

Package: rsparse (via r-universe)

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Type Package

Title Statistical Learning on Sparse Matrices

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Description Implements many algorithms for statistical learning on sparse matrices - matrix factorizations, matrix completion, elastic net regressions, factorization machines. Also 'rsparse' enhances 'Matrix' package by providing methods for multithreaded <sparse, dense> matrix products and native slicing of the sparse matrices in Compressed Sparse Row (CSR) format. List of the algorithms for regression problems: 1) Elastic Net regression via Follow The Proximally-Regularized Leader (FTRL) Stochastic Gradient Descent (SGD), as per McMahan et al, (<doi:10.1145/2487575.2488200>) 2) Factorization Machines via SGD, as per Rendle (2010, <doi:10.1109/ICDM.2010.127>) List of algorithms for matrix factorization and matrix completion: 1) Weighted Regularized Matrix Factorization (WRMF) via Alternating Least Squares (ALS) - paper by Hu, Koren, Volinsky (2008, <doi:10.1109/ICDM.2008.22>) 2) Maximum-Margin Matrix Factorization via ALS, paper by Rennie, Srebro (2005, <doi:10.1145/1102351.1102441>) 3) Fast Truncated Singular Value Decomposition (SVD), Soft-Thresholded SVD, Soft-Impute matrix completion via ALS - paper by Hastie, Mazumder et al. (2014, <doi:10.48550/arXiv.1410.2596>) 4) Linear-Flow matrix factorization, from 'Practical linear models for large-scale one-class collaborative filtering' by Sedhain, Bui, Kawale et al (2016, ISBN:978-1-57735-770-4) 5) GlobalVectors (GloVe) matrix factorization via SGD, paper by Pennington, Socher, Manning (2014, <<https://aclanthology.org/D14-1162/>>) Package is reasonably fast and memory efficient - it allows to work with large datasets - millions of rows and millions of columns. This is particularly useful for practitioners working on recommender systems.

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ByteCompile true

Depends R (>= 3.6.0), methods, Matrix (>= 1.3)

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Suggests testthat, covr

StagedInstall TRUE

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BugReports <https://github.com/dselivanov/rsparse/issues>

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detect_number_omp_threads

Detects number of OpenMP threads in the system

Description

Detects number of OpenMP threads in the system respecting environment variables such as OMP_NUM_THREADS and OMP_THREAD_LIMIT

Usage

detect_number_omp_threads()

FactorizationMachine *Second order Factorization Machines*

Description

Creates second order Factorization Machines model

Methods

Public methods:

- [FactorizationMachine\\$new\(\)](#)
- [FactorizationMachine\\$partial_fit\(\)](#)
- [FactorizationMachine\\$fit\(\)](#)
- [FactorizationMachine\\$predict\(\)](#)
- [FactorizationMachine\\$clone\(\)](#)

Method new(): creates Creates second order Factorization Machines model

Usage:

```
FactorizationMachine$new(
  learning_rate_w = 0.2,
  rank = 4,
  lambda_w = 0,
  lambda_v = 0,
  family = c("binomial", "gaussian"),
  intercept = TRUE,
  learning_rate_v = learning_rate_w
)
```

Arguments:

learning_rate_w learning rate for features intercatations

rank dimension of the latent dimensions which models features interactions

lambda_w regularization for features interactions
 lambda_v regularization for features
 family one of "binomial", "gaussian"
 intercept logical, indicates whether or not include intercept to the model
 learning_rate_v learning rate for features

Method partial_fit(): fits/updates model

Usage:

FactorizationMachine\$partial_fit(x, y, weights = rep(1, length(y)), ...)

Arguments:

x input sparse matrix. Native format is Matrix::RsparseMatrix. If x is in different format, model will try to convert it to RsparseMatrix with as(x, "RsparseMatrix"). Dimensions should be (n_samples, n_features)
 y vector of targets
 weights numeric vector of length 'n_samples'. Defines how to amplify SGD updates for each sample. May be useful for highly unbalanced problems.
 ... not used at the moment

Method fit(): shorthand for applying 'partial_fit' 'n_iter' times

Usage:

FactorizationMachine\$fit(x, y, weights = rep(1, length(y)), n_iter = 1L, ...)

