

Package ‘clustree’

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Type Package

Title Visualise Clusterings at Different Resolutions

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Description Deciding what resolution to use can be a difficult question when approaching a clustering analysis. One way to approach this problem is to look at how samples move as the number of clusters increases. This package allows you to produce clustering trees, a visualisation for interrogating clusterings as resolution increases.

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URL <https://github.com/lazappi/clustree>

BugReports <https://github.com/lazappi/clustree/issues>

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clustree-package	<i>Clustree</i>
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Description

Deciding what resolution to use can be a difficult question when approaching a clustering analysis. One way to approach this problem is to look at how samples move as the number of clusters increases. This package allows you to produce clustering trees, a visualisation for interrogating clusterings as resolution increases.

add_node_points	<i>Add node points</i>
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Description

Add node points to a clustering tree plot with the specified aesthetics.

Usage

```
add_node_points(node_colour, node_size, node_alpha, allowed)
```

Arguments

node_colour	either a value indicating a colour to use for all nodes or the name of a metadata column to colour nodes by
node_size	either a numeric value giving the size of all nodes or the name of a metadata column to use for node sizes
node_alpha	either a numeric value giving the alpha of all nodes or the name of a metadata column to use for node transparency
allowed	vector of allowed node attributes to use as aesthetics

aggr_metadata	<i>Aggregate metadata</i>
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Description

Aggregate a metadata column to get a summarized value for a cluster node

Usage

```
aggr_metadata(node_data, col_name, col_aggr, metadata, is_cluster)
```

Arguments

node_data	data.frame containing information about a set of cluster nodes
col_name	the name of the metadata column to aggregate
col_aggr	string naming a function used to aggregate the column
metadata	data.frame providing metadata on samples
is_cluster	logical vector indicating which rows of metadata are in the node to be summarized

Value

data.frame with aggregated data

 assert_colour_node_aes

Assert colour node aesthetics

Description

Raise error if an incorrect set of colour node parameters has been supplied.

Usage

```
assert_colour_node_aes(node_aes_name, prefix, metadata, node_aes, node_aes_aggr,
  min, max)
```

Arguments

node_aes_name	name of the node aesthetic to check
prefix	string indicating columns containing clustering information
metadata	data.frame containing metadata on each sample that can be used as node aesthetics
node_aes	value of the node aesthetic to check
node_aes_aggr	aggregation function associated with the node aesthetic
min	minimum numeric value allowed
max	maximum numeric value allowed

 assert_node_aes

Assert node aesthetics

Description

Raise error if an incorrect set of node parameters has been supplied.

Usage

```
assert_node_aes(node_aes_name, prefix, metadata, node_aes, node_aes_aggr)
```

Arguments

node_aes_name	name of the node aesthetic to check
prefix	string indicating columns containing clustering information
metadata	data.frame containing metadata on each sample that can be used as node aesthetics
node_aes	value of the node aesthetic to check
node_aes_aggr	aggregation function associated with the node aesthetic

 assert_numeric_node_aes

Assert numeric node aesthetics

Description

Raise error if an incorrect set of numeric node parameters has been supplied.

Usage

```
assert_numeric_node_aes(node_aes_name, prefix, metadata, node_aes,
    node_aes_aggr, min, max)
```

Arguments

node_aes_name	name of the node aesthetic to check
prefix	string indicating columns containing clustering information
metadata	data.frame containing metadata on each sample that can be used as node aesthetics
node_aes	value of the node aesthetic to check
node_aes_aggr	aggregation function associated with the node aesthetic
min	minimum numeric value allowed
max	maximum numeric value allowed

build_tree_graph

Build tree graph

Description

Build a tree graph from a set of clusterings, metadata and associated aesthetics

Usage

```
build_tree_graph(clusterings, prefix, count_filter, prop_filter, metadata,
    node_aes_list)
```

Arguments

clusterings	numeric matrix containing clustering information, each column contains clustering at a separate resolution
prefix	string indicating columns containing clustering information
count_filter	count threshold for filtering edges in the clustering graph
prop_filter	in proportion threshold for filtering edges in the clustering graph
metadata	data.frame containing metadata on each sample that can be used as node aesthetics
node_aes_list	nested list containing node aesthetics

Value

`tidygraph::tbl_graph` object containing the tree graph

`calc_sc3_stability` *Calculate SC3 stability*

Description

Calculate the SC3 stability index for every cluster at every resolution in a set of clusterings. The index varies from 0 to 1, where 1 suggests that a cluster is more stable across resolutions. See `calc_sc3_stability_cluster()` for more details.

