Package 'netrankr'

September 18, 2018

Type Package

Title Analyzing Partial Rankings in Networks

Version 0.2.1

Description Implements methods for centrality related analyses of networks. While the package includes the possibility to build more than 20 indices, its main focus lies on index-free assessment of centrality via partial rankings obtained by neighborhood-inclusion or positional dominance. These partial rankings can be analyzed with different methods, including probabilistic methods like computing expected node ranks and relative rank probabilities (how likely is it that a node is more central than another?). The methodology is described in depth in the vignettes and in Schoch (2018) <doi:10.1016/j.socnet.2017.12.003>.

URL https://schochastics.github.io/netrankr

 $\pmb{BugReports} \ \text{https://github.com/schochastics/netrankr/issues}$

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LazyData true

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LinkingTo Rcpp,RcppArmadillo

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VignetteBuilder knitr

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R topics documented:

 $aggregate_positions$

Quantification of (indirect) relations

Description

Function to aggregate positions defined via indirect relations to construct centrality scores.

Usage

Index

```
aggregate_positions(tau_x, type = "sum")
```

Arguments

tau_x Numeric matrix containing indirect relations calculated with indirect_relations.
type String indicating the type of aggregation to be used. See Details for options.

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Details

The predefined functions are mainly wrappers around base R functions. type='sum', for instance, is equivalent to rowSums(). A non-base functions is type='invsum' which calculates the inverse of type='sum'. type='self' is mostly useful for walk based relations, e.g. to count closed walks. Other self explanatory options are type='mean', type='min', type='max' and type='prod'.

Value

Scores for the index defined by the indirect relation tau_x and the used aggregation type.

Author(s)

David Schoch

See Also

indirect_relations, transform_relations

```
library(igraph)
library(magrittr)
g <- graph.empty(n=11,directed = FALSE)</pre>
g \leftarrow add_edges(g,c(1,11,2,4,3,5,3,11,4,8,5,9,5,11,6,7,6,8,
                   6,10,6,11,7,9,7,10,7,11,8,9,8,10,9,10))
g %>% indirect_relations(type='adjacency') %>%
 aggregate_positions(type='sum')
#closeness centrality
g %>% indirect_relations(type='dist_sp') %>%
  aggregate_positions(type='invsum')
#betweenness centrality
g %>% indirect_relations(type='depend_sp') %>%
  aggregate_positions(type='sum')
#eigenvector centrality
g %>% indirect_relations(type='walks',FUN=walks_limit_prop) %>%
  aggregate_positions(type='sum')
#subgraph centrality
g %>% indirect_relations(type='walks',FUN=walks_exp) %>%
  aggregate_positions(type='self')
```

approx_rank_expected Approximation of expected ranks

Description

Implements a variety of functions to approximate expected ranks for partial rankings.

Usage

```
approx_rank_expected(P, method = "lpom")
```

Arguments

P A partial ranking as matrix object calculated with neighborhood_inclusion or

positional_dominance.

method String indicating which method to be used. see Details.

Details

The method parameter can be set to

lpom local partial order model

glpom extension of the local partial order model.

loof1 based on a connection with relative rank probabilities.

loof2 extension of the previous method.

Which of the above methods performs best depends on the structure and size of the partial ranking. See vignette("benchmarks",package="netrankr") for more details.

Value

A vector containing approximated expected ranks.

Author(s)

David Schoch

References

Brüggemann R., Simon, U., and Mey,S, 2005. Estimation of averaged ranks by extended local partial order models. *MATCH Commun. Math. Comput. Chem.*, 54:489-518.

Brüggemann, R. and Carlsen, L., 2011. An improved estimation of averaged ranks of partial orders. *MATCH Commun. Math. Comput. Chem.*, 65(2):383-414.

De Loof, L., De Baets, B., and De Meyer, H., 2011. Approximation of Average Ranks in Posets. *MATCH Commun. Math. Comput. Chem.*, 66:219-229.

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See Also

```
approx_rank_relative, exact_rank_prob, mcmc_rank_prob
```

Examples

```
P <- matrix(c(0,0,1,1,1,0,0,0,1,0,0,0,0,0,1,rep(0,10)),5,5,byrow=TRUE)
#Exact result
exact_rank_prob(P)$expected.rank
approx_rank_expected(P,method = 'lpom')
approx_rank_expected(P,method = 'glpom')</pre>
```

approx_rank_relative Approximation of relative rank probabilities

Description

Approximate relative rank probabilities P(rk(u) < rk(v)). In a network context, P(rk(u) < rk(v)) is the probability that u is less central than v, given the partial ranking P.

Usage

```
approx_rank_relative(P, iterative = TRUE, num.iter = 10)
```

Arguments

P A partial ranking as matrix object calculated with neighborhood_inclusion or

positional_dominance.

iterative Logical scalar if iterative approximation should be used.

num.iter Number of iterations to be used. defaults to 10 (see Details).