Arguments:

x input sparse matrix. Native format is Matrix::RsparseMatrix. If x is in different format, model will try to convert it to RsparseMatrix with as(x, "RsparseMatrix"). Dimensions should be (n_samples, n_features)
 y vector of targets
 weights numeric vector of length 'n_samples'. Defines how to amplify SGD updates for each sample. May be useful for highly unbalanced problems.
 n_iter number of SGD epochs
 ... not used at the moment

Method predict(): makes predictions based on fitted model

Usage:

FactorizationMachine\$predict(x, ...)

Arguments:

x input sparse matrix of shape (n_samples, n_features)
 ... not used at the moment

Method clone(): The objects of this class are cloneable with this method.

Usage:

FactorizationMachine\$clone(deep = FALSE)

Arguments:

deep Whether to make a deep clone.

Examples

```
# Factorization Machines can fit XOR function!
x = rbind(
  c(0, 0),
  c(0, 1),
  c(1, 0),
  c(1, 1)
)
y = c(0, 1, 1, 0)

x = as(x, "RsparseMatrix")
fm = FactorizationMachine$new(learning_rate_w = 10, rank = 2, lambda_w = 0,
  lambda_v = 0, family = 'binomial', intercept = TRUE)
res = fm$fit(x, y, n_iter = 100)
preds = fm$predict(x)
all(preds[c(1, 4)] < 0.01)
all(preds[c(2, 3)] > 0.99)
```

FTRL

Logistic regression model with FTRL proximal SGD solver.

Description

Creates 'Follow the Regularized Leader' model. Only logistic regression implemented at the moment.

Methods

Public methods:

- [FTRL\\$new\(\)](#)
- [FTRL\\$partial_fit\(\)](#)
- [FTRL\\$fit\(\)](#)
- [FTRL\\$predict\(\)](#)
- [FTRL\\$coef\(\)](#)
- [FTRL\\$clone\(\)](#)

Method `new()`: creates a model

Usage:

```
FTRL$new(
  learning_rate = 0.1,
  learning_rate_decay = 0.5,
  lambda = 0,
  l1_ratio = 1,
  dropout = 0,
  family = c("binomial")
)
```

Arguments:

learning_rate learning rate

learning_rate_decay learning rate which controls decay. Please refer to FTRL proximal paper for details. Usually convergence does not heavily depend on this parameter, so default value 0.5 is safe.

lambda regularization parameter

l1_ratio controls L1 vs L2 penalty mixing. 1 = Lasso regression, 0 = Ridge regression. Elastic net is in between

dropout dropout - percentage of random features to exclude from each sample. Acts as regularization.

family a description of the error distribution and link function to be used in the model. Only binomial (logistic regression) is implemented at the moment.

Method partial_fit(): fits model to the data

Usage:

```
FTRL$partial_fit(x, y, weights = rep(1, length(y)), ...)
```

Arguments:

x input sparse matrix. Native format is Matrix::RsparseMatrix. If x is in different format, model will try to convert it to RsparseMatrix with as(x, "RsparseMatrix"). Dimensions should be (n_samples, n_features)

y vector of targets

weights numeric vector of length 'n_samples'. Defines how to amplify SGD updates for each sample. May be useful for highly unbalanced problems.

... not used at the moment

Method fit(): shorthand for applying 'partial_fit' 'n_iter' times

Usage:

```
FTRL$fit(x, y, weights = rep(1, length(y)), n_iter = 1L, ...)
```

Arguments:

x input sparse matrix. Native format is Matrix::RsparseMatrix. If x is in different format, model will try to convert it to RsparseMatrix with as(x, "RsparseMatrix"). Dimensions should be (n_samples, n_features)

y vector of targets

weights numeric vector of length 'n_samples'. Defines how to amplify SGD updates for each sample. May be useful for highly unbalanced problems.

n_iter number of SGD epochs

... not used at the moment

Method predict(): makes predictions based on fitted model

Usage:

```
FTRL$predict(x, ...)
```

Arguments:

x input matrix

... not used at the moment

Method `coef()`: returns coefficients of the regression model

Usage:

```
FTRL$coef()
```

Method `clone()`: The objects of this class are cloneable with this method.

Usage:

```
FTRL$clone(deep = FALSE)
```

Arguments:

`deep` Whether to make a deep clone.