Usage

```
calc_sc3_stability(clusterings)
```

Arguments

`clusterings` numeric matrix containing clustering information, each column contains clustering at a separate resolution

Value

matrix with stability score for each cluster

`calc_sc3_stability_cluster`
Calculate single SC3 stability

Description

Calculate the SC3 stability index for a single cluster in a set of clusterings. The index varies from 0 to 1, where 1 suggests that a cluster is more stable across resolutions.

Usage

```
calc_sc3_stability_cluster(clusterings, res, cluster)
```

Arguments

`clusterings` numeric matrix containing clustering information, each column contains clustering at a separate resolution

`res` resolution of the cluster to calculate stability for

`cluster` index of the cluster to calculate stability for

Details

This index was originally introduced in the SC3 package for clustering single-cell RNA-seq data. Clusters are awarded increased stability if they share the same samples as a cluster at another resolution and penalised at higher resolutions. We use a slightly different notation to describe the score but the results are the same:

$$s(c_{k,i}) = \frac{1}{\text{size}(L) + 1} \sum_{l \in L} \sum_{j \in N_l} \frac{\text{size}(c_{k,i} \cap c_{l,j})}{\text{size}(c_{l,j}) * \text{size}(N_l)^2}$$

Where:

- $c_{\{x, y\}}$ is cluster y at resolution x
- k is the resolution of the cluster we want to score
- i is the index of the cluster we want to score
- L is the set of all resolutions except k
- l is a resolution in L
- N_l is the set of clusters at resolution l that share samples with $c_{\{k, i\}}$
- j is a cluster in N_l

Value

SC3 stability index

See Also

The documentation for the `calculate_stability` function in the SC3 package

clustree	<i>Plot a clustering tree</i>
----------	-------------------------------

Description

Creates a plot of a clustering tree showing the relationship between clusterings at different resolutions.

Usage

```
clustree(x, ...)
```

```
## S3 method for class 'matrix'
clustree(x, prefix, suffix = NULL, metadata = NULL,
  count_filter = 0, prop_filter = 0.1, layout = c("tree", "sugiyama"),
  use_core_edges = TRUE, highlight_core = FALSE, node_colour = prefix,
  node_colour_aggr = NULL, node_size = "size", node_size_aggr = NULL,
  node_size_range = c(4, 15), node_alpha = 1, node_alpha_aggr = NULL,
```

```

node_text_size = 3, scale_node_text = FALSE, node_text_colour = "black",
edge_width = 1.5, edge_arrow = TRUE, edge_arrow_ends = c("last",
"first", "both"), return = c("plot", "graph", "layout"), ...)

## S3 method for class 'data.frame'
clustree(x, prefix, ...)

## S3 method for class 'SingleCellExperiment'
clustree(x, prefix, exprs = "counts", ...)

## S3 method for class 'seurat'
clustree(x, prefix = "res.", exprs = c("data", "raw.data",
"scale.data"), ...)

```

Arguments

<code>x</code>	object containing clustering data
<code>...</code>	extra parameters passed to other methods
<code>prefix</code>	string indicating columns containing clustering information
<code>suffix</code>	string at the end of column names containing clustering information
<code>metadata</code>	data.frame containing metadata on each sample that can be used as node aesthetics
<code>count_filter</code>	count threshold for filtering edges in the clustering graph
<code>prop_filter</code>	in proportion threshold for filtering edges in the clustering graph
<code>layout</code>	string specifying the "tree" or "sugiyama" layout, see igraph::layout_as_tree() and igraph::layout_with_sugiyama() for details
<code>use_core_edges</code>	logical, whether to only use core tree (edges with maximum in proportion for a node) when creating the graph layout, all (unfiltered) edges will still be displayed
<code>highlight_core</code>	logical, whether to increase the edge width of the core network to make it easier to see
<code>node_colour</code>	either a value indicating a colour to use for all nodes or the name of a metadata column to colour nodes by
<code>node_colour_aggr</code>	if <code>node_colour</code> is a column name than a string giving the name of a function to aggregate that column for samples in each cluster
<code>node_size</code>	either a numeric value giving the size of all nodes or the name of a metadata column to use for node sizes
<code>node_size_aggr</code>	if <code>node_size</code> is a column name than a string giving the name of a function to aggregate that column for samples in each cluster
<code>node_size_range</code>	numeric vector of length two giving the maximum and minimum point size for plotting nodes
<code>node_alpha</code>	either a numeric value giving the alpha of all nodes or the name of a metadata column to use for node transparency