Details

The iterative approach generally gives better approximations than the non iterative, if only slightly. The default number of iterations is based on the observation, that the approximation does not improve significantly beyond this value. This observation, however, is based on very small networks such that increasing it for large network may yield better results. See vignette("benchmarks", package="netrankr") for more details.

Value

a matrix containing approximation of mutual rank probabilities. relative.rank[i,j] is the probability that i is ranked lower than j

Author(s)

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References

De Loof, K. and De Baets, B and De Meyer, H., 2008. Properties of mutual rank probabilities in partially ordered sets. In *Multicriteria Ordering and Ranking: Partial Orders, Ambiguities and Applied Issues*, 145-165.

See Also

```
approx_rank_expected, exact_rank_prob, mcmc_rank_prob
```

Examples

```
P \leftarrow \mathsf{matrix}(c(0,0,1,1,1,0,0,0,1,0,0,0,0,0,1,\mathsf{rep}(0,10)),5,5,\mathsf{byrow=TRUE}) P \mathsf{approx\_rank\_relative}(P,\mathsf{iterative} = \mathsf{FALSE}) \mathsf{approx\_rank\_relative}(P,\mathsf{iterative} = \mathsf{TRUE})
```

comparable_pairs

Comparable pairs in a partial ranking

Description

Calculates the fraction of comparable pairs in a partial ranking. This fraction is identical to the density of the induced undirected graph of a partial ranking.

Usage

```
comparable_pairs(P)
```

Arguments

Ρ

A partial ranking as matrix object calculated with neighborhood_inclusion or positional_dominance.

Value

Fraction of comparable pairs in P.

Author(s)

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Examples

```
library(igraph)
g <- sample_gnp(100,0.1)
P <- neighborhood_inclusion(g)
comparable_pairs(P)
# All pairs of vertices are comparable in a threshold graph
tg <- threshold_graph(100,0.3)
P <- neighborhood_inclusion(g)
comparable_pairs(P)</pre>
```

compare_ranks

Count occurrences of pairs in rankings

Description

Counts the number of concordant, discordant and (left/right) ties between two rankings.

Usage

```
compare_ranks(x, y)
```

Arguments

x A numeric vector.

y A numeric vector with the same length as x.

Details

Explicitly calculating the number of occurring cases is more robust than using correlation indices as given in the cor function. Especially left and right ties can significantly alter correlations.

Value

A list containing

```
concordant number of concordant pairs: x[i] > x[j] and y[i] > y[j] discordant number of discordant pairs: x[i] > x[j] and y[i] < y[j] ties number of tied pairs: x[i] == x[j] and y[i] == y[j] left number of left ties: x[i] == x[j] and y[i] != y[j] right number of right ties: x[i] != x[j] and y[i] == y[j]
```

Author(s)

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Examples

```
library(igraph)
tg <- threshold_graph(100,0.2)
compare_ranks(degree(tg),closeness(tg)) #only concordant pairs
compare_ranks(degree(tg),betweenness(tg)) #no discordant pairs
## Rank Correlation
cor(degree(tg),closeness(tg),method='kendall') #1
cor(degree(tg),betweenness(tg),method='kendall') #not 1, although no discordant pairs</pre>
```

dominance_graph

Partial ranking as directed graph

Description

Turns a partial ranking into a directed graph. An edge (u,v) is present if P[u,v]=1, meaning that u is dominated by v.

Usage

```
dominance_graph(P)
```

Arguments

Ρ

A partial ranking as matrix object calculated with neighborhood_inclusion or positional_dominance.

Value

Directed graph as an igraph object.

Author(s)

David Schoch

```
library(igraph)
g <- threshold_graph(20,0.1)
P <- neighborhood_inclusion(g)
d <- dominance_graph(P)
## Not run: plot(d)

# to reduce overplotting use transitive reduction
P <- transitive_reduction(P)
d <- dominance_graph(P)
## Not run: plot(d)</pre>
```

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exact_rank_prob Probabilistic centrality rankings

Description

Performs a complete and exact rank analysis of a given partial ranking. This includes rank probabilities, relative rank probabilities and expected ranks.

Usage

```
exact_rank_prob(P, only.results = T, verbose = F, force = F)
```

Arguments

P A partial ranking as matrix object calculated with neighborhood_inclusion or

positional_dominance.

only.results Logical. return only results (default) or additionally the ideal tree and lattice if

FALSE.

verbose Logical. should diagnostics be printed. Defaults to FALSE.

force Logical. If FALSE (default), stops the analysis if the partial ranking has more

than 40 elements and less than 0.4 comparable pairs. Only change if you know

what you are doing.

