

Package ‘codyn’

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Description

Univariate and multivariate temporal and spatial diversity indices, rank abundance curves, and community stability metrics for ecologists.

Details

The functions in `codyn` implement metrics that are either explicitly temporal and include the option to calculate them over multiple replicates, or spatial and include the option to calculate them over multiple time points. Functions fall into five categories: static diversity indices, temporal diversity indices, spatial diversity indices, rank abundance curves, and community stability metrics. The diversity indices in `codyn` are temporal and spatial analogs to traditional diversity indices. Specifically, `codyn` includes functions to calculate community richness, evenness and diversity at a given point in space and time. In addition, `codyn` contains functions to calculate species turnover, mean rank shifts, and lags in community similarity between two time points. For the components of rank abundance curves, the shape of the rank abundance curve, species abundances and multivariate metrics of community composition, `codyn` contains functions to calculate these metrics either between time points or a single time point between two paired replicates. The community stability metrics in `codyn` calculate overall stability and patterns of species covariance and synchrony over time. Finally, `codyn` contains vignettes that describe methods and reproduce figures from published papers to help users contextualize and apply functions to their own data. Work on this package was supported by National Science Foundation grant #1262458 to the National Center for Ecological Analysis and Synthesis (NCEAS), the University of Wisconsin, and the University of New Mexico, and a SESYNC Synthesis Postdoctoral Fellowship to MLA.

Functions

- `turnover`: Calculates species turnover between time periods
- `rank_shift`: Calculates the mean relative change in species rank abundances
- `rate_change`: Calculates the rate change in a community over time
- `rate_change_interval`: Produces a data frame containing differences in species composition between samples at increasing time intervals
- `community_stability`: Calculates community stability over time
- `variance_ratio`: Computes the ratio of the variance of aggregate species abundances in a community
- `synchrony`: Calculates the degree synchrony in species abundances
- `temporal_torus_translation`: Calculates a test statistic on a null ecological community created via cyclic shifts. `confint` provides mean and confidence intervals of this null distribution
- `community_structure`: Calculates richness and evenness (using specified metric) for a replicate

- **community_diversity**: Calculates diversity (using specified metric) for a replicate
- **RAC_change**: Calculates changes in species richness, evenness, species' ranks, gain, and losses for a replicate over time
- **abundance_change**: For each species in a replicate, calculates changes in abundance over time
- **curve_change**: Calculates changes in the shape of the RAC curve for each replicate over time
- **multivariate_change**: Calculates changes in community composition and dispersion over time
- **RAC_difference**: Calculates differences in species richness, evenness, species' ranks, shared species between paired samples at a single point in time
- **abundance_difference**: Calculates differences in abundance for each species in paired samples at a single point in time
- **curve_difference**: Calculates differences in the shape of the RAC between paired samples at a single point in time
- **multivariate_difference**: Calculates differences in community composition and dispersion of all replicates between treatments at a single point in time

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abundance_change

Species Abundance Changes

Description

Calculates the abundance change for species in a replicate between two time points. Changes are on abundance values provided, if relative data is used, then changes in relative abundance will be calculated.

Usage

```
abundance_change(df, time.var, species.var, abundance.var,
  replicate.var = NULL, reference.time = NULL)
```

Arguments

<code>df</code>	A data frame containing time, species, and abundance columns and an optional columns of replicates.
<code>time.var</code>	The name of the time column.
<code>species.var</code>	The name of the species column.
<code>abundance.var</code>	The name of the abundance column.
<code>replicate.var</code>	The name of the optional replicate column. If specified, replicate must be unique within the dataset and cannot be nested within treatments or blocks.
<code>reference.time</code>	The name of the optional time point that all other time points should be compared to (e.g. the first year of data). If not specified, each comparison is between consecutive time points (the first and second year, second and third year, etc.)

Value

The `abundance_change` function returns a data frame with a subset of the following columns:

- `replicate.var`: A column with the specified `replicate.var`, if it is specified.
- `time.var`: A column with the specified `time.var` and a second column, with '2' appended to the name. Time is subtracted from `time2`
- `species.var`: A column with the specified `species.var`.
- `change`: A numeric column of the change in abundance between time points. A positive value occurs when a species increases in abundance over time, and a negative value when a species decreases in abundance over time.

References

Avolio et al. Submitted

Examples

```
data(pplots)
# Without replicates
df <- subset(pplots, plot == 25)
abundance_change(df = df,
  species.var = "species",
  abundance.var = "relative_cover",
  time.var = "year")

# With replicates
df <- subset(pplots, year < 2004 & plot %in% c(6, 25, 32))
abundance_change(df = df,
  species.var = "species",
  abundance.var = "relative_cover",
```

```

      replicate.var = "plot",
      time.var = "year")

# With reference year
df <- subset(pplots, year < 2005 & plot %in% c(6, 25, 32))
abundance_change(df = df,
  species.var = "species",
  abundance.var = "relative_cover",
  replicate.var = "plot",
  time.var = "year",
  reference.time = 2002)

```

abundance_difference *Abundance Differences*

Description

Calculates the abundance difference for species between two samples. Differences are on abundance values provided, if relative data is used, then differences in relative abundance will be calculated. There are three ways differences can be calculated. 1) Between treatments within a block (note: block.var and treatment.var need to be specified). 2) Between treatments, pooling all replicates into a single species pool (note: pool = TRUE, treatment.var needs to be specified, and block.var = NULL). 3) All pairwise combinations between all replicates (note: block.var = NULL, pool = FALSE and specifying treatment.var is optional. If treatment.var is specified, the treatment that each replicate belongs to will also be listed in the output).

Usage

```

abundance_difference(df, time.var = NULL, species.var, abundance.var,
  replicate.var, treatment.var = NULL, pool = FALSE,
  block.var = NULL, reference.treatment = NULL)

```

Arguments

df	A data frame containing a species, abundance, and replicate columns and optional time, treatment, and block columns.
time.var	The name of the optional time column.
species.var	The name of the species column.
abundance.var	The name of the abundance column.
replicate.var	The name of the replicate column. Replicate identifiers must be unique within the dataset and cannot be nested within treatments or blocks.
treatment.var	The name of the optional treatment column.
pool	An argument to allow abundance values to be pooled within a treatment. The default value is "FALSE", a value of "TRUE" averages abundance of each species within a treatment at a given time point.
block.var	The name of the optional block column.

reference.treatment

The name of the optional treatment that all other treatments will be compared to (e.g. only controls will be compared to all other treatments). If not specified all pairwise treatment comparisons will be made.

Value

The `abundance_difference` function returns a data frame with a subset of the following columns:

- `species.var`: A column that has same name and type as the `species.var` column.
- `difference`: A numeric column of the abundance differences between the two samples being compared (replicates or treatments). A numeric column of the change in abundance between consecutive timepoints. A positive value occurs when a species has greater abundance in `replicate.var2` than in `replicate.var` and/or in `treatment.var2` than in `treatment.var`.
- `replicate.var`: A column that has same name and type as the `replicate.var` column, represents the first replicate being compared. Note, a replicate column will be returned only when `pool = FALSE` or `block.var = NULL`.
- `replicate.var2`: A column that has the same type as the `replicate.var` column, and is named `replicate.var` with a 2 appended to it, represents the second replicate being compared. Note, a `replicate.var` column will be returned only when `pool = FALSE` and `block.var = NULL`.
- `time.var`: A column that has the same name and type as the `time.var` column, if `time.var` is specified.
- `treatment.var`: A column that has same name and type as the `treatment.var` column, represents the first treatment being compared. A `treatment.var` column will be returned when `pool = TRUE`, `block.var` is specified, or `treatment.var` is specified.
- `treatment.var2`: A column that has the same type as the `treatment.var` column, and is named `treatment.var` with a 2 appended to it, represents the second treatment being compared. A `treatment.var` column will be returned when `pool = TRUE`, `block.var` is specified, or `treatment.var` is specified.
- `block.var`: A column that has same name and type as the `block.var` column, if `block.var` is specified.