Examples

```
library(rsparse)
library(Matrix)
i = sample(1000, 1000 * 100, TRUE)
j = sample(1000, 1000 * 100, TRUE)
y = sample(c(0, 1), 1000, TRUE)
x = sample(c(-1, 1), 1000 * 100, TRUE)
odd = seq(1, 99, 2)
x[i %in% which(y == 1) & j %in% odd] = 1
x = sparseMatrix(i = i, j = j, x = x, dims = c(1000, 1000), repr="R")

ftrl = FTRL$new(learning_rate = 0.01, learning_rate_decay = 0.1,
lambda = 10, l1_ratio = 1, dropout = 0)
ftrl$partial_fit(x, y)

w = ftrl$coef()
head(w)
sum(w != 0)
p = ftrl$predict(x)
```

GloVe

Global Vectors

Description

Creates Global Vectors matrix factorization model

Public fields

`components` represents context embeddings

`bias_i` bias term *i* as per paper

`bias_j` bias term *j* as per paper

`shuffle` logical = FALSE by default. Whether to perform shuffling before each SGD iteration. Generally shuffling is a good practice for SGD.

Methods

Public methods:

- `GloVe$new()`
- `GloVe$fit_transform()`
- `GloVe$get_history()`
- `GloVe$clone()`

Method `new()`: Creates GloVe model object

Usage:

```
GloVe$new(
  rank,
  x_max,
  learning_rate = 0.15,
  alpha = 0.75,
  lambda = 0,
  shuffle = FALSE,
  init = list(w_i = NULL, b_i = NULL, w_j = NULL, b_j = NULL)
)
```

Arguments:

`rank` desired dimension for the latent vectors

`x_max` integer maximum number of co-occurrences to use in the weighting function

`learning_rate` numeric learning rate for SGD. I do not recommend that you modify this parameter, since AdaGrad will quickly adjust it to optimal

`alpha` numeric = 0.75 the alpha in weighting function formula: $f(x) = 1 \text{ if } x > x_{max}; \text{ else } (x/x_{max})^{\alpha}$

`lambda` numeric = 0.0 regularization parameter

`shuffle` see shuffle field

`init` list(`w_i` = NULL, `b_i` = NULL, `w_j` = NULL, `b_j` = NULL) initialization for embeddings (`w_i`, `w_j`) and biases (`b_i`, `b_j`). `w_i`, `w_j` - numeric matrices, should have #rows = rank, #columns = expected number of rows (`w_i`) / columns(`w_j`) in the input matrix. `b_i`, `b_j` = numeric vectors, should have length of #expected number of rows(`b_i`) / columns(`b_j`) in input matrix

Method `fit_transform()`: fits model and returns embeddings

Usage:

```
GloVe$fit_transform(
  x,
  n_iter = 10L,
  convergence_tol = -1,
  n_threads = getOption("rsparse_omp_threads", 1L),
  ...
)
```

Arguments:

`x` An input term co-occurrence matrix. Preferably in `dgTMatrix` format

`n_iter` integer number of SGD iterations

convergence_tol numeric = -1 defines early stopping strategy. Stop fitting when one of two following conditions will be satisfied: (a) passed all iterations (b) $\text{cost_previous_iter} / \text{cost_current_iter} - 1 < \text{convergence_tol}$.

n_threads number of threads to use

... not used at the moment

Method get_history(): returns value of the loss function for each epoch

Usage:

GloVe\$get_history()

Method clone(): The objects of this class are cloneable with this method.

Usage:

GloVe\$clone(deep = FALSE)

Arguments:

deep Whether to make a deep clone.

References

<http://nlp.stanford.edu/projects/glove/>

Examples

```
data('movielens100k')
co_occurrence = crossprod(movielens100k)
glove_model = GloVe$new(rank = 4, x_max = 10, learning_rate = .25)
embeddings = glove_model$fit_transform(co_occurrence, n_iter = 2, n_threads = 1)
embeddings = embeddings + t(glove_model$components) # embeddings + context embeddings
identical(dim(embeddings), c(ncol(movielens100k), 10L))
```

LinearFlow

Linear-Flow model for one-class collaborative filtering

Description

Creates *Linear-Flow* model described in [Practical Linear Models for Large-Scale One-Class Collaborative Filtering](#). The goal is to find item-item (or user-user) similarity matrix which is **low-rank and has small Frobenius norm**. Such double regularization allows to better control the generalization error of the model. Idea of the method is somewhat similar to **Sparse Linear Methods(SLIM)** but scales to large datasets much better.

