

Package ‘PlackettLuce’

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Type Package

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Description Functions to prepare rankings data and fit the Plackett-Luce model jointly attributed to Plackett (1975) <[doi:10.2307/2346567](https://doi.org/10.2307/2346567)> and Luce (1959, ISBN:0486441369). The standard Plackett-Luce model is generalized to accommodate ties of any order in the ranking. Partial rankings, in which only a subset of items are ranked in each ranking, are also accommodated in the implementation. Disconnected/weakly connected networks implied by the rankings may be handled by adding pseudo-rankings with a hypothetical item. Optionally, a multivariate normal prior may be set on the log-worth parameters and ranker reliabilities may be incorporated as proposed by Raman and Joachims (2014) <[doi:10.1145/2623330.2623654](https://doi.org/10.1145/2623330.2623654)>. Maximum a posteriori estimation is used when priors are set. Methods are provided to estimate standard errors or quasi-standard errors for inference as well as to fit Plackett-Luce trees. See the package website or vignette for further details.

License GPL-3

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Author Heather Turner [aut, cre] (ORCID:

<<https://orcid.org/0000-0002-1256-3375>>),

Ioannis Kosmidis [aut] (ORCID: <<https://orcid.org/0000-0003-1556-0302>>),

David Firth [aut] (ORCID: <<https://orcid.org/0000-0003-0302-2312>>),

Jacob van Etten [ctb] (ORCID: <<https://orcid.org/0000-0001-7554-2558>>)

Maintainer Heather Turner <ht@heatherturner.net>

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`adjacency`*Create an Adjacency Matrix for a set of Rankings*

Description

Convert a set of rankings to an adjacency matrix summarising wins and losses between pairs of items.

Usage

```
adjacency(object, weights = NULL, ...)
```

Arguments

<code>object</code>	a <code>rankings</code> object, or an object that can be coerced by <code>as.rankings</code> .
<code>weights</code>	an optional vector of weights for the rankings.
<code>...</code>	further arguments passed to/from methods.

Details

For a "rankings" object based on N items, the adjacency matrix is an N by N matrix, with element (i, j) being the number of times item i wins over item j. For example, in the ranking {1} > {3, 4} > {2}, item 1 wins over items 2, 3, and 4, and items 3 and 4 win over item 2.

If `weights` is specified, the values in the adjacency matrix are the weighted counts.

Value

An N by N matrix, where N is the number of items that can be ranked.

Examples

```
X <- matrix(c(2, 1, 2, 1, 2,
             3, 2, 0, 0, 1,
             1, 0, 2, 2, 3), nrow = 3, byrow = TRUE)
X <- as.rankings(X)
adjacency(X)

adjacency(X, weights = c(1, 1, 2))
```

 aggregate

 Aggregate Rankings

Description

Aggregate rankings, returning an "aggregated_rankings" object of the unique rankings and their frequencies. The frequencies can be extracted via the function `freq()`.

Usage

```
## S3 method for class 'rankings'
aggregate(x, freq = NULL, ...)

as.aggregated_rankings(x, ...)

## S3 method for class 'aggregated_rankings'
x[i, j, ..., drop = FALSE, as.aggregated_rankings = TRUE]

freq(x)
```

Arguments

<code>x</code>	A "rankings" object for <code>aggregate()</code> ; an object that can be coerced to a "aggregated_rankings" object for <code>as.aggregated_rankings()</code> , otherwise an "aggregated_rankings" object.
<code>freq</code>	A vector of frequencies for rankings that have been previously aggregated.
<code>...</code>	Additional arguments, currently unused.
<code>i</code>	indices specifying rankings to extract, as for <code>[]</code> .
<code>j</code>	indices specifying items to extract, as for <code>[]</code> .
<code>drop</code>	if TRUE return single row/column matrices as a vector.
<code>as.aggregated_rankings</code>	if TRUE create an "aggregated_rankings" object from the indexed rankings. Otherwise index the underlying matrix of ranks and return in a data frame with the corresponding frequencies.

Value

A data frame of class "aggregated_rankings", with columns

<code>ranking</code>	A "rankings" object of the unique rankings.
<code>freq</code>	The corresponding frequencies.

Methods are available for `rbind()` and `as.matrix()`.

See Also

`preflib()` for an object that can be coerced to an "aggregated_rankings" object.

Examples

```

# create a rankings object with duplicated rankings
R <- matrix(c(1, 2, 0, 0,
             0, 1, 2, 3,
             2, 1, 1, 0,
             1, 2, 0, 0,
             2, 1, 1, 0,
             1, 0, 3, 2), nrow = 6, byrow = TRUE)
colnames(R) <- c("apple", "banana", "orange", "pear")
R <- as.rankings(R)

# aggregate the rankings
A <- aggregate(R)

# subsetting applies to the rankings, e.g. first two unique rankings
A[1:2]

# (partial) rankings of items 2 to 4 only
A[, 2:4]

# convert to a matrix
as.matrix(A)

# frequencies are automatically used as weights by PlackettLuce()
mod <- PlackettLuce(A)
mod$weights

```

beans

Preferred Bean Varieties in Nicaragua

Description

This is a subset of data from trials of bean varieties (*Phaseolus vulgaris* L.) in Nicaragua over five growing seasons. Farmers were asked to try three varieties of bean from a total of ten varieties and to rank them in order of preference. In addition, for each variety the farmers were asked to compare each trial variety to the local variety and state whether they considered it to be better or worse.

Usage

```
beans
```

Format

A data frame with 842 records and 14 variables:

variety_a The name of variety A in the comparison.

variety_b The name of variety B in the comparison.

variety_c The name of variety C in the comparison.

best The variety the farmer ranked in first place ("A", "B" or "C").

worst The variety the farmer ranked in last place ("A", "B" or "C").

var_a How the farmer ranked variety A compared to the local variety ("Worse" or "Better").

var_b How the farmer ranked variety B compared to the local variety ("Worse" or "Better").

var_c How the farmer ranked variety C compared to the local variety ("Worse" or "Better").

season A factor specifying the growing season ("Po - 15", "Ap - 15", "Pr - 16", "Po - 16", "Ap - 16").

year The year of planting.

maxTN The maximum temperature at night during the vegetative cycle (degrees Celsius).

lon The geographic coordinate longitude (X axis) for where the plot was established.

lat The geographic coordinate latitude (Y axis) for where the plot was established.

planting_date A Date, specifying the start date of planting the trial.

Details

There are three crop seasons in Central America:

Primera May - August.

Postrera September - October.

Apante November - January.

Beans can be planted near the beginning of each season, though are most commonly planted in the Postrera or Apante seasons.

Source

van Etten, J. et al. (2019) *PNAS*, **116** (10), 4194–4199, [doi:10.1073/pnas.1813720116](https://doi.org/10.1073/pnas.1813720116).

Examples

```
# Consider the best and worst rankings. These give the variety the
# farmer thought was best or worst, coded as A, B or C for the
# first, second or third variety assigned to the farmer
# respectively.
data(bean)
head(bean[c("best", "worst")], 2)

# Fill in the missing item
bean$middle <- complete(bean[c("best", "worst")],
                        items = c("A", "B", "C"))
head(bean[c("best", "middle", "worst")], 2)

# This gives an ordering of the three varieties the farmer was
# given. The names of these varieties are stored in separate
# columns
varieties <- bean[c("variety_a", "variety_b", "variety_c")]
head(varieties, 2)
```

```

# Use these names to decode the orderings of order 3
order3 <- decode(beans[c("best", "middle", "worst")],
  items = beans[c("variety_a", "variety_b", "variety_c")],
  code = c("A", "B", "C"))

# Now consider the paired comparisons against the local variety
head(beans[c("var_a", "var_b", "var_c")], 2)

# Convert these results to a vector and get the corresponding trial variety
outcome <- unlist(beans[c("var_a", "var_b", "var_c")])
trial_variety <- unlist(beans[c("variety_a", "variety_b", "variety_c")])

# Create a data frame of the implied orderings of order 2
order2 <- data.frame(Winner = ifelse(outcome == "Worse",
  "Local", trial_variety),
  Loser = ifelse(outcome == "Worse",
  trial_variety, "Local"),
  stringsAsFactors = FALSE, row.names = NULL)

head(order2, 2)

# Finally combine the rankings of order 2 and order 3
R <- rbind(as.rankings(order3, input = "orderings"),
  as.rankings(order2, input = "orderings"))

head(R)
tail(R)

```

 choices

Choices Object

Description

Convert a set of rankings to a list of choices, alternatives, and rankings. The choices and the corresponding alternatives make up the exchangeable part of the Plackett-Luce with ties.

