

Package ‘NetworkDistance’

January 20, 2025

Type Package

Title Distance Measures for Networks

Version 0.3.4

Description Network is a prevalent form of data structure in many fields. As an object of analysis, many distance or metric measures have been proposed to define the concept of similarity between two networks. We provide a number of distance measures for networks. See Jurman et al (2011) <[doi:10.3233/978-1-60750-692-8-227](https://doi.org/10.3233/978-1-60750-692-8-227)> for an overview on spectral class of inter-graph distance measures.

License MIT + file LICENSE

Encoding UTF-8

Depends R (>= 3.0.0)

Imports Matrix, Rcpp, Rdpack, RSpectra, doParallel, foreach, graphon, parallel, stats, igraph, network, pracma, utils

LinkingTo Rcpp, RcppArmadillo

Suggests graphics, knitr, rmarkdown

RdMacros Rdpack

RoxygenNote 7.1.1

VignetteBuilder knitr

NeedsCompilation yes

Author Kisung You [aut, cre] (<<https://orcid.org/0000-0002-8584-459X>>)

Maintainer Kisung You <kisungyou@outlook.com>

Repository CRAN

Date/Publication 2021-08-21 15:00:08 UTC

Contents

graph20	2
nd.centrality	3
nd.csd	4
nd.dsd	5

nd.edd	6
nd.extremal	7
nd.gdd	8
nd.graphon	9
nd.hamming	10
nd.him	11
nd.moments	12
nd.nfd	13
nd.wsd	14
NetworkDistance	15

Index	17
--------------	-----------

graph20	<i>20 adjacency matrices from Erdős–Rényi models</i>
---------	--

Description

Simulated list of 20 adjacency matrices of 28 nodes. First 10 are from Erdős–Rényi model with $p = 0.9$, and the latter 10 are generated using $p = 0.5$. Each element in the list is of size (28×28) , symmetric, having values in 0 or 1, and every diagonal element is set as 0 in accordance with no self-loop assumption.

Usage

```
data(graph20)
```

Format

A list of 20 adjacency matrices of size (28×28) .

Details

Below is the code used to generate *graph20*:

```
require(stats)
graph20 = list()
for (i in 1:10){ # type-1 adjacency matrices
  rbin = rbinom(784,1,0.9)
  mat = matrix(rbin, nrow=28)
  matout = mat*t(mat)
  diag(matout) = 0
  graph20[[i]]=matout
}
for (i in 11:20){ # type-2 adjacency matrices
  rbin = rbinom(784,1,0.5)
  mat = matrix(rbin, nrow=28)
  matout = mat*t(mat)
}
```

```

    diag(matout) = 0
    graph20[[i]]=matout
  }

```

nd centrality

Centrality Distance

Description

Centrality is a core concept in studying the topological structure of complex networks, which can be either defined for each node or edge. `nd centrality` offers 3 distance measures on node-defined centralities. See this [Wikipedia page](#) for more on network/graph centrality.

Usage

```

nd centrality(
  A,
  out.dist = TRUE,
  mode = c("Degree", "Close", "Between"),
  directed = FALSE
)

```

Arguments

<code>A</code>	a list of length N containing $(M \times M)$ adjacency matrices.
<code>out.dist</code>	a logical; TRUE for computed distance matrix as a <code>dist</code> object.
<code>mode</code>	type of node centrality definitions to be used.
<code>directed</code>	a logical; FALSE as symmetric, undirected graph.

Value

a named list containing

D an $(N \times N)$ matrix or `dist` object containing pairwise distance measures.

features an $(N \times M)$ matrix where rows are node centralities for each graph.

References

Roy M, Schmid S, Trédan G (2014). "Modeling and Measuring Graph Similarity: The Case for Centrality Distance." In *FOMC 2014, 10th ACM International Workshop on Foundations of Mobile Computing*, 53.

Examples

```
## load example data
data(graph20)

## use 3 types of centrality measures
out1 <- nd.centrality(graph20, out.dist=FALSE,mode="Degree")
out2 <- nd.centrality(graph20, out.dist=FALSE,mode="Close")
out3 <- nd.centrality(graph20, out.dist=FALSE,mode="Between")

## visualize
opar = par(no.readonly=TRUE)
par(mfrow=c(1,3), pty="s")
image(out1$D[,20:1], main="Degree", col=gray(0:32/32), axes=FALSE)
image(out2$D[,20:1], main="Close", col=gray(0:32/32), axes=FALSE)
image(out3$D[,20:1], main="Between", col=gray(0:32/32), axes=FALSE)
par(opar)
```

nd.csd

L₂ Distance of Continuous Spectral Densities

Description

The method employs spectral density of eigenvalues from Laplacian in that for each, we have corresponding spectral density $\rho(w)$ as a sum of narrow Lorentz distributions with bandwidth parameter. Since it involves integration of a function over the non-compact domain, it may blow up to infinity and the code automatically aborts the process.