Super class

`rsparse::MatrixFactorizationRecommender` -> LinearFlow

Public fields

v right singular vector of the user-item matrix. Size is n_items * rank. In the paper this matrix is called v

Methods

Public methods:

- `LinearFlow$new()`
- `LinearFlow$fit_transform()`
- `LinearFlow$transform()`
- `LinearFlow$cross_validate_lambda()`
- `LinearFlow$clone()`

Method `new()`: creates Linear-Flow model with rank latent factors.

Usage:

```
LinearFlow$new(
  rank = 8L,
  lambda = 0,
  init = NULL,
  preprocess = identity,
  solve_right_singular_vectors = c("soft_impute", "svd")
)
```

Arguments:

`rank` size of the latent dimension

`lambda` regularization parameter

`init` initialization of the orthogonal basis.

`preprocess` `identity()` by default. User specified function which will be applied to user-item interaction matrix before running matrix factorization (also applied during inference time before making predictions). For example we may want to normalize each row of user-item matrix to have 1 norm. Or apply `log1p()` to discount large counts.

`solve_right_singular_vectors` type of the solver for initialization of the orthogonal basis. Original paper uses SVD. See paper for details.

Method `fit_transform()`: performs matrix factorization

Usage:

```
LinearFlow$fit_transform(x, ...)
```

Arguments:

`x` input matrix

`...` not used at the moment

Method `transform()`: calculates user embeddings for the new input

Usage:

```
LinearFlow$transform(x, ...)
```

Arguments:

`x` input matrix

`...` not used at the moment

Method `cross_validate_lambda()`: performs fast tuning of the parameter 'lambda' with warm re-starts

Usage:

```
LinearFlow$cross_validate_lambda(
  x,
  x_train,
  x_test,
  lambda = "auto@10",
  metric = "map@10",
  not_recommend = x_train,
  ...
)
```

Arguments:

`x` input user-item interactions matrix. Model performs matrix factorization based only on this matrix

`x_train` user-item interactions matrix. Model recommends items based on this matrix. Usually should be different from 'x' to avoid overfitting

`x_test` target user-item interactions. Model will evaluate predictions against this matrix, 'x_test' should be treated as future interactions.

`lambda` numeric vector - sequence of regularization parameters. Supports special value like 'auto@10'. This will automatically find a sequence of lambda of length 10. This is recommended way to check for 'lambda'.

`metric` a metric against which model will be evaluated for top-k recommendations. Currently only map@k and ndcg@k are supported (k can be any integer)

`not_recommend` matrix same shape as 'x_train'. Specifies which items to not recommend for each user.

... not used at the moment

Method `clone()`: The objects of this class are cloneable with this method.

Usage:

```
LinearFlow$clone(deep = FALSE)
```

Arguments:

`deep` Whether to make a deep clone.

References

- <http://www.bkveton.com/docs/ijcai2016.pdf>

Examples

```
data("movielens100k")
train = movielens100k[1:900, ]
cv = movielens100k[901:nrow(movielens100k), ]
model = LinearFlow$new(
  rank = 10, lambda = 0,
  solve_right_singular_vectors = "svd"
)
user_emb = model$fit_transform(train)
preds = model$predict(cv, k = 10)
```

Description

ap_k calculates **Average Precision at K** (ap@k). Please refer to [Information retrieval wikipedia article](#)

ndcg_k() calculates **Normalized Discounted Cumulative Gain at K** (ndcg@k). Please refer to [Discounted cumulative gain](#)

Usage

```
ap_k(predictions, actual, ...)
```

```
ndcg_k(predictions, actual, ...)
```

Arguments

predictions	matrix of predictions. Predctions can be defined 2 ways: <ol style="list-style-type: none"> 1. predictions = integer matrix with item indices (correspond to column numbers in actual) 2. predictions = character matrix with item identifiers (characters which correspond to colnames(actual)) which has attribute "indices" (integer matrix with item indices which correspond to column numbers in actual).
actual	sparse Matrix of relevant items. Each non-zero entry considered as relevant item. Value of the each non-zero entry considered as relevance for calculation of ndcg@k. It should inherit from Matrix::sparseMatrix. Internally Matrix::RsparseMatrix is used.
...	other arguments (not used at the moment)

Examples

```
predictions = matrix(
  c(5L, 7L, 9L, 2L),
  nrow = 1
)
actual = matrix(
  c(0, 0, 0, 0, 1, 0, 1, 0, 1, 0),
  nrow = 1
)
actual = as(actual, "RsparseMatrix")
identical(rsparse::ap_k(predictions, actual), 1)
```

`movielens100k`*MovieLens 100K Dataset*

Description

This data set consists of:

1. 100,000 ratings (1-5) from 943 users on 1682 movies.
2. Each user has rated at least 20 movies.

