha = 0.05)
```

Arguments

model	a model object.
method	the method for getting the individual p-values. Can be either "summary" for the p-values of the coefficient estimates, or "anova" for the p-values of the variables themselves (see Details).
alpha	the p-value above which a variable is removed.

Details

Stepwise variable selection is a common procedure for simplifying models. It maximizes predictive efficiency in an objective and reproducible way, and is useful when the individual importance of the predictors is not known a priori (Hosmer & Lemeshow, 2000). The step R function performs such procedure using an information criterion (AIC) to select the variables, but it often leaves variables that are not significant in the model. Such variables can be subsequently removed with a manual stepwise procedure (e.g. Crawley 2007, p. 442; Barbosa & Real 2010, 2012; Estrada & Arroyo 2012). The modelTrim function performs such removal automatically until all remaining variables are significant. It can also be applied to a full model (i.e., without previous use of the step function), as it serves as a backward stepwise selection procedure based on the significance of the coefficients (if method = "summary", the default) or on the significance of the variables themselves (if method = "anova", better when there are categorical variables in the model).

Value

The input model object after removal of non-significant variables.

Author(s)

A. Marcia Barbosa

References

Barbosa A.M. & Real R. (2010) Favourable areas for expansion and reintroduction of Iberian lynx accounting for distribution trends and genetic diversity of the European rabbit. Wildlife Biology in Practice 6: 34-47

Barbosa A.M. & Real R. (2012) Applying fuzzy logic to comparative distribution modelling: a case study with two sympatric amphibians. The Scientific World Journal, Article ID 428206

Crawley M.J. (2007) The R Book. John Wiley & Sons, Chichester (UK)

Estrada A. & Arroyo B. (2012) Occurrence vs abundance models: Differences between species with varying aggregation patterns. Biological Conservation, 152: 37-45

Hosmer D. W. & Lemeshow S. (2000) Applied Logistic Regression (2nd ed). John Wiley and Sons, New York

See Also

step

Examples

```
# load sample data:

data(rotif.env)

names(rotif.env)


# build a stepwise model of a species' occurrence based on some of the variables:

mod <- with(rotif.env, step(glm(Abrigh ~ Area + Altitude + AltitudeRange +
HabitatDiversity + HumanPopulation, family = binomial)))


# examine the model:

summary(mod)  # contains non-significant variables


# use modelTrim to get rid of non-significan effects:

mod <- modelTrim(mod)