References

Avolio et al. Submitted

Examples

```
data(pplots)
# With block and no time
df <- subset(pplots, year == 2002 & block < 3)
abundance_difference(df = df,
  species.var = "species",
  abundance.var = "relative_cover",
  treatment.var = "treatment",
  block.var = "block",
  replicate.var = "plot")
```

```
# With blocks and time
df <- subset(pplots, year < 2004 & block < 3)
abundance_difference(df = df,
                    species.var = "species",
                    abundance.var = "relative_cover",
                    treatment.var = "treatment",
                    block.var = "block",
                    replicate.var = "plot",
                    time.var = "year")

# With blocks, time and reference treatment
df <- subset(pplots, year < 2004 & block < 3)
abundance_difference(df = df,
                    species.var = "species",
                    abundance.var = "relative_cover",
                    treatment.var = "treatment",
                    block.var = "block",
                    replicate.var = "plot",
                    time.var = "year",
                    reference.treatment = "N1P0")

# Pooling by treatment with time
df <- subset(pplots, year < 2004)
abundance_difference(df = df,
                    species.var = "species",
                    abundance.var = "relative_cover",
                    treatment.var = "treatment",
                    pool = TRUE,
                    replicate.var = "plot",
                    time.var = "year")

# All pairwise replicates with treatment
df <- subset(pplots, year < 2004 & plot %in% c(21, 25, 32))
abundance_difference(df = df,
                    species.var = "species",
                    abundance.var = "relative_cover",
                    replicate.var = "plot",
                    time.var = "year",
                    treatment.var = "treatment")

# All pairwise replicates without treatment
df <- subset(pplots, year < 2004 & plot %in% c(21, 25, 32))
abundance_difference(df = df,
                    species.var = "species",
                    abundance.var = "relative_cover",
                    replicate.var = "plot",
                    time.var = "year")
```


Description

A dataset of tallgrass prairie plant composition at one annually burned and one unburned site over time at the Konza Prairie LTER, Manhattan Kansas (Collins et al. 2008).

Usage

```
data(collins08)
```

Format

A data frame with 2058 rows and 4 variables

Details

A data frame containing a column for replicate, year, species and abundance:

- replicate: A factor column of spatial replicates with two levels ("annually burned" and "unburned")
- year: An integer column of sampling time points
- species: A factor column of species sampled
- abundance: A numeric column of abundance values

Source

Collins, Scott L., Katharine N. Suding, Elsa E. Cleland, Michael Batty, Steven C. Pennings, Katherine L. Gross, James B. Grace, Laura Gough, Joe E. Fargione, and Christopher M. Clark. (2008) "Rank clocks and plant community dynamics." *Ecology* 89, no. 12: 3534-41.

community_diversity *Community Diversity*

Description

Calculates Shannon's or Inverse Simpson's diversity of a community, but only one measure of diversity can be calculated at a time and must be specified.

Usage

```
community_diversity(df, time.var = NULL, abundance.var,  
  replicate.var = NULL, metric = c("Shannon", "InverseSimpson"))
```

Arguments

<code>df</code>	A data frame containing species and abundance columns and optional columns of time and/or replicate.
<code>time.var</code>	The name of the optional time column
<code>abundance.var</code>	The name of the abundance column
<code>replicate.var</code>	The name of the optional replicate column. If specified, replicate must be unique within the dataset and cannot be nested within treatments or blocks.
<code>metric</code>	The diversity measure to return: <ul style="list-style-type: none"> • "Shannon": The default measure, calculates Shannon's diversity. • "InverseSimpson": Calculates inverse of Simpson's diversity.

Value

The `community_diversity` function returns a data frame with the following attributes:

- `time.var`: A column that has the same name and type as the `time.var` column, if `time.var` is specified.
- `replicate.var`: A column that has same name and type as the `replicate.var` column, if `replicate.var` is specified.
- `Shannon`: A numeric column of Shannon's diversity if `metric = "Shannon"`
- `InverseSimpson`: A numeric column of the inverse of Simpson's diversity if `metric = "InverseSimpson"`

References

Magurran, A.E. 2004. *Measuring Biological Diversity*. Blackwell Publishing, Malden MA, USA.

Examples

```
data(pplots)
#Example with both time and replicates
df <- subset(pplots, plot == 25 | plot == 6)
community_diversity(df,
  time.var="year",
  replicate.var = "plot",
  abundance.var = "relative_cover") # for Shannon's diversity measure

df <- subset(pplots, plot == 25 | plot == 6)
community_diversity(df,
  time.var="year",
  replicate.var = "plot",
  abundance.var = "relative_cover",
  metric = "InverseSimpson") # for Inverse of Simpson's diversity measure

#Example with no replicates
df <- subset(pplots, plot == 25)
community_diversity(df,
  time.var="year",
```

```
abundance.var = "relative_cover") # for Shannon's diversity measure

#Example with no time or replicate
df <- subset(pplots, plot == 25 & year == 2002)
community_diversity(df,
  abundance.var = "relative_cover") # for Shannon's diversity measure
```

community_stability *Community Stability*

Description

Calculates the stability of the overall community over time as the temporal mean / temporal standard deviation of aggregate species abundances (Tilman 1999).

Usage

```
community_stability(df, time.var, abundance.var, replicate.var = NA)
```

Arguments

df	A data frame containing time, species and abundance columns and an optional column of replicates
time.var	The name of the time column
abundance.var	The name of the abundance column
replicate.var	The name of the optional replicate column

Details

The input data frame needs to contain columns for time, species and abundance; time.var, species.var and abundance.var are used to indicate which columns contain those variables. If multiple replicates are included in the data frame, that column should be specified with replicate.var. Each replicate should reflect a single experimental unit - there should be a single community represented within each time point and replicate.

Value

The community_stability function returns a numeric stability value unless a replication column is specified in the input data frame. If replication is specified, the function returns a data frame with the following columns:

- stability: A numeric column with the stability values.
- replicate.var: A column that shares the same name and type as the replicate.var column in the input data frame.

References

Tilman, D. "The Ecological Consequences of Changes in Biodiversity: A Search for General Principles." *Ecology* 80, no. 5 (July 1999): 1455-74. doi:10.1890/0012-9658(1999)080[1455:TECOCI]2.0.CO;2.

Examples

```
data(knz_001d)
community_stability(knz_001d[knz_001d$subplot=="A_1", ],
  time.var = "year",
  abundance.var = "abundance") # for one subplot
community_stability(knz_001d,
  time.var = "year",
  abundance.var = "abundance",
  replicate.var = "subplot") # across all subplots
```

community_structure *Community Structure*

Description

Calculates species richness and evenness of a community. Evenness may be calculated as Simpson's (1/D/S), EQ, or Evar, but only one metric of evenness can be calculated at a time and must be specified.