Usage

```
choices(rankings, names = FALSE)
```

Arguments

rankings	a " rankings " object, or an object that can be coerced by <code>as.rankings</code> .
names	logical: if TRUE use the object names in the returned "choices" object, else use object indices.

Value

A data frame of class "choices" with elements:

choices	A list where each element represents the set of items chosen for a single rank in the ranking.
---------	--

alternatives A list where each element represents the set of items to choose from for a single rank in the ranking.

ranking A list where each element represents the ranking that the choice belongs to.

The list stores the number of choices and the names of the objects as the attributes "nchoices" and "objects" respectively.

Examples

```
R <- matrix(c(1, 2, 0, 0,
             4, 1, 2, 3,
             2, 1, 1, 1,
             1, 2, 3, 0,
             2, 1, 1, 0,
             1, 0, 3, 2), nrow = 6, byrow = TRUE)
colnames(R) <- c("apple", "banana", "orange", "pear")

actual_choices <- choices(R, names = TRUE)
actual_choices[1:6,]

coded_choices <- choices(R, names = FALSE)
coded_choices[1:2,]
as.data.frame(coded_choices)[1:2,]
attr(coded_choices, "objects")
```

complete

Complete Orderings with the Missing Redundant Rank

Description

Given orderings with one rank missing, complete the ordering by assigning the remaining item(s) to the final rank.

Usage

```
complete(orderings, items)
```

Arguments

orderings A data frame of orderings with one rank missing.

items A vector of item names.

Value

A vector of the missing items, which will be a list if there are any ties.

Examples

```
# Orderings of 3 items, when only the best and worst are recorded
orderings <- data.frame(best = c("A", "B", "A"),
                        worst = c("C", "C", NA))
orderings$middle <- complete(orderings, items = c("A", "B", "C"))
```

connectivity

Check Connectivity of Rankings

Description

Check the connectivity of the network underlying a set of rankings.

Usage

```
connectivity(x, verbose = TRUE)
```

Arguments

x	an adjacency matrix as returned by adjacency , a "rankings" object, or an object that can be coerced by <code>as.rankings</code> .
verbose	logical, if TRUE, a message is given if the network is not strongly connected.

Details

Ranked items are connected in a directed graph according to the implied wins and losses between pairs of items. The wins and losses can be summarised as an adjacency matrix using [adjacency](#). From this adjacency matrix, the graph is inferred and it is checked for connectivity. A message is given if the network is not strongly connected, i.e. with at least one win and one loss between all partitions of the network into two groups. Features of clusters in the network are returned - if the network is strongly connected, all items belong to the same cluster.

Value

A list with elements

membership	a labelled vector of indices specifying membership of clusters in the network of items
csize	the sizes of clusters in the network of items
no	the number of clusters in the network of items

Examples

```

## weakly connected network:
## one win between two clusters
X <- matrix(c(1, 2, 0, 0,
              2, 1, 3, 0,
              0, 0, 1, 2,
              0, 0, 2, 1), ncol = 4, byrow = TRUE)
X <- as.rankings(X)
res <- connectivity(X)
res$membership
## keep items in cluster 1
na.omit(X[,res$membership == 1])

## two weakly connected items:
## item 1 always loses; item 4 only wins against item 1
X <- matrix(c(4, 1, 2, 3,
              0, 2, 1, 3), nr = 2, byrow = TRUE)
X <- as.rankings(X)
res <- connectivity(X)
res$membership

## item 1 always wins; item 4 always loses
X <- matrix(c(1, 2, 3, 4,
              1, 3, 2, 4), nr = 2, byrow = TRUE)
res <- connectivity(as.rankings(X))
res$membership

## all in separate clusters: always 1 > 2 > 3 > 4
## also miscoded rankings and redundant ranking
X <- matrix(c(1, 2, 3, 4,
              1, 0, 2, 3,
              1, 1, 2, 0,
              1, 0, 3, 4,
              2, 2, 0, 4,
              0, 0, 3, 0,
              2, 4, 0, 0), ncol = 4, byrow = TRUE)
res <- connectivity(as.rankings(X))
res$membership

```

 decode

Decode Orderings using a Key to Item Names

Description

Decode orderings by replacing numeric or character coded values with item names.

Usage

```
decode(orderings, items, code = NULL)
```

Arguments

orderings	A data frame of coded orderings.
items	A data frame of the items in each ranking, or a vector of common items.
code	(Optional) a vector giving the key to the code. If missing, <code>names(items)</code> is used for a character code, while <code>seq(items)</code> is used for a numeric code.

Value

A data frame with the coded values replaced by the item names.

Examples

```
# orderings of up to 3 items coded as A, B, C
orderings <- data.frame(Rank1 = c("A", "B"),
                       Rank2 = c("C", "A"),
                       Rank3 = c("B", NA),
                       stringsAsFactors = FALSE)
items <- data.frame(A = c("banana", "apple"),
                   B = c("orange", "pear"),
                   C = c("apple", NA),
                   stringsAsFactors = FALSE)
decode(orderings, items)

# orderings with ties of up to 3 items, coded 1:3
orderings <- data.frame(Rank1 = c(1, 3),
                       Rank2 = I(list(c(2, 3), 2)),
                       Rank3 = c(NA, 1),
                       stringsAsFactors = FALSE)
items <- data.frame(A = c("banana", "apple"),
                   B = c("orange", "pear"),
                   C = c("apple", "orange"),
                   stringsAsFactors = FALSE)
decode(orderings, items)

# same items in each comparison
items <- c(A = "banana", B = "orange", C = "pear")
decode(orderings, items)
```

fitted.PlackettLuce *Fitted Probabilities for PlackettLuce Objects*

Description

Fitted probabilities for all choice/alternative combinations in the data.

Usage

```
## S3 method for class 'PlackettLuce'
fitted(object, aggregate = TRUE, free = TRUE, ...)

## S3 method for class 'pltree'
fitted(object, aggregate = TRUE, free = TRUE, ...)
```

Arguments

<code>object</code>	an object as returned by PlackettLuce or pltree .
<code>aggregate</code>	logical; if TRUE observations of the same choice from the same set of alternatives are aggregated.
<code>free</code>	logical; if TRUE only free choices are included, i.e. choices of one item from a set of one item are excluded.
<code>...</code>	further arguments, currently ignored.

Value

A list with the following components

<code>choices</code>	The selected item(s).
<code>alternatives</code>	The set of item(s) that the choice was made from.
<code>ranking</code>	The ranking(s) including this choice.
<code>n</code>	The weighted count of rankings including this choice (equal to the ranking weight if <code>aggregate = FALSE</code>).
<code>fitted</code>	The fitted probability of making this choice.

If `object` was a "pltree" object, the list has an additional element, `node`, specifying which node the ranking corresponds to.

See Also

[choices](#)

Examples

```
R <- matrix(c(1, 2, 0, 0,
             4, 1, 2, 3,
             2, 1, 1, 1,
             1, 2, 3, 0,
             2, 1, 1, 0,
             1, 0, 3, 2), nrow = 6, byrow = TRUE)
colnames(R) <- c("apple", "banana", "orange", "pear")

mod <- PlackettLuce(R)
fit <- fitted(mod)
fit
```

group	<i>Group Rankings</i>
-------	-----------------------

Description

Create an object of class "grouped_rankings" which associates a group index with an object of class "rankings". This allows the rankings to be linked to covariates with group-specific values as the basis for model-based recursive partitioning, see [pltree](#).

Usage

```
group(x, index, ...)

as.grouped_rankings(x, ...)

## S3 method for class 'paircomp'
as.grouped_rankings(x, ...)

## S3 method for class 'grouped_rankings'
x[i, j, ..., drop = TRUE, as.grouped_rankings = TRUE]

## S3 method for class 'grouped_rankings'
format(x, max = 2L, width = 20L, ...)
```

Arguments

x	a "rankings" object for group(); an object that can be coerced to a "grouped_rankings" object for as.grouped_rankings(), otherwise a "rankings" object.
index	a numeric vector of length equal to the number of rankings specifying the subject for each ranking.
...	additional arguments passed on to as.rankings by grouped_rankings or as.grouped_rankings; unused by format.
i	indices specifying groups to extract, may be any data type accepted by [] .
j	indices specifying items to extract, as for [] .
drop	if TRUE return single row/column matrices as a vector.
as.grouped_rankings	if TRUE return a grouped_rankings object, otherwise return a matrix/vector.
max	the maximum number of rankings to format per subject.
width	the maximum width in number of characters to format each ranking.