summary(mod)  # only significant variables now
```

---

multConvert   *Multiple conversion*

Description

This function can simultaneously convert multiple columns of a matrix or data frame.

Usage

multConvert(data, conversion, cols = 1:ncol(data))

Arguments

data	A matrix or data frame containing columns that need to be converted
conversion	the conversion to apply, e.g. as.factor or a custom-made fucntion
cols	the columns of data to convert

Details

Sometimes we need to change the data type (class, mode) of a variable in R. There are various possible conversions, performed by functions like as.integer, as.factor or as.character. If we need to perform the same conversion on a number of variables (columns) in a data frame, we can convert them all simultaneously using this function. By default it converts all the columns in the data frame, but you can specify just a few of them. multConvert can also be used to apply other kinds of transformations - for example, if you need to divide some of your columns by 100, just write a function to do this and then use multConvert to apply this function to any group of columns.

Value

The input data with the specified columns converted as asked.

Author(s)

A. Marcia Barbosa

Examples

data(rotif.env)

str(rotif.env)

# convert the first 4 columns to character:
converted.rotif.env <- multConvert(data = rotif.env, conversion = as.character, cols = 1:4)

str(converted.rotif.env)

names(rotif.env)

divide some columns by 100:

div100 <- function(x) x / 100

rotif.env.cent <- multConvert(data = rotif.env, conversion = div100, cols = c(6:10, 12:17))

head(rotif.env.cent)

---

multGLM   *Multiple GLMs*

Description

This function calculates generalized linear models for a set of (species) presence/absence data in a data frame.

Usage

multGLM(data, sp.cols, var.cols, id.col = NULL, family = "binomial",
test.sample = 0, FDR = FALSE, step = TRUE, trace = 0, start = "null.model",
direction = "both", Y.prediction = FALSE, P.prediction = TRUE,
Favourability = TRUE, group.preds = TRUE, trim = TRUE, ...)

Arguments

data	a data frame in wide format (see splist2presabs) containing, in separate columns, your species' binary (0/1) occurrence data and the predictor variables.
sp.cols	index numbers of the columns containing the species data to be modelled.
var.cols	index numbers of the columns containing the predictor variables to be used for modelling.
id.col	(optional) index number of column containing the row identifiers (if defined, it will be included in the output predictions data frame).
family	argument to be passed to the glm function; only 'binomial' is implemented in multGLM so far.
test.sample	a subset of data to set aside for subsequent model testing. Can be a value between 0 and 1 for a proportion of the data to choose randomly (e.g. 0.2 for 20%), or an integer number for a particular number of cases to choose randomly among the records in data, or a vector of integers for the index numbers of the particular rows to set aside, or "Huberty" for his rule of thumb based on the number of variables (Huberty 1994, Fielding & Bell 1997).
FDR	logical, whether to do a preliminary exclusion of variables based on their bivariate relationship with the response and the false discovery rate (see FDR).
step	logical, whether to use the step function to perform a stepwise variable selection based on AIC.
trace	if positive, information is printed during the running of step. Larger values may give more detailed information.
start	character, whether to start with the 'null.model' (so AIC variable selection starts forward) or with the 'full.model' (so selection starts backward). Used only if step = TRUE.

direction	argument to be passed to step specifying the direction of variable selection ('forward', 'backward' or 'both'). Used only if $step = TRUE$.
Y.prediction	logical, whether to include output predictions in the scale of the predictor variables ($type =$ "link" in predict.glm).
P.prediction	logical, whether to include output predictions in the scale of the response variable, i.e. probability ($type =$ "response" in predict.glm).
Favourability	logical, whether to apply the Favourability function to extract the effect of prevalence on probability (Real et al. 2006) and include its results in the output.
group.preds	logical, whether to group together predictions of similar type (Y, P or F) in the output predictions table (e.g. if FALSE: sp1_Y, sp1_P, sp1_F, sp2_Y, sp2_P, sp2_F; if TRUE: sp1_Y, , sp2_Y, sp1_P, sp2_P, sp1_F, sp2_F).
trim	logical, whether to trim non-significant variables off the models using the modelTrim function; can be used whether or not step is TRUE; works as a backward variable elimination procedure based on significance.
...	additional arguments to be passed to modelTrim.

Details

This function automatically calculates binomial GLMs for one or more species (or other binary variables) in a data frame. The function can optionally perform stepwise variable selection (and it does so by default) instead of forcing all variables into the models, starting from either the null model (the default, so selection starts forward) or from the full model (so selection starts backward) and using Akaike's information criterion (AIC) as a variable selection criterion. Instead or subsequently, it can also perform stepwise removal of non-significant variables from the models using the modelTrim function. There is also an optional preliminary selection of the variables with a significant bivariate relationship with the response, based on the false discovery rate (FDR), but note that some variables can be significant in a multivariate model even if they would not have been selected by FDR. Favourability is also calculated. By default, all data are used in model training, but you can define an optional test.sample to be reserved for model testing afterwards. You may also want to do a previous check for multicollinearity among variables, e.g. the variance inflation factor (VIF).

The multGLM function will create a list of the resulting models (each with the name of the corresponding species column) and a data frame with their predictions (Y, P and/or F, all of which are optional). If you plan on representing these predictions in a GIS based on .dbf tables, remember that dbf only allows up to 10 characters in column names; multGLM predictions will add 2 characters (_Y, _P and/or _F) to each of your species column names, so use species names/codes with up to 8 characters in the data set that you are modelling. You can create (sub)species name abbreviations with the spCodes function.

Value

This function returns a list with the following components:

predictions	a data frame with the model predictions (if either of Y.prediction, P.prediction, or Favourability are TRUE).
models	a list of the resulting model objects.

Author(s)

A. Marcia Barbosa

References

Fielding A.H. & Bell J.F. (1997) A review of methods for the assessment of prediction errors in conservation presence/absence models. Environmental Conservation 24: 38-49

Huberty C.J. (1994) Applied Discriminant Analysis. Wiley, New York, 466 pp. Schaafsma W. & van Vark G.N. (1979) Classification and discrimination problems with applications. Part IIa. Statistica Neerlandica 33: 91-126

Real R., Barbosa A.M. & Vargas J.M. (2006) Obtaining environmental favourability functions from logistic regression. Environmental and Ecological Statistics 13: 237-245.

See Also

glm, Fav, step, modelTrim, multicol

Examples