Usage

```
nd.csd(A, out.dist = TRUE, bandwidth = 1)
```

Arguments

A	a list of length N containing $(M \times M)$ adjacency matrices.
out.dist	a logical; TRUE for computed distance matrix as a <code>dist</code> object.
bandwidth	common bandwidth of positive real number.

Value

a named list containing

D an $(N \times N)$ matrix or `dist` object containing pairwise distance measures.

spectra an $(N \times M - 1)$ matrix where each row is top- $M - 1$ vibrational spectra.

References

Ipsen M, Mikhailov AS (2002). “Evolutionary reconstruction of networks.” *Physical Review E*, **66**(4). ISSN 1063-651X, 1095-3787.

Examples

```
## load example data
data(graph20)

## compute distance matrix
output = nd.csd(graph20, out.dist=FALSE, bandwidth=1.0)

## visualize
opar = par(no.readonly=TRUE)
par(pty="s")
image(output$D[,20:1], main="two group case", axes=FALSE, col=gray(0:32/32))
par(opar)
```

nd.dsd

Discrete Spectral Distance

Description

Discrete Spectral Distance (DSD) is defined as the Euclidean distance between the spectra of various matrices, such as adjacency matrix A ("Adj"), (unnormalized) Laplacian matrix $L = D - A$ ("Lap"), signless Laplacian matrix $|L| = D + A$ ("SLap"), or normalized Laplacian matrix $\tilde{L} = D^{-1/2} L D^{-1/2}$.

Usage

```
nd.dsd(A, out.dist = TRUE, type = c("Lap", "SLap", "NLap", "Adj"))
```

Arguments

A a list of length N containing $(M \times M)$ adjacency matrices.
out.dist a logical; TRUE for computed distance matrix as a `dist` object.
type type of target structure. One of "Lap", "SLap", "NLap", "Adj" as defined above.

Value

a named list containing

D an $(N \times N)$ matrix or `dist` object containing pairwise distance measures.

spectra an $(N \times M - 1)$ matrix where each row is top- $M - 1$ vibrational spectra.

References

Wilson RC, Zhu P (2008). “A study of graph spectra for comparing graphs and trees.” *Pattern Recognition*, **41**(9), 2833–2841. ISSN 00313203.

Examples

```
## load example data and extract only a few
data(graph20)
gr.small = graph20[c(1:5,11:15)]

## compute distance matrix
output <- nd.dsd(gr.small, out.dist=FALSE)

## visualize
opar <- par(no.readonly=TRUE)
par(pty="s")
image(output$D[,10:1], main="two group case", axes=FALSE, col=gray(0:32/32))
par(opar)
```

nd.edd

Edge Difference Distance

Description

It is of the most simplest form that Edge Difference Distance (EDD) takes two adjacency matrices and takes Frobenius norm of their differences.

Usage

```
nd.edd(A, out.dist = TRUE)
```

Arguments

A a list of length N containing $(M \times M)$ adjacency matrices.
out.dist a logical; TRUE for computed distance matrix as a `dist` object.

Value

a named list containing

D an $(N \times N)$ matrix or `dist` object containing pairwise distance measures.

References

Hammond DK, Gur Y, Johnson CR (2013). “Graph Diffusion Distance: A Difference Measure for Weighted Graphs Based on the Graph Laplacian Exponential Kernel.” In *Proceedings of the IEEE global conference on information and signal processing (GlobalSIP'13)*, 419–422.

Examples

```
## load example data
data(graph20)

## compute distance matrix
output = nd.edd(graph20, out.dist=FALSE)

## visualize
opar <- par(no.readonly=TRUE)
par(pty="s")
image(output$D[,20:1], main="two group case", axes=FALSE, col=gray(0:32/32))
par(opar)
```

nd.extremal

*Extremal distance with top-k eigenvalues***Description**

Extremal distance (nd.extremal) is a type of spectral distance measures on two graphs' graph Laplacian,

$$L := D - A$$

where A is an adjacency matrix and $D_{ii} = \sum_j A_{ij}$. It takes top- k eigenvalues from graph Laplacian matrices and take normalized sum of squared differences as metric. Note that it is 1. *non-negative*, 2. *separated*, 3. *symmetric*, and satisfies 4. *triangle inequality* in that it is indeed a metric.