Usage

```
community_structure(df, time.var = NULL, abundance.var,
  replicate.var = NULL, metric = c("Evar", "SimpsonEvenness", "EQ"))
```

Arguments

df	A data frame containing species and abundance columns and optional columns of time and/or replicate.
time.var	The name of the optional time column
abundance.var	The name of the abundance column
replicate.var	The name of the optional replicate column. If specified, replicate must be unique within the dataset and cannot be nested within treatments or blocks.
metric	The measure of evenness to return: <ul style="list-style-type: none"> • "Evar": The default measure, calculates evenness as Evar from Smith and Wilson 1996 • "SimpsonEvenness": Calculates Simpson's evenness • "EQ": Calculates evenness as EQ from Smith and Wilson 1996

Value

The `community_structure` function returns a data frame with the following attributes:

- `time.var`: A column that has the same name and type as the `time.var` column, if `time.var` is specified.
- `replicate.var`: A column that has same name and type as the `replicate.var` column, if specified.
- `richness`: A numeric column of species richness
- `Evar`: A numeric column of Evar if `evenness = "Evar"`
- `EQ`: A numeric column of EQ if `evenness = "EQ"`
- `SimpsonEvenness`: A numeric column of Simpson's evenness if `evenness = "SimpsonEvenness"`

References

Smith, B. and Wilson, J. B. 1996. A consumer's guide to evenness indices. *Oikos* 76: 70-82.

Examples

```
data(pplots)
#Example with both time and replicates
df <- subset(pplots, plot == 25 | plot == 6)
community_structure(df,
                    time.var="year",
                    replicate.var = "plot",
                    abundance.var = "relative_cover") # for Evar evenness measure

df <- subset(pplots, plot == 25 | plot == 6)
community_structure(df,
                    time.var="year",
                    replicate.var = "plot",
                    abundance.var = "relative_cover",
                    metric = "SimpsonEvenness") # for Simpson's evenness measure

#Example with no replicates
df <- subset(pplots, plot == 25)
community_structure(df,
                    time.var="year",
                    abundance.var = "relative_cover",
                    metric = "EQ") # for EQ evenness measure

#Example with only a single time point and no replicates
df <- subset(pplots, plot == 25 & year == 2002)
community_structure(df,
                    abundance.var = "relative_cover") # for Evar evenness measure
```

confint.cyclic_shift *Confidence Intervals from a Cyclic Shift Permutation*

Description

Calculates confidence intervals for the S3 object produced by `cyclic_shift`

Usage

```
## S3 method for class 'cyclic_shift'
confint(object, parm = "out", level = 0.95, ...)
```

Arguments

<code>object</code>	An object of class <code>cyclic_shift</code>
<code>parm</code>	which parameter is to be given a confidence interval. At present there is only one option: the mean of the null distribution. Defaults to "out", referring to the null distribution in objects of class <code>cyclic_shift</code> .
<code>level</code>	the confidence level required.
<code>...</code>	further arguments to <code>quantile</code>

Value

A dataframe with the following columns:

- `lowerCI`: A numeric column with the lowest confidence interval value.
- `upperCI`: A numeric column with the highest confidence interval value.
- `nullMean`: A numeric column with the average value of the specified test statistic when calculated on a null community.

References

Hallett, Lauren M., Joanna S. Hsu, Elsa E. Cleland, Scott L. Collins, Timothy L. Dickson, Emily C. Farrer, Laureano A. Gherardi, et al. "Biotic Mechanisms of Community Stability Shift along a Precipitation Gradient." *Ecology* 95, no. 6 (2014): 1693-1700.

Harms, Kyle E., Richard Condit, Stephen P. Hubbell, and Robin B. Foster. "Habitat Associations of Trees and Shrubs in a 50-Ha Neotropical Forest Plot." *Journal of Ecology* 89, no. 6 (2001): 947-59.

Examples

```
# Calculate a covariance matrix on a null community
data(knz_001d)
a1_cyclic <- cyclic_shift(subset(knz_001d, subplot == "A_1"),
  time.var = "year",
  species.var = "species",
  abundance.var = "abundance",
  FUN = cov,
```

```

                                bootnumber = 10)

# Return CI on a1_cyclic
confint(a1_cyclic)

```

curve_change	<i>Curve Change</i>
--------------	---------------------

Description

Calculates the area difference between two rank abundance curves between two time periods. If replicate is specified, it must be measured in both time points, otherwise it will be dropped for that time period comparison.

Usage

```
curve_change(df, time.var, species.var, abundance.var,
             replicate.var = NULL, reference.time = NULL)
```

Arguments

df	A data frame containing time, species, and abundance columns and an optional columns of replicates.
time.var	The name of the time column.
species.var	The name of the species column.
abundance.var	The name of the abundance column.
replicate.var	The name of the optional replicate column. If specified, replicate must be unique within the dataset and cannot be nested within treatments or blocks.
reference.time	The name of the optional time point that all other time points should be compared to (e.g. the first year of data). If not specified, each comparison is between consecutive time points (the first and second year, second and third year, etc.)

Value

The curve_change function returns a data frame with the following attributes:

- time.var: A column with the specified time.var and a second column, with '2' appended to the name. Time is subtracted from time2.
- curve_change: A numeric column of the change in curves between time points.
- replicate.var: A column that has same name and type as the replicate.var column, if specified.

References

Avolio et al. Submitted

Examples

```

data(pplots)
# Without replicates
df <- subset(pplots, plot == 25)
curve_change(df = df,
             species.var = "species",
             abundance.var = "relative_cover",
             time.var = "year")

# With replicates
df <- subset(pplots, year < 2004 & plot %in% c(6, 25, 32))
curve_change(df = df,
             species.var = "species",
             abundance.var = "relative_cover",
             replicate.var = "plot",
             time.var = "year")

# With reference year
df <- subset(pplots, year < 2005 & plot %in% c(6, 25, 32))
curve_change(df = df,
             species.var = "species",
             abundance.var = "relative_cover",
             replicate.var = "plot",
             time.var = "year",
             reference.time = 2002)

```

curve_difference	<i>Curve Difference</i>
------------------	-------------------------

Description

Calculates the area difference between two rank abundance curves. There are three ways differences can be calculated. 1) Between all treatments within a block (note: block.var and treatment.var need to be specified. 2) Between treatments, pooling all replicates into a single species pool (note: pool = TRUE, treatment.var needs to be specified, and block.var = NULL. 3) All pairwise combinations between all replicates (note: block.var = NULL, pool = FALSE and specifying treatment.var is optional. If treatment.var is specified, the treatment that each replicate belongs to will also be listed in the output).

Usage

```

curve_difference(df, time.var = NULL, species.var, abundance.var,
               replicate.var, treatment.var = NULL, pool = FALSE,
               block.var = NULL, reference.treatment = NULL)

```

Arguments

df A data frame containing a species, abundance, and replicate columns and optional time, treatment, and block columns.

time.var	The name of the optional time column.
species.var	The name of the species column.
abundance.var	The name of the abundance column.
replicate.var	The name of the replicate column. Replicate identifiers must be unique within the dataset and cannot be nested within treatments or blocks.
treatment.var	The name of the optional treatment column.
pool	An argument to allow abundance values to be pooled within a treatment. The default value is "FALSE", a value of "TRUE" averages abundance of each species within a treatment at a given time point.
block.var	The name of the optional block column.
reference.treatment	The name of the optional treatment that all other treatments will be compared to (e.g. only controls will be compared to all other treatments). If not specified all pairwise treatment comparisons will be made.