```
data(rotif.env)

names(rotif.env)


# make models for 3 of the species in rotif.env:

mods <- multGLM(rotif.env, sp.cols = 45:47, var.cols = 5:17, id.col = 1,
step = TRUE, FDR = TRUE, trim = TRUE)

names(mods)

head(mods$predictions)

names(mods$models)

mods$models[[1]]

mods$models[["Ttetra"]]
```

---

multicol	*Analyse multicollinearity in a dataset, including VIF*

Description

This function analyses multicollinearity in a set of variables, including the R-squared, tolerance and variance inflation factor (VIF).

Usage

```
multicol(vars)
```

Arguments

vars   A matrix or data frame containing the numeric variables for which to calculate multicollinearity. Only the 'independent' (predictor, explanatory, right hand side) variables should be entered, as the result obtained for each variable depends on all the other variables present in the analysed data set.

Details

Testing collinearity among covariates is a recommended step of data exploration before applying a statistical model (Zuur et al. 2010). The multicol function calculates the degree of multicollinearity in a set of numeric variables, using three closely related measures: R squared (the coefficient of determination of a linear regression of each predictor variable on all other predictor variables, i.e., the amount of variation in each variable that is accounted for by other variables in the dataset); tolerance (1 - R squared), i.e. the amount of variation in each variable that is not included in the remaining variables; and the variance inflation factor: VIF = 1 / (1 - R squared), which, in a linear model with these variables as predictors, reflects the degree to which the variance of an estimated regression coefficient is increased due only to the correlations among covariates (Marquardt 1970; Mansfield & Helms 1982).

Value

The function returns a matrix with one row per analysed variable, the names of the variables as row names, and 3 columns: R-squared, Tolerance and VIF.

Author(s)

A. Marcia Barbosa

References

Marquardt D.W. (1970) Generalized inverses, ridge regression, biased linear estimation, and non-linear estimation. Technometrics 12: 591-612.

Mansfield E.R. & Helms B.P. (1982) Detecting multicollinearity. The American Statistician 36: 158-160.

Zuur A.F., Ieno E.N. & Elphick C.S. (2010) A protocol for data exploration to avoid common statistical problems. Methods in Ecology and Evolution 1: 3-14.

Examples

```
data(rotif.env)
names(rotif.env)

# calculate multicollinearity among the predictor variables:
multicol(rotif.env[ , 5:17])
```

```
# more examples uing R datasets:
multicol(trees)

# you'll get a warning and some NA results if any of the variables is not numeric:
multicol(OrchardSprays)

# so define the subset of numeric 'vars' to calculate 'multicol' for:
multicol(OrchardSprays[ , 1:3])
```

---

multTSA   *Trend Surface Analysis for multiple species*

Description

This function performs trend surface analysis for multiple species at a time. It converts categorical presence-absence (1-0) data into continuous surfaces denoting the spatial trend in species' occurrence patterns.

Usage