Usage

```
nd.extremal(A, out.dist = TRUE, k = ceiling(nrow(A)/5))
```

Arguments

A a list of length N containing adjacency matrices.
out.dist a logical; TRUE for computed distance matrix as a dist object.
k the number of largest eigenvalues to be used.

Value

a named list containing

D an $(N \times N)$ matrix or dist object containing pairwise distance measures.

spectra an $(N \times k)$ matrix where each row is top- k Laplacian eigenvalues.

References

Jakobson D, Rivin I (2002). "Extremal metrics on graphs I." *Forum Mathematicum*, **14**(1). ISSN 0933-7741, 1435-5337.

Examples

```
## load data
data(graph20)

## compute distance matrix
output = nd.extremal(graph20, out.dist=FALSE, k=2)

## visualize
opar = par(no.readonly=TRUE)
par(pty="s")
image(output$D[,20:1], main="two group case", col=gray(0:32/32), axes=FALSE)
par(opar)
```

nd.gdd

Graph Diffusion Distance

Description

Graph Diffusion Distance (nd.gdd) quantifies the difference between two weighted graphs of same size. It takes an idea from heat diffusion process on graphs via graph Laplacian exponential kernel matrices. For a given adjacency matrix A , the graph Laplacian is defined as

$$L := D - A$$

where $D_{ii} = \sum_j A_{ij}$. For two adjacency matrices A_1 and A_2 , GDD is defined as

$$d_{gdd}(A_1, A_2) = \max_t \sqrt{\|\exp(-tL_1) - \exp(-tL_2)\|_F^2}$$

where $\exp(\cdot)$ is matrix exponential, $\|\cdot\|_F$ a Frobenius norm, and L_1 and L_2 Laplacian matrices corresponding to A_1 and A_2 , respectively.

Usage

```
nd.gdd(A, out.dist = TRUE, vect = seq(from = 0.1, to = 1, length.out = 10))
```

Arguments

<code>A</code>	a list of length N containing adjacency matrices.
<code>out.dist</code>	a logical; TRUE for computed distance matrix as a dist object.
<code>vect</code>	a vector of parameters t whose values will be used.

Value

a named list containing

D an $(N \times N)$ matrix or dist object containing pairwise distance measures.

maxt an $(N \times N)$ matrix whose entries are maximizer of the cost function.

References

Hammond DK, Gur Y, Johnson CR (2013). “Graph Diffusion Distance: A Difference Measure for Weighted Graphs Based on the Graph Laplacian Exponential Kernel.” In *Proceedings of the IEEE global conference on information and signal processing (GlobalSIP'13)*, 419–422.

Examples

```
## load data and extract a subset
data(graph20)
gr.small = graph20[c(1:5,11:15)]

## compute distance matrix
output = nd.gdd(gr.small, out.dist=FALSE)

## visualize
opar = par(no.readonly=TRUE)
par(pty="s")
image(output$D[,10:1], main="two group case", col=gray((0:32)/32), axes=FALSE)
par(opar)
```

nd.graphon

Graphon Estimates Distance

Description

Graphon is a symmetric measurable function

$$W : [0, 1]^2 \rightarrow [0, 1]$$

that is considered to be a generating model for an observed network. `nd.graphon` computes distances between networks based on the estimated graphons of each network. Estimation methods are taken from **graphon** package. For more details, see the function links below.

Usage

```
nd.graphon(
  A,
  out.dist = TRUE,
  method = c("completion", "LG", "nbdsmooth", "SBA", "USVT"),
  ...
)
```

Arguments

<code>A</code>	a list of length N containing $(M \times M)$ adjacency matrices.
<code>out.dist</code>	a logical; TRUE for computed distance matrix as a <code>dist</code> object.
<code>method</code>	type of graphon estimation methods to be used.

... extra parameters to be passed onto graphon estimation functions. See also [est.completion](#), [est.LG](#), [est.nbdsmooth](#), [est.SBA](#), and [est.USVT](#) for details.

Value

a named list containing

D an $(N \times N)$ matrix or `dist` object containing pairwise distance measures.

References

Mukherjee SS, Sarkar P, Lin L (2017). “On clustering network-valued data.” In Guyon I, Luxburg UV, Bengio S, Wallach H, Fergus R, Vishwanathan S, Garnett R (eds.), *Advances in neural information processing systems 30*, 7071–7081. Curran Associates, Inc.

