Package 'scBubbletree'

September 16, 2024

Description scBubbletree is a quantitative method for visual exploration of scRNA-seq data. It preserves biologically meaningful properties of scRNA-seq data, such as local and global cell distances, as well as the density distribution of cells across the sample. scBubbletree is scalable and avoids the overplotting problem, and is able to visualize diverse cell attributes derived from multiomic single-cell experiments. Importantly, Importantly, scBubbletree is easy to use and to integrate with popular approaches for scRNA-seq data analysis. **License** GPL-3 + file LICENSE **Depends** R (>= 4.2.0) Imports reshape2, future, future.apply, ape, scales, Seurat, ggplot2, ggtree, patchwork, proxy, methods, stats, base, utils Suggests BiocStyle, knitr, testthat, cluster, SingleCellExperiment **Encoding UTF-8** NeedsCompilation no biocViews Visualization, Clustering, SingleCell, Transcriptomics, RNASeq BugReports https://github.com/snaketron/scBubbletree/issues URL https://github.com/snaketron/scBubbletree **SystemRequirements** Python (>= 3.6), leidenalg (>= 0.8.2) RoxygenNote 6.1.1 VignetteBuilder knitr git_url https://git.bioconductor.org/packages/scBubbletree git branch RELEASE 3 19 git_last_commit 1efa88b git_last_commit_date 2024-04-30

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Author Simo Kitanovski [aut, cre]

Maintainer Simo Kitanovski <simokitanovski@gmail.com>

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Description

Method for quantitative visualization of single cell RNA-seq data

Details

This package contains functions for clustering, hierarchical grouping of clusters and visualization of scRNA-seq data.

Author(s)

Authors and maintainers:

• Simo Kitanovski <simokitanovski@uni-due.de>(ORCID)

See Also

Useful links:

- https://github.com/snaketron/scBubbletree
- Report bugs at https://github.com/snaketron/scBubbletree/issues

d_500

d_500 Dataset: 500 PBMCs

Description

d_500 is a list with 3 elements:

- 1. A = numeric matrix A^500x15 with n=500 rows for PBMCs and f=15 principal components.
- 2. f = character vector f of length 500. Each element in f represents the predicted cell type of a specific cell.
- 3. fs = numeric matrix containing normalized gene expressions of 12 marker genes in 500 cells.

Usage

```
data("d_500", package = "scBubbletree")
```

Format

Format of d_500: list

Details

This data is a sample drawn from a larger dataset of 2,700 PBMCs. The original dataset was processed as described in vignette (accessed 23, Sep, 2022):

https://satijalab.org/seurat/articles/multimodal_reference_mapping.html

See R script inst/script/get_d_500.R to see how this dataset was created.

Source

https://satijalab.org/seurat/articles/multimodal_reference_mapping.html

```
data("d_500", package = "scBubbletree")
A <- d_500$A
base::dim(A)

f <- d_500$f
base::table(f)

fs <- d_500$fs
base::dim(fs)</pre>
```

d_ccl

d_ccl

Dataset: scRNA-seq data of 3,918 cells from 5 adenocarcinoma cell lines

Description

d_ccl is a list with 3 elements:

- 1. A = numeric matrix with n=3,918 rows for cells and f=15 principal components
- 2. m = data.frame meta data
- 3. e = numeric matrix containing normalized gene expressions of 5 marker genes

Usage

```
data("d_ccl", package = "scBubbletree")
```

Format

Format of d_ccl: list

Details

d_ccl is a scRNA-seq dataset containing a mixture of 3,918 cells from five human lung adenocarcinoma cell lines (HCC827, H1975, A549, H838 and H2228). The dataset is available here:

https://github.com/LuyiTian/sc_mixology/blob/master/data/ sincell_with_class_5cl.RData

The library has been prepared with 10x Chromium platform and sequenced with Illumina NextSeq 500 platform. Raw data has been processed with Cellranger. The tool demuxlet has been used to predict the identity of each cell based on known genetic differences between the different cell lines.

See R script inst/script/get_d_ccl.R to see how this dataset was created.