Value

The `curve_difference` function returns a data frame with the following attributes:

- `curve_diff`: A numeric column of the area difference in curves between the two samples being compared (replicates or treatments).
- `replicate.var`: A column that has same name and type as the `replicate.var` column, represents the first replicate being compared. Note, a replicate column will be returned only when `pool` is `FALSE` or `block.var = NULL`.
- `replicate.var2`: A column that has the same type as the `replicate.var` column, and is named `replicate.var` with a 2 appended to it, represents the second replicate being compared. Note, a `replicate.var` column will be returned only when `pool` is `FALSE` and `block.var = NULL`.
- `time.var`: A column that has the same name and type as the `time.var` column, if `time.var` is specified.
- `treatment.var`: A column that has same name and type as the `treatment.var` column, represents the first treatment being compared. A `treatment.var` column will be returned when `pool` is `TRUE` or `block.var` is specified, or `treatment.var` is specified.
- `treatment.var2`: A column that has the same type as the `treatment.var` column, and is named `treatment.var` with a 2 appended to it, represents the second treatment being compared. A `treatment.var` column will be returned when `pool` is `TRUE` or `block.var` is specified, or `treatment.var` is specified.
- `block.var`: A column that has same name and type as the `block.var` column, if `block.var` is specified.

References

Avolio et al. Submitted

Examples

```
data(pplots)
# With block and no time
df <- subset(pplots, year == 2002 & block < 3)
curve_difference(df = df,
  species.var = "species",
  abundance.var = "relative_cover",
  treatment.var = "treatment",
  block.var = "block",
  replicate.var = "plot")

# With blocks and time
df <- subset(pplots, year < 2004 & block < 3)
curve_difference(df = df,
  species.var = "species",
  abundance.var = "relative_cover",
  treatment.var = "treatment",
  block.var = "block",
  replicate.var = "plot",
  time.var = "year")

# With blocks, time, and reference treatment
df <- subset(pplots, year < 2004 & block < 3)
curve_difference(df = df,
  species.var = "species",
  abundance.var = "relative_cover",
  treatment.var = "treatment",
  block.var = "block",
  replicate.var = "plot",
  time.var = "year",
  reference.treatment = "N1P0")

# Pooling by treatment with time
df <- subset(pplots, year < 2004)
curve_difference(df = df,
  species.var = "species",
  abundance.var = "relative_cover",
  treatment.var = "treatment",
  pool = TRUE,
  replicate.var = "plot",
  time.var = "year")

# All pairwise replicates with treatment
df <- subset(pplots, year < 2004 & plot %in% c(21, 25, 32))
curve_difference(df = df,
  species.var = "species",
  abundance.var = "relative_cover",
  replicate.var = "plot",
  time.var = "year",
  treatment.var = "treatment")

# All pairwise replicates without treatment
```

```
df <- subset(pplots, year < 2004 & plot %in% c(21, 25, 32))
curve_difference(df = df,
  species.var = "species",
  abundance.var = "relative_cover",
  replicate.var = "plot",
  time.var = "year")
```

cyclic_shift

Cyclic Shift Permutations

Description

Performs a user-specified function on a null ecological community created via cyclic shift permutations (Harms et al. 2001, Hallett et al. 2014). The null community is formed by randomly selected different starting years for the time series of each species. This generates a null community matrix in which species abundances vary independently but within-species autocorrelation is maintained. The user-specified function must require a species x time input.

Usage

```
cyclic_shift(df, time.var, species.var, abundance.var,
  replicate.var = NA, FUN, bootnumber)
```

Arguments

df	A data frame containing time, species and abundance columns and an optional column of replicates
time.var	The name of the time column
species.var	The name of the species column
abundance.var	The name of the abundance column
replicate.var	The name of the replicate column. Defaults to NA, indicating no replicates (i.e., data are from a single plot).
FUN	A function to calculate on the null community
bootnumber	Number of null simulations to run

Details

The input data frame needs to contain columns for time, species and abundance; time.var, species.var and abundance.var are used to indicate which columns contain those variables.

Value

The cyclic_shift function returns an S3 object of class "cyclic_shift" and parameter "out". The length of the "out" parameter is the number of null iterations as specified by bootnumber. If multiple replicates are specified, null values are averaged among replicates for each iteration, but a different cyclic shift permutation is applied to each replicate within an iteration.

References

Hallett, Lauren M., Joanna S. Hsu, Elsa E. Cleland, Scott L. Collins, Timothy L. Dickson, Emily C. Farrer, Laureano A. Gherardi, et al. "Biotic Mechanisms of Community Stability Shift along a Precipitation Gradient." *Ecology* 95, no. 6 (2014): 1693-1700.

Harms, Kyle E., Richard Condit, Stephen P. Hubbell, and Robin B. Foster. "Habitat Associations of Trees and Shrubs in a 50-Ha Neotropical Forest Plot." *Journal of Ecology* 89, no. 6 (2001): 947-59.

Examples

```
# Calculate a covariance matrix on a null community
data(knz_001d)
a1_cyclic <- cyclic_shift(subset(knz_001d, subplot == "A_1"),
                          time.var = "year",
                          species.var = "species",
                          abundance.var = "abundance",
                          FUN = cov,
                          bootnumber = 10)
```

knz_001d

Data from Konza Prairie, watershed 001d

Description

Plant composition within multiple replicates at an annually burned tallgrass prairie site in the Konza Prairie LTER, Manhattan KS (Watershed 001d).

Usage

```
data(knz_001d)
```

Format

A data frame with 8768 rows and 4 variables

Details

A data frame containing a column for species, year, subplot and abundance:

- species: A factor column of species sampled
- year: An integer column of sampling time points
- subplot: A factor column of spatial replicates with 20 levels
- abundance: A numeric column of abundance values

Source

Konza Prairie LTER Dataset ID: PVC02, watershed 1D

Collins, S. L. (2000) Disturbance frequency and community stability in native tallgrass prairie. *American Naturalist* 155:311-325.

mean_rank_shift	<i>Mean Rank Shifts</i>
-----------------	-------------------------

Description

A measure of the relative change in species rank abundances, which indicates shifts in relative abundances over time (Collins et al. 2008). Mean rank shifts are calculated as the average difference in species' ranks between consecutive time periods, among species that are present across the entire time series.

Usage

```
mean_rank_shift(df, time.var, species.var, abundance.var,  
  replicate.var = as.character(NA))
```

Arguments

df	A data frame containing time, species and abundance columns and an optional column of replicates
time.var	The name of the time column
species.var	The name of the species column
abundance.var	The name of the abundance column
replicate.var	The name of the optional replicate column

Details

The input data frame needs to contain columns for time, species and abundance; time.var, species.var and abundance.var are used to indicate which columns contain those variables. If multiple replicates are included in the data frame, that column should be specified with replicate.var. Each replicate should reflect a single experimental unit - there must be a single abundance value per species within each time point and replicate.

Value

mean_rank_shift returns a data frame with the following columns:

- time.var_pair: A factor column that returns the two time periods being compared, separated by a dash. The name of this column is the same as the time.var column in the input dataframe followed by "_pair".
- MRS: A numeric column with the mean rank shift values.
- replicate.var: A column that has same name and type as the replicate.var column, if replication is specified.

multivariate_change *Using dissimilarity-based measures to calculate changes in composition and dispersion*

Description

Calculates the changes in composition and dispersion based off a Bray-Curtis dissimilarity matrix. Composition change is the pairwise distance between centroids of compared time periods and ranges from 0-1, where identical communities give 0 and completely different communities give 1. Dispersion change is the difference between time periods in the dispersion of replicates, i.e. the average distance between a replicate and its centroid.