Value

An object of class "grouped_rankings", which is a vector of of group IDs with the following attributes:

rankings	The "rankings" object.
index	An index match each ranking to each group ID.
R	A matrix with items ordered from last to first place, for each ranking.
S	The rankings matrix with the ranks replaced by the size of the chosen set for free choices and zero for forced choices.
id	A list with elements of the adjacency matrix that are incremented by each ranking.

See Also[pltree](#)**Examples**

```
# ungrouped rankings (5 rankings, 4 items)
R <- as.rankings(matrix(c(1, 2, 0, 0,
                        0, 2, 1, 0,
                        0, 0, 1, 2,
                        2, 1, 0, 0,
                        0, 1, 2, 3), ncol = 4, byrow = TRUE))

length(R)
R

# group rankings (first three in group 1, next two in group 2)
G <- group(R, c(1, 1, 1, 2, 2))
length(G)

## by default up to 2 rankings are shown per group, "..." indicates if
## there are further rankings
G
print(G, max = 1)

## select rankings from group 1
G[1,]

## exclude item 3 from ranking
G[, -3]

## rankings from group 2, excluding item 3
## - note group 2 becomes the first group
G[2, -3]

## index underlying rankings without creating new grouped_rankings object
G[2, -3, as.grouped_rankings = FALSE]
```

itempar.PlackettLuce *Extract Item Parameters of Plackett-Luce Models*

Description

Methods for `itempar` to extract the item parameters (worth or log-worth) from a Plackett-Luce model or tree. In the case of a tree, item parameters are extracted for each terminal node.

Usage

```
## S3 method for class 'PlackettLuce'  
itempar(object, ref = NULL, alias = TRUE, vcov = TRUE, log = FALSE, ...)  
  
## S3 method for class 'pltree'  
itempar(object, ...)  
  
## S3 method for class 'PLADMM'  
itempar(object, ref = NULL, alias = TRUE, vcov = TRUE, log = FALSE, ...)
```

Arguments

<code>object</code>	a fitted model object as returned by <code>PlackettLuce</code> , <code>pladmm</code> , or <code>pltree</code> .
<code>ref</code>	a vector of labels or position indices of item parameters which should be used as restriction/for normalization. If <code>NULL</code> (the default), all items are used with a zero sum (<code>log = TRUE</code>) or unit sum (<code>log = FALSE</code>) constraint.
<code>alias</code>	logical. If <code>TRUE</code> (the default), the aliased parameter is included in the return vector (and in the variance-covariance matrix if <code>vcov = TRUE</code>). If <code>FALSE</code> , it is removed. If the restriction given in <code>ref</code> depends on several parameters, the first parameter of the restriction specified is (arbitrarily) chosen to be removed if <code>alias</code> is <code>FALSE</code> .
<code>vcov</code>	logical. If <code>TRUE</code> (the default), the (transformed) variance-covariance matrix of the item parameters is attached as attribute <code>vcov</code> . If <code>FALSE</code> , a NA-matrix is attached.
<code>log</code>	logical. Whether to return log-abilities (<code>TRUE</code>) or abilities (<code>FALSE</code>).
<code>...</code>	further arguments which are currently not used.

Value

An object of class "itempar", see `itempar`.

Examples

```
R <- matrix(c(1, 2, 0, 0,  
             4, 1, 2, 3,  
             2, 1, 1, 1,  
             1, 2, 3, 0,
```

```

      2, 1, 1, 0,
      1, 0, 3, 2), nrow = 6, byrow = TRUE)
colnames(R) <- c("apple", "banana", "orange", "pear")

mod <- PlackettLuce(R)
coef(mod)

# equivalent to default coefficients, i.e. log abilities
itempar(mod, ref= 1, log = TRUE)

# abilities, normalized so abilities for apple and pear sum to 1
itempar(mod, ref = 1:2)

```

nascar

Results from 2002 NASCAR Season

Description

This is an example dataset from *Hunter 2004* recording the results of 36 car races in the 2002 NASCAR season in the United States. Each record is an ordering of the drivers according to their finishing position.

Usage

```
nascar
```

Format

A matrix with 36 rows corresponding to the races and 43 columns corresponding to the positions. The columns contain the ID for the driver that came first to last place respectively. The "drivers" attribute contains the names of the 87 drivers.

References

Hunter, D. R. (2004) MM algorithms for generalized Bradley-Terry models. *The Annals of Statistics*, **32**(1), 384–406.

Examples

```

# convert orderings to rankings
nascar[1:2, ]
R <- as.rankings(nascar, input = "orderings",
                items = attr(nascar, "drivers"))
R[1:2, 1:4, as.rankings = FALSE]
format(R[1:2], width = 60)

# fit model as in Hunter 2004, excluding drivers that only lose
keep <- seq_len(83)

```

```

R2 <- R[, keep]
mod <- PlackettLuce(R2, npseudo = 0)

# show coefficients as in Table 2 of Hunter 2004
avRank <- apply(R, 2, function(x) mean(x[x > 0]))
coefs <- round(coef(mod)[order(avRank[keep])], 2)
head(coefs, 3)
tail(coefs, 3)

```

PlackettLuce

Fit a Plackett-Luce Model

Description

Fit a Plackett-Luce model to a set of rankings. The rankings may be partial (each ranking completely ranks a subset of the items) and include ties of arbitrary order.

Usage

```

PlackettLuce(
  rankings,
  npseudo = 0.5,
  normal = NULL,
  gamma = NULL,
  adherence = NULL,
  weights = freq(rankings),
  na.action = getOption("na.action"),
  start = NULL,
  method = c("iterative scaling", "BFGS", "L-BFGS"),
  epsilon = 1e-07,
  steffensen = 0.1,
  maxit = c(500, 10),
  trace = FALSE,
  verbose = TRUE,
  ...
)

```

Arguments

- | | |
|----------|---|
| rankings | a " rankings " object, or an object that can be coerced by <code>as.rankings</code> . An " aggregated_rankings " object can be used to specify rankings and weights simultaneously. A " grouped_rankings " object should be used when estimating adherence for rankers with multiple rankings per ranker. |
| npseudo | when using pseudodata: the number of wins and losses to add between each object and a hypothetical reference object. |
| normal | a optional list with elements named <code>mu</code> and <code>Sigma</code> specifying the mean and covariance matrix of a multivariate normal prior on the <i>log</i> worths. |

<code>gamma</code>	a optional list with elements named <code>shape</code> and <code>rate</code> specifying parameters of a gamma prior on adherence parameters for each ranker (use <code>grouped_rankings</code> to group multiple rankings by ranker). The short-cut <code>TRUE</code> may be used to specify a <code>Gamma(10, 10)</code> prior. If <code>NULL</code> (or <code>FALSE</code>), adherence is fixed to adherence for all rankers.
<code>adherence</code>	an optional vector of adherence values for each ranker. If missing, adherence is fixed to 1 for all rankers. If <code>gamma</code> is not <code>NULL</code> , this specifies the starting values for the adherence.
<code>weights</code>	an optional vector of weights for each ranking.
<code>na.action</code>	a function to handle any missing rankings, see <code>na.omit()</code> .
<code>start</code>	starting values for the worth parameters and the tie parameters on the raw scale (worth parameters need not be scaled to sum to 1). If <code>normal</code> is specified, <code>exp(normal\$mu)</code> is used as starting values for the worth parameters. Coefficients from a previous fit can be passed as the result of a call to <code>coef.PlackettLuce</code> , or the <code>coefficients</code> element of a "PlackettLuce" object.
<code>method</code>	the method to be used for fitting: "iterative scaling" (iterative scaling to sequentially update the parameter values), "BFGS" (the BFGS optimisation algorithm through the <code>optim</code> interface), "L-BFGS" (the limited-memory BFGS optimisation algorithm as implemented in the <code>lbfgs</code> package). Iterative scaling is used by default, unless a prior is specified by <code>normal</code> or <code>gamma</code> , in which case the default is "BFGS".
<code>epsilon</code>	the maximum absolute difference between the observed and expected sufficient statistics for the ability parameters at convergence.
<code>steffensen</code>	a threshold defined as for <code>epsilon</code> after which to apply Steffensen acceleration to the iterative scaling updates.
<code>maxit</code>	a vector specifying the maximum number of iterations. If <code>gamma</code> is <code>NULL</code> , only the first element is used and specifies the maximum number of iterations of the algorithm specified by <code>method</code> . If <code>gamma</code> is not <code>NULL</code> , a second element may be supplied to specify the maximum number of iterations of an alternating algorithm, where the adherence parameters are updated alternately with the other parameters. The default is to use 10 outer iterations.
<code>trace</code>	logical, if <code>TRUE</code> show trace of iterations.
<code>verbose</code>	logical, if <code>TRUE</code> show messages from validity checks on the rankings.
<code>...</code>	additional arguments passed to <code>optim</code> or <code>lbfgs</code> . In particular the convergence tolerance may be adjusted using e.g. <code>control = list(reltol = 1e-10)</code> .