Source

https://github.com/LuyiTian/sc_mixology/blob/master/data/ sincell_with_class_5cl.RData

References

Tian, Luyi, et al. "Benchmarking single cell RNA-sequencing analysis pipelines using mixture control experiments." Nature methods 16.6 (2019): 479-487

```
data("d_ccl", package = "scBubbletree")
A <- d_ccl$A
base::dim(A)

m <- d_ccl$m
utils::head(m)</pre>
```

```
e <- d_ccl$e
base::dim(e)</pre>
```

get_bubbletree_dummy

Build bubbletree given matrix A and vector cs of externally generated cluster IDs

Description

get_bubbletree_dummy takes two main inputs:

- 1. numeric matrix $A^{n \times f}$, which represents a low-dimensional projection (obtained e.g. by PCA) of the original high-dimensional scRNA-seq data, with n rows as cells and f columns as low-dimension features.
- 2. vector cs of cluster IDs of each cell

The function get_bubbletree_dummy performs one main operation. It organizes the bubbles (defined by cs) in a hierarchical dendrogram (bubbletree) which represents the hierarchical relationships between the clusters (bubbles).

Usage

Arguments

X	numeric matrix ($A^{n \times f}$ with n cells, and f low-dimensional projections of the
	original single cell RNA-seq dataset)

cs vector, cluster IDs

integer, number of bootstrap iterations to perform in order to generate bubbletree. If B=0, cluster centroids are used to compute inter-cluster distances and N_eff is ignored, i.e. all cells are used to compute centroids.

N_eff integer, number of cells to draw randomly from each cluster when computing inter-cluster distances. Maximum available number of cells are used for clusters that contain lower number of cells than N_eff

hclust_distance

distance measure to be used: euclidean (default) or manhattan, see documenta-

tion of stats::dist

hclust_method the agglomeration method to be used, default = average. See documentation of

stats::hclust

cores integer, number of PC cores for parallel execution

round_digits integer, number of decimal places to keep when showing the relative frequency

of cells in each bubble

show_simple_count

logical, if show_simple_count=T, cell counts in each bubble will be divided by 1,000 to improve readability. This is only useful for samples that are composed

of millions of cells.

verbose logical, progress messages

Details

This function is similar to get_bubbletree_kmeans and get_bubbletree_graph. It skips the clustering step. See documentation of get_bubbletree_kmeans and get_bubbletree_graph.

Value

A input x matrix k number of clusters

km NULL

ph boot_ph: bootstrap dendrograms H_b ; main_ph: consensus dendrogram \hat{H}

pair_dist inter-cluster distances used to generate the dendrograms

cluster cluster assignments of each cell input_par list of all input parameters tree ggtree bubbletree object

tree_meta meta-data associated with the bubbletree

Author(s)

Simo Kitanovski <simo.kitanovski@uni-due.de>

See Also

get_k, get_r, get_bubbletree_kmeans, get_bubbletree_graph, get_gini, get_gini_k, get_num_tiles, get_num_violins, get_cat_tiles, d_500

```
# input data
data("d_500", package = "scBubbletree")
A <- d_500$A</pre>
```

get_bubbletree_graph 7

get_bubbletree_graph Louvain clustering and hierarchical grouping of k' clusters (bubbles)

Description

get_bubbletree_graph takes two main inputs:

- 1. numeric matrix $A^{n \times f}$, which represents a low-dimensional projection (obtained e.g. by PCA) of the original high-dimensional scRNA-seq data, with n rows as cells and f columns as low-dimension features.
- 2. clustering resolution r

The function get_bubbletree_graph performs two main operations. First, it performs Louvain clustering to identify groups (bubbles) of transcriptionally similar cells; second, it organizes the bubbles in a hierarchical dendrogram (bubbletree) which adequatly represents inter-cluster relationships.

Usage

Arguments

- x numeric matrix $(A^{n \times f})$ with n cells, and f low-dimensional projections of the original single cell RNA-seq dataset)
- r number, clustering resolution

B integer, number of bootstrap iterations to perform in order to generate bubble-

tree. If B=0, cluster centroids are used to compute inter-cluster distances and

N_eff is ignored, i.e. all cells are used to compute centroids.