<code>node_alpha_aggr</code>	if <code>node_aggr</code> is a column name than a string giving the name of a function to aggregate that column for samples in each cluster
<code>node_text_size</code>	numeric value giving the size of node labels if <code>scale_node_text</code> is FALSE
<code>scale_node_text</code>	logical indicating whether to scale node labels along with the node size
<code>node_text_colour</code>	colour value for node labels
<code>edge_width</code>	numeric value giving the width of plotted edges
<code>edge_arrow</code>	logical indicating whether to add an arrow to edges
<code>edge_arrow_ends</code>	string indicating which ends of the line to draw arrow heads if <code>edge_arrow</code> is TRUE, one of "last", "first", or "both"
<code>return</code>	string specifying what to return, either "plot" (a ggplot object), "graph" (a <code>tbl_graph</code> object) or "layout" (a <code>ggraph</code> layout object)
<code>exprs</code>	source of gene expression information to use as node aesthetics, for <code>SingleCellExperiment</code> objects it must be a name in <code>assayNames(x)</code> , for a <code>seurat</code> object it must be one of <code>data</code> , <code>raw.data</code> or <code>scale.data</code>

Details

Data sources

Plotting a clustering tree requires information about which cluster each sample has been assigned to at different resolutions. This information can be supplied in various forms, as a matrix, `data.frame` or more specialised object. In all cases the object provided must contain numeric columns with the naming structure PXS where P is a prefix indicating that the column contains clustering information, X is a numeric value indicating the clustering resolution and S is any additional suffix to be removed. For `SingleCellExperiment` objects this information must be in the `colData` slot and for `Seurat` objects it must be in the `meta.data` slot. For all objects except matrices any additional columns can be used as aesthetics, for matrices an additional metadata `data.frame` can be supplied if required.

Filtering

Edges in the graph can be filtered by adjusting the `count_filter` and `prop_filter` parameters. The `count_filter` removes any edges that represent less than that number of samples, while the `prop_filter` removes edges that represent less than that proportion of cells in the node it points towards.

Node aesthetics

The aesthetics of the plotted nodes can be controlled in various ways. By default the colour indicates the clustering resolution, the size indicates the number of samples in that cluster and the transparency is set to 100. Each of these can be set to a specific value or linked to a supplied metadata column. For a `SingleCellExperiment` or `Seurat` object the names of genes can also be used. If a metadata column is used than an aggregation function must also be supplied to combine the samples in each cluster. This function must take a vector of values and return a single value.

Layout

The clustering tree can be displayed using either the Reingold-Tilford tree layout algorithm or the Sugiyama layout algorithm for layered directed acyclic graphs. These layouts were selected

as they are the algorithms available in the `igraph` package designed for trees. The Reingold-Tilford algorithm places children below their parents while the Sugiyama places nodes in layers while trying to minimise the number of crossing edges. See `igraph::layout_as_tree()` and `igraph::layout_with_sugiyama()` for more details. When `use_core_edges` is `TRUE` (default) only the core tree of the maximum proportion edges for each node are used for constructing the layout. This can often lead to more attractive layouts where the core tree is more visible.

Value

a `ggplot` object (default), a `tbl_graph` object or a `ggraph` layout object depending on the value of `return`

Examples

```
data(iris_clusts)
clustree(iris_clusts, prefix = "K")
```

clustree_overlay	<i>Overlay a clustering tree</i>
------------------	----------------------------------

Description

Creates a plot of a clustering tree overlaid on a scatter plot of individual samples.