Details

The function derives rank probabilities from a given partial ranking (for instance returned by neighborhood_inclusion or positional_dominance). This includes the calculation of expected ranks, (relative) rank probabilities and the number of possible rankings. Note that the set of rankings grows exponentially in the number of elements and the exact calculation becomes infeasible quite quickly and approximations need to be used. See vignette("benchmarks") for guidelines and approx_rank_relative, approx_rank_expected, and mcmc_rank_prob for approximative methods.

Value

lin.ext	Number of possible rankings that extend P.
mse	Array giving the equivalence classes of P.
rank.prob	Matrix containing rank probabilities: $rank.prob[u,k]$ is the probability that u has rank k .
relative.rank	Matrix containing relative rank probabilities: $relative.rank[u,v]$ is the probability that u is ranked lower than v .
expected.rank	Expected ranks of nodes in any centrality ranking.
rank.spread	Standard deviation of the ranking probabilities.
topo.order	Random ranking used to build the lattice of ideals (if only.results = FALSE).
tree	Adjacency list (incoming) of the tree of ideals (if only.results = FALSE).

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lattice Adjacency list (incoming) of the lattice of ideals (if only.results = FALSE).

ideals List of order ideals (if only.results = FALSE).

In all cases, higher numerical ranks imply a higher position in the ranking. That is, the lowest ranked node has rank 1.

Author(s)

David Schoch, Julian Müller

References

De Loof, K. 2009. Efficient computation of rank probabilities in posets. *Phd thesis*, Ghent University.

De Loof, K., De Meyer, H. and De Baets, B., 2006. Exploiting the lattice of ideals representation of a poset. *Fundamenta Informaticae*, 71(2,3):309-321.

See Also

```
approx_rank_relative, approx_rank_expected, mcmc_rank_prob
```

Examples

```
P <- matrix(c(0,0,1,1,1,0,0,0,1,0,0,0,0,1,rep(0,10)),5,5,byrow=TRUE)
P
res <- exact_rank_prob(P)

#a warning is displayed if only one ranking is possible
tg <- threshold_graph(20,0.2)
P <- neighborhood_inclusion(tg)
res <- exact_rank_prob(P)</pre>
```

florentine_m

Florentine family marriage network

Description

Florentine family marriage network

Usage

florentine_m

Format

An igraph object containing marriage links of florentine families

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References

Padgett, J.F. and Ansell, C.K., 1993. Robust Action and the Rise of the Medici, 1400-1434. *American Journal of Sociology*, **98**(6), 1259-1319.

get_rankings

Rankings that extend a partial ranking

Description

Returns all possible rankings that extend a partial ranking.

Usage

```
get_rankings(data, force = F)
```

Arguments

data List as returned by exact_rank_prob when run with only.results = FALSE force Logical scalar. Stops function if the number of rankings is too large. Only

change to TRUE if you know what you are doing

Details

The ith row of the matrix contains the rank of node i in all possible rankings that are in accordance with the partial ranking P. The lowest rank possible is associated with 1.

Value

A matrix containing ranks of nodes in all possible rankings.

Author(s)

David Schoch

```
P \leftarrow \mathsf{matrix}(\mathsf{c}(\emptyset,\emptyset,1,1,1,\emptyset,\emptyset,\emptyset,1,\emptyset,\emptyset,\emptyset,\emptyset,1,\mathsf{rep}(\emptyset,10)),5,5,\mathsf{byrow=TRUE}) P \mathsf{res} \leftarrow \mathsf{exact\_rank\_prob}(P,\mathsf{only}.\mathsf{results} = \mathsf{FALSE}) \mathsf{get\_rankings}(\mathsf{res})
```

hyperbolic_index

hyperbolic_index

Hyperbolic (centrality) index

Description

The hyperbolic index is an index that considers all closed walks of even or odd length on induced neighborhoods of a vertex.

Usage

```
hyperbolic_index(g, type = "odd")
```

Arguments

g igraph object.

type string. 'even' if only even length walks should be considered. 'odd' (Default) if

only odd length walks should be used.

Details

The hyperbolic index is an illustrative index that should not be used for any serious analysis. Its purpose is to show that with enough mathematical trickery, any desired result can be obtained when centrality indices are used.

Value

A vector containing centrality scores.

Author(s)

David Schoch

```
library(igraph)
g <- graph.empty(n=11,directed = FALSE)
g <- add_edges(g,c(1,11,2,4,3,5,3,11,4,8,5,9,5,11,6,7,6,8,6,10,6,11,7,9,7,10,7,11,8,9,8,10,9,10))
hyperbolic_index(g,type = "odd")
hyperbolic_index(g,type = "even")</pre>
```

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Description

This shiny gadget can be used to build centrality indices based on specific indirect relations, transformations and aggregation functions.

Usage

```
index_builder()
```

Value

code to calculate the specified index.

indirect_relations Indirect relations in a network

Description

Derive indirect relations for a given network. Observed relations, like presents or absence of a relation, are commonly not the center of analysis, but are transformed in a new set of indirect relation like shortest path distances among nodes. These transformations are usually an implicit step when centrality indices are used. Making this step explicit gives more possibilities, for example calculating partial centrality rankings with positional_dominance.

