

# Package ‘netrankr’

September 9, 2020

**Type** Package

**Title** Analyzing Partial Rankings in Networks

**Version** 0.3.0

**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://github.com/schochastics/netrankr/>

**BugReports** <https://github.com/schochastics/netrankr/issues>

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aggregate_positions	<i>Quantification of (indirect) relations</i>
---------------------	---

---

**Description**

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

**Usage**

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

**Arguments**

tau_x	Numeric matrix containing indirect relations calculated with <a href="#">indirect_relations</a> .
type	String indicating the type of aggregation to be used. See Details for options.

## 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](#)

## Examples

```
library(igraph)
library(magrittr)

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))

#degree
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 <a href="#">neighborhood_inclusion</a> or <a href="#">positional_dominance</a> .
method	String indicating which method to be used. see Details.

### Details

The *method* parameter can be set to

**lpom** local partial order model

**glopom** 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.

**See Also**

[approx\\_rank\\_relative](#), [exact\\_rank\\_prob](#), [mcmc\\_rank\\_prob](#)

**Examples**

```
P <- matrix(c(0,0,1,1,1,0,0,0,1,0,0,0,0,1,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')
```

---

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**

<code>P</code>	A partial ranking as matrix object calculated with <a href="#">neighborhood_inclusion</a> or <a href="#">positional_dominance</a> .
<code>iterative</code>	Logical scalar if iterative approximation should be used.
<code>num.iter</code>	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)**

David Schoch

## 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 <- matrix(c(0,0,1,1,1,0,0,0,1,0,0,0,0,0,1,rep(0,10)),5,5,byrow=TRUE)
P
approx_rank_relative(P,iterative = FALSE)
approx_rank_relative(P,iterative = TRUE)
```

---

comparable\_pairs

*Comparable pairs in a partial order*

---

## Description

Calculates the fraction of comparable pairs in a partial order.

## Usage

```
comparable_pairs(P)
```

## Arguments

P                    A partial order as matrix object, e.g. calculated with [neighborhood\\_inclusion](#) or [positional\\_dominance](#).

## Value

Fraction of comparable pairs in P.

## Author(s)

David Schoch

## See Also

[incomparable\\_pairs](#)

**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)

```

---

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)**

David Schoch

**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
```

---

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**

P                    A partial ranking as matrix object calculated with [neighborhood\\_inclusion](#) or [positional\\_dominance](#).

**Value**

Directed graph as an igraph object.

**Author(s)**

David Schoch

**Examples**

```
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)
```



---

exact_rank_prob	<i>Probabilistic centrality rankings</i>
-----------------	--

---

### 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 <a href="#">neighborhood_inclusion</a> or <a href="#">positional_dominance</a> .
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: <code>rank.prob[u,k]</code> is the probability that u has rank k.
relative.rank	Matrix containing relative rank probabilities: <code>relative.rank[u,v]</code> 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 <code>only.results = FALSE</code> ).
tree	Adjacency list (incoming) of the tree of ideals (if <code>only.results = FALSE</code> ).

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)
```

---

florentine\_m

*Florentine family marriage network*

---

### Description

Florentine family marriage network

### Usage

```
florentine_m
```

### Format

An igraph object containing marriage links of florentine families

## 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	<i>Rankings that extend a partial ranking</i>
--------------	---

---

## Description

Returns all possible rankings that extend a partial ranking.

## Usage

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

## Arguments

data	List as returned by <a href="#">exact_rank_prob</a> when run with <code>only.results = FALSE</code>
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  $i$ th 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

## 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)
P
res <- exact_rank_prob(P,only.results = FALSE)
get_rankings(res)
```

---

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

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

---

incomparable\_pairs     *Incomparable pairs in a partial order*

---

### Description

Calculates the fraction of incomparable pairs in a partial order.