Usage

```
multivariate_change(df, time.var, species.var, abundance.var,
  replicate.var, treatment.var = NULL, reference.time = NULL)
```

Arguments

df	A data frame containing time, species, abundance and replicate columns and an optional column of treatment.
time.var	The name of the time column.
species.var	The name of the species column.
abundance.var	The name of the abundance column.
replicate.var	The name of the replicate column. Replicate identifiers must be unique within the dataset and cannot be nested within treatments or blocks.
treatment.var	The name of the optional treatment column.
reference.time	The name of the optional time point that all other time points should be compared to (e.g. the first year of data). If not specified, each comparison is between consecutive time points (the first and second year, second and third year, etc.)

Value

The multivariate_change function returns a data frame with the following attributes:

- time.var: A column with the specified time.var and a second column, with '2' appended to the name. Time is subtracted from time2 for dispersion change.
- composition_change: A numeric column that is the distance between the centroids of two time points, or NA if a real distance could not be calculated.
- dispersion_change: A numeric column that is the difference in the average dispersion of the replicates around the centroid for the two time periods. A negative value indicates replicates are converging over time (there is less dispersion at time period 2 than time period 1) and a positive value indicates replicates are diverging over time (there is more dispersion at time period 2 than time period 1).
- treatment.var: A column that has same name and type as the treatment.var column, if treatment.var is specified.

References

Avolio et al. 2015; Avolio et al. Submitted, Marti Anderson et al. 2006.

Examples

```
data(pplots)
# With treatment
multivariate_change(pplots,
  time.var="year",
  replicate.var = "plot",
  treatment.var = "treatment",
  species.var = "species",
  abundance.var = "relative_cover")
# In each year there are 6 replicates and there are 4 years of data for 3
# time comparisons, thus 24 total observations in each treatment.

# With treatment and reference year
multivariate_change(pplots,
  time.var="year",
  replicate.var = "plot",
  treatment.var = "treatment",
  species.var = "species",
  abundance.var = "relative_cover",
  reference.time = 2002)
# In each year there are 6 replicates and there are 4 years of data for 3
# time comparisons, thus 24 total observations in each treatment.

# Without treatment
df <- subset(pplots, treatment == "N1P0")
multivariate_change(df,
  time.var="year",
  replicate.var = "plot",
  species.var = "species",
  abundance.var = "relative_cover")
# In each year there are 6 replicates and there are 4 years of data for 3
# time comparisons, thus 24 total observations.
```

multivariate_difference

Using dissimilarity-based measures to calculate differences in composition and dispersion between pairs of treatments at a single time point

Description

Calculates the difference in composition and dispersion between treatments based off a Bray-Curtis dissimilarity matrix at a single point in time. Composition difference is the pairwise distance between centroids of compared treatments and ranges from 0-1, where identical communities give 0

and completely different communities give 1. Dispersion difference is the difference between treatments in the dispersion of replicates, i.e. the average distance between a replicate and its centroid.

Usage

```
multivariate_difference(df, time.var = NULL, species.var, abundance.var,
  replicate.var, treatment.var, reference.treatment = NULL)
```

Arguments

<code>df</code>	A data frame containing a species, abundance, replicate, and treatment columns and optional time column.
<code>time.var</code>	The name of the optional time column.
<code>species.var</code>	The name of the species column.
<code>abundance.var</code>	The name of the abundance column.
<code>replicate.var</code>	The name of the replicate column. Replicate identifiers must be unique within the dataset and cannot be nested within treatments or blocks.
<code>treatment.var</code>	The name of the treatment column.
<code>reference.treatment</code>	The name of the optional treatment that all other treatments will be compared to (e.g. only controls will be compared to all other treatments). If not specified all pairwise treatment comparisons will be made.

Value

The `multivariate_difference` function returns a data frame with the following attributes:

- `treatment.var`: A column that has same name and type as the `treatment.var` column, if `treatment.var` is specified.
- `treatment.var2`: A column that has the same type as the `treatment.var` column, and is named `treatment.var` with a 2 appended to it.
- `composition_diff`: A numeric column that is the euclidean distance between the centroids of two treatments at a single point in time.
- `abs_dispersion_diff`: A numeric column that is the absolute value of the difference in the average dispersion of the replicates around the centroid for the two treatments.
- `trt_greater_disp`: A column that has same type as the `treatment.var` column, and specifies which of the two treatments has greater dispersion.
- `time.var`: A characteristic column that has the same name and type as the `time.var` column, if specified.

References

Avolio et al. Submitted, Avolio et al. 2015, Marti Anderson et al. 2006

Examples

```
data(pplots)
# Without time
df <- subset(pplots, year == 2002)
multivariate_difference(df,
  replicate.var = "plot",
  treatment.var = "treatment",
  species.var = "species",
  abundance.var = "relative_cover")
# There are 6 replicates for each of three treatments, thus 18 total
# observations.

# Without time and with reference treatment
df <- subset(pplots, year == 2002)
multivariate_difference(df,
  replicate.var = "plot",
  treatment.var = "treatment",
  species.var = "species",
  abundance.var = "relative_cover",
  reference.treatment = "N1P0")
# There are 6 replicates for each of three treatments, thus 18 total
# observations.

# With time
multivariate_difference(pplots,
  time.var = "year",
  replicate.var = "plot",
  species.var = "species",
  abundance.var = "relative_cover",
  treatment.var = "treatment")
# In each year there are 6 replicates for each of three treatments, for a
# total of 18 observations.
```

pplots

Phosphorus plots data from Avolio et al. 2014

Description

A dataset of tallgrass prairie plant composition in a nitrogen and phosphorus addition experiment at Konza Prairie, Manhattan Kansas (Avolio et al. 2014). This dataset is a subset of the full dataset.

Usage

```
data(pplots)
```

Format

A data frame with 1232 rows and 6 variables

Details

A data frame containing a column for replicate, year, species, abundance, block and treatment :

- plot: An integer column of spatial replicates with 18 levels (6-48)
- year: An integer column of sampling time points
- species: A factor column of species sampled
- relative_cover: A numeric column of relative cover values
- block: An integer column of dummy blocking variable, grouping treatment plots into blocks
- treatment: A factor column of nitrogen and phosphorus treatments applied to the plots

Source

Avolio, ML, Koerner, S, La Pierre, K, Wilcox, K, Wilson, GTW, Smith, MD, Collins, S. 2014. Changes in plant community composition, not diversity, to a decade of nitrogen and phosphorus additions drive changes in aboveground productivity in a tallgrass prairie. *Journal of Ecology*. 102: 1649-1660.

RAC_change

Rank Abundance Curve Changes

Description

Calculates change of the five aspects of rank abundance curves (richness, evenness, rank, species gains, and species losses) for a replicate between two time points.

Usage

```
RAC_change(df, time.var, species.var, abundance.var,
           replicate.var = NULL, reference.time = NULL)
```

Arguments

df	A data frame containing time, species, and abundance columns and an optional columns of replicates.
time.var	The name of the time column.
species.var	The name of the species column.
abundance.var	The name of the abundance column.
replicate.var	The name of the optional replicate column. If specified, replicate must be unique within the dataset and cannot be nested within treatments or blocks.
reference.time	The name of the optional time point that all other time points should be compared to (e.g. the first year of data). If not specified, each comparison is between consecutive time points (the first and second year, second and third year, etc.)