Usage

```
indirect_relations(g, type = "dist_sp", lfparam = NULL, dwparam = NULL,
 netflowmode = "", rspxparam = NULL, FUN = identity, ...)
```

Arguments

g	igraph object. The network for which relations should be derived.
type	String giving the relation to be calculated. See Details for options.
lfparam	Numeric parameter. Only used if type = "dist_lf".
dwparam	Numeric parameter. Only used if type = "dist_walk".
netflowmode	String, one of raw, frac, or norm. Only used if type = "depend_netflow".
rspxparam	Numeric parameter. Only used if type = "depend_rsps" or type = "depend_rspn".
FUN	A function that allows the transformation of relations. See Details.

Additional arguments passed to FUN.

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Details

The type parameter has the following options.

'adjacency' returns the adjacency matrix of the network.

'weights' returns the weighted adjacency matrix of the network if an edge attribute 'weight' is present.

'dist_sp' returns shortest path distances between all pairs of nodes.

'depend_sp' returns dyadic dependencies

$$\delta(u, s) = \sum_{t \in V} \frac{\sigma(s, t|u)}{\sigma(s, t)}$$

where $\sigma(s,t|u)$ is the number of shortest paths from s to t that include u and $\sigma(s,t)$ is the total number of shortest (s,t)-paths. This relation is used for betweenness-like centrality indices.

'walks' returns walk counts between pairs of nodes, usually they are weighted decreasingly in their lengths or other properties which can be done by adding a function in FUN. See transform_relations for options.

'dist_resist' returns the resistance distance between all pairs of nodes.

'dist_lf' returns a logarithmic forest distance $d_{\alpha}(s,t)$. The logarithmic forest distances form a one-parametric family of distances, converging to shortest path distances as $\alpha - > 0$ and to the resistance distance as $\alpha - > \infty$. See (Chebotarev, 2011) for more details. The parameter 1fparam can be used to tune α .

'dist_walk' returns the walk distance $d_{\alpha}^{W}(s,t)$ between nodes. The walk distances form a one-parametric family of distances, converging to shortest path distances as $\alpha - > 0$ and to longest walk distances for $\alpha - > \infty$. Walk distances contain the logarithmic forest distances as a special case. See (Chebotarev, 2012) for more details.

'dist_rwalk' returns the expected length of a random walk between two nodes. For more details see (Noh and Rieger, 2004)

'depend_netflow' returns dependencies based on network flow (See Freeman et al.,1991). If netflowmode="raw", the function returns

$$\delta(u,s) = \sum_{t \in V} f(s,t,G) - f(s,t,G-v)$$

where f(s,t,G) is the maximum flow from s to t in G and f(s,t,G-v) in G without the node v. For netflowmode="frac" it returns dependencies in the form, similar to shortest path dependencies:

$$\delta(u,s) = \sum_{t \in V} \frac{f(s,t,G) - f(s,t,G-v)}{f(s,t,G)}$$

'depend_curflow' returns pairwise dependencies based on current flow. The relation is based on the same idea as 'depend_sp' and 'depend_netflow'. However, instead of considering shortest paths or network flow, the current flow (or equivalent: random walks) between nodes are of interest. See (Newman, 2005) for details.

'depend_exp' returns pairwise dependencies based on 'communicability':

$$\delta(u,s) = \sum_{t \in V} \frac{exp(A)_{st} - exp(A + E(u))_{st}}{exp(A)_{st}},$$

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where E(u) has nonzeros only in row and column u, and in this row and column has -1 if A has +1. See (Estrada et al., 2009) for additional details.

'depend_rsps'. Simple randomized shortest path dependencies. The simple RSP dependency of a node u with respect to absorbing paths from s to t, is defined as the expected number of visits through u over all s-t-walks. The parameter rspxparam is the "inverse temperature parameter". If it converges to infinity, only shortest paths are considered and the expected number of visits to a node on a shortest path approaches the probability of following that particular path. When the parameter converges to zero, then the dependencies converge to the expected number of visits to a node over all absorbing walks with respect to the unbiased random walk probabilities. This means for undirected networks, that the relations converge to adjacency. See (Kivimäki et al., 2016) for details.

'depend_rspn' Net randomized shortest path dependencies. The parameter rspxparam is the "inverse temperature parameter". The asymptotic for the infinity case are the same as for 'depend_rsps'. If the parameter approaches zero, then it converges to 'depend_curflow'. The net randomized shortest path dependencies are closely related to the random walk interpretation of current flows. See (Kivimäki et al., 2016) for technical details.

The function FUN is used to transform the indirect relation. See transform_relations for predefined functions and additional help.

Value

A matrix containing indirect relations in a network.