Examples

```
## load example data
data(graph20)

## compute USVT-based distance
output <- nd.graphon(graph20, out.dist=FALSE, method="usvt")

## visualize
opar = par(no.readonly=TRUE)
par(pty="s")
image(output$D[,20:1], main="USVT", col=gray(0:32/32), axes=FALSE)
par(opar)
```

nd.hamming

Hamming Distance

Description

Hamming Distance is the count of discrepancy between two binary networks for each edge. Therefore, if used with non-binary networks, it might return a warning message and distorted results. It was originally designed to compare two strings of equal length, see [Wikipedia page](#) for more detailed introduction.

Usage

```
nd.hamming(A, out.dist = TRUE)
```

Arguments

A a list of length N containing adjacency matrices.
out.dist a logical; TRUE for computed distance matrix as a `dist` object.

Value

a named list containing

D an $(N \times N)$ matrix or dist object containing pairwise distance measures.

References

Hamming RW (1950). "Error Detecting and Error Correcting Codes." *Bell System Technical Journal*, **29**(2), 147–160. ISSN 00058580.

Examples

```
## load example data and extract only a few
data(graph20)
gr.small = graph20[c(1:5,11:15)]

## compute distance matrix
output = nd.hamming(gr.small, out.dist=FALSE)

## visualize
opar = par(no.readonly=TRUE)
par(pty="s")
image(output$D[,10:1], main="two group case", axes=FALSE, col=gray(0:32/32))
par(opar)
```

nd.him

HIM Distance

Description

Hamming-Ipsen-Mikhailov (HIM) combines the local Hamming edit distance and the global Ipsen-Mikhailov distance to merge information at each scale. For Ipsen-Mikhailov distance, it is provided as `nd.csd` in our package for consistency. Given a parameter ξ (ξ), it is defined as

$$HIM_{\xi}(A, B) = \sqrt{H^2(A, B) + \xi \cdot IM^2(A, B)} / \sqrt{1 + \xi}$$

where H and IM stand for Hamming and I-M distance, respectively.

Usage

```
nd.him(A, out.dist = TRUE, xi = 1, ntest = 10)
```

Arguments

<code>A</code>	a list of length N containing $(M \times M)$ adjacency matrices.
<code>out.dist</code>	a logical; TRUE for computed distance matrix as a dist object.
<code>xi</code>	a parameter to control balance between two distances.
<code>ntest</code>	the number of searching over <code>nd.csd</code> parameter.

Value

a named list containing

D an $(N \times N)$ matrix or dist object containing pairwise distance measures.

References

Jurman G, Visintainer R, Filosi M, Riccadonna S, Furlanello C (2015). “The HIM glocal metric and kernel for network comparison and classification.” In *2015 IEEE International Conference on Data Science and Advanced Analytics (DSAA)*, 1–10. ISBN 978-1-4673-8272-4.

See Also

[nd.hamming](#), [nd.csd](#)

Examples

```
## load example data
data(graph20)

## compute distance matrix
output = nd.him(graph20, out.dist=FALSE)

## visualize
opar = par(no.readonly=TRUE)
par(pty="s")
image(output$D[,20:1], main="two group case", axes=FALSE, col=gray(0:32/32))
par(opar)
```

nd.moments

Log Moments Distance

Description

For a graph with an adjacency matrix A , *graph moment* is defined as

$$\rho_m(A) = \text{tr}(A/n)^m$$

where n is the number of vertices and m is an order of the moment. nd.moments computes pairwise distances based on log of graph moments from $m = 1$ to $m = k$.

Usage

```
nd.moments(
  A,
  k = 3,
  metric = c("euclidean", "maximum", "manhattan", "canberra", "binary", "minkowski"),
  out.dist = TRUE
)
```

Arguments

A	a list of length N containing $(M \times M)$ adjacency matrices.
k	the integer order of moments. If k is too large, it may incur numerical overflow.
metric	type of distance measures for log-moment features. See dist for more details.
out.dist	a logical; TRUE for computed distance matrix as a <code>dist</code> object.

Value

a named list containing

D an $(N \times N)$ matrix or `dist` object containing pairwise distance measures.

References

Mukherjee SS, Sarkar P, Lin L (2017). “On clustering network-valued data.” In Guyon I, Luxburg UV, Bengio S, Wallach H, Fergus R, Vishwanathan S, Garnett R (eds.), *Advances in neural information processing systems 30*, 7071–7081. Curran Associates, Inc.