Value

The RAC_change function returns a data frame with a subset of the following columns:

- replicate.var: A column that has same name and type as the replicate.var column, if replicate.var is specified.
- time.var: A column with the specified time.var and a second column, with '2' appended to the name. Time is subtracted from time2.
- richness_change: A numeric column that is the change in richness between the two time periods for a replicate divided by the total number of unique species in both time periods. A positive value occurs when there is an increase in species richness over time, and a negative value when there is a decreases in species richness over time.
- evenness_change: A numeric column that is the change in evenness(measured with Evar) between the two time periods for a replicate. A positive value occurs when evenness increases over time, and a negative value when evenness decreases in over time.
- rank_change: A numeric column that is the absolute value of the average change in species ranks between the two time periods for a replicate divided by the total number of unique species in both time periods. Species that are not present in both time periods are given the S+1 rank in the sample it is absent in, where S is the number of species in that sample.
- gains: A numeric column of the number of species that are present at time period 2 that were not present at time period 1 for a replicate divided by the total number of unique species in both time periods. This is equivalent to the turnover function with metric = "appearances".
- losses: A numeric column of the number of species that are not present at time period 2 but were present at time period 1 for a replicate divided by the total number of unique species in both time periods. This is equivalent to the turnover function with metric = "disappearance".

References

Avolio et al. Submitted

Examples

```
data(pplots)
# Without replicates
df <- subset(pplots, plot == 25)
RAC_change(df = df,
           species.var = "species",
           abundance.var = "relative_cover",
           time.var = "year")

# With replicates
df <- subset(pplots, year < 2004 & plot %in% c(6, 25, 32))
RAC_change(df = df,
           species.var = "species",
           abundance.var = "relative_cover",
           replicate.var = "plot",
           time.var = "year")

# With reference year
```

```
df <- subset(pplots, year < 2005 & plot %in% c(6, 25, 32))
RAC_change(df = df,
           species.var = "species",
           abundance.var = "relative_cover",
           replicate.var = "plot",
           time.var = "year",
           reference.time = 2002)
```

RAC_difference

Rank Abundance Curve Differences

Description

Calculates differences between two samples for four comparable aspects of rank abundance curves (richness, evenness, rank, species composition). There are three ways differences can be calculated. 1) Between treatments within a block (note: block.var and treatment.var need to be specified). 2) Between treatments, pooling all replicates into a single species pool (note: pool = TRUE, treatment.var needs to be specified, and block.var will be NULL). 3) All pairwise combinations between all replicates (note: block.var = NULL, pool = FALSE and specifying treatment.var is optional. If treatment.var is specified, the treatment that each replicate belongs to will also be listed in the output).

Usage

```
RAC_difference(df, time.var = NULL, species.var, abundance.var,
              replicate.var, treatment.var = NULL, pool = FALSE,
              block.var = NULL, reference.treatment = NULL)
```

Arguments

df	A data frame containing a species, abundance, and replicate columns and optional time, treatment, and block columns.
time.var	The name of the optional time column.
species.var	The name of the species column.
abundance.var	The name of the abundance column.
replicate.var	The name of the replicate column. Replicate identifiers must be unique within the dataset and cannot be nested within treatments or blocks.
treatment.var	The name of the optional treatment column.
pool	An argument to allow abundance values to be pooled within a treatment. The default value is "FALSE", a value of "TRUE" averages abundance of each species within a treatment at a given time point.
block.var	The name of the optional block column.
reference.treatment	The name of the optional treatment that all other treatments will be compared to (e.g. only controls will be compared to all other treatments). If not specified all pairwise treatment comparisons will be made.

Value

The RAC_difference function returns a data frame with the following attributes:

- `time.var`: A column that has the same name and type as the `time.var` column, if `time.var` is specified.
- `block.var`: A column that has same name and type as the `block.var` column, if `block.var` is specified.
- `replicate.var`: A column that has same name and type as the `replicate.var` column, represents the first replicate being compared. Note, a replicate column will be returned only when `pool` is `FALSE` or `block.var = NULL`.
- `replicate.var2`: A column that has the same type as the `replicate.var` column, and is named `replicate.var` with a 2 appended to it, represents the second replicate being compared. Note, a `replicate.var` column will be returned only when `pool` is `FALSE` and `block.var = NULL`.
- `treatment.var`: A column that has the same name and type as the `treatment.var` column, represents the first treatment being compared. A `treatment.var` column will be returned when `pool` is `TRUE` or `block.var` is present, or `treatment.var` is specified.
- `treatment.var2`: A column that has the same type as the `treatment.var` column, and is named `treatment.var` with a 2 appended to it, represents the second treatment being compared. A `treatment.var` column will be returned when `pool` is `TRUE` or `block.var` is present, or `treatment.var` is specified.
- `richness_diff`: A numeric column that is the difference between the compared samples (treatments or replicates) in species richness divided by the total number of unique species in both samples. A positive value occurs when there is greater species richness in `replicate.var2` than `replicate.var` or `treatment.var2` than `treatment.var`.
- `evenness_diff`: A numeric column of the difference between the compared samples (treatments or replicates) in evenness (measured by Evar). A positive value occurs when there is greater evenness in `replicate.var2` than `replicate.var` or `treatment.var2` than `treatment.var`.
- `rank_diff`: A numeric column of the absolute value of average difference between the compared samples (treatments or replicates) in species' ranks divided by the total number of unique species in both samples. Species that are not present in both samples are given the $S+1$ rank in the sample it is absent in, where S is the number of species in that sample.
- `species_diff`: A numeric column of the number of species that are different between the compared samples (treatments or replicates) divided by the total number of species in both samples. This is equivalent to the Jaccard Index.

References

Avolio et al. Submitted

Examples

```
data(pplots)
# With block and no time
df <- subset(pplots, year == 2002 & block < 3)
RAC_difference(df = df,
              species.var = "species",
```

```

        abundance.var = "relative_cover",
        treatment.var = 'treatment',
        block.var = "block",
        replicate.var = "plot")

# With blocks and time
df <- subset(pplots, year < 2004 & block < 3)
RAC_difference(df = df,
  species.var = "species",
  abundance.var = "relative_cover",
  treatment.var = 'treatment',
  block.var = "block",
  replicate.var = "plot",
  time.var = "year")

# With blocks, time and reference treatment
df <- subset(pplots, year < 2004 & block < 3)
RAC_difference(df = df,
  species.var = "species",
  abundance.var = "relative_cover",
  treatment.var = 'treatment',
  block.var = "block",
  replicate.var = "plot",
  time.var = "year",
  reference.treatment = "N1P0")

# Pooling by treatment with time
df <- subset(pplots, year < 2004)
RAC_difference(df = df,
  species.var = "species",
  abundance.var = "relative_cover",
  treatment.var = 'treatment',
  pool = TRUE,
  replicate.var = "plot",
  time.var = "year")

# All pairwise replicates with treatment
df <- subset(pplots, year < 2004 & plot %in% c(21, 25, 32))
RAC_difference(df = df,
  species.var = "species",
  abundance.var = "relative_cover",
  replicate.var = "plot",
  time.var = "year",
  treatment.var = "treatment")

# All pairwise replicates without treatment
df <- subset(pplots, year < 2004 & plot %in% c(21, 25, 32))
RAC_difference(df = df,
  species.var = "species",
  abundance.var = "relative_cover",
  replicate.var = "plot",
  time.var = "year")

```

rank_shift	<i>Mean Rank Shifts</i>
------------	-------------------------

Description

A measure of the relative change in species rank abundances, which indicates shifts in relative abundances over time (Collins et al. 2008). Mean rank shifts are calculated as the average difference in species' ranks between consecutive time periods, among species that are present across the entire time series.