Author(s)

David Schoch

References

Chebotarev, P., 2012. The walk distances in graphs. *Discrete Applied Mathematics*, 160(10), pp.1484-1500.

Chebotarev, P., 2011. A class of graph-geodetic distances generalizing the shortest-path and the resistance distances. *Discrete Applied Mathematics* 159,295-302.

Noh, J.D. and Rieger, H., 2004. Random walks on complex networks. *Physical Review Letters*, 92(11), p.118701.

Freeman, L.C., Borgatti, S.P., and White, D.R., 1991. Centrality in Valued Graphs: A Measure of Betweenness Based on Network Flow. *Social Networks* 13(2), 141-154.

Newman, M.E., 2005. A measure of betweenness centrality based on random walks. *Social Networks*, 27(1), pp.39-54.

Estrada, E., Higham, D.J., and Hatano, N., 2009. Communicability betweenness in complex networks. *Physica A* 388,764-774.

Kivimäki, I., Lebichot, B., Saramäki, J., and Saerens, M., 2016. Two betweenness centrality measures based on Randomized Shortest Paths *Scientific Reports* 6: 19668

See Also

aggregate_positions to build centrality indices, positional_dominance to derive dominance relations

is_preserved

Examples

```
library(igraph)
g <- graph.empty(n=11,directed = FALSE)
g <- add_edges(g,c(1,11,2,4,3,5,3,11,4,8,5,9,5,11,6,7,6,8,6,10,6,11,7,9,7,10,7,11,8,9,8,10,9,10))

#shortest path distances
D <- indirect_relations(g,type = "dist_sp")

#inverted shortest path distances
D <- indirect_relations(g,type = "dist_sp", FUN = dist_inv)
#shortes path dependencies (used for betweenness)
D <- indirect_relations(g,type = "depend_sp")

#walks attenuated exponentially by their length
W <- indirect_relations(g,type = "walks",FUN = walks_exp)</pre>
```

is_preserved

Check preservation

Description

Checks if a partial ranking is preserved in the ranking induced by scores.

Usage

```
is_preserved(P, scores)
```

Arguments

P A partial ranking as matrix object calculated with neighborhood_inclusion or

positional_dominance.

scores Numeric vector containing the scores of a centrality index.

Details

In order for a score vector to preserve a partial ranking, the following condition must be fulfilled: P[u,v]==1 & scores[i] <= scores[j].

Value

Logical scaler whether scores preserves the relations in P.

Author(s)

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Examples

```
library(igraph)
# standard measures of centrality preserve the neighborhood inclusion preorder
g <- graph.empty(n=11,directed = FALSE)
g <- add_edges(g,c(1,11,2,4,3,5,3,11,4,8,5,9,5,11,6,7,6,8,6,10,6,11,7,9,7,10,7,11,8,9,8,10,9,10))
P<-neighborhood_inclusion(g)

is_preserved(P,degree(g))
is_preserved(P,betweenness(g))
is_preserved(P,closeness(g))</pre>
```

majorization_gap

Majorization gap

Description

Calculates the (normalized) majorization gap of an undirected graph. The majorization gap indicates how far the degree sequence of a graph is from a degree sequence of a threshold_graph.

Usage

```
majorization_gap(g, norm = TRUE)
```

Arguments

g An igraph object

norm True (Default) if the normalized majorization gap should be returned.

Details

The distance is measured by the number of *reverse unit transformations* necessary to turn the degree sequence into a threshold sequence. First, the *corrected conjugated degree sequence* d' is calculated from the degree sequence d as follows:

$$d'_k = |\{i : i < k \land d_i \ge k - 1\}| + |\{i : i > k \land d_i \ge k\}|.$$

the majorization gap is then defined as

$$1/2\sum_{k=1}^{n} \max\{d_k' - d_k, 0\}$$

The higher the value, the further away is a graph to be a threshold graph.

Value

Majorization gap of an undirected graph.

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Author(s)

David Schoch

References

Schoch, D., Valente, T. W. and Brandes, U., 2017. Correlations among centrality indices and a class of uniquely ranked graphs. *Social Networks* **50**, 46–54.

Arikati, S.R. and Peled, U.N., 1994. Degree sequences and majorization. *Linear Algebra and its Applications*, **199**, 179-211.

Examples

```
library(igraph)
g <- graph.star(5,'undirected')
majorization_gap(g) #0 since star graphs are threshold graphs

g <- sample_gnp(100,0.15)
majorization_gap(g,norm=TRUE) #fraction of reverse unit transformation
majorization_gap(g,norm=FALSE) #number of reverse unit transformation</pre>
```

mcmc_rank_prob

Estimate rank probabilities with Markov Chains

Description

Performs a probabilistic rank analysis based on an almost uniform sample of possible rankings that preserve a partial ranking.