Value

An object of class "PlackettLuce", which is a list containing the following elements:

<code>call</code>	The matched call.
<code>coefficients</code>	The model coefficients.
<code>loglik</code>	The maximized log-likelihood.
<code>null.loglik</code>	The maximized log-likelihood for the null model (all alternatives including ties have equal probability).

<code>df.residual</code>	The residual degrees of freedom.
<code>df.null</code>	The residual degrees of freedom for the null model.
<code>rank</code>	The rank of the model.
<code>logposterior</code>	If a prior was specified, the maximised log posterior.
<code>gamma</code>	If a gamma prior was specified, the list of parameters.
<code>normal</code>	If a normal prior was specified, the list of parameters.
<code>iter</code>	The number of iterations run.
<code>rankings</code>	The rankings passed to <code>rankings</code> , converted to a "rankings" object if necessary.
<code>weights</code>	The weights applied to each ranking in the fitting.
<code>adherence</code>	The fixed or estimated adherence per ranker.
<code>ranker</code>	The ranker index mapping rankings to rankers (the "index" attribute of rankings if specified as a "grouped_rankings" object.)
<code>ties</code>	The observed tie orders corresponding to the estimated tie parameters.
<code>conv</code>	The convergence code: 0 for successful convergence; 1 if reached <code>maxit</code> (outer) iterations without convergence; 2 if Steffensen acceleration cause log-likelihood to increase; negative number if L-BFGS algorithm failed for other reason.

Model definition

A single ranking is given by

$$R = \{C_1, C_2, \dots, C_J\}$$

where the items in set C_1 are ranked higher than (better than) the items in C_2 , and so on. If there are multiple objects in set C_j these items are tied in the ranking.

For a set of items S , let

$$f(S) = \delta_{|S|} \left(\prod_{i \in S} \alpha_i \right)^{\frac{1}{|S|}}$$

where $|S|$ is the cardinality (size) of the set, δ_n is a parameter related to the prevalence of ties of order n (with $\delta_1 \equiv 1$), and α_i is a parameter representing the worth of item i . Then under an extension of the Plackett-Luce model allowing ties up to order D , the probability of the ranking R is given by

$$\prod_{j=1}^J \frac{f(C_j)}{\sum_{k=1}^{\min(D_j, D)} \sum_{S \in \binom{A_j}{k}} f(S)}$$

where D_j is the cardinality of A_j , the set of alternatives from which C_j is chosen, and $\binom{A_j}{k}$ is all the possible choices of k items from A_j . The value of D can be set to the maximum number of tied items observed in the data, so that $\delta_n = 0$ for $n > D$.

When the worth parameters are constrained to sum to one, they represent the probability that the corresponding item comes first in a ranking of all items, given that first place is not tied.

The 2-way tie prevalence parameter δ_2 is related to the probability that two items *of equal worth* tie for first place, given that the first place is not a 3-way or higher tie. Specifically, that probability is $\delta_2 / (2 + \delta_2)$.

The 3-way and higher tie-prevalence parameters are similarly interpretable, in terms of tie probabilities among equal-worth items.

When intermediate tie orders are not observed (e.g. ties of order 2 and order 4 are observed, but no ties of order 3), the maximum likelihood estimate of the corresponding tie prevalence parameters is zero, so these parameters are excluded from the model.

Pseudo-rankings

In order for the maximum likelihood estimate of an object's worth to be defined, the network of rankings must be strongly connected. This means that in every possible partition of the objects into two nonempty subsets, some object in the second set is ranked higher than some object in the first set at least once.

If the network of rankings is not strongly connected then pseudo-rankings may be used to connect the network. This approach posits a hypothetical object with log-worth 0 and adds npseudo wins and npseudo losses to the set of rankings.

The parameter npseudo is the prior strength. With npseudo = 0 the MLE is the posterior mode. As npseudo approaches infinity the log-worth estimates all shrink towards 0. The default, npseudo = 0.5, is sufficient to connect the network and has a weak shrinkage effect. Even for networks that are already connected, adding pseudo-rankings typically reduces both the bias and variance of the estimators of the worth parameters.

Incorporating prior information on log-worths

Prior information can be incorporated by using normal to specify a multivariate normal prior on the log-worths. The log-worths are then estimated by maximum a posteriori (MAP) estimation. Model summaries (deviance, AIC, standard errors) are based on the log-likelihood evaluated at the MAP estimates, resulting in a finite sample bias that should disappear as the number of rankings increases. Inference based on these model summaries is valid as long as the prior is considered fixed and not tuned as part of the model.

Incorporating a prior is an alternative method of penalization, therefore npseudo is set to zero when a prior is specified.

Incorporating ranker adherence parameters

When rankings come from different rankers, the model can be extended to allow for varying reliability of the rankers, as proposed by Raman and Joachims (2014). In particular, replacing $f(S)$ by

$$h(S) = \delta_{|S|} \left(\prod_{i \in S} \alpha_i \right)^{\frac{\eta_g}{|S|}}$$

where $\eta_g > 0$ is the adherence parameter for ranker g . In the standard model, all rankers are assumed to have equal reliability, so $\eta_g = 1$ for all rankers. Higher $\eta_g = 1$ increases the distance between item worths, giving greater weight to the ranker's choice. Conversely, lower $\eta_g = 1$ shrinks the item worths towards equality so the ranker's choice is less relevant.

The adherence parameters are not estimable by maximum likelihood, since for given item worths the maximum likelihood estimate of adherence would be infinity for rankers that give rankings consistent with the items ordered by worth and zero for all other rankers. Therefore it is essential

to include a prior on the adherence parameters when these are estimated rather than fixed. Setting `gamma = TRUE` specifies the default $\Gamma(10, 10)$ prior, which has a mean of 1 and a probability of 0.99 that the adherence is between 0.37 and 2. Alternative parameters can be specified by a list with elements `shape` and `rate`. Setting `scale` and `rate` to a common value θ specifies a mean of 1; $\theta \geq 2$ will give low prior probability to near-zero adherence; as θ increases the density becomes more concentrated (and more symmetrical) about 1.

Since the number of adherence parameters will typically be large and it is assumed the worth and tie parameters are of primary interest, the adherence parameters are not included in model summaries, but are included in the returned object.

Controlling the fit

For models without priors, using `nspseudo = 0` will use standard maximum likelihood, if the network is connected (and throw an error otherwise).

The fitting algorithm is set by the `method` argument. The default method `"iterative scaling"` is a slow but reliable approach. In addition, this has the most control on the accuracy of the final fit, since convergence is determined by direct comparison of the observed and expected values of the sufficient statistics for the worth parameters, rather than a tolerance on change in the log-likelihood.

The `"iterative scaling"` algorithm is slow because it is a first order method (does not use derivatives of the likelihood). From a set of starting values that are 'close enough' to the final solution, the algorithm can be accelerated using [Steffensen's method](#). `PlackettLuce` attempts to apply Steffensen's acceleration when all differences between the observed and expected values of the sufficient statistics are less than `steffensen`. This is an ad-hoc rule defining 'close enough' and in some cases the acceleration may produce negative worth parameters or decrease the log-likelihood. `PlackettLuce` will only apply the update when it makes an improvement.

The `"BFGS"` and `"L-BFGS"` algorithms are second order methods, therefore can be quicker than the default method. Control parameters can be passed on to `optim` or `lbfgs`.

For models with priors, the iterative scaling method cannot be used, so BFGS is used by default.

