# **Importing Libraries**

## In [2]:

```
### Run this cell to install surprise library that we will use in exercise 3
# ! pip install scikit-surprise
```

## In [1]:

```
import matplotlib.pyplot as plt
from scipy import sparse
from sklearn.model_selection import train_test_split
from sklearn.decomposition import NMF
import numpy as np
import pandas as pd
import math
from surprise import SVD
from surprise import Dataset
from surprise import accuracy
from surprise.model_selection import train_test_split
from surprise.model_selection import GridSearchCV
from surprise.model_selection import cross_validate
```

### In [2]:

```
import warnings
warnings.filterwarnings("ignore")
```

# **Exercise 2**

In this task, matrix factorization has been implemented using Stochastic Gradient Descent.

# **Utility Functions**

```
In [17]:
```

```
def read_data():
    movieLdf = pd.read_csv("u.data", delimiter="\t", header=None)
    movieLdf = movieLdf.drop(movieLdf.columns[3], axis=1)
    return movieLdf
```

```
In [18]:
```

```
def lossFn(data, p, q, lamda):
    loss = 0
    for x in data:
        a = (x[2]-predict(p,q,x[0]-1,x[1]-1))*(x[2]-predict(p,q,x[0]-1,x[1]-1))
        b = lamda * (np.dot(p[x[0]-1].T, p[x[0]-1]) + np.dot(q[x[1]-1].T, q[x[1]-1]))
        loss = loss + ( a + b )
    return loss
```

```
In [19]:
```

```
def predict(p, q, user, item):
    return np.dot(p[user].T, q[item])
```

Following function has been created for Stochastic gradient descent algorithm (Algorithm LearnLatentFactors) as described on slide 29:

## In [20]:

```
def trainMF(data, valid, nUsers, nItems, learnR, lamda, factors, epochs, eps):
   p = np.random.normal(size=((nUsers, factors)))
   q = np.random.normal(size=((nItems, factors)))
   trainRMSE=[]
   validRMSE=[]
   for i in range(epochs):
        np.random.shuffle(data)
        prevFnV = lossFn(data, p, q, lamda)
        for x in data:
            1 = x[2] - predict(p, q, x[0]-1, x[1]-1)
            p[x[0]-1] = p[x[0]-1] + learnR * (l*q[x[1]-1] - lamda*p[x[0]-1])
            q[x[1]-1] = q[x[0]-1] + learnR * (l*p[x[0]-1] - lamda*q[x[1]-1])
        trainRMSE.append(np.sqrt(np.mean(np.square(data[:,2] -\
                        [predict(p, q, x[0]-1, x[1]-1) for x in data]))))
        if valid is not None:
            validRMSE.append(np.sqrt(np.mean(np.square(valid[:,2] -\
                            [predict(p, q, x[0]-1, x[1]-1) for x in valid]))))
        nextFnV = lossFn(data, p, q, lamda)
        if(np.abs(prevFnV-nextFnV) < eps):</pre>
            break;
   return p, q, trainRMSE, validRMSE
```

### The following function is used for K-fold split:

## In [21]:

```
def cross_validation_split(dataset, folds=3):
    dataset_split = list()
    dataset_copy = np.copy(dataset)
    fold_size = math.ceil(len(dataset) / folds)
    for i in range(folds):
        if (i < folds-1):
            dataset_split.append(dataset_copy[i*fold_size:(i+1)*fold_size])
        else:
            dataset_split.append(dataset_copy[i*fold_size:len(dataset)])
    return dataset_split</pre>
```

## Function to perform training using LearnLatentFactors Algorithm

### In [22]:

## **Main Function**

```
In [24]:
```

```
movieLdf = read_data() ### read data
```

```
In [25]:
```

```
### Split the data into train and test set. The train set is further divided into train and
trainData = movieLdf.values[0:int(0.9*100000)]
testData = movieLdf.values[int(0.9*100000):]
```

Hyper paramters optimization to get best paramters that give use minimum rmse on validation set