```
multTSA(data, sp.cols, coord.cols, id.col = NULL, degree = 3, step = TRUE,
Favourability = FALSE, suffix = "_TS", save.models = FALSE)
```

Arguments

data	a matrix or data frame containing, at least, two columns with spatial coordinates, and one column per species containing their presence (1) and absence (0) data, with localities in rows.
sp.cols	names or index numbers of the columns containing the species presences and absences in data. Must contain only zeros (0) for absences and ones (1) for presences.
coord.cols	names or index numbers of the columns containing the spatial coordinates in data (x and y, or longitude and latitude, in this order!).
id.col	optionally, the name or index number of a column (to be included in the output) containing locality identifiers in data.
degree	the degree of the spatial polynomial to use (see Details). The default is 3.
step	logical value indicating whether the regression of presence-absence on the spatial polynomial should do a stepwise inclusion of the polynomial terms (using the step function with default settings, namely backward AIC selection), rather than forcing all terms into the equation. The default is TRUE.
Favourability	logical value indicating whether the probability values obtained from the regression should be converted to favourability, so that they are more directly comparable among species with different prevalence (see Real et al. 2006, Acevedo & Real 2012). The default is FALSE.

suffix	character indicating the suffix to add to the trend surface columns in the resulting data frame. The default is "_TS".
save.models	logical value indicating whether the models obtained from the regressions should be saved in the results. The default is FALSE.

Details

Trend Surface Analysis is a way to model the spatial structure in species' distributions by regressing occurrence data on the spatial coordinates x and y, for a linear trend, or on polynomial terms of these coordinates (x^2, y^2, x*y, etc.), for curvilinear trends (Legendre & Legendre, 1998; Borcard et al., 2011). Second- and third-degree polynomials are often used. multTSA allows specifying the degree of the spatial polynomial to use. By default, it uses a 3rd-degree polynomial and performs stepwise AIC selection of the polynomial terms to include.

Value

A matrix or data frame containing the identifier column (if provided in id.col) and one column per species containing the value predicted by the trend surface analysis.

Author(s)

A. Marcia Barbosa

References

Acevedo P. & Real R. (2012) Favourability: concept, distinctive characteristics and potential usefulness. Naturwissenschaften 99: 515-522

Borcard D., Gillet F. & Legendre P. (2011) Numerical Ecology with R. Springer, New York.

Legendre P. & Legendre L. (1998) Numerical Ecology. Elsevier, Amsterdam.

Real R., Barbosa A.M. & Vargas J.M. (2006) Obtaining environmental favourability functions from logistic regression. Environmental and Ecological Statistics 13: 237-245

See Also

distPres, poly

Examples

```
data(rotif.env)

head(rotif.env)

names(rotif.env)

tsa <- multTSA(rotif.env, sp.cols = 18:47, coord.cols = c("Longitude", "Latitude"),
id.col = 1)

head(tsa)
```

---

percentTestData   *Percent test data*

Description

Based on the work of Schaafsma & van Vark (1979), Huberty (1994) provided a heuristic ("rule of thumb") for determining an adequate proportion of data to set aside for testing species presence/absence models, based on the number of predictor variables that are used (Fielding & Bell 1997). The percentTestData function calculates this proportion as a percentage.

Usage

percentTestData(nvar)

Arguments

nvar   the number of variables in the model.

Value

A numeric value of the percentage of data to leave out of the model for further model testing.

Author(s)

A. Marcia Barbosa

References

Huberty C.J. (1994) Applied Discriminant Analysis. Wiley, New York, 466 pp.

Schaafsma W. & van Vark G.N. (1979) Classification and discrimination problems with applications. Part IIa. Statistica Neerlandica 33: 91-126

Fielding A.H. & Bell J.F. (1997) A review of methods for the assessment of prediction errors in conservation presence/absence models. Environmental Conservation 24: 38-49

See Also

multGLM

Examples

```
# say you're building a model with 15 variables:

percentTestData(15)

# the result tells you that 21% is an appropriate percentage of data
# to set aside for testing your model, so train it with 79% of the data
```

---

rotif.env	*Rotifers and environmental variables on TDWG level 4 regions of the world*

Description

These data were extracted from a database of monogonont rotifer species presence records on the geographical units used by the Biodiversity Information Standards (formerly Taxonomic Database Working Group, TDWG; base maps available from www.kew.org/science-conservation/research-data/ resources/gis-unit/tdwg-world) and a few environmental (including human and spatial) variables on the same spatial units. The original data were compiled and published by Fontaneto et al. (2012) in long (narrow, stacked) format. Here they are presented in wide or unstacked format (presence-absence table, obtained with the splist2presabs function), reduced to the species recorded in at least 100 (roughly one third) different TDWG level 4 units, and with abbreviations of the species' names (obtained with the spCodes function). Mind that this is not a complete picture of these species' distributions, due to insufficient sampling in many regions.

Usage