Usage

```
mcmc_rank_prob(P, rp = nrow(P)^3)
```

Arguments

P P A partial ranking as matrix object calculated with neighborhood_inclusion or

positional_dominance.

rp Integer indicating the number of samples to be drawn.

Details

This function can be used instead of exact_rank_prob if the number of elements in P is too large for an exact computation. As a rule of thumb, the number of samples should be at least cubic in the number of elements in P. See vignette("benchmarks",package="netrankr") for guidelines and benchmark results.

Value

```
expected.rank Estimated expected ranks of nodes

relative.rank Matrix containing estimated relative rank probabilities: relative.rank[u,v] is the probability that u is ranked lower than v.
```

Author(s)

David Schoch

References

```
Bubley, R. and Dyer, M., 1999. Faster random generation of linear extensions. Discrete Mathematics, 201(1):81-88
```

See Also

```
exact_rank_prob, approx_rank_relative, approx_rank_expected
```

Examples

```
## Not run:
data("florentine_m")
P <- neighborhood_inclusion(florentine_m)
res <- exact_rank_prob(P)
mcmc <- mcmc_rank_prob(P,rp = vcount(g)^3)

# mean absolute error (expected ranks)
mean(abs(res$expected.rank-mcmc$expected.rank))
## End(Not run)</pre>
```

neighborhood_inclusion

Neighborhood-inclusion preorder

Description

Calculates the neighborhood-inclusion preorder of an undirected graph.

Usage

```
neighborhood_inclusion(g)
```

Arguments

g An igraph object

Details

Neighborhood-inclusion is defined as

$$N(u) \subseteq N[v]$$

where N(u) is the neighborhood of u and $N[v] = N(v) \cup \{v\}$ is the closed neighborhood of v. $N(u) \subseteq N[v]$ implies that $c(u) \le c(v)$, where c is a centrality index based on a specific path algebra. Indices falling into this category are closeness (and variants), betweenness (and variants) as well as many walk-based indices (eigenvector and subgraph centrality, total communicability,...).

Value

The neighborhood-inclusion preorder of g as matrix object. P[u, v]=1 if $N(u) \subseteq N[v]$

Author(s)

David Schoch

References

Schoch, D. and Brandes, U., 2016. Re-conceptualizing centrality in social networks. *European Journal of Applied Mathematics* 27(6), 971-985.

Brandes, U. Heine, M., Müller, J. and Ortmann, M., 2017. Positional Dominance: Concepts and Algorithms. *Conference on Algorithms and Discrete Applied Mathematics*, 60-71.

See Also

positional_dominance, exact_rank_prob

```
library(igraph)
#the neighborhood inclusion preorder of a star graph is complete
g <- graph.star(5,'undirected')</pre>
P <- neighborhood_inclusion(g)</pre>
comparable_pairs(P)
#the same holds for threshold graphs
tg <- threshold_graph(50,0.1)</pre>
P <- neighborhood_inclusion(tg)</pre>
comparable_pairs(P)
#standard centrality indices preserve neighborhood-inclusion
g <- graph.empty(n=11,directed = FALSE)</pre>
g <- add_edges(g,c(1,11,2,4,3,5,3,11,4,8,5,9,5,11,6,7,6,8,
                    6,10,6,11,7,9,7,10,7,11,8,9,8,10,9,10))
P <- neighborhood_inclusion(g)</pre>
is_preserved(P,degree(g))
is_preserved(P,closeness(g))
is_preserved(P,betweenness(g))
```

netrankr 21

netrankr

netrankr: An R package for centrality and partial rankings in networks

Description

netrankr provides several functions to analyze partial rankings for network centrality. The main focus lies on methods that do not necessarily rely on indices like degree, betweenness or closeness. However, the package also provides more than 20 indices, which can be constructed via a Rstudio addin.

The package follows the philosophy, that centrality can be decomposed in a series of micro steps. Starting from a network, indirect_relations can be derived which can either be aggregated into an index with aggregate_positions, or alternatively turned into a partial ranking with positional_dominance. The partial ranking can then be further analyzed with exact_rank_prob, to obtain probabilistic centrality rankings.

Details

Some features of the package are:

- Working with the neighborhood inclusion preorder. This forms the bases for any centrality
 analysis on undirected and unweighted graphs. More details can be found in the dedicated
 vignette: vignette("neighborhood_inclusion", package = "netrankr")
- Constructing graphs with a unique centrality ranking. This class of graphs, known as threshold
 graphs, can be used to benchmark centrality indices, since they only allow for one ranking of
 the nodes. For more details consult the vignette: vignette("threshold_graph", package = "netrankr")
- Probabilistic centrality. Why apply a handful of indices and choosing the one that fits best, when it is possible to analyze **all** centrality rankings at once? The package includes several function to calculate rank probabilities of nodes in a network. These include expected ranks and relative rank probabilities (how likely is it that a node is more central than another?) Consult vignette("probabilistic_cent", package = "netrankr") for more info.