Usage

```
rank_shift(df, time.var, species.var, abundance.var,  
           replicate.var = as.character(NA))
```

Arguments

df	A data frame containing time, species and abundance columns and an optional column of replicates
time.var	The name of the time column
species.var	The name of the species column
abundance.var	The name of the abundance column
replicate.var	The name of the optional replicate column

Details

The input data frame needs to contain columns for time, species and abundance; time.var, species.var and abundance.var are used to indicate which columns contain those variables. If multiple replicates are included in the data frame, that column should be specified with replicate.var. Each replicate should reflect a single experimental unit - there must be a single abundance value per species within each time point and replicate.

Value

rank_shift returns a data frame with the following columns:

- time.var_pair: A factor column that returns the two time periods being compared, separated by a dash. The name of this column is the same as the time.var column in the input dataframe followed by "_pair".
- MRS: A numeric column with the mean rank shift values.
- replicate.var: A column that has same name and type as the replicate.var column, if replication is specified.

References

Collins, Scott L., Katharine N. Suding, Elsa E. Cleland, Michael Batty, Steven C. Pennings, Katherine L. Gross, James B. Grace, Laura Gough, Joe E. Fargione, and Christopher M. Clark. (2008) "Rank clocks and plant community dynamics." *Ecology* 89, no. 12: 3534-41.

Examples

```
# Calculate mean rank shifts within replicates
data(knz_001d)

myoutput <- rank_shift(knz_001d,
                       time.var = "year",
                       species.var = "species",
                       abundance.var = "abundance",
                       replicate.var = "subplot")

# Calculate mean rank shifts for a data frame with no replication

myoutput_singlerep <- rank_shift(subset(knz_001d, subplot=="A_1"),
                                 time.var = "year",
                                 species.var = "species",
                                 abundance.var = "abundance")
```

rate_change

Rate of community change over successive time intervals

Description

Calculates the slope of the differences in species composition within a community over increasing time intervals, which provides a measure of the rate of directional change in community composition. Differences in species composition are characterized by Euclidean distances, which are calculated on pair-wise communities across the entire time series. For example, a data set with six time intervals will have distance values for five one-year time lags (year 1 vs year 2, year 2 vs year 3 ...), four two-year time lags (year 1 vs year 3, year 2 vs year 4 ...) and so forth. These distance values are regressed against the time lag interval. The slope of the regression line is reported as an indication of the rate and direction of compositional change in the community.

Usage

```
rate_change(df, time.var, species.var, abundance.var, replicate.var = NA)
```

Arguments

df	A data frame containing time, species and abundance columns and an optional column of replicates
time.var	The name of the time column
species.var	The name of the species column
abundance.var	The name of the abundance column
replicate.var	The name of the optional replicate column

Details

The input data frame needs to contain columns for time, species and abundance; `time.var`, `species.var` and `abundance.var` are used to indicate which columns contain those variables. If multiple replicates are included in the data frame, that column should be specified with `replicate.var`. Each replicate should reflect a single experimental unit - there must be a single abundance value per species within each time point and replicate.

The `rate_change` function uses linear regression to relate Euclidean distances to time lag intervals. It is recommended that fit of this relationship be verified using `rate_change_interval`, which returns the full set of community distance values and associated time lag intervals.

Value

The `rate_change` function returns a numeric rate change value unless a replication column is specified in the input data frame. If replication is specified, the function returns a data frame with the following attributes:

- `rate_change`: A numeric column with the synchrony values.
- `replicate.var`: A column that shares the same name and type as the `replicate.var` column in the input data frame.

References

Collins, S. L., Micheli, F. and Hartt, L. 2000. A method to determine rates and patterns of variability in ecological communities. - *Oikos* 91: 285-293.

Examples

```
data(knz_001d)
rate_change(knz_001d[knz_001d$subplot=="A_1", ],
            time.var = "year",
            species.var = "species",
            abundance.var = "abundance") # for one subplot

rate_change(knz_001d,
            time.var = "year",
            species.var = "species",
            abundance.var = "abundance",
            replicate.var = "subplot") # across all subplots
```

rate_change_interval *Differences in community composition over successive time lag intervals*

Description

Calculates the differences in species composition within a community over increasing time intervals. Differences in species composition are characterized by Euclidean distances, which are calculated on pair-wise communities across the entire time series. For example, a data set with 6 time intervals will have distance values for five one-year time lags (year 1 vs year 2, year 2 vs year 3 ...), 4 two-year time lags (year 1 vs year 3, year 2 vs year 4 ...) and so forth. Returns the full set of community distance values and associated time lag intervals.

Usage

```
rate_change_interval(df, time.var, species.var, abundance.var,
  replicate.var = NA)
```

Arguments

df	A data frame containing time, species and abundance columns and an optional column of replicates
time.var	The name of the time column
species.var	The name of the species column
abundance.var	The name of the abundance column
replicate.var	The name of the optional replicate column

Value

The `rate_change_interval` function returns a data frame with the following attributes:

- `interval`: A numeric column containing the interval length between time periods.
- `distance`: A numeric column containing the Euclidean distances.
- `replicate.var`: A column that shares the same name and type as the `replicate.var` column in the input data frame.

The input data frame needs to contain columns for time, species and abundance; `time.var`, `species.var` and `abundance.var` are used to indicate which columns contain those variables. If multiple replicates are included in the data frame, that column should be specified with `replicate.var`. Each replicate should reflect a single experimental unit - there must be a single abundance value per species within each time point and replicate.

References

Collins, S. L., Micheli, F. and Hartt, L. 2000. A method to determine rates and patterns of variability in ecological communities. - *Oikos* 91: 285-293.

Examples

```
data(knz_001d)
rate_change_interval(knz_001d[knz_001d$subplot=="A_1", ],
  time.var = "year",
  species.var = "species",
```

```

abundance.var = "abundance") # for one subplot

rate_change_interval(knz_001d,
  time.var = "year",
  species.var = "species",
  abundance.var = "abundance",
  replicate.var = "subplot") # across all subplots

```

synchrony	<i>Species synchrony</i>
-----------	--------------------------

Description

Calculates the degree synchrony in species abundances within a community over time. Includes the option for two different synchrony metrics. The first, developed by Loreau and de Mazancourt (2008), compares the variance of the aggregated community with the variance of individual components. The second, developed by Gross et al. (2014), compares the average correlation of each individual species with the rest of the aggregated community.

Usage

```

synchrony(df, time.var, species.var, abundance.var, metric = "Loreau",
  replicate.var = NA)

```

Arguments

<code>df</code>	A data frame containing time, species and abundance columns and an optional column of replicates
<code>time.var</code>	The name of the time column
<code>species.var</code>	The name of the species column
<code>abundance.var</code>	The name of the abundance column
<code>metric</code>	The synchrony metric to return: <ul style="list-style-type: none"> • "Loreau": The default metric, calculates synchrony following Loreau and de Mazancourt (2008). • "Gross": Calculates synchrony following Gross et al. (2014).
<code>replicate.var</code>	The name of the optional replicate column

Details

The input data frame needs to contain columns for time, species and abundance; `time.var`, `species.var` and `abundance.var` are used to indicate which columns contain those variables. If multiple replicates are included in the data frame, that column should be specified with `replicate.var`. Each replicate should reflect a single experimental unit - there must be a single abundance value per species within each time point and replicate.

Value

The synchrony function returns a numeric synchrony value unless a replication column is specified in the input data frame. If replication is specified, the function returns a data frame with the following attributes:

- synchrony: A numeric column with the synchrony values.
- replicate.var: A column that shares the same name and type as the replicate.var column in the input data frame.

