Package 'lmds'

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| Title Landmark Multi-Dimensional Scaling |
| Version 0.1.0 |
| Description A fast dimensionality reduction method scaleable to large numbers of samples. Landmark Multi-Dimensional Scaling (LMDS) is an extension of classical Torgerson MDS, but rather than calculating a complete distance matrix between all pairs of samples, only the distances between a set of landmarks and the samples are calculated. |
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| <pre>BugReports https://github.com/dynverse/lmds/issues</pre> |
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| R topics documented: |
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| cmdscale_landmarks | Perform MDS on landmarks and project other samples to the same space |
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Description

Perform MDS on landmarks and project other samples to the same space

Usage

```
cmdscale_landmarks(dist_2lm, ndim = 3, rescale = TRUE, ...)
```

Arguments

| dist_2lm | Distance matrix between the landmarks and all the samples in original dataset |
|----------|---|
| ndim | The number of dimensions |
| rescale | Whether or not to rescale the final dimensionality reduction (recommended) |
| | Extra params to pass to irlba::irlba() |

Value

The dimensionality reduction in the form of a ncol(dist_2lm) by ndim matrix.

Examples

```
library(Matrix)
x <- as.matrix(iris[,1:4])
dist_2lm <- select_landmarks(x)
cmdscale_landmarks(dist_2lm)</pre>
```

lmds Landmark MDS

Description

A fast dimensionality reduction method scaleable to large numbers of samples. Landmark Multi-Dimensional Scaling (LMDS) is an extension of classical 'Torgerson MDS', but rather than calculating a complete distance matrix between all pairs of samples, only the distances between a set of landmarks and the samples are calculated.

A fast dimensionality reduction method scaleable to large numbers of samples. Landmark Multi-Dimensional Scaling (LMDS) is an extension of classical Torgerson MDS′, but rather than calculating a complete distance matrix between all pairs of samples, only the distances between a set of landmarks and the samples are calculated.

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Usage

```
lmds(x, ndim = 3, distance_method = c("euclidean", "pearson",
    "spearman", "cosine", "manhattan"), landmark_method = c("sample"),
    num_landmarks = 500)
```

Arguments

x A matrix, optionally sparse.

ndim The number of dimensions
distance_method The distance metric to use. Options are "euclidean" (default), "pearson", "spearman", "cosine", "manhattan".

landmark_method The landmark selection method to use. Options are "sample" (default).

num_landmarks The number of landmarks to use,

Value

The dimensionality reduction in the form of a nrow(x) by ndim matrix.

Examples

```
library(Matrix)
x <- Matrix::rsparsematrix(1000, 1000, .01)
lmds(x, ndim = 3)</pre>
```

select_landmarks

Select landmarks from dataset

Description

In addition, the distances between the landmarks and all samples are calculated.

Usage

```
select_landmarks(x, distance_method = c("euclidean", "pearson",
   "spearman", "cosine", "manhattan"), landmark_method = c("sample"),
   num_landmarks = 500)
```

Arguments

x A matrix, optionally sparse.

distance_method

The distance metric to use. Options are "euclidean" (default), "pearson", "spearman", "cosine", "manhattan".

landmark_method

The landmark selection method to use. Options are "sample" (default).

num_landmarks

The number of landmarks to use,

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Value

The distance matrix between the landmarks and all samples. In addition, an attribute "landmark_ix" denotes the indices of landmarks that were sampled.

Examples

```
library(Matrix)
x <- Matrix::rsparsematrix(1000, 1000, .01)
select_landmarks(x)</pre>
```

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