MovieLens data sets were collected by the GroupLens Research Project at the University of Minnesota.

Usage

```
data("movielens100k")
```

Format

A sparse column-compressed matrix (`Matrix::dgCMatrix`) with 943 rows and 1682 columns.

1. rows are users
2. columns are movies
3. values are ratings

Source

<https://en.wikipedia.org/wiki/MovieLens#Datasets>

`PureSVD`*PureSVD recommender model decomposition*

Description

Creates PureSVD recommender model. Solver is based on Soft-SVD which is very similar to truncated SVD but optionally adds regularization based on nuclear norm.

Super class

```
rsparse::MatrixFactorizationRecommender -> PureSVD
```

Methods

Public methods:

- `PureSVD$new()`
- `PureSVD$fit_transform()`
- `PureSVD$transform()`
- `PureSVD$clone()`

Method `new()`: create PureSVD model

Usage:

```
PureSVD$new(
  rank = 10L,
  lambda = 0,
  init = NULL,
  preprocess = identity,
  method = c("svd", "impute"),
  ...
)
```

Arguments:

`rank` size of the latent dimension

`lambda` regularization parameter

`init` initialization of item embeddings

`preprocess` `identity()` by default. User specified function which will be applied to user-item interaction matrix before running matrix factorization (also applied during inference time before making predictions). For example we may want to normalize each row of user-item matrix to have 1 norm. Or apply `log1p()` to discount large counts.

`method` type of the solver for initialization of the orthogonal basis. Original paper uses SVD. See paper for details.

... not used at the moment

Method `fit_transform()`: performs matrix factorization

Usage:

```
PureSVD$fit_transform(x, n_iter = 100L, convergence_tol = 0.001, ...)
```

Arguments:

`x` input sparse user-item matrix(of class `dgCMatrix`)

`n_iter` maximum number of iterations

`convergence_tol` `numeric = -Inf` defines early stopping strategy. Stops fitting when one of two following conditions will be satisfied: (a) passed all iterations (b) relative change of Frobenious norm of the two consequent solution is less then provided `convergence_tol`.

... not used at the moment

Method `transform()`: calculates user embeddings for the new input

Usage:

```
PureSVD$transform(x, ...)
```

Arguments:

x input matrix
 ... not used at the moment

Method clone(): The objects of this class are cloneable with this method.

Usage:

```
PureSVD$clone(deep = FALSE)
```

Arguments:

deep Whether to make a deep clone.

Examples

```
data('movielens100k')
i_train = sample(nrow(movielens100k), 900)
i_test = setdiff(seq_len(nrow(movielens100k)), i_train)
train = movielens100k[i_train, ]
test = movielens100k[i_test, ]
rank = 32
lambda = 0
model = PureSVD$new(rank = rank, lambda = lambda)
user_emb = model$fit_transform(sign(test), n_iter = 100, convergence_tol = 0.00001)
item_emb = model$components
preds = model$predict(sign(test), k = 1500, not_recommend = NULL)
mean(ap_k(preds, actual = test))
```

ScaleNormalize

Re-scales input matrix proportionally to item popularity

Description

scales input user-item interaction matrix as per eq (16) from the paper. Usage of such rescaled matrix with [PureSVD] model will be equal to running PureSVD on the scaled cosine-based inter-item similarity matrix.

Public fields

norm which norm model should make equal to one
 scale how to rescale norm vector

Methods**Public methods:**

- [ScaleNormalize\\$new\(\)](#)
- [ScaleNormalize\\$fit\(\)](#)
- [ScaleNormalize\\$transform\(\)](#)
- [ScaleNormalize\\$fit_transform\(\)](#)

- [ScaleNormalize\\$clone\(\)](#)

Method new(): creates model

Usage:

```
ScaleNormalize$new(scale = 0.5, norm = 2, target = c("rows", "columns"))
```

Arguments:

scale numeric, how to rescale norm vector

norm numeric, which norm model should make equal to one

target character, defines whether rows or columns should be rescaled

Method fit(): fits the modes

Usage:

```
ScaleNormalize$fit(x)
```

Arguments:

x input sparse matrix

Method transform(): transforms new matrix

Usage:

```
ScaleNormalize$transform(x)
```

Arguments:

x input sparse matrix

Method fit_transform(): fits the model and transforms input

Usage:

```
ScaleNormalize$fit_transform(x)
```

Arguments:

x input sparse matrix

Method clone(): The objects of this class are cloneable with this method.