Note

As the maximum tie order increases, the number of possible choices for each rank increases rapidly, particularly when the total number of items is high. This means that the model will be slower to fit with higher D . In addition, due to the current implementation of the `vcov()` method, computation of the standard errors (as by `summary()`) can take almost as long as the model fit and may even become infeasible due to memory limits. As a rule of thumb, for > 10 items and > 1000 rankings, we recommend `PlackettLuce()` for ties up to order 4. For higher order ties, a rank-ordered logit model, see `ROlogit::rologit()` or generalized Mallows Model as in `BayesMallows::compute_mallows()` may be more suitable, as they do not model tied events explicitly.

References

Raman, K. and Joachims, T. (2014) Methods for Ordinal Peer Grading. [arXiv:1404.3656](#).

See Also

Handling rankings: [rankings](#), [aggregate](#), [group](#), [choices](#), [adjacency](#), [connectivity](#).

Inspect fitted Plackett-Luce models: `coef`, `deviance`, `fitted`, `itempar`, `logLik`, `print`, `qvcalc`, `summary`, `vcov`.

Fit Plackett-Luce tree: `pltree`.

Example data sets: `beans`, `nascar`, `pudding`, `preflib`.

Vignette: `vignette("Overview", package = "PlackettLuce")`.

Examples

```
# Six partial rankings of four objects, 1 is top rank, e.g
# first ranking: item 1, item 2
# second ranking: item 2, item 3, item 4, item 1
# third ranking: items 2, 3, 4 tie for first place, item 1 second
R <- matrix(c(1, 2, 0, 0,
             4, 1, 2, 3,
             2, 1, 1, 1,
             1, 2, 3, 0,
             2, 1, 1, 0,
             1, 0, 3, 2), nrow = 6, byrow = TRUE)
colnames(R) <- c("apple", "banana", "orange", "pear")

# create rankings object
R <- as.rankings(R)

# Standard maximum likelihood estimates
mod_mle <- PlackettLuce(R, npseudo = 0)
coef(mod_mle)

# Fit with default settings
mod <- PlackettLuce(R)
# log-worths are shrunk towards zero
coef(mod)

# independent N(0, 9) priors on log-worths, as in Raman and Joachims
prior <- list(mu = rep(0, ncol(R)),
             Sigma = diag(rep(9, ncol(R))))
mod_normal <- PlackettLuce(rankings = R, normal = prior)
# slightly weaker shrinkage effect vs pseudo-rankings,
# with less effect on tie parameters (but note small number of rankings here)
coef(mod_normal)

# estimate adherence assuming every ranking is from a separate ranker
mod_separate <- PlackettLuce(rankings = R, normal = prior, gamma = TRUE)
coef(mod_separate)
# gives more weight to rankers 4 & 6 which rank apple first,
# so worth of apple increased relative to banana
mod_separate$adherence

# estimate adherence based on grouped rankings
# - assume two rankings from each ranker
G <- group(R, rep(1:3, each = 2))
mod_grouped <- PlackettLuce(rankings = G, normal = prior, gamma = TRUE)
```

```
coef(mod_grouped)
# first ranker is least consistent so down-weighted
mod_grouped$adherence
```

pladmm

Fit a Plackett-Luce Model with Linear Predictor for Log-worth

Description

Fit a Plackett-Luce model where the log-worth is predicted by a linear function of covariates. The rankings may be partial (each ranking completely ranks a subset of the items), but ties are not supported.

Usage

```
pladmm(
  rankings,
  formula,
  data = NULL,
  weights = freq(rankings),
  start = NULL,
  contrasts = NULL,
  rho = 1,
  n_iter = 500,
  rtol = 1e-04
)
```

Arguments

rankings	a " rankings " object, or an object that can be coerced by <code>as.rankings</code> . An " aggregated_rankings " object can be used to specify rankings and weights simultaneously.
formula	a formula specifying the linear model for log-worth.
data	a data frame containing the variables in the model.
weights	weights for the rankings.
start	starting values for the coefficients.
contrasts	an optional list specifying contrasts for the factors in formula. See the <code>contrasts.arg</code> of model.matrix() .
rho	the penalty parameter in the penalized likelihood, see details.
n_iter	the maximum number of iterations (also for inner loops).
rtol	the convergence tolerance (also for inner loops)

Details

The log-worth is modelled as a linear function of item covariates:

$$\log \alpha_i = \beta_0 + \beta_1 x_{i1} + \dots + \beta_p x_{ip}$$

where β_0 is fixed by the constraint that $\sum_i \alpha_i = 1$.

The parameters are estimated using an Alternating Directions Method of Multipliers (ADMM) algorithm proposed by Yildiz (2020). ADMM alternates between estimating the worths α_i and the linear coefficients β_k , encapsulating them in a quadratic penalty on the likelihood:

$$L(\boldsymbol{\beta}, \boldsymbol{\alpha}, \mathbf{u}) = \mathcal{L}(\mathcal{D}|\boldsymbol{\alpha}) + \frac{\rho}{2} \|\mathbf{X}\boldsymbol{\beta} - \log \boldsymbol{\alpha} + \mathbf{u}\|_2^2 - \frac{\rho}{2} \|\mathbf{u}\|_2^2$$

where \mathbf{u} is a dual variable that imposes the equality constraints (so that $\log \boldsymbol{\alpha}$ converges to $\mathbf{X}\boldsymbol{\beta}$).

Note

This is a prototype function and the user interface is planned to change in upcoming versions of PlackettLuce.

References

Yildiz, I., Dy, J., Erdogmus, D., Kalpathy-Cramer, J., Ostmo, S., Campbell, J. P., Chiang, M. F. and Ioannidis, S. (2020) Fast and Accurate Ranking Regression In Proceedings of the Twenty Third International Conference on Artificial Intelligence and Statistics, **108**, 77—88.

Examples

```
# data.frame of rankings for salad dressings A B C D
# 1 = most tart, 4 = least tart
salad[1:3,]

# create data frame of corresponding features
# (acetic and gluconic acid concentrations in salad dressings)
features <- data.frame(salad = LETTERS[1:4],
                      acetic = c(0.5, 0.5, 1, 0),
                      gluconic = c(0, 10, 0, 10))

# fit Plackett-Luce model based on covariates
res_PLADMM <- pladmm(salad, ~ acetic + gluconic, data = features, rho = 8)
## coefficients
coef(res_PLADMM)
## worth
res_PLADMM$pi
## worth as predicted by linear function
res_PLADMM$tilde_pi
## equivalent to
drop(exp(res_PLADMM$x %*% coef(res_PLADMM)))
```

 pltree

Plackett-Luce Trees

Description

Recursive partitioning based on Plackett-Luce models.

Usage

```
pltree(formula, data, worth, na.action, cluster, ref = NULL, ...)
```

Arguments

formula	A symbolic description of the model to be fitted, of the form $y \sim x_1 + \dots + x_n$ where y should be an object of class <code>grouped_rankings</code> and x_1, \dots, x_n are used as partitioning variables.
data	An optional data object containing the variables in the model. Either a data frame of variables in formula or a list of length 2 giving data frames for variables in formula and in worth.
worth	A optional formula specifying a linear model for log-worth. If NULL, worth is estimated separately for each item with <code>PlackettLuce()</code> . Otherwise, the model in each node of the tree is fitted with <code>pladmm()</code> .
na.action	how NAs are treated for variables in formula, applied to the underlying rankings.
cluster	an optional vector of cluster IDs to be employed for clustered covariances in the parameter stability tests, see <code>mob</code> .
ref	an integer or character string specifying the reference item (for which log ability will be set to zero). If NULL the first item is used.
...	additional arguments, passed to <code>PlackettLuce</code> or <code>pladmm()</code> .

Details

Plackett-Luce trees are an application of model-based recursive partitioning (implemented in `mob`) to Plackett-Luce models for rankings. The partitioning is based on ranking covariates, e.g. attributes of the judge making the ranking, or conditions under which the ranking is made. The response should be a `grouped_rankings` object that groups rankings with common covariate values. This may be included in a data frame alongside the covariates.

Most arguments of `PlackettLuce` can be passed on by `pltree`. However, Plackett-Luce tree with fixed adherence are not implemented. Arguably it makes more sense to estimate adherence or reliability within the nodes of the Plackett-Luce tree.

Various methods are provided for "pltree" objects, most of them inherited from "modelparty" objects (e.g. `print`, `summary`), or "bttree" objects (`plot`). The `plot` method employs the `node_btplot` panel-generating function. The See Also section gives details of separately documented methods.