### Usage

```
incomparable_pairs(P)
```

### Arguments

P                    A partial order as matrix object, e.g. calculated with [neighborhood\\_inclusion](#) or [positional\\_dominance](#).

### Value

Fraction of incomparable pairs in P.

### Author(s)

David Schoch

### See Also

[comparable\\_pairs](#)

### 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)
```

---

index_builder	<i>Centrality Index Builder</i>
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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	<i>Indirect relations in a network</i>
--------------------	--

---

### 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**

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

**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  $\alpha \rightarrow 0$  and to the resistance distance as  $\alpha \rightarrow \infty$ . See (Chebotarev, 2011) for more details. The parameter `lfparam` can be used to tune  $\alpha$ .

`'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  $\alpha \rightarrow 0$  and to longest walk distances for  $\alpha \rightarrow \infty$ . 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':

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

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

### 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)
```

---

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 <a href="#">neighborhood_inclusion</a> or <a href="#">positional_dominance</a> .
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:  
 $P[u,v]=1 \ \& \ scores[i] \leq scores[j]$ .

**Value**

Logical scalar whether scores preserves the relations in P.

**Author(s)**

David Schoch

**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))
```

---

majorization_gap	<i>Majorization gap</i>
------------------	-------------------------

---

**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:

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

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.

## 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
```

## 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)
```

---

`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  $c(u) \leq c(v)$ , 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 Ortman, M., 2017. Positional Dominance: Concepts and Algorithms. *Conference on Algorithms and Discrete Applied Mathematics*, 60-71.

**See Also**

[positional\\_dominance](#), [exact\\_rank\\_prob](#)

## Examples

```

library(igraph)
#the neighborhood inclusion preorder of a star graph is complete
g <- graph.star(5, 'undirected')
P <- neighborhood_inclusion(g)
comparable_pairs(P)

#the same holds for threshold graphs
tg <- threshold_graph(50,0.1)
P <- neighborhood_inclusion(tg)
comparable_pairs(P)

#standard centrality indices preserve neighborhood-inclusion
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,closeness(g))
is_preserved(P,betweenness(g))

```

---

 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')`

---

plot\_rank\_intervals *Plot rank intervals*

---

### Description

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

### Usage

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

### Arguments

<code>P</code>	A partial ranking as matrix object calculated with <a href="#">neighborhood_inclusion</a> or <a href="#">positional_dominance</a> .
<code>cent.df</code>	A data frame containing centrality scores of indices (optional). See Details.
<code>ties.method</code>	String specifying how ties are treated in the base <a href="#">rank</a> 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](#)

**Examples**

```

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)

```

---

positional\_dominance    *Generalized Dominance in Graphs*

---

**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 <a href="#">indirect_relations</a> .
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](#)

**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))
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
```

---

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**

P 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.

**Author(s)**

David Schoch

**See Also**

[plot\\_rank\\_intervals](#), [exact\\_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)
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

**Examples**

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

---

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, bseq)
```

**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.
bseq	(0,1)-vector a binary sequence that produces a threshold graph. See details

**Details**

Either n and p must be specified or bseq. 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. If bseq is given instead, a threshold graph is constructed from that sequence. 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](#)

**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")
```

---

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)

```

**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:

`dist_2pow` returns  $2^{-x}$

`dist_inv` returns  $1/x$

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

`dist_powd` returns  $\alpha^x$  where  $\alpha$  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  $\sum_{k=0}^{\infty} \frac{A^{2k}}{(2k)!}$

`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)

David Schoch

---

transitive\_reduction *Transitive Reduction*

---

### Description

Calculates the transitive reduction of a partial ranking.

### Usage

```
transitive_reduction(P)
```

### Arguments

P A partial ranking as matrix object calculated with [neighborhood\\_inclusion](#) or [positional\\_dominance](#).

### Value

transitive reduction of P

### Author(s)

David Schoch

### Examples

```
library(igraph)

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

R <- transitive_reduction(P)
sum(R)
```

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