N_eff integer, number of cells to draw randomly from each cluster when computing

inter-cluster distances. Maximum available number of cells are used for clusters

that contain lower number of cells than N eff

n_start, iter_max

parameters for Louvain clustering, see documentation of function FindClusters,

R-package Seurat

algorithm character, four clustering algorithms: 'original', 'LMR', 'SLM' and 'Leiden',

see documentation of function FindClusters, R-package Seurat

knn_k integer, defines k for the k-nearest neighbor algorithm, see documentation of

function FindClusters, R-package Seurat

hclust_method the agglomeration method to be used, default = average. See documentation of

stats::hclust

hclust_distance

distance measure to be used: euclidean (default) or manhattan, see documenta-

tion of stats::dist

cores integer, number of PC cores for parallel execution

round_digits integer, number of decimal places to keep when showing the relative frequency

of cells in each bubble

show_simple_count

logical, if show_simple_count=T, cell counts in each bubble will be divided by

1,000 to improve readability. This is only useful for samples that are composed

of millions of cells.

verbose logical, progress messages

Details

For Louvain clustering get_bubbletree_graph uses the function FindClusters implemented in R-package Seurat. For additional information on the clustering procedure see the documentation of FindClusters. To organize the resulting clusters in a hierarchical dendrogram the algorithm performs the following steps:

- 1. In bootrap iteration b from 1:B
- 2. draw up to N_{eff} number of cells at random from each cluster without replacement
- 3. compute Euclidean distances (in space $A^{n \times f}$) between all pairs of cells in cluster i and cluster j
- 4. compute mean Euclidean distance between cluster i and j and populate inter-cluster distance matrix $D_b^{k \times k}$
- 5. perform hierarchical clustering with average linkage based on $D_b^{k \times k}$ to generate dendrogram H_b
- 6. compute average distance matrix \hat{D} and use is as input to build consensus hierarchical dendrogram (\hat{H} ; bubbletree) with average linkage
- 7. quantify branch robustness in \hat{H} count how many times each branch is found among bootrap dendrograms (H_b
- 8. visualize the bubbletree (\hat{H}) with R-package ggtree

get_bubbletree_graph 9

Special case: If B=0, then cluster centroids are used to compute inter-cluster distances and N_eff is ignored, i.e. all cells are used to compute centroids. This leads to computational efficiency, however, by doing so we lose information about the robustness of branches.

Value

A	input x matrix
k	number of clusters
r	clustering resolution
ph	boot_ph: bootstrap dendrograms H_b ; main_ph: consensus dendrogram \hat{H}
pair_dist	inter-cluster distances used to generate the dendrograms
cluster	cluster assignments of each cell
input_par	list of all input parameters
tree	ggtree bubbletree object
tree_meta	meta-data associated with the bubbletree

Author(s)

Simo Kitanovski <simo.kitanovski@uni-due.de>

See Also

get_k, get_bubbletree_dummy, get_bubbletree_kmeans, get_gini, get_gini_k, d_500, get_num_tiles, get_num_violins, get_cat_tiles

Examples

b\$tree

get_bubbletree_kmeans k-means clustering and hierarchical grouping of k clusters (bubbles)

Description

get_bubble_kmeans takes two main inputs:

1. numeric matrix $A^{n \times f}$, which represents a low-dimensional projection (obtained e.g. by PCA) of the original high-dimensional scRNA-seq data, with n rows as cells and f columns as low-dimension features.

2. number k of clusters

The function get_bubble_kmeans performs two main operations. First, it performs k-means clustering to identify groups (bubbles) of transcriptionally similar cells; second, it organizes the bubbles in a hierarchical dendrogram (bubbletree) which adequatly represents inter-cluster relationships.

Usage

Arguments

N_eff

X	numeric matrix ($A^{n \times f}$ with n cells, and f low-dimensional projections of the
	original single cell RNA-seq dataset)

k integer, number of clusters

B integer, number of bootstrap iterations to perform in order to generate bubbletree. If B=0, cluster centroids are used to compute inter-cluster distances and N_eff is ignored, i.e. all cells are used to compute centroids.

integer, number of cells to draw randomly from each cluster when computing inter-cluster distances. Maximum available number of cells are used for clusters that contain lower number of cells than N_eff

n_start, iter_max, kmeans_algorithm

parameters for k-means clustering, see documentation of function k-means, R-package stats

hclust_distance

distance measure to be used: euclidean (default) or manhattan, see documenta-

tion of stats::dist

hclust_method the agglomeration method to be used, default = average. See documentation of

stats::hclust

cores integer, number of PC cores for parallel execution

round_digits integer, number of decimal places to keep when showing the relative frequency

of cells in each bubble

show_simple_count

logical, if show_simple_count=T, cell counts in each bubble will be divided by 1,000 to improve readability. This is only useful for samples that are composed

of millions of cells.