References

Gross, Kevin, Bradley J. Cardinale, Jeremy W. Fox, Andrew Gonzalez, Michel Loreau, H. Wayne Polley, Peter B. Reich, and Jasper van Ruijven. (2014) "Species richness and the temporal stability of biomass production: A new analysis of recent biodiversity experiments." *The American Naturalist* 183, no. 1: 1-12. doi:10.1086/673915.

Loreau, Michel, and Claire de Mazancourt. (2008) "Species synchrony and its drivers: Neutral and nonneutral community dynamics in fluctuating environments." *The American Naturalist* 172, no. 2: E48-66. doi:10.1086/589746.

Examples

```
data(knz_001d)
synchrony(knz_001d[knz_001d$subplot=="A_1",],
          time.var = "year",
          species.var = "species",
          abundance.var = "abundance") # for one subplot

## Not run:
synchrony(knz_001d,
          time.var = "year",
          species.var = "species",
          abundance.var = "abundance",
          replicate.var = "subplot") # across all subplots

synchrony(knz_001d,
          time.var = "year",
          species.var = "species",
          abundance.var = "abundance",
          replicate.var = "subplot",
          metric="Gross") # With Gross et al. (2014) metric.

## End(Not run)
```

Description

Computes species turnover between time periods as the proportion of species either gained or lost relative to the total number of species observed across both time periods. Includes an option to compute turnover as just the proportion of species gained (i.e., "appearances") or lost (i.e., "disappearances").

Usage

```
turnover(df, time.var, species.var, abundance.var, replicate.var = NA,
         metric = "total")
```

Arguments

<code>df</code>	A data frame containing time, species and abundance columns and an optional column of replicates
<code>time.var</code>	The name of the time column
<code>species.var</code>	The name of the species column
<code>abundance.var</code>	The name of the abundance column
<code>replicate.var</code>	The name of the optional replicate column
<code>metric</code>	The turnover metric to return: <ul style="list-style-type: none"> • <code>total</code>: The default metric, calculates summed appearances and disappearances relative to total species richness across both time periods. • <code>appearance</code>: Calculates the number of species that appeared in the second time period relative to total species richness across both time periods. • <code>disappearance</code>: Calculates the number of species that disappeared in the second time period relative to total species richness across both time periods.

Details

The input data frame needs to contain columns for time, species and abundance; `time.var`, `species.var` and `abundance.var` are used to indicate which columns contain those variables. If multiple replicates are included in the data frame, that column should be specified with `replicate.var`. Each replicate should reflect a single experimental unit - there must be a single abundance value per species within each time point and replicate.

Value

The turnover function returns a data frame with the following attributes:

- `turnover`: A numeric column with the turnover values. The name of this column is the same as the specified metric (default is "total").
- `time.var`: A column containing the second time point; the name and type of this column is the same as the `time.var` column in the input dataframe.
- `replicate.var`: A column that has same name and type as the `replicate.var` column, if replication is specified.

References

Cleland, Elsa E., Scott L. Collins, Timothy L. Dickson, Emily C. Farrer, Katherine L. Gross, Laureano A. Gherardi, Lauren M. Hallett, et al. (2013) "Sensitivity of grassland plant community composition to spatial vs. temporal variation in precipitation." *Ecology* 94, no. 8: 1687-96.

Examples

```
data(knz_001d)

# Calculate relative total turnover within replicates
total.res <- turnover(df=knz_001d,
  time.var = "year",
  species.var = "species",
  abundance.var = "abundance",
  replicate.var="subplot")

# Calculate relative species appearances within replicates
appear.res <- turnover(df=knz_001d,
  time.var = "year",
  species.var = "species",
  abundance.var = "abundance",
  replicate.var="subplot",
  metric="appearance")

# Calculate relative species disappearances within replicates
disappear.res <- turnover(df=knz_001d,
  time.var = "year",
  species.var = "species",
  abundance.var = "abundance",
  replicate.var="subplot",
  metric="disappearance")
```

variance_ratio

Variance Ratio

Description

Computes the ratio of the variance of aggregate species abundances in a community to the sum of the variances of individual, component species. A variance ratio = 1 indicates that species do not covary, a variance ratio > 1 indicates predominately positive covariance among species and a variance ratio < 1 indicates predominately negative covariance (Schluter 1984).

Includes a cyclic shift null modeling option to test if the variance ratio significantly differs from 1. The null community is created by randomly selecting different starting points for each species' time series, which generates a community in which species abundances vary independently but within-species autocorrelation is maintained (Hallett et al. 2014). This randomization is repeated a user-specific number of times and confidence intervals are reported for the resultant null distribution of variance ratios. If the dataframe includes multiple replicates, the variance ratios for the actual and null communities are averaged within each iteration unless specified otherwise.

Usage

```
variance_ratio(df, time.var, species.var, abundance.var, bootnumber,
  replicate.var = NA, average.replicates = TRUE, level = 0.95, li,
  ui)
```

Arguments

df	A data frame containing time, species and abundance columns and an optional column of replicates
time.var	The name of the time column
species.var	The name of the species column
abundance.var	The name of the abundance column
bootnumber	The number of null model iterations used to calculate confidence intervals
replicate.var	The name of the (optional) replicate column
average.replicates	If true returns the variance ratio and CIs averaged
level	The confidence level for the null mean
li	(deprecated) lower confidence interval
ui	(deprecated) upper confidence interval across replicates; if false returns the variance ratio and CI for each replicate. Defaults to true.

Details

The input data frame needs to contain columns for time, species and abundance; time.var, species.var and abundance.var are used to indicate which columns contain those variables. If multiple replicates are included in the data frame, that column should be specified with replicate.var. Each replicate should reflect a single experimental unit - there must be a single abundance value per species within each time point and replicate.

Null model confidence intervals default to the standard lowest 2.5% and upper 97.5% of the null distribution, typically these do not need to be change, but they can be user-modified to set more stringent CIs. @references Hallett, Lauren M., Joanna S. Hsu, Elsa E. Cleland, Scott L. Collins, Timothy L. Dickson, Emily C. Farrer, Laureano A. Gherardi, et al. (2014) "Biotic Mechanisms of Community Stability Shift along a Precipitation Gradient." *Ecology* 95, no. 6: 1693-1700. doi: 10.1890/13-0895.1

Schluter, Dolph. (1984) "A Variance Test for Detecting Species Associations, with Some Example Applications." *Ecology* 65, no. 3: 998-1005. doi:10.2307/1938071.

Value

The variance_ratio function returns a data frame with the following attributes:

- VR: A numeric column with the actual variance ratio value.
- lowerCI: A numeric column with the lowest confidence interval value.
- upperCI: A numeric column with the highest confidence interval value.
- nullmean: A numeric column with the average null variance ratio value.
- replicate.var: A column that has same name and type as the replicate.var column, if replication is specified.

Examples

```
data(knz_001d)

# Calculate the variance ratio and CIs averaged within replicates
# Here the null model is repeated once, for final use it is recommended to set a
# large bootnumber (eg, 10000)

res_averagedreplicates <- variance_ratio(knz_001d,
    time.var = "year",
    species.var = "species",
    abundance.var = "abundance",
    bootnumber = 1,
    replicate = "subplot")

#Calculate the variance ratio and CIs for each replicate

res_withinreplicates <- variance_ratio(knz_001d,
    time.var = "year",
    species.var = "species",
    abundance.var = "abundance",
    bootnumber = 1,
    replicate = "subplot",
    average.replicates = FALSE)
```


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