Value

An object of class "pltree" inheriting from "btree" and "modelparty".

See Also

[btree](#) For fitting Bradley-Terry trees (equivalent to the Plackett-Luce model for paired comparisons without ties).

coef, vcov, AIC and predict methods are documented on [pltree-summaries](#).

[itempar](#), extracts the abilities or item parameters in each node of the tree using `itempar.PlackettLuce`.

[fitted](#), computes probabilities for the observed choices based on the full tree.

Examples

```
# Bradley-Terry example

if (require(psychotree)){
  ## Germany's Next Topmodel 2007 data
  data("Topmodel2007", package = "psychotree")
  ## convert paircomp object to grouped rankings
  R <- as.grouped_rankings(Topmodel2007$preference)
  ## rankings are grouped by judge
  print(R[1:2,], max = 4)
  ## Topmodel2007[, -1] gives covariate values for each judge
  print(Topmodel2007[1:2, -1])

  ## fit partition model based on all variables except preference
  ## set npseudo = 0 as all judges rank all models
  tm_tree <- pltree(R ~ ., data = Topmodel2007[, -1], minsize = 5,
                   npseudo = 0)

  ## plot shows abilities constrained to sum to 1
  plot(tm_tree, abbreviate = 1, yscale = c(0, 0.5))
  ## instead show log-abilities with Anja as reference (need to used index)
  plot(tm_tree, abbreviate = 1, worth = FALSE, ref = 6,
       yscale = c(-1.5, 2.2))

  ## log-abilities, zero sum contrast
  itempar(tm_tree, log = TRUE)
}
```

pltree-summaries

Plackett-Luce Tree Summaries

Description

Obtain the coefficients, variance-covariance matrix, AIC, or predictions from a Plackett-Luce tree fitted by `pltree()`.

Usage

```
## S3 method for class 'pltree'
coef(object, node = NULL, drop = TRUE, ...)

## S3 method for class 'pltree'
vcov(object, node = nodeids(object, terminal = TRUE), ...)

## S3 method for class 'pltree'
AIC(object, newdata = NULL, ...)

## S3 method for class 'pltree'
predict(
  object,
  newdata = NULL,
  type = c("itempar", "rank", "best", "node"),
  ...
)
```

Arguments

object	a fitted model object of class "pltree".
node	a vector of node ids specifying the nodes to summarise, by default the ids of the terminal nodes.
drop	if TRUE return the coefficients as a vector when only one node is selected.
...	additional arguments passed to <code>itempar</code> by <code>predict</code> , and to <code>model.frame</code> by <code>AIC</code> .
newdata	an optional data frame to use instead of the original data. For <code>AIC</code> this must include the response variable.
type	the type of prediction to return for each group, one of: "itempar" to give the result of <code>itempar</code> (by default the fitted probability of each item being ranked first out of all objects), "rank" the corresponding rank, "best" the topped ranked item, or "node" the node of the tree the group belongs to.

Details

`AIC` computes $-2L + 2p$ where L is the joint likelihood of the observed rankings under the tree model and p is the degrees of freedom used to fit the tree model.

Examples

```
data(beans)
# fit tree based on pairwise comparisons with variety B
pairB <- data.frame(Winner = ifelse(beans$var_b == "Worse",
                                   "Local", beans$variety_b),
                   Loser = ifelse(beans$var_b == "Worse",
                                   beans$variety_b, "Local"),
                   stringsAsFactors = FALSE, row.names = NULL)
beans$G <- as.rankings(pairB, input = "orderings",
```

```

        index = rep(seq(nrow(beans)), 1))

mod <- pltree(G ~ ., data = beans[c("G", "maxTN")])

coef(mod, node = 3)
AIC(mod)

# treat first row from each year as new data
newdata <- beans[!duplicated(beans$year),]

## fitted probabilities
predict(mod, newdata)

## fitted log-abilities, with Local as reference
predict(mod, newdata, log = TRUE, ref = "Local")

## variety ranks
predict(mod, newdata, type = "rank")

## top ranked variety
predict(mod, newdata, type = "best")

## node the trial belongs to
predict(mod, newdata, type = "node")

```

preflib

Read Preflib Election Data Files

Description

Read orderings from .soc, .soi, .toc or .toi file types storing election data as defined by [{PrefLib}: A Library for Preferences](#).

Usage

```
read.soc(file)
```

```
read.soi(file)
```

```
read.toc(file)
```

```
read.toi(file)
```

```
## S3 method for class 'preflib'
as.aggregated_rankings(x, ...)
```

Arguments

file An election data file, conventionally with extension .soc, .soi, .toc or .toi according to data type.

- x An object of class "preflib".
- ... Additional arguments passed to `as.rankings()`: freq, input or items will be ignored with a warning as they are set automatically.

Details

The file types supported are

- .soc** Strict Orders - Complete List
- .soi** Strict Orders - Incomplete List
- .toc** Orders with Ties - Complete List
- .toi** Orders with Ties - Incomplete List

Note that the file types do not distinguish between types of incomplete orderings, i.e. whether they are a complete ranking of a subset of items (as supported by `PlackettLuce()`) or top- n rankings of n items from the full set of items (not currently supported by `PlackettLuce()`).

The numerically coded orderings and their frequencies are read into a data frame, storing the item names as an attribute. The `as.aggregated_rankings` method converts these to an "aggregated_rankings" object with the items labelled by the item names.

A Preflib file may be corrupt, in the sense that the ordered items do not match the named items. In this case, the file can be read in as a data frame (with a warning) using the corresponding `read.*` function, but `as.aggregated_rankings` will throw an error.

Value

A data frame of class "preflib" with first column Freq, giving the frequency of the ranking in that row, and remaining columns Rank 1, ..., Rank r giving the items ranked from first to last place in that ranking. Ties are represented by vector elements in list columns. The data frame has an attribute "items" giving the labels corresponding to each item number.

Note

The Netflix and cities datasets used in the examples are from Bennet and Lanning (2007) and Caragiannis et al (2017) respectively. These data sets require a citation for re-use.

References

- Mattei, N. and Walsh, T. (2013) PrefLib: A Library of Preference Data. *Proceedings of Third International Conference on Algorithmic Decision Theory (ADT 2013)*. Lecture Notes in Artificial Intelligence, Springer.
- Caragiannis, I., Chatzigeorgiou, X., Krimpas, G. A., and Voudouris, A. A. (2017) Optimizing positional scoring rules for rank aggregation. In *Proceedings of the 31st AAAI Conference on Artificial Intelligence*.
- Bennett, J. and Lanning, S. (2007) The Netflix Prize. *Proceedings of The KDD Cup and Workshops*.

Examples

```

# strict complete orderings of four films on Netflix
netflix <- read.soc(system.file("extdata", "netflix.soc",
                              package = "PlackettLuce"))

head(netflix)
attr(netflix, "items")

head(as.aggregated_rankings(netflix))

# strict incomplete orderings of 6 random cities from 36 in total
cities <- read.soi(system.file("extdata", "cities.soi",
                              package = "PlackettLuce"))

# complete orderings with ties of 30 skaters
skaters <- read.toc(system.file("extdata", "skaters.toc",
                                package = "PlackettLuce"))

# incomplete orderings with ties: most important qualities for success
# from 20 in total
qualities <- read.toi(system.file("extdata", "education_qualities.toi",
                                  package = "PlackettLuce"))

# alternatively read from a url
# - can take a little while depending on speed of internet connection

## Not run:
# incomplete orderings with ties: most important qualities for success
# from 20 in total
preflib <- "https://raw.githubusercontent.com/PrefLib/PrefLib-Data/main/datasets"
qualities2 <- read.toi(file.path(preflib, "00032%20-%20education/00032-00000007.toi"))
all.equal(qualities, qualities2)

## End(Not run)

```

pudding

Paired Comparisons of Chocolate Pudding

Description

This is an example dataset from *Davidson (1970)* comprising paired comparisons of chocolate pudding, with six brands in total. The responses include tied outcomes, i.e. no preference.

Usage

```
pudding
```

Format

A data frame with 15 records and 6 variables:

- i The first brand in the comparison.
- j The second brand in the comparison.
- r_{ij} The frequency of paired comparisons of brand i and brand j.
- w_{ij} The frequency of preferences for i over j.
- w_{ji} The frequency of preferences for j over i.
- t_{ij} The frequency of no preference between i and j.

References

Davidson, R. R. (1970). On extending the Bradley-Terry model to accommodate ties in paired comparison experiments. *Journal of the American Statistical Association*, **65**, 317–328.