verbose logical, progress messages

Details

For k-means clustering get_bubble_kmeans uses the function kmeans implemented in R-package stats (version 4.2.0). For additional information on the clustering procedure see the documentation of kmeans. To organize the resulting clusters in a hierarchical dendrogram the algorithm performs the following steps:

1. In bootrap iteration b from 1:B

2. draw up to N_{eff} number of cells at random from each cluster without replacement

3. compute Euclidean distances (in space $A^{n \times f}$) between pairs of cells in cluster i and cluster j

4. compute mean Euclidean distance between cluster i and j and populate inter-cluster distance matrix $D_b^{k \times k}$

5. perform hierarchical clustering with average linkage based on $D_b^{k \times k}$ to generate dendrogram H_b

6. compute average distance matrix \hat{D} and use is as input to build consensus hierarchical dendrogram (\hat{H} ; bubbletree) with average linkage

7. quantify branch robustness in \hat{H} count how many times each branch is found among bootrap dendrograms (H_b

8. visualize the bubbletree (\hat{H}) with R-package ggtree

Special case: If B=0, then cluster centroids are used to compute inter-cluster distances and N_eff is ignored, i.e. all cells are used to compute centroids. This leads to computational efficiency, however, by doing so we lose information about the robustness of branches.

Value

A input matrix x k number of clusters

km k-means clustering results identical to those generated by function k-means from

R-package stats

ph boot_ph: bootstrap dendrograms H_b ; main_ph: consensus dendrogram \hat{H}

pair_dist inter-cluster distances used to generate the dendrograms

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```
cluster cluster assignments of each cell
input_par list of all input parameters
tree ggtree bubbletree object
tree_meta meta-data associated with the bubbletree
```

Author(s)

Simo Kitanovski <simo.kitanovski@uni-due.de>

See Also

```
get_k, get_bubbletree_dummy, get_bubbletree_graph, get_gini, get_gini_k, d_500, get_num_tiles, get_num_violins, get_cat_tiles
```

Examples

get_cat_tiles

b\$tree

Visualization of categorical cell features using tile plots

Description

get_cat_tiles creates tile plot to visualize the relative frequency of categorical cell features between and within the bubbles of a bubbletree

Usage

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Arguments

btd bubbletree object

f character vector, categorical cell features

integrate_vertical

logical, if integrate_vertical=TRUE: relative frequency of the features is shown in each bubble, if integrate_vertical=FALSE: relative frequencies of the features

is shown within each bubble

round_digits integer, number of decimal places to keep when showing the relative frequency

of cells in each bubble

tile_text_size integer, size of tile labels

x_axis_name character, x-axis title

rotate_x_axis_labels

logical, should the x-axis labels be shown horizontally (rotate_x_axis_labels =

FALSE) or vertically (rotate_x_axis_labels = TRUE)

tile_bw logical, tile grayscale (tile_bw = TRUE) vs. color (tile_bw = FALSE, default)

Details

get_cat_tiles uses two main inputs:

- 1. bubbletree object
- 2. character vector of categorical cell features.

The order of the cells used to generat the bubbletree (input 1.) should correspond to the order of cells in the vector of categorical cell features (input 2.)

This function computes:

- 1. with integrate_vertical=T: relative frequencies of each feature across the different bubbles
- 2. with integrate_vertical=F: within-bubble relative frequencies (composition) of different features

Value

plot ggplot2, tile plot

table data.frame, raw data used to generate the plot

Author(s)

Simo Kitanovski <simo.kitanovski@uni-due.de>

See Also

get_k, get_r get_bubbletree_dummy, get_bubbletree_kmeans, get_bubbletree_graph, get_gini, get_gini_k, get_num_tile, get_num_violins, d_500

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Examples

