Package ‘NetworkDistance’

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Title Distance Measures for Networks
Version 0.3.2
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> for an overview on spectral class of inter-graph distance measures.

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Simulated list of 20 adjacency matrices of 28 nodes. First 10 are from Erdős–Rényi model with \( p = 0.8 \), and the latter 10 are generated using \( p = 0.2 \). Each element in the list is of size \((28 \times 28)\), symmetric, having values in 0 or 1, and every diagonal element is set as 0 in accordance with no self-loop assumption.

```r
require(stats)
data(graph20)
```

A list of 20 adjacency matrices of size \((28 \times 28)\).

Below is the code used to generate `graph20`:

```r
require(stats)
graph20 = list()
for (i in 1:10){ # type-1 adjacency matrices
  rbin = rbinom(784,1,p=0.8)
  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,p=0.2)
  mat = matrix(rbin, nrow=28)
  matout = mat*t(mat)
  diag(matout) = 0
  graph20[[i]]=matout
}```
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

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

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.
- **mode**
  type of node centrality definitions to be used.
- **directed**
  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

Examples

```r
## Not run:
## 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
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)

## End(Not run)
```

nd.csd

$L_2$ 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 dist 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.
## 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 $\tilde{L} = D^{-1/2}LD^{-1/2}$.

**Usage**

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

**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 "Adj", "Lap", "SLap", "NLap" 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**

Examples

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

## Compute Distance Matrix and Visualize
## Not run:
output <- nd.dsd(gr.small, out.dist=FALSE)
opar <- par(pty="s")
image(output$D[,1:1], main="two group case", axes=FALSE, col=gray(0:32/32))
par(opar)

## End(Not run)
```

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**

```r
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**

Examples

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

## Compute Distance Matrix and Visualize
output = nd.edd(graph20, out.dist=FALSE)
opar = 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**

```r
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; \( \text{TRUE} \) for computed distance matrix as a \text{dist} object.
- **k** the number of largest eigenvalues to be used.

**Value**

a named list containing

- **D** an \((N \times N)\) matrix or \text{dist} object containing pairwise distance measures.
- **spectra** an \((N \times k)\) matrix where each row is top-\( k \) Laplacian eigenvalues.

**References**

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

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

Arguments

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


Examples

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

## Compute Distance Matrix and Visualize
output = nd.gdd(gr.small, out.dist=FALSE)
par(pty="s")
image(output$D[,10:1], main="two group case", 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

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

Arguments

- **A**: a list of length \( N \) containing adjacency matrices.
- **out.dist**: a logical; \( \text{TRUE} \) for computed distance matrix as a \text{dist} object.

Value

A named list containing

- **D**: an \((N \times N)\) matrix or \text{dist} object containing pairwise distance measures.

References

Examples

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

## compute distance and visualize
output = nd.hamming(gr.small, out.dist=FALSE)
opar = 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 \( \xi \) (\( xi \)), 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

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

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.
- **xi**: a parameter to control balance between two distances.
- **ntest**: the number of searching over nd.csd parameter.

Value

A named list containing

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

References

nd.nfd

See Also

nd.hamming, nd.csd

Examples

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

## compute distance matrix and visualize
output = nd.him(graph20, out.dist=FALSE)
image(output$D[,20:1], main="two group case", axes=FALSE, col=gray(0:32/32))

## End(Not run)

---

**nd.nfd**  
*Network Flow Distance*

**Description**

Network Flow Distance

**Usage**

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

**Arguments**

- **A**  
  a list of length $N$ containing adjacency matrices.
- **order**  
  the order of Laplacian; currently only 0 and 1 are supported.
- **out.dist**  
  a logical; TRUE for computed distance matrix as a `dist` object.
- **vect**  
  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

```r
## 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)

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)
## End(Not run)
```

### `nd.wsd`

**Distance with Weighted Spectral Distribution**

**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**

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

**Arguments**

- `A` a list of length `N` containing `(M x 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 x N)` matrix or `dist` object containing pairwise distance measures.
- `spectra` an `(N x M)` matrix of rows being eigenvalues for each graph.

**References**

NetworkDistance

**Examples**

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

## compute distance matrix and visualize
output = nd.wsd(gr.small, out.dist=FALSE, K=10)
opar = par(pty="s")
image(output$D[,10:1], main="two group case", axes=FALSE, col=gray(0:32/32))
par(opar)
```

---

**NetworkDistance**  
*Distance Measures for Networks*

**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.
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