The package provides several additional vignettes that explain the functionality of netrankr and its conceptual ideas. See browseVignettes(package = 'netrankr') or the online manual.

Description

Compute rank intervals (minimal and maximal possible rank) and visualize them with ggplot.

Usage

```
plot_rank_intervals(P, cent.df, ties.method = "min")
```

22 plot_rank_intervals

Arguments

cent.df

P	A partial ranking as matrix object calculated with neighborhood_inclusion or
	positional_dominance.

A data frame containing centrality scores of indices (optional). See Details.

ties.method String specifying how ties are treated in the base rank function.

Details

If a data frame of centrality scores is added, the respective ranks of nodes are shown in the intervals. Note that some points might fall outside of the intervals depending how ties are treated.

Value

```
a ggplot object.
```

Author(s)

David Schoch

See Also

rank_intervals

```
library(igraph)
library(ggplot2)
g <- graph.empty(n=11,directed = FALSE)
g <- add_edges(g,c(1,11,2,4,3,5,3,11,4,8,5,9,5,11,6,7,6,8,6,10,6,11,7,9,7,10,7,11,8,9,8,10,9,10))
P <- neighborhood_inclusion(g)
## Not run: plot_rank_intervals(P)

#adding index based rankings
cent_scores <- data.frame(
   degree = degree(g),
   betweenness = round(betweenness(g),4),
   closeness = round(closeness(g),4),
   eigenvector = round(eigen_centrality(g)$vector,4))
## Not run: plot_rank_intervals(P,cent.df=cent_scores)</pre>
```

positional_dominance 23

Description

generalized dominance relations.

Usage

```
positional_dominance(A, type = "one-mode", map = FALSE, benefit = TRUE)
```

Arguments

A Matrix containing attributes or relations, for instance calculated by indirect_relations.

type A string which is either 'one-mode' (Default) if A is a regular one-mode network

or 'two-mode' if A is a general data matrix.

map Logical scalar, whether rows can be sorted or not (Default). See Details.

benefit Logical scalar, whether the attributes or relations are benefit or cost variables.

Details

Positional dominance is a generalization of neighborhood-inclusion for arbitrary network data. In the default case, it checks for all pairs u,v if $A_{ut} \geq A_{vt}$ holds for all t if benefit = TRUE or $A_{ut} \leq A_{vt}$ holds for all t if benefit = FALSE. This form of dominance is referred to as dominance under total heterogeneity. If map=TRUE, the rows of A are sorted decreasingly (benefit = TRUE) or increasingly (benefit = FALSE) and then the dominance condition is checked. This second form of dominance is referred to as dominance under total homogeneity, while the first is called dominance under total heterogeneity.

Value

Dominance relations as matrix object. An entry [u,v] is 1 if u is dominated by v.

Author(s)

David Schoch

References

Brandes, U., 2016. Network positions. *Methodological Innovations* 9, 2059799116630650.

Schoch, D. and Brandes, U., 2016. Re-conceptualizing centrality in social networks. *European Journal of Applied Mathematics* 27(6), 971-985.

See Also

neighborhood_inclusion, indirect_relations, exact_rank_prob

24 rank_intervals

Examples

```
library(igraph)

g <- graph.empty(n=11,directed = FALSE)

g <- add_edges(g,c(1,11,2,4,3,5,3,11,4,8,5,9,5,11,6,7,6,8,6,0,6,10,6,11,7,9,7,10,7,11,8,9,8,10,9,10))
P<-neighborhood_inclusion(g)
comparable_pairs(P)

# positional dominance under total heterogeneity
dist <- indirect_relations(g,type = "dist_sp")
D <- positional_dominance(dist,map = FALSE,benefit = FALSE)
comparable_pairs(D)

# positional dominance under total homogeneity
D_map <- positional_dominance(dist,map = TRUE,benefit = FALSE)
comparable_pairs(D_map) #more comparables than D</pre>
```

rank_intervals

Rank interval of nodes

Description

Calculate the maximal and minimal rank possible for each node in any ranking that is in accordance with the partial ranking P.

Usage

```
rank_intervals(P)
```

Arguments

Ρ

A partial ranking as matrix object calculated with neighborhood_inclusion or positional_dominance.

Details

Note that the returned mid_point is not the same as the expected rank, for instance computed with exact_rank_prob. It is simply the average of min_rank and max_rank. For exact rank probabilities use exact_rank_prob.

Value

A data frame with the minimal, maximal rank of each node together with the mid point of the two extrema.

spectral_gap 25

Author(s)

David Schoch

See Also

```
plot_rank_intervals, exact_rank_prob
```

Examples

```
 P \leftarrow \texttt{matrix}(c(\emptyset,\emptyset,1,1,1,0,\emptyset,\emptyset,1,\emptyset,\emptyset,\emptyset,0,1,rep(\emptyset,10)),5,5,byrow=TRUE) \\ rank\_intervals(P)
```

spectral_gap

Spectral gap of a graph

Description

The spectral (or eigen) gap of a graph is the absolute difference between the biggest and second biggest eigenvalue of the adjacency matrix. To compare spectral gaps across networks, the fraction can be used.