```
# input data
data("d_500", package = "scBubbletree")
A <- d_500$A
f <- d_500$f
b \leftarrow get_bubbletree_graph(x = A,
                           N_{eff} = 100
g_v <- get_cat_tiles(btd = b,</pre>
                      f = f,
                      integrate_vertical = TRUE,
                      round_digits = 2,
                      tile_text_size = 3,
                      x_axis_name = "Feature",
                      rotate_x_axis_labels = TRUE)
g_h <- get_cat_tiles(btd = b,</pre>
                     integrate_vertical = FALSE,
                     round_digits = 2,
                     tile_text_size = 3,
                     x_axis_name = "Feature",
                     rotate_x_axis_labels = TRUE)
b$tree|g_v$plot|g_h$plot
```

get_gini

Gini impurity index computed for a clustering solution and a vector of categorical cell feature labels

Description

How well is a set of categorical feature labels (e.g. cell type predictions) partitioned accross the different clusters of a clustering solution? We can assess this using the Gini impurity index (see details below).

Inputs are two equal-sized vectors:

- 1) clusters IDs
- 2) labels

Output:

- 1) cluster-specific purity -> Gini impurity (GI) index
- 2) clustering solution impurity -> Weighted Gini impurity (WGI) index

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Usage

```
get_gini(labels, clusters)
```

Arguments

labels character or numeric vector of labels
clusters character or numeric vector of cluster IDs

Details

To quantify the purity of a cluster (or bubble) i with n_i number of cells, each of which carries one of L possible labels (e.g. cell type), we computed the Gini impurity index:

$$GI_i = \sum_{j=1}^{L} \pi_{ij} (1 - \pi_{ij}),$$

with π_{ij} as the relative frequency of label j in cluster i. In homogeneous ('pure') clusters most cells carry a distinct label. Hence, the π 's are close to either 1 or 0, and GI takes on a small value close to zero. In 'impure' clusters cells carry a mixture of different labels. In this case most π are far from either 1 or 0, and GI diverges from 0 and approaches 1. If the relative frequencies of the different labels in cluster i are equal to the (background) relative frequencies of the labels in the sample, then cluster i is completely 'impure'.

To compute the overall Gini impurity of a bubbletree, which represents a clustering solution with k bubbles, we estimated the weighted Gini impurity (WGI) by computing the weighted (by the cluster size) average of the GIs:

$$WGI = \sum_{i=1}^{k} GI_i n_i / n,$$

with n_i as the number of cells in cluster i and $n = \sum_i n_i$.

Value

gi Gini impurity of each bubble

wgi Weighted Gini impurity index of the bubbletree

Author(s)

Simo Kitanovski <simo.kitanovski@uni-due.de>

See Also

get_k, get_r, get_bubbletree_kmeans, get_bubbletree_dummy, get_bubbletree_graph, get_gini_k, d_500

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get_gini_k	Gini impurity index computed for a list of clustering solutions obtained
	by functions get_k or get_r and a vector of categorical cell feature labels

Description

Given The Gini impurity (GI) index allows us to quantitatively evaluate how well a set of labels (categorical features) are split across a set of bubbles. We have a completely perfect split (GI = 0) when each bubble is 'pure', i.e. each bubble contains labels coming from distinct a class. In contrast to this, we have completely imperfect split (GI = 1) when the relative frequency distribution of the labels in each bubble is identical to the background relative frequency distribution of the labels.

Cell type predictions are a type of categorical features that are often used to evaluate the goodness of the clustering. get_gini_k takes as input: 1) a vector of labels for each cell (e.g. cell types) and 2) object returned by function get_k or get_r. Then it computes for each k or r the cluster purity and weightred gini impurity of each clustering solution mean GI, which is another way of finding an optimal clustering resolution.

Usage

```
get_gini_k(labels, obj)
```

Arguments

labels	character/factor vector of labels
obj	object returned by functions get_k or get_r

Details

To quantify the purity of a cluster (or bubble) i with n_i number of cells, each of which carries one of L possible labels (e.g. cell type), we computed the Gini impurity index:

$$GI_i = \sum_{j=1}^{L} \pi_{ij} (1 - \pi_{ij}),$$

with π_{ij} as the relative frequency of label j in cluster i. In homogeneous ('pure') clusters most cells carry a distinct label. Hence, the π 's are close to either 1 or 0, and GI takes on a small value close to zero. In 'impure' clusters cells carry a mixture of different labels. In this case most π are far from either 1 or 0, and GI diverges from 0 and approaches 1. If the relative frequencies of the different labels in cluster i are equal to the (background) relative frequencies of the labels in the sample, then cluster i is completely 'impure'.

To compute the overall Gini impurity of a bubbletree, which represents a clustering solution with k bubbles, we estimated the weighted Gini impurity (WGI) by computing the weighted (by the cluster size) average of the GIs:

$$WGI = \sum_{i=1}^{k} GI_i n_i / n,$$

with n_i as the number of cells in cluster i and $n = \sum_i n_i$.

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Value