Usage

```
clustree_overlay(x, ...)

## S3 method for class 'matrix'
clustree_overlay(x, prefix, metadata, x_value, y_value,
  suffix = NULL, count_filter = 0, prop_filter = 0.1,
  node_colour = prefix, node_colour_aggr = NULL, node_size = "size",
  node_size_aggr = NULL, node_size_range = c(4, 15), node_alpha = 1,
  node_alpha_aggr = NULL, edge_width = 1, use_colour = c("edges",
  "points"), alt_colour = "black", point_size = 3, point_alpha = 0.2,
  point_shape = 18, label_nodes = FALSE, label_size = 3,
  plot_sides = FALSE, side_point_jitter = 0.45, side_point_offset = 1,
  ...)

## S3 method for class 'data.frame'
clustree_overlay(x, prefix, ...)

## S3 method for class 'SingleCellExperiment'
clustree_overlay(x, prefix, x_value, y_value,
  exprs = "counts", red_dim = NULL, ...)

## S3 method for class 'seurat'
```

```
clustree_overlay(x, x_value, y_value, prefix = "res.",
  exprs = c("data", "raw.data", "scale.data"), red_dim = NULL, ...)
```

Arguments

<code>x</code>	object containing clustering data
<code>...</code>	extra parameters passed to other methods
<code>prefix</code>	string indicating columns containing clustering information
<code>metadata</code>	data.frame containing metadata on each sample that can be used as node aesthetics
<code>x_value</code>	numeric metadata column to use as the x axis
<code>y_value</code>	numeric metadata column to use as the y axis
<code>suffix</code>	string at the end of column names containing clustering information
<code>count_filter</code>	count threshold for filtering edges in the clustering graph
<code>prop_filter</code>	in proportion threshold for filtering edges in the clustering graph
<code>node_colour</code>	either a value indicating a colour to use for all nodes or the name of a metadata column to colour nodes by
<code>node_colour_aggr</code>	if <code>node_colour</code> is a column name than a string giving the name of a function to aggregate that column for samples in each cluster
<code>node_size</code>	either a numeric value giving the size of all nodes or the name of a metadata column to use for node sizes
<code>node_size_aggr</code>	if <code>node_size</code> is a column name than a string giving the name of a function to aggregate that column for samples in each cluster
<code>node_size_range</code>	numeric vector of length two giving the maximum and minimum point size for plotting nodes
<code>node_alpha</code>	either a numeric value giving the alpha of all nodes or the name of a metadata column to use for node transparency
<code>node_alpha_aggr</code>	if <code>node_aggr</code> is a column name than a string giving the name of a function to aggregate that column for samples in each cluster
<code>edge_width</code>	numeric value giving the width of plotted edges
<code>use_colour</code>	one of "edges" or "points" specifying which element to apply the colour aesthetic to
<code>alt_colour</code>	colour value to be used for edges or points (whichever is NOT given by <code>use_colour</code>)
<code>point_size</code>	numeric value giving the size of sample points
<code>point_alpha</code>	numeric value giving the alpha of sample points
<code>point_shape</code>	numeric value giving the shape of sample points
<code>label_nodes</code>	logical value indicating whether to add labels to clustering graph nodes
<code>label_size</code>	numeric value giving the size of node labels is <code>label_nodes</code> is TRUE
<code>plot_sides</code>	logical value indicating whether to produce side on plots

side_point_jitter	numeric value giving the y-direction spread of points in side plots
side_point_offset	numeric value giving the y-direction offset for points in side plots
exprs	source of gene expression information to use as node aesthetics, for <code>SingleCellExperiment</code> objects it must be a name in <code>assayNames(x)</code> , for a <code>Seurat</code> object it must be one of <code>data</code> , <code>raw.data</code> or <code>scale.data</code>
red_dim	dimensionality reduction to use as a source for <code>x_value</code> and <code>y_value</code>

Details

Data sources

Plotting a clustering tree requires information about which cluster each sample has been assigned to at different resolutions. This information can be supplied in various forms, as a matrix, data.frame or more specialised object. In all cases the object provided must contain numeric columns with the naming structure PXS where P is a prefix indicating that the column contains clustering information, X is a numeric value indicating the clustering resolution and S is any additional suffix to be removed. For `SingleCellExperiment` objects this information must be in the `colData` slot and for `Seurat` objects it must be in the `meta.data` slot. For all objects except matrices any additional columns can be used as aesthetics.