```
data(rotif.env)
```

Format

A data frame with 291 observations on the following 47 variables.

TDWG4 a factor with 291 levels indicating the abbreviation code of each TDWG4 region

LEVEL_NAME a factor with 291 levels indicating the name of each TDWG4 region

REGION_NAME a factor with 47 levels indicating the name of the main geographical region to which each TDWG4 level belongs

CONTINENT a factor with 9 levels indicating the continent to which each TDWG4 level belongs

Area a numeric vector

Altitude a numeric vector

AltitudeRange a numeric vector

HabitatDiversity a numeric vector

HumanPopulation a numeric vector

Latitude a numeric vector

Longitude a numeric vector

Precipitation a numeric vector

PrecipitationSeasonality a numeric vector

TemperatureAnnualRange a numeric vector

Temperature a numeric vector

TemperatureSeasonality a numeric vector

UrbanArea a numeric vector

Abrigh a numeric vector

Afissa a numeric vector

Apriod a numeric vector

Bangul a numeric vector

Bcalyc a numeric vector

Bplica a numeric vector

Bquadr a numeric vector

Burceo a numeric vector

Cgibba a numeric vector

Edilat a numeric vector

Flongi a numeric vector

Kcochl a numeric vector

Kquadr a numeric vector

Ktropi a numeric vector

Lbulla a numeric vector

Lclost a numeric vector

Lhamat a numeric vector

Lluna a numeric vector

Llunar a numeric vector

Lovali a numeric vector

Lpatel a numeric vector

Lquadr a numeric vector

Mventr a numeric vector

Ppatul a numeric vector

Pquadr a numeric vector

Pvulga a numeric vector

Specti a numeric vector

Tpatin a numeric vector

Tsimil a numeric vector

Ttetra a numeric vector

Source

Fontaneto D., Barbosa A.M., Segers H. & Pautasso M. (2012) The 'rotiferologist' effect and other global correlates of species richness in monogonont rotifers. Ecography, 35: 174-182.

Examples

```
data(rotif.env)

head(rotif.env)
```

---

rotifers	*Rotifer species on TDWG level 4 regions of the world*

Description

These data were extracted from a database of monogonont rotifer species records on the geographical units used by the Biodiversity Information Standards (formerly Taxonomic Database Working Group, TDWG; base maps available at www.kew.org/science-conservation/research-data/resources/gis-unit/tdwg-world). The original data were compiled and published by Fontaneto et al. (2012) for all TDWG levels. Here they are reduced to the TDWG - level 4 units and to the species recorded in at least 100 (roughly one third) of these units. Mind that this is not a complete picture of these species' distributions, due to insufficient sampling in many regions.

Usage

```
data("rotifers")
```

Format

A data frame with 3865 observations on the following 2 variables.

TDWG4 a factor with 274 levels corresponding to the code names of the TDWG level 4 regions in which the records were taken

species a factor with 30 levels corresponding to the names of the (sub)species recorded in at least 100 different TDWG level 4 regions

Source

Fontaneto D., Barbosa A.M., Segers H. & Pautasso M. (2012) The 'rotiferologist' effect and other global correlates of species richness in monogonont rotifers. Ecography, 35: 174-182.

Examples

```
data(rotifers)

head(rotifers, 10)
```

---

simMat	*Pair-wise (fuzzy) similarity matrix*

Description

simMat takes a matrix or data frame containing species occurrence data or regional species composition, either categorical (0 or 1) or fuzzy (between 0 and 1), and uses the fuzSim function to calculate a square matrix of pair-wise similarities between them, using a fuzzy logic version (Barbosa, 2015) of the specified similarity index.

Usage

    simMat(data, method)

Arguments

data   a matrix or data frame containing (optionally fuzzy) species presence-absence
                    data (in wide format, i.e. one column per species), with 1 meaning presence,
                    0 meaning absence, and values in between for fuzzy presence (or the degree to
                    which each locality belongs to the set of species presences; see Zadeh, 1965).
                    Fuzzy presence-absence can be obtained, for example, with multGLM, distPres
                    or multTSA. These data can also be transposed for comparing regional species
                    compositions.

method   the similarity index whose fuzzy version to use. See fuzSim for available op-
                    tions.

Details

The fuzzy versions of species occurrence data and of the similarity indices (Barbosa, 2015) intro-
duce tolerance for small spatial differences in species' occurrence localities and may compensate
under-sampling and geo-referencing errors. The results of simMat can be used for determining
chorotypes or biotic regions using the **RMACOQUI** package (Olivero et al. 2011).

Value

This function returns a square matrix of pair-wise similarity among the species distributions (columns)
in data. Similarity is calculated with the fuzzy version of the index specified in method, which
yields traditional binary similarity if the data are binary (0 or 1), or fuzzy similarity if the data are
fuzzy (between 0 and 1) (Barbosa, 2015).

Author(s)

A. Marcia Barbosa

References

Barbosa A.M. (2015) fuzzySim: applying fuzzy logic to binary similarity indices in ecology. Meth-
ods in Ecology and Evolution, in press.

Olivero, J., Real, R. & Marquez, A.L. (2011) Fuzzy chorotypes as a conceptual tool to improve
insight into biogeographic patterns. Systematic Biology, 60: 645-60.

See Also

fuzSim

Examples