```
 \begin{array}{ll} {\tt gi\_summary} & {\tt GI \ for \ each \ bubble \ of \ a \ clustering \ solution \ with \ clustering \ resolution \ k \ or \ r} \\ {\tt wgi\_summary} & {\tt WGI \ for \ each \ clustering \ solution \ with \ clustering \ resolution \ k \ or \ r} \\ \end{array}
```

Author(s)

Simo Kitanovski <simo.kitanovski@uni-due.de>

See Also

```
get_k, get_r, get_gini, get_bubbletree_kmeans, get_bubbletree_graph, get_bubbletree_dummy, d_500, get_num_tiles, get_num_violins, get_cat_tiles
```

Examples

```
# input data
data("d_500", package = "scBubbletree")
A <- d_500$A
f <- d_500 f
b_k \leftarrow get_k(x = A,
           ks = 1:5,
           B_gap = 5,
           n_start = 100,
           iter_max = 200,
            kmeans_algorithm = "MacQueen",
           cores = 1)
b_r \leftarrow get_r(x = A,
            rs = c(0.1, 0.5, 1),
            B_gap = 5,
            n_start = 20,
            iter_max = 100,
            algorithm = "original",
            cores = 1)
get_gini_k(labels = f, obj = b_k)
get_gini_k(labels = f, obj = b_r)
```

get_k

Finding optimal number k of clusters

Description

To perform k-means clustering we must specify a number k of clusters. Data-driven metrics, such as the Gap statistic or the within-cluster sum of squares (WCSS), can be used to infer appropriate k from the data. get_k computes the Gap statistic and WCSS for a number of clusters ks.

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Usage

Arguments

x numeric matrix A^{nxf} with n cells, and f low-dimensional projections

ks integer vector, k values to consider

B_gap integer, number of Monte Carlo ("bootstrap") samples taken when computing

the Gap statistic (see documentation of function clusGap, R-package cluster)

 $n_start, iter_max, kmeans_algorithm$

parameters for k-means clustering, see documentation of function k-means, R-

package stats

cores integer, number of PC cores for parallel execution

verbose logical, progress messages

Details

To compute the Gap statistic get_k adapts the algorithm of function clustGap from R-package cluster (version 2.1.3). For k-means clustering get_k uses the function kmeans implemented in R-package stats (version 4.2.0). For additional information see the respective documentations.

Value

```
boot_obj The results: k-means clustering solutions, the Gap statistic and WCSS gap_stats_summary, wcss_stats_summary
```

main results; Gap statistic and WCSS estimates. Means, standard errors and 95% confidence intervals are provided for each \boldsymbol{k}

gap_stats, wcss_stats

intermediate results; Gap statistic and WCSS estimates for each \boldsymbol{k} and bootstrap iteration \boldsymbol{b}

Author(s)

Simo Kitanovski <simo.kitanovski@uni-due.de>

See Also

get_r, get_bubbletree_dummy, get_bubbletree_graph, get_bubbletree_kmeans, get_gini, get_gini_k, d_500, get_num_tiles, get_num_violins, get_cat_tiles

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Examples

get_num_tiles

Visualization of numeric cell features using tile plots

Description

get_num_tiles creates tile plot to visualize a summary (e.g. mean, median or sum) of a numeric cell feature (e.g. gene expression of a specific gene) in each bubble of a bubbletree

Usage

Arguments

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```
rotate_x_axis_labels
```

logical, should the x-axis labels be shown horizontally (rotate_x_axis_labels = FALSE) or vertically (rotate_x_axis_labels = TRUE)

tile_bw logical, tile grayscale (tile_bw = TRUE) vs. color (tile_bw = FALSE, default)

Details

get_num_tiles uses two main inputs:

- 1. bubbletree object
- 2. numeric vector or matrix of numeric cell features.