Filtering

Edges in the graph can be filtered by adjusting the `count_filter` and `prop_filter` parameters. The `count_filter` removes any edges that represent less than that number of samples, while the `prop_filter` removes edges that represent less than that proportion of cells in the node it points towards.

Node aesthetics

The aesthetics of the plotted nodes can be controlled in various ways. By default the colour indicates the clustering resolution, the size indicates the number of samples in that cluster and the transparency is set to 100. Each of these can be set to a specific value or linked to a supplied metadata column. For a `SingleCellExperiment` or `Seurat` object the names of genes can also be used. If a metadata column is used then an aggregation function must also be supplied to combine the samples in each cluster. This function must take a vector of values and return a single value.

Colour aesthetic

The colour aesthetic can be applied to either edges or sample points by setting `use_colour`. If "edges" is selected edges will be coloured according to the clustering resolution they originate at. If "points" is selected they will be coloured according to the cluster they are assigned to at the highest resolution.

Dimensionality reductions

For `SingleCellExperiment` and `Seurat` objects precomputed dimensionality reductions can be used for x or y aesthetics. To do so `red_dim` must be set to the name of a dimensionality reduction in `reducedDimNames(x)` (for a `SingleCellExperiment`) or `x@dr` (for a `Seurat` object). `x_value` and `y_value` can then be set to `red_dimX` when `red_dim` matches the `red_dim` argument and X is the column of the dimensionality reduction to use.

Value

a ggplot object if plot_sides is FALSE or a list of ggplot objects if plot_sides is TRUE

Examples

```
data(iris_clusts)
clustree_overlay(iris_clusts, prefix = "K", x_value = "PC1", y_value = "PC2")
```

get_tree_edges	<i>Get tree edges</i>
----------------	-----------------------

Description

Extract the edges from a set of clusterings

Usage

```
get_tree_edges(clusterings, prefix)
```

Arguments

clusterings	numeric matrix containing clustering information, each column contains clustering at a separate resolution
prefix	string indicating columns containing clustering information

Value

data.frame containing edge information

get_tree_nodes	<i>Get tree nodes</i>
----------------	-----------------------

Description

Extract the nodes from a set of clusterings and add relevant attributes

Usage

```
get_tree_nodes(clusterings, prefix, metadata, node_aes_list)
```

Arguments

clusterings	numeric matrix containing clustering information, each column contains clustering at a separate resolution
prefix	string indicating columns containing clustering information
metadata	data.frame containing metadata on each sample that can be used as node aesthetics
node_aes_list	nested list containing node aesthetics

Value

data.frame containing node information

iris_clusts	<i>Clustered Iris dataset</i>
-------------	-------------------------------

Description

Iris dataset clustered using k-means with a range of values of k

Usage

```
iris_clusts
```

Format

iris_clusts is a data.frame containing the normal iris dataset with additional columns holding k-means clusterings at different values of k and the first two principal components

Source

```
set.seed(1)
iris_mat <- as.matrix(iris[1:4])
iris_km <- sapply(1:5, function(x) {
  km <- kmeans(iris_mat, centers = x, iter.max = 100, nstart = 10)
  km$cluster
})
colnames(iris_km) <- paste0("K", 1:5)
iris_clusts <- cbind(iris, iris_km)
iris_pca <- prcomp(iris_clusts[1:4])
iris_clusts$PC1 <- iris_pca$x[, 1]
iris_clusts$PC2 <- iris_pca$x[, 2]
```

overlay_node_points *Overlay node points*

Description

Overlay clustering tree nodes on a scatter plot with the specified aesthetics.

Usage

```
overlay_node_points(nodes, x_value, y_value, node_colour, node_size, node_alpha)
```

Arguments

nodes	data.frame describing nodes
x_value	column of nodes to use for the x position
y_value	column of nodes to use for the y position
node_colour	either a value indicating a colour to use for all nodes or the name of a metadata column to colour nodes by
node_size	either a numeric value giving the size of all nodes or the name of a metadata column to use for node sizes
node_alpha	either a numeric value giving the alpha of all nodes or the name of a metadata column to use for node transparency

plot_overlay_side *Plot overlay side*

Description

Plot the side view of a clustree overlay plot. If the ordinary plot shows the tree from above this plot shows it from the side, highlighting either the x or y dimension and the clustering resolution.