Examples

```
# create orderings for each set of paired comparisons

# wins for brand i and wins for brand j
i_wins <- data.frame(Winner = pudding$i, Loser = pudding$j)
j_wins <- data.frame(Winner = pudding$j, Loser = pudding$i)

# ties: use an array list (easier with R >= 3.6.0)
if (getRversion() < "3.6.0"){
  n <- nrow(pudding)
  ties <- data.frame(Winner = array(split(pudding[c("i", "j")], 1:n), n),
                    Loser = rep(NA, 15))
} else {
  ties <- data.frame(Winner = asplit(pudding[c("i", "j")], 1),
                    Loser = rep(NA, 15))
}
head(ties, 2)

# convert to rankings
R <- as.rankings(rbind(i_wins, j_wins, ties),
                input = "orderings")
head(R, 2)
tail(R, 2)

# define weights as frequencies of each ranking
w <- unlist(pudding[c("w_ij", "w_ji", "t_ij")])

# fit Plackett-Luce model: limit iterations to match paper
mod <- PlackettLuce(R, npseudo = 0, weights = w, maxit = 7)
```

Description

A method for `qvcalc` to compute a set of quasi variances (and corresponding quasi standard errors) for estimated item parameters from a Plackett-Luce model.

Usage

```
## S3 method for class 'PlackettLuce'
qvcalc(object, ref = 1L, ...)
```

Arguments

<code>object</code>	a "PlackettLuce" object as returned by <code>PlackettLuce</code> .
<code>ref</code>	An integer or character string specifying the reference item (for which log worth will be set to zero). If NULL the sum of the log worth parameters is set to zero.
<code>...</code>	additional arguments, currently ignored..

Details

For details of the method see Firth (2000), Firth (2003) or Firth and de Menezes (2004). Quasi variances generalize and improve the accuracy of "floating absolute risk" (Easton et al., 1991). This device for economical model summary was first suggested by Ridout (1989).

Ordinarily the quasi variances are positive and so their square roots (the quasi standard errors) exist and can be used in plots, etc.

Value

A list of class "qv", with components

<code>covmat</code>	The full variance-covariance matrix for the item parameters.
<code>qvframe</code>	A data frame with variables <code>estimate</code> , <code>SE</code> , <code>quasiSE</code> and <code>quasiVar</code> , the last two being a quasi standard error and quasi-variance for each parameter.
<code>dispersion</code>	NULL (dispersion is fixed to 1).
<code>relerrs</code>	Relative errors for approximating the standard errors of all simple contrasts.
<code>factorname</code>	NULL (not required for this method).
<code>coef.indices</code>	NULL (not required for this method).
<code>modelcall</code>	The call to <code>PlackettLuce</code> to fit the model from which the item parameters were estimated.

References

Easton, D. F, Peto, J. and Babiker, A. G. A. G. (1991) Floating absolute risk: an alternative to relative risk in survival and case-control analysis avoiding an arbitrary reference group. *Statistics in Medicine* **10**, 1025–1035.

Firth, D. (2000) Quasi-variances in Xlisp-Stat and on the web. *Journal of Statistical Software* **5.4**, 1–13. At <https://www.jstatsoft.org>

Firth, D. (2003) Overcoming the reference category problem in the presentation of statistical models. *Sociological Methodology* **33**, 1–18.

Firth, D. and de Menezes, R. X. (2004) Quasi-variances. *Biometrika* **91**, 65–80.

Menezes, R. X. de (1999) More useful standard errors for group and factor effects in generalized linear models. *D.Phil. Thesis*, Department of Statistics, University of Oxford.

Ridout, M.S. (1989). Summarizing the results of fitting generalized linear models to data from designed experiments. In: *Statistical Modelling: Proceedings of GLIM89 and the 4th International Workshop on Statistical Modelling held in Trento, Italy, July 17–21, 1989* (A. Decarli et al., eds.), pp 262–269. New York: Springer.

See Also

[worstErrors](#), [plot.qv](#).

Examples

```
# Six partial rankings of four objects, 1 is top rank, e.g
# first ranking: item 1, item 2
# second ranking: item 2, item 3, item 4, item 1
# third ranking: items 2, 3, 4 tie for first place, item 1 second
R <- matrix(c(1, 2, 0, 0,
              4, 1, 2, 3,
              2, 1, 1, 1,
              1, 2, 3, 0,
              2, 1, 1, 0,
              1, 0, 3, 2), nrow = 6, byrow = TRUE)
colnames(R) <- c("apple", "banana", "orange", "pear")

mod <- PlackettLuce(R)
qv <- qvcalc(mod)
qv
plot(qv)
```

rankings

Rankings Object

Description

Create a "rankings" object from data or convert a matrix of rankings or ordered items to a "rankings" object.

Usage

```
rankings(data, id, item, rank, aggregate = FALSE, verbose = TRUE, ...)
```

```
as.rankings(x, ..., verbose = TRUE)
```

```
## Default S3 method:
```

```

as.rankings(
  x,
  input = c("rankings", "orderings"),
  freq = NULL,
  index = NULL,
  aggregate = FALSE,
  items = NULL,
  labels = NULL,
  ...,
  verbose = TRUE
)

## S3 method for class 'grouped_rankings'
as.rankings(x, ..., verbose = TRUE)

## S3 method for class 'matrix'
as.rankings(
  x,
  input = c("rankings", "orderings"),
  freq = NULL,
  index = NULL,
  aggregate = FALSE,
  items = NULL,
  labels = NULL,
  ...,
  verbose = TRUE
)

## S3 method for class 'rankings'
x[i, j, ..., drop = TRUE, as.rankings = TRUE]

## S3 method for class 'rankings'
format(x, width = 40L, ...)

```

Arguments

<code>data</code>	a data frame with columns specified by <code>id</code> , <code>item</code> and <code>rank</code> .
<code>id</code>	an index of data specifying the column containing ranking IDs.
<code>item</code>	an index of data specifying the column containing item IDs,
<code>rank</code>	an index of data specifying the column containing item ranks.
<code>aggregate</code>	if TRUE, aggregate the rankings via <code>aggregate()</code> before returning.
<code>verbose</code>	logical; if TRUE print messages when changes are made to rankings data.
<code>...</code>	further arguments passed to/from methods.
<code>x</code>	for <code>as.rankings</code> , a matrix with one column per item and one row per ranking, or an object that can be coerced to such as matrix; for <code>[</code> and <code>format</code> , a "rankings" object.

<code>input</code>	for <code>as.rankings</code> , whether rows in the input matrix contain numeric "rankings" (dense, standard/modified competition or fractional rankings) or "orderings", i.e. the items ordered by rank.
<code>freq</code>	an optional column index (number, character or logical) specifying a column of <code>x</code> that holds ranking frequencies, or a vector of ranking frequencies. If provided, an "aggregated_rankings" object will be returned.
<code>index</code>	an optional column index (number, character or logical) specifying a column of <code>x</code> that holds a grouping index, or a numeric vector to for grouping. If provided, the rankings will be grouped by <code>group()</code> before returning.
<code>items</code>	for <code>input = "orderings"</code> , a character vector specifying the full set of items. Values in <code>x</code> are matched to this by value (if character) or position (if numeric). Use <code>decode()</code> for orderings requiring more complex decoding.
<code>labels</code>	for <code>input = "orderings"</code> an optional vector of labels for the items, corresponding to the sorted unique values of <code>x</code> .
<code>i</code>	indices specifying rankings to extract, as for <code>[]</code> .
<code>j</code>	indices specifying items to extract, as for <code>[]</code> .
<code>drop</code>	if TRUE return single row/column matrices as a vector.
<code>as.rankings</code>	if TRUE return a rankings object, otherwise return a matrix/vector.
<code>width</code>	the width in number of characters to format each ranking - rankings that are too wide will be truncated.

Details

Each ranking in the input data will be converted to a dense ranking, which rank items from 1 (first place) to n_r (last place). Items not ranked should have a rank of 0 or NA. Tied items are given the same rank with no rank skipped. For example `{1, 0, 2, 1}`, ranks the first and fourth items in first place and the third item in second place; the second item is unranked.

Records in data with missing `id` or `item` are dropped. Duplicated items in the rankings are resolved if possible: redundant or inconsistent ranks are set to NA. Rankings with only 1 item are set to NA (rankings with zero items are automatically treated as NA). Any issues causing records to be removed or recoded produce a message if `verbose = TRUE`.