Examples

```
## load example data
data(graph20)

## compute distance based on different k's.
out3 <- nd.moments(graph20, k=3, out.dist=FALSE)
out5 <- nd.moments(graph20, k=5, out.dist=FALSE)
out7 <- nd.moments(graph20, k=7, out.dist=FALSE)
out9 <- nd.moments(graph20, k=9, out.dist=FALSE)

## visualize
opar = par(no.readonly=TRUE)
par(mfrow=c(2,2), pty="s")
image(out3$D[,20:1], col=gray(0:32/32), axes=FALSE, main="k=3")
image(out5$D[,20:1], col=gray(0:32/32), axes=FALSE, main="k=5")
image(out7$D[,20:1], col=gray(0:32/32), axes=FALSE, main="k=7")
image(out9$D[,20:1], col=gray(0:32/32), axes=FALSE, main="k=9")
par(opar)
```

nd.nfd

Network Flow Distance

Description

Network Flow Distance

Usage

```
nd.nfd(
  A,
  order = 0,
  out.dist = TRUE,
  vect = seq(from = 0, to = 10, length.out = 1000)
)
```

Arguments

<code>A</code>	a list of length N containing adjacency matrices.
<code>order</code>	the order of Laplacian; currently only 0 and 1 are supported.
<code>out.dist</code>	a logical; TRUE for computed distance matrix as a <code>dist</code> object.
<code>vect</code>	a vector of parameters t whose values will be used.

Value

a named list containing

D an $(N \times N)$ matrix or `dist` object containing pairwise distance measures.

Examples

```
## Not run:
## load example data
data(graph20)

# compute two diffusion-based distances and visualize
out1 = nd.gdd(graph20, out.dist=FALSE)
out2 = nd.nfd(graph20, out.dist=FALSE)

# visualize
opar <- par(no.readonly=TRUE)
par(mfrow=c(1,2), pty="s")
image(out1$D[,20:1], col=gray((0:32)/32), main="nd.gdd", axes=FALSE)
image(out2$D[,20:1], col=gray((0:32)/32), main="nd.nfd", axes=FALSE)
par(opar)

## End(Not run)
```

Description

Normalized Laplacian matrix contains topological information of a corresponding network via its spectrum. `nd.wsd` adopts weighted spectral distribution of eigenvalues and brings about a metric via binning strategy.

Usage

```
nd.wsd(A, out.dist = TRUE, K = 50, wN = 4)
```

Arguments

A a list of length N containing $(M \times M)$ adjacency matrices.

out.dist a logical; TRUE for computed distance matrix as a `dist` object.

K the number of bins for the spectrum interval $[0, 2]$.

wN a decaying exponent; default is 4 set by authors.

Value

a named list containing

D an $(N \times N)$ matrix or `dist` object containing pairwise distance measures.

spectra an $(N \times M)$ matrix of rows being eigenvalues for each graph.

References

Fay D, Haddadi H, Thomason A, Moore AW, Mortier R, Jamakovic A, Uhlig S, Rio M (2010). “Weighted Spectral Distribution for Internet Topology Analysis: Theory and Applications.” *IEEE/ACM Transactions on Networking*, **18**(1), 164–176. ISSN 1063-6692, 1558-2566.

Examples

```
## load example data and extract a few
data(graph20)
gr.small = graph20[c(1:5, 11:15)]

## compute distance matrix
output = nd.wsd(gr.small, out.dist=FALSE, K=10)

## visualize
opar = par(no.readonly=TRUE)
par(pty="s")
image(output$D[,10:1], main="two group case", axes=FALSE, col=gray(0:32/32))
par(opar)
```

Description

Network has gathered much attention from many disciplines, as many of real data can be well represented in the relational form. The concept of distance - or, metric - between two networks is the starting point for inference on population of networks. **NetworkDistance** package provides a not-so-comprehensive collection of distance measures for measuring dissimilarity between two network objects. Data should be supplied as *adjacency* matrices, where we support three formats of data representation; matrix object in **R** base, network class from **network** package, and igraph class from **igraph** package.

Index

* datasets

graph20, 2

dist, 13

est.completion, 10

est.LG, 10

est.nbdsmooth, 10

est.SBA, 10

est.USVT, 10

graph20, 2

nd.central, 3

nd.csd, 4, 11, 12

nd.dsd, 5

nd.edd, 6

nd.extremal, 7

nd.gdd, 8

nd.graphon, 9

nd.hamming, 10, 12

nd.him, 11

nd.moments, 12

nd.nfd, 13

nd.wsd, 14

NetworkDistance, 15