Usage:

```
ScaleNormalize$clone(deep = FALSE)
```

Arguments:

deep Whether to make a deep clone.

References

See [EigenRec: Generalizing PureSVD for Effective and Efficient Top-N Recommendations](#) for details.

`soft_impute`*SoftImpute/SoftSVD matrix factorization*

Description

Fit SoftImpute/SoftSVD via fast alternating least squares. Based on the paper by Trevor Hastie, Rahul Mazumder, Jason D. Lee, Reza Zadeh by "Matrix Completion and Low-Rank SVD via Fast Alternating Least Squares" - <http://arxiv.org/pdf/1410.2596>

Usage

```
soft_impute(  
  x,  
  rank = 10L,  
  lambda = 0,  
  n_iter = 100L,  
  convergence_tol = 0.001,  
  init = NULL,  
  final_svd = TRUE  
)
```

```
soft_svd(  
  x,  
  rank = 10L,  
  lambda = 0,  
  n_iter = 100L,  
  convergence_tol = 0.001,  
  init = NULL,  
  final_svd = TRUE  
)
```

Arguments

<code>x</code>	sparse matrix. Both CSR <code>dgRMatrix</code> and CSC <code>dgCMatrix</code> are supported. CSR matrix is preferred because in this case algorithm will benefit from multithreaded CSR * dense matrix products (if OpenMP is supported on your platform). On many-cores machines this reduces fitting time significantly.
<code>rank</code>	maximum rank of the low-rank solution.
<code>lambda</code>	regularization parameter for the nuclear norm
<code>n_iter</code>	maximum number of iterations of the algorithms
<code>convergence_tol</code>	convergence tolerance. Internally functions keeps track of the relative change of the Frobenious norm of the two consequent iterations. If the change is less than <code>convergence_tol</code> then the process is considered as converged and function returns result.

<code>init</code>	<code>svd</code> like object with <code>u</code> , <code>v</code> , <code>d</code> components to initialize algorithm. Algorithm benefit from warm starts. <code>init</code> could be rank up rank of the maximum allowed rank. If <code>init</code> has rank less than max rank it will be padded automatically.
<code>final_svd</code>	logical whether need to make final preprocessing with SVD. This is not necessary but cleans up rank nicely - highly recommended to leave it TRUE.

Value

`svd`-like object - `list(u, v, d)`. `u`, `v`, `d` components represent left, right singular vectors and singular values.

Examples

```
set.seed(42)
data('movielens100k')
k = 10
seq_k = seq_len(k)
m = movielens100k[1:100, 1:200]
svd_ground_true = svd(m)
svd_soft_svd = soft_svd(m, rank = k, n_iter = 100, convergence_tol = 1e-6)
m_restored_svd = svd_ground_true$u[, seq_k] %*%
  diag(x = svd_ground_true$d[seq_k]) %*%
  t(svd_ground_true$v[, seq_k])
m_restored_soft_svd = svd_soft_svd$u %*%
  diag(x = svd_soft_svd$d) %*%
  t(svd_soft_svd$v)
all.equal(m_restored_svd, m_restored_soft_svd, tolerance = 1e-1)
```

WRMF

*Weighted Regularized Matrix Factorization for collaborative filtering***Description**

Creates a matrix factorization model which is solved through Alternating Least Squares (Weighted ALS for implicit feedback). For implicit feedback see "Collaborative Filtering for Implicit Feedback Datasets" (Hu, Koren, Volinsky). For explicit feedback it corresponds to the classic model for rating matrix decomposition with MSE error. These two algorithms are proven to work well in recommender systems.