For `as.rankings` with `input = "orderings"`, unused ranks may be filled with zeroes for numeric `x` or NA. It is only necessary to have as many columns as ranks that are used.

The method for `[]` will return a reduced rankings object by default, recoding as dense rankings and setting invalid rankings to NA as necessary. To extract rows and/or columns of the rankings as a matrix or vector, set `as.rankings = FALSE`, see examples.

Value

By default, a "rankings" object, which is a matrix of dense rankings with methods for several generics including `aggregate`, `[]`, `format`, `rbind()` and `as.matrix()`.

If the object is created with `aggregate = TRUE`, or ranking frequencies are specified via `freq`, the rankings are post-processed to create an "aggregated_rankings" object.

If a group index is specified via `index`, the (possibly aggregated) rankings are post-processed to create a "grouped_rankings" object.

Examples

```

# create rankings from data in long form

# example long form data
x <- data.frame(ranking = c(rep(1:4, each = 4), 5, 5, 5),
               letter = c(LETTERS[c(1:3, 3, 1:4, 2:5, 1:2, 1)], NA,
                          LETTERS[3:5]),
               rank = c(4:1, rep(NA, 4), 3:4, NA, NA, 1, 3, 4, 2, 2, 2, 3))

# ranking 1 has different rank for same item, but order of items unambiguous
# all ranks are missing in ranking 2
# some ranks are missing in ranking 3
# ranking 4 has inconsistent ranks for two items and a rank with missing item
# ranking 5 is fine - an example of a tie
split(x, x$ranking)

# fix issues when creating rankings object
rankings(x, id = "ranking", item = "letter", rank = "rank")

# convert existing matrix of rankings

R <- matrix(c(1, 2, 0, 0,
             4, 1, 2, 3,
             2, 1, 1, 1,
             1, 2, 3, 0,
             2, 1, 1, 0,
             1, 0, 3, 2), nrow = 6, byrow = TRUE)
colnames(R) <- c("apple", "banana", "orange", "pear")
R <- as.rankings(R)

# first three rankings
R[1:3,]

# exclude pear from the rankings
R[, -4]

# extract rankings 2 and 3 as numeric matrix
R[2:3, , as.rankings = FALSE]

# same as
as.matrix(R)[2:3,]

# extract rankings for item 1 as a vector
R[,1, as.rankings = FALSE]

```

Description

Full rankings of four salad dressings collected by Vargo (1989) and published in Critchlow and Fligner (1991). The salad dressings are formulated with varying percentages of acetic and gluconic acid as follows: A = (0.5, 0), B = (0.5, 10.0), C = (1.0, 0), and D = (0, 10.0). The salad dressings are ranked by tartness from 1 (most tart) to 4 (least tart).

Usage

```
salad
```

Format

A data frame with 32 records across 4 variables: A, B, C, and D, representing the four salad dressings.

Source

Critchlow, D. E. and Fligner, M. A. (1991) Paired comparison, triple comparison, and ranking experiments as generalized linear models, and their implementation on GLIM. *Psychometrika* **56**(3), 517—533.

References

Vargo, M. D. (1989) Microbiological spoilage of a moderate acid food system using a dairy-based salad dressing model. Unpublished masters thesis, Ohio State University, Department of Food Science and Nutrition, Columbus, OH, USA.

Examples

```
# create data frame of acetic and gluconic acid concentrations
features <- data.frame(salad = LETTERS[1:4],
                      acetic = c(0.5, 0.5, 1, 0),
                      gluconic = c(0, 10, 0, 10))

# fit Plackett-Luce model based on covariates
res_PLADMM <- pladmm(salad, ~ acetic + gluconic, data = features, rho = 8)
# worth (relative tartness) as predicted by linear function of features
itempar(res_PLADMM)
```

simulate.PlackettLuce *Simulate from PlackettLuce fitted objects*

Description

Simulate from PlackettLuce fitted objects

Usage

```
## S3 method for class 'PlackettLuce'
simulate(
  object,
  nsim = 1L,
  seed = NULL,
  multinomial = FALSE,
  max_combinations = 20000,
  ...
)
```

Arguments

<code>object</code>	an object representing a fitted model.
<code>nsim</code>	number of response vectors to simulate. Defaults to 1.
<code>seed</code>	an object specifying if and how the random number generator should be initialised. Either <code>NULL</code> or an integer that will be used in a call to <code>set.seed</code> before simulating the rankings. If set, the value is saved as the <code>seed</code> attribute of the returned value. The default, <code>NULL</code> , will not change the random generator state, and return <code>Random.seed</code> as the <code>seed</code> attribute.
<code>multinomial</code>	use multinomial sampling anyway? Default is <code>FALSE</code> . see Details.
<code>max_combinations</code>	a positive number. Default is 20000. See Details.
<code>...</code>	additional optional arguments.

Details

If `multinomial` is `FALSE` (default) and there are no tie parameters in the object (i.e. `length(object$ties) == 1`), then rankings are sampled by ordering exponential random variates with rate 1 scaled by the estimated item-worth parameters `object$coefficients` (see, Diaconis, 1988, Chapter 9D for details).

In all other cases, the current implementation uses direct multinomial sampling, and will throw an error if there are more than `max_combinations` combinations of items that the sampler has to decide from. This is a hard-coded exit to prevent issues relating to the creation of massive objects in memory.

If `length(object$ties) > 1` the user's setting for `multinomial` is ignored and `simulate.PlackettLuce` operates as if `multinomial` is `TRUE`.

Value

A data frame of `rankings` objects of the same dimension as `object$rankings`.

References

Diaconis (1988). *Group Representations in Probability and Statistics*. Institute of Mathematical Statistics Lecture Notes 11. Hayward, CA.

Examples

```

R <- matrix(c(1, 2, 0, 0,
             4, 1, 2, 3,
             2, 1, 1, 1,
             1, 2, 3, 0,
             2, 1, 1, 0,
             1, 0, 3, 2), nrow = 6, byrow = TRUE)
colnames(R) <- c("apple", "banana", "orange", "pear")
mod <- PlackettLuce(R)
simulate(mod, 5)

s1 <- simulate(mod, 3, seed = 112)
s2 <- simulate(mod, 2, seed = 112)

identical(s1[1:2], s2[1:2])

```

summaries

*Plackett-Luce Model Summaries***Description**

Obtain the coefficients, model summary or coefficient variance-covariance matrix for a model fitted by `PlackettLuce`.

Usage

```

## S3 method for class 'PlackettLuce'
coef(object, ref = 1L, log = TRUE, type = "all", ...)

## S3 method for class 'PlackettLuce'
summary(object, ref = 1L, ...)

## S3 method for class 'PlackettLuce'
vcov(object, ref = 1L, type = c("expected", "observed"), ...)

```

Arguments

<code>object</code>	An object of class "PlackettLuce" as returned by <code>PlackettLuce</code> .
<code>ref</code>	An integer or character string specifying the reference item (for which log worth will be set to zero). If <code>NULL</code> the sum of the log worth parameters is set to zero.
<code>log</code>	A logical indicating whether to return parameters on the log scale with the item specified by <code>ref</code> set to zero.
<code>type</code>	For <code>coef</code> , the type of coefficients to return: one of "ties", "worth" or "all". For <code>vcov</code> , the type of Fisher information to base the estimation on: either "expected" or "observed".
<code>...</code>	additional arguments, passed to <code>vcov</code> by <code>summary</code> .

Details

By default, parameters are returned on the log scale, as most suited for inference. If `log = FALSE`, the worth parameters are returned, constrained to sum to one so that they represent the probability that the corresponding item comes first in a ranking of all items, given that first place is not tied.

The variance-covariance matrix is returned for the worth and tie parameters on the log scale, with the reference as specified by `ref`. For models estimated by maximum likelihood, the variance-covariance is the inverse of the Fisher information of the log-likelihood.

For models with a normal or gamma prior, the variance-covariance is based on the Fisher information of the log-posterior. When adherence parameters have been estimated, the log-posterior is not linear in the parameters. In this case there is a difference between the expected and observed Fisher information. By default, `vcov` will return the variance-covariance based on the expected information, but `type` gives to option to use the observed information instead. For large samples, the difference between these options should be small. Note that the estimation of the adherence parameters is accounted for in the computation of the variance-covariance matrix, but only the sub-matrix corresponding to the worth and tie parameters is estimated.

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