```
# load and look at the rotif.env presence-absence data:

data(rotif.env)

head(rotif.env)

names(rotif.env)


# build a matrix of similarity among these binary data
# using e.g. Jaccard's index:

bin.sim.mat <- simMat(rotif.env[ , 18:47], method = "Jaccard")

head(bin.sim.mat)


# calculate a fuzzy version of the presence-absence data
# based on inverse distance to presences:

rotifers.invd <- distPres(rotif.env, sp.cols = 18:47,
coord.cols = c("Longitude", "Latitude"), id.col = 1, suffix = ".d", p = 1, inv = TRUE)

head(rotifers.invd)


# build a matrix of fuzzy similarity among these fuzzy distribution data,
# using the fuzzy version of Jaccard's index:

fuz.sim.mat <- simMat(rotifers.invd[ , -1], method = "Jaccard")

head(fuz.sim.mat)


# plot the similarity matrices as colours:

image(x = 1:ncol(bin.sim.mat), y = 1:nrow(bin.sim.mat), z = bin.sim.mat,
col = rev(heat.colors(256)), xlab = "", ylab = "", axes = FALSE,
main = "Binary similarity")
axis(side = 1, at = 1:ncol(bin.sim.mat), tick = FALSE,
labels = colnames(bin.sim.mat), las = 2)
axis(side = 2, at = 1:nrow(bin.sim.mat), tick = FALSE,
labels = rownames(bin.sim.mat), las = 2)

image(x = 1:ncol(fuz.sim.mat), y = 1:nrow(fuz.sim.mat), z = fuz.sim.mat,
col = rev(heat.colors(256)), xlab = "", ylab = "", axes = FALSE,
main = "Fuzzy similarity")
axis(side = 1, at = 1:ncol(fuz.sim.mat), tick = FALSE,
labels = colnames(fuz.sim.mat), las = 2, cex = 0.5)
axis(side = 2, at = 1:nrow(fuz.sim.mat), tick = FALSE,
```

```
labels = rownames(fuz.sim.mat), las = 2)


# plot a UPGMA dendrogram from each similarity matrix:

plot(hclust(as.dist(1 - bin.sim.mat), method = "average"),
main = "Binary cluster dendrogram")

plot(hclust(as.dist(1 - fuz.sim.mat), method = "average"),
main = "Fuzzy cluster dendrogram")


# you can get fuzzy chorotypes from these similarity matrices
# (or fuzzy biotic regions if your \code{\link{transpose}} \code{data},
# so that localities are in columns and species in rows)
# using the \pkg{RMACOQUI} package (Olivero et al. 2011)
```

---

spCodes   *Obtain unique abbreviations of species names*

Description

This function takes a vector of species names and converts them to abbreviated species codes containing the specified numbers of characters from the genus, the specific and optionally also the subspecific name. Separators can be specified by the user. The function checks that the resulting codes are unique.

Usage

```
spCodes(species, nchar.gen = 3, nchar.sp = 3, nchar.ssp = 0, sep.species = " ",
sep.spcode = "")
```

Arguments

species	a character vector containig the species names to be abbreviated.
nchar.gen	the number of characters from the genus name to be included in the resulting species code.
nchar.sp	the number of characters from the specific name to be included in the resulting species code.
nchar.ssp	optionally, the number of characters from the subspecific name to be included in the resulting species code. Set it to 0 if you have subspecific names in species but do not want them included in the resulting species codes.
sep.species	the character that separates genus, specific and subspecific names in species. The default is a white space.
sep.spcode	the character you want separating genus and species abbreviations in the resulting species codes. The default is an empty character (no separator).

Value

This function returns a character vector containing the species codes resulting from the abbreviation. If the numbers of characters specified do not make for unique codes, an error message is displayed showing which species names caused it, so that you can try again with different nchar.gen, nchar.sp and/or nchar.ssp.

Author(s)

A. Marcia Barbosa

See Also

substr, strsplit

Examples