Usage

```
plot_overlay_side(nodes, edges, points, prefix, side_value, graph_attr,
  node_size_range, edge_width, use_colour, alt_colour, point_size, point_alpha,
  point_shape, label_nodes, label_size, y_jitter, y_offset)
```

Arguments

nodes	data.frame describing nodes
edges	data.frame describing edges
points	data.frame describing points
prefix	string indicating columns containing clustering information
side_value	string giving the metadata column to use for the x axis
graph_attr	list describing graph attributes
node_size_range	numeric vector of length two giving the maximum and minimum point size for plotting nodes
edge_width	numeric value giving the width of plotted edges
use_colour	one of "edges" or "points" specifying which element to apply the colour aesthetic to
alt_colour	colour value to be used for edges or points (whichever is NOT given by use_colour)
point_size	numeric value giving the size of sample points
point_alpha	numeric value giving the alpha of sample points
point_shape	numeric value giving the shape of sample points
label_nodes	logical value indicating whether to add labels to clustering graph nodes
label_size	numeric value giving the size of node labels is label_nodes is TRUE
y_jitter	numeric value giving the y-direction spread of points in side plots
y_offset	numeric value giving the y-direction offset for points in side plots

Value

RETURN_DESCRIPTION

sc_example	<i>Simulated scRNA-seq dataset</i>
------------	------------------------------------

Description

A simulated scRNA-seq dataset generated using the `splatter` package and clustered using the SC3 and Seurat packages.

Usage

```
sc_example
```

Format

`sc_example` is a list holding a simulated scRNA-seq dataset. Items in the list included the simulated counts, normalised log counts, tSNE dimensionality reduction and cell assignments from SC3 and Seurat clustering.

Source

```

# Simulation
library("splatter") # Version 1.2.1

sim <- splatSimulate(batchCells = 200, nGenes = 10000,
                    group.prob = c(0.4, 0.2, 0.2, 0.15, 0.05),
                    de.prob = c(0.1, 0.2, 0.05, 0.1, 0.05),
                    method = "groups", seed = 1)
sim_counts <- counts(sim)[1:1000, ]

# SC3 Clustering
library("SC3") # Version 1.7.6
library("scater") # Version 1.6.2

sim_sc3 <- SingleCellExperiment(assays = list(counts = sim_counts))
rowData(sim_sc3)$feature_symbol <- rownames(sim_counts)
sim_sc3 <- normalise(sim_sc3)
sim_sc3 <- sc3(sim_sc3, ks = 1:8, biology = FALSE, n_cores = 1)
sim_sc3 <- runTSNE(sim_sc3)

# Seurat Clustering
library("Seurat") # Version 2.2.0

sim_seurat <- CreateSeuratObject(sim_counts)
sim_seurat <- NormalizeData(sim_seurat, display.progress = FALSE)
sim_seurat <- FindVariableGenes(sim_seurat, do.plot = FALSE,
                               display.progress = FALSE)
sim_seurat <- ScaleData(sim_seurat, display.progress = FALSE)
sim_seurat <- RunPCA(sim_seurat, do.print = FALSE)
sim_seurat <- FindClusters(sim_seurat, dims.use = 1:6,
                           resolution = seq(0, 1, 0.1),
                           print.output = FALSE)

sc_example <- list(counts = counts(sim_sc3),
                  tsne = reducedDim(sim_sc3),
                  sc3_clusters = colData(sim_sc3),
                  seurat_clusters = sim_seurat@meta.data)

```

store_node_aes

Store node aesthetics

Description

Store the names of node attributes to use as aesthetics as graph attributes

Usage

```
store_node_aes(graph, node_aes_list, metadata)
```

Arguments

graph	graph to store attributes in
node_aes_list	nested list containing node aesthetics
metadata	data.frame containing metadata that can be used as aesthetics

Value

graph with additional attributes

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