Super class

`rsparse::MatrixFactorizationRecommender` -> WRMF

Methods**Public methods:**

- `WRMF$new()`
- `WRMF$fit_transform()`

- `WRMF$transform()`
- `WRMF$clone()`

Method `new()`: creates WRMF model

Usage:

```
WRMF$new(
  rank = 10L,
  lambda = 0,
  dynamic_lambda = TRUE,
  init = NULL,
  preprocess = identity,
  feedback = c("implicit", "explicit"),
  solver = c("conjugate_gradient", "cholesky", "nnls"),
  with_user_item_bias = FALSE,
  with_global_bias = FALSE,
  cg_steps = 3L,
  precision = c("double", "float"),
  ...
)
```

Arguments:

`rank` size of the latent dimension

`lambda` regularization parameter

`dynamic_lambda` whether 'lambda' is to be scaled according to the number

`init` initialization of item embeddings

`preprocess` `identity()` by default. User specified function which will be applied to user-item interaction matrix before running matrix factorization (also applied during inference time before making predictions). For example we may want to normalize each row of user-item matrix to have 1 norm. Or apply `log1p()` to discount large counts. This corresponds to the "confidence" function from "Collaborative Filtering for Implicit Feedback Datasets" paper. Note that it will not automatically add +1 to the weights of the positive entries.

`feedback` character - feedback type - one of `c("implicit", "explicit")`

`solver` character - solver name. One of `c("conjugate_gradient", "cholesky", "nnls")`. Usually approximate "conjugate_gradient" is significantly faster and solution is on par with "cholesky". "nnls" performs non-negative matrix factorization (NNMF) - restricts user and item embeddings to be non-negative.

`with_user_item_bias` bool controls if model should calculate user and item biases. At the moment only implemented for "explicit" feedback.

`with_global_bias` bool controls if model should calculate global biases (mean). At the moment only implemented for "explicit" feedback.

`cg_steps` integer > 0 - max number of internal steps in conjugate gradient (if "conjugate_gradient" solver used). `cg_steps = 3` by default. Controls precision of linear equation solution at the each ALS step. Usually no need to tune this parameter

`precision` one of `c("double", "float")`. Should embedding matrices be numeric or float (from float package). The latter is usually 2x faster and consumes less RAM. BUT float matrices are not "base" objects. Use carefully.

... not used at the moment

Method `fit_transform()`: fits the model

Usage:

```
WRMF$fit_transform(
  x,
  n_iter = 10L,
  convergence_tol = ifelse(private$feedback == "implicit", 0.005, 0.001),
  ...
)
```

Arguments:

`x` input matrix (preferably matrix in CSC format -‘CsparseMatrix’
`n_iter` max number of ALS iterations
`convergence_tol` convergence tolerance checked between iterations
 ... not used at the moment

Method `transform()`: create user embeddings for new input

Usage:

```
WRMF$transform(x, ...)
```

Arguments:

`x` user-item interaction matrix (preferrably as ‘dgRMatrix’)
 ... not used at the moment

Method `clone()`: The objects of this class are cloneable with this method.

Usage:

```
WRMF$clone(deep = FALSE)
```

Arguments:

`deep` Whether to make a deep clone.

References

- Hu, Yifan, Yehuda Koren, and Chris Volinsky. "Collaborative filtering for implicit feedback datasets." 2008 Eighth IEEE International Conference on Data Mining. Ieee, 2008.
- <https://math.stackexchange.com/questions/1072451/analytic-solution-for-matrix-factorization-us-1073170#1073170>
- <http://activisiongamescience.github.io/2016/01/11/Implicit-Recommender-Systems-Biased-Matrix-F>
- <https://jessesw.com/Rec-System/>
- <http://www.benfrederickson.com/matrix-factorization/>
- <http://www.benfrederickson.com/fast-implicit-matrix-factorization/>
- Franc, Vojtech, Vaclav Hlavac, and Mirko Navara. "Sequential coordinate-wise algorithm for the non-negative least squares problem." International Conference on Computer Analysis of Images and Patterns. Springer, Berlin, Heidelberg, 2005.
- Zhou, Yunhong, et al. "Large-scale parallel collaborative filtering for the netflix prize." International conference on algorithmic applications in management. Springer, Berlin, Heidelberg, 2008.

Examples

```
data('movielens100k')
train = movielens100k[1:900, ]
cv = movielens100k[901:nrow(movielens100k), ]
model = WRMF$new(rank = 5, lambda = 0, feedback = 'implicit')
user_emb = model$fit_transform(train, n_iter = 5, convergence_tol = -1)
item_emb = model$components
preds = model$predict(cv, k = 10, not_recommnd = cv)
```

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