The order of the cells used to generat the bubbletree (input 1.) should correspond to the order of cells in the vector/matrix of numeric cell features (input 2.)

This function computes summaries of numeric cell feature in each bubble: 1. mean = mean of feature 2. median = median of feature 3. sum = sum of feature 4. pct nonzero = sum of cells with feature > 0.5. pct zero = sum of cells with feature = 0

Important note: NA and NULL values are omitted.

Value

plot ggplot2, tile plot

table data.frame, raw data used to generate the plot

Author(s)

Simo Kitanovski <simo.kitanovski@uni-due.de>

See Also

get_k, get_r get_bubbletree_dummy, get_bubbletree_kmeans, get_bubbletree_graph, get_gini, get_gini_k, get_cat_tile, get_num_violins, d_500, d_ccl

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```
rotate_x_axis_labels = TRUE)
```

b\$tree|g\$plot

get_num_violins

Visualization of numeric cell features using violin plots

Description

get_num_violins creates violin plot to visualize the distribution of of numeric cell features (e.g. gene expressions) in each bubble of a bubbletree

Usage

Arguments

btd bubbletree object

fs numeric vector or matrix, numeric cell features

x_axis_name character, x-axis title

rotate_x_axis_labels

logical, should the x-axis labels be shown horizontally (rotate_x_axis_labels =

FALSE) or vertically (rotate_x_axis_labels = TRUE)

Details

get_num_violins uses two main inputs:

- 1. bubbletree object
- 2. numeric vector or matrix of numeric cell features.

The order of the cells used to generat the bubbletree (input 1.) should correspond to the order of cells in the vector/matrix of numeric cell features (input 2.)

This function visualizes densities of numeric cell feature in the different bubble.

Value

plot ggplot2, violin plot

table data.frame, raw data used to generate the plot

Author(s)

Simo Kitanovski <simo.kitanovski@uni-due.de>

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See Also

get_k, get_r get_bubbletree_dummy, get_bubbletree_kmeans, get_bubbletree_graph, get_gini, get_gini_k, get_cat_tile, get_num_tiles, d_500

Examples

get_r

Finding optimal clustering resulution r and number of communities k'

Description

To perform Louvain clustering we must specify a clustering resulution r. Data-driven metrics, such as the Gap statistic or the within-cluster sum of squares (WCSS) can be used to infer appropriate r from the data. get_r computes the Gap statistic and WCSS for a vector of clustering resolutions rs.

Usage

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Arguments

	X	numeric matrix A^{nxf} with n cells, and f low-dimensional projections
	rs	number vector, r values to consider
	B_gap	integer, number of Monte Carlo ("bootstrap") samples taken when computing the Gap statistic (see documentation of function clusGap, R-package cluster)
n_start, iter_max		
		parameters for Louvain clustering, see documentation of function FindClusters, R-package Seurat
	algorithm	character, four clustering algorithms: 'original', 'LMR', 'SLM' and 'Leiden', see documentation of function FindClusters, R-package Seurat
	knn_k	integer, defines k for the k -nearest neighbor algorithm, see documentation of function FindClusters, R -package Seurat
	cores	integer, number of PC cores for parallel execution
	verbose	logical, progress messages

Details

To compute the Gap statistic get_r adapts the algorithm of function clustGap from R-package cluster (version 2.1.3). For Louvain clustering get_r uses the function FindClusters implemented in Rpackage Seurat. For additional information see the respective documentations.

Value

```
boot_obj
                 The results: k-means clustering solutions, the Gap statistic and WCSS
gap_stats_summary, wcss_stats_summary
                 main results; Gap statistic and WCSS estimates. Means, standard errors and
                  95% confidence intervals are provided for each r and k'
gap_stats, wcss_stats
                  intermediate results; Gap statistic and WCSS estimates for each r and k' and
                  bootstrap iteration b
```

Author(s)

Simo Kitanovski <simo.kitanovski@uni-due.de>

See Also

get_k, get_bubbletree_dummy, get_bubbletree_graph, get_bubbletree_kmeans, get_gini, get_gini_k, d_500, get_num_tiles, get_num_violins, get_cat_tiles, d_ccl

```
# input data
data("d_500", package = "scBubbletree")
A <- d_500$A
```

get_r

b\$gap_stats_summary

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