Usage

```
spectral_gap(g, method = "frac")
```

Arguments

g igraph object

method A string, either "frac" or "abs"

Details

The spectral gap is bounded between 0 and 1 if method="frac". The closer the value to one, the bigger the gap.

Value

Numeric value

Author(s)

David Schoch

```
#The fractional spectral gap of a threshold graph is usually close to 1 g <- threshold_graph(50,0.3) spectral_gap(g,method = "frac")
```

26 threshold_graph

threshold_graph

Random threshold graphs

Description

Constructs a random threshold graph. A threshold graph is a graph where the neighborhood inclusion preorder is complete.

Usage

```
threshold_graph(n, p)
```

Arguments

n The number of vertices in the graph.

p The probability of inserting dominating vertices. Equates approximately to the density of the graph. See Details.

Details

Threshold graphs can be constructed with a binary sequence. For each 0, an isolated vertex is inserted and for each 1, a vertex is inserted that connects to all previously inserted vertices. The probability of inserting a dominating vertices is controlled with parameter p. An important property of threshold graphs is, that all centrality indices induce the same ranking.

Value

A threshold graph as igraph object

Author(s)

David Schoch

References

Mahadev, N. and Peled, U. N., 1995. Threshold graphs and related topics.

Schoch, D., Valente, T. W. and Brandes, U., 2017. Correlations among centrality indices and a class of uniquely ranked graphs. *Social Networks* 50, 46–54.

See Also

neighborhood_inclusion, positional_dominance

transform_relations 27

Examples

```
library(igraph)
g <- threshold_graph(10,0.3)
## Not run:
plot(g)

# star graphs and complete graphs are threshold graphs
complete <- threshold_graph(10,1) #complete graph
plot(complete)

star <- threshold_graph(10,0) #star graph
plot(star)

## End(Not run)

# centrality scores are perfectly rank correlated
cor(degree(g),closeness(g),method = "kendall")</pre>
```

transform_relations

Transform indirect relations

Description

Mostly wrapper functions that can be used in conjunction with indirect_relations to fine tune indirect relations.

Usage

```
dist_2pow(x)
dist_inv(x)
dist_dpow(x, alpha = 1)
dist_powd(x, alpha = 0.5)
walks_limit_prop(x)
walks_exp(x, alpha = 1)
walks_exp_even(x, alpha = 1)
walks_exp_odd(x, alpha = 1)
walks_attenuated(x, alpha = 1/max(x) * 0.99)
walks_uptok(x, alpha = 1, k = 3)
```

28 transform_relations

Arguments

x Matrix of relations.

alpha Potential weighting factor.

k For walk counts up to a certain length.

Details

The predefined functions follow the naming scheme relation_transformation. Predefined functions walks_* are thus best used with type="walks" in indirect_relations. Theoretically, however, any transformation can be used with any relation. The results might, however, not be interpretable.

The following functions are implemented so far:

 $\operatorname{dist_2pow\ returns\ }2^{-x}$

 $dist_inv returns 1/x$

 $\operatorname{dist_dpow}$ returns $x^{-\alpha}$ where α should be chosen greater than 0.

dist_powd returns α^x where α should be chosen between 0 and 1.

walks_limit_prop returns the limit proportion of walks between pairs of nodes. Calculating row-Sums of this relation will result in the principle eigenvector of the network.

walks_exp returns $\sum_{k=0}^{\infty} \frac{A^k}{k!}$

walks_exp_even returns $\sum_{k=0}^{\infty} rac{A^{2k}}{(2k)!}$

walks_exp_odd returns $\sum_{k=0}^{\infty} \frac{A^{2k+1}}{(2k+1)!}$

walks_attenuated returns $\sum_{k=0}^{\infty} \alpha^k A^k$

walks_uptok returns $\sum_{j=0}^k \alpha^j A^j$

Walk based transformation are defined on the eigen decomposition of the adjacency matrix using the fact that

$$f(A) = X f(\Lambda) X^T.$$

Care has to be taken when using user defined functions.

Value

Transformed relations as matrix

Author(s)

transitive_reduction 29

transitive_reduction Transitive Reduction

Description

Calculates the transitive reduction of a partial ranking.

Usage

```
transitive_reduction(P)
```

Arguments

Ρ

A partial ranking as matrix object calculated with neighborhood_inclusion or positional_dominance.

Value

transitive reduction of P

Author(s)

David Schoch

```
library(igraph)

g <- threshold_graph(100,0.1)
P <- neighborhood_inclusion(g)
sum(P)

R <- transitive_reduction(P)
sum(R)</pre>
```

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