```
data(rotifers)

head(rotifers)

## add a column to 'rotifers' with shorter versions of the species names:

## Not run:
rotifers$spcode <- spCodes(rotifers$species, sep.species = "_", nchar.gen = 1,
nchar.sp = 4, nchar.ssp = 0, sep.spcode = ".")

# this produces an error due to resulting species codes not being unique

## End(Not run)

rotifers$spcode <- spCodes(rotifers$species, sep.species = "_", nchar.gen = 1,
nchar.sp = 5, nchar.ssp = 0, sep.spcode = ".")

# with a larger number of characters from the specific name,
# resulting codes are now unique

## check out the result:
head(rotifers)
```

---

splist2presabs   *Convert a species list to a presence-absence table*

Description

This function takes a locality+species dataset in long (stacked) format, i.e., a matrix or data frame containing localities in one column and their recorded species in another column, and converts them to a presence-absence table (wide format) suitable for mapping and for computing distributional similarities (see e.g. simMat). Try out the Examples below for an illustration.

Usage

$$\mathrm{splist2presabs(data,\ sites.col,\ sp.col,\ keep.n = FALSE)}$$

Arguments

data	a matrix or data frame with localities in one column and species in another column. Type $\mathrm{data(rotifers);\ head(rotifers)}$ for an example.
sites.col	the name or index number of the column containing the localities in data.
sp.col	the name or index number of the column containing the species names or codes in data.
keep.n	logical value indicating whether to get in the resulting table the number of times each species appears in each locality; if FALSE (the default), only presence (1) or absence (0) is recorded.

Value

A data frame containing the localities in the first column and then one column per species indicating their presence (or their number of records if $\mathrm{keep.n = TRUE}$) and absence. Type $\mathrm{data(rotif.env);\ head(rotif.env[,18:47])}$ for an example.

Author(s)

A. Marcia Barbosa

See Also

table

Examples

$\mathrm{data(rotifers)}$

$\mathrm{head(rotifers)}$

$\mathrm{rotifers.presabs <-\ splist2presabs(rotifers,\ sites.col = "TDWG4",}$
$\mathrm{sp.col = "species",\ keep.n = FALSE)}$

$\mathrm{head(rotifers.presabs)}$

---

timer	*Timer*

Description

Reporting of time elapsed since a given start time. This function is used internally by other functions in the package.

Usage

```
timer(start.time)
```

Arguments

start.time   A date-time object of class POSIXct, e.g. as given by Sys.time.

Value

The function returns a message informing of the time elapsed since the input start.time.

Author(s)

A. Marcia Barbosa

See Also

Sys.time, proc.time, difftime

Examples

```
# get starting time:
start <- Sys.time()

# do some random analysis:
sapply(rnorm(50000), function(x) x*5)

# see how long it took:
timer(start)
```

---

transpose	*Transpose (part of) a matrix or dataframe*

Description

This function transposes (a specified part of) a matrix or data frame, optionally using one of its columns as column names for the transposed result. It can be useful for turning a species presence-absence table into a regional species composition table.

Usage

```
transpose(data, sp.cols = 1:ncol(data), reg.names = NULL)
```

Arguments

data   a matrix or data frame containing the species occurrence data to transpose.

sp.cols   names or index numbers of the columns containing the species occurrences in data which are meant to be transposed.

reg.names   name or index number of the column in data containing the region names, to be used as column names in the transposed result.

Value

This function returns the transposed sp.cols of data, with the column specified in reg.names as column names.

Author(s)

A. Marcia Barbosa

See Also

t

Examples

data(rotif.env)

head(rotif.env)

names(rotif.env)

rotif.reg <- transpose(rotif.env, sp.cols = 18:47, reg.names = 1)

head(rotif.reg)

Index