### In [26]:

```
n_factors = [20, 40]
listAlpha = [0.005, 0.003]
listLamda = [0.01, 0.1]

listRMSE = np.zeros((3, 3, 3))

for i , factors in enumerate(n_factors):
    for j, alpha in enumerate(listAlpha):
        for k, lamda in enumerate(listLamda):
            print("n_Factors=", factors, " alpha=", alpha, " lambda=", lamda)
            listRMSE[i, j, k] = kFoldCV(trainData, factors, alpha, lamda, 3, 10) #Using onl
            print("RMSE on validation set = ", listRMSE[i, j, k])
```

```
n Factors= 20 alpha= 0.005 lambda= 0.01
RMSE on validation set = 1.0999056337461874
n Factors= 20 alpha= 0.005 lambda= 0.1
RMSE on validation set = 1.069648250590203
n_Factors= 20 alpha= 0.003 lambda= 0.01
RMSE on validation set = 1.0661329946623537
n Factors= 20 alpha= 0.003 lambda= 0.1
RMSE on validation set = 1.046282108508331
n Factors= 40 alpha= 0.005 lambda= 0.01
RMSE on validation set = 1.10578560665887
n Factors= 40 alpha= 0.005 lambda= 0.1
RMSE on validation set = 1.0865546917210298
n Factors= 40 alpha= 0.003 lambda= 0.01
RMSE on validation set = 1.0721369953113398
n_Factors= 40 alpha= 0.003 lambda= 0.1
RMSE on validation set = 1.065423065430813
```

Above results shows that we got minimum RMSE on validation set with n\_Factor=20, alpha= 0.003 and lambda= 0.1. So we will use these hyper paramters.

```
In [27]:
```

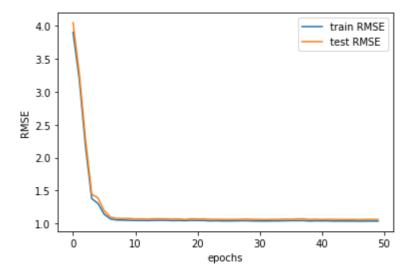
```
p, q, listTrainRMSE, listTestRMSE = trainMF(trainData, testData, 943, 1682, 0.003, 0.1, 20,
```

## In [28]:

```
plt.plot(listTrainRMSE, label="train RMSE")
plt.plot(listTestRMSE, label="test RMSE")
plt.xlabel("epochs")
plt.ylabel("RMSE")
plt.legend()
```

# Out[28]:

<matplotlib.legend.Legend at 0x2c916bc1788>



#### RMSE on test set

```
In [78]:
```

```
listTestRMSE[0] ###Rmse on test set
```

## Out[78]:

2.937676496510638

# **Exercise 3**

# **Library 1 Surprise:**

Surprise is a Python scikit building and analyzing recommender systems.

Users can use both built-in datasets (Movielens, Jester), and their own custom datasets. Provide various ready-to-use prediction algorithms such as baseline algorithms, neighborhood methods, matrix factorization-based (SVD, PMF, SVD++, NMF), and many others. Also, various similarity measures (cosine, MSD, pearson...) are built-in. Provide tools to evaluate, analyse and compare the algorithms performance. Cross-validation procedures can be run very easily using powerful CV iterators (inspired by scikit-learn excellent tools), as well as exhaustive search over a set of parameters. The name SurPRISE (roughly:)) stands for Simple Python Recommendation System Engine.

We used SVD to solve the problem

Singular value decomposition takes a rectangular matrix of gene expression data (defined as A, where A is a n x p matrix) in which the n rows represents the genes, and the p columns represents the experimental conditions. The SVD theorem states:

```
A(nxp)= U(nxn) * S(nxp) * V.T(pxp)

Where

U.T*U = I(nxn)

V.T*V = I(px)p (i.e. U and V are orthogonal)
```

Where the columns of U are the left singular vectors (gene coefficient vectors); S (the same dimensions as A) has singular values and is diagonal (mode amplitudes); and V.T has rows that are the right singular vectors (expression level vectors). The SVD represents an expansion of the original data in a coordinate system where the covariance matrix is diagonal.

Calculating the SVD consists of finding the eigenvalues and eigenvectors of AA.T and A.TA. The eigenvectors of A.TA make up the columns of V, the eigenvectors of AA.T make up the columns of U. Also, the singular values in S are square roots of eigenvalues from AA.T or A.TA. The singular values are the diagonal entries of the S matrix and are arranged in descending order. The singular values are always real numbers. If the matrix A is a real matrix, then U and V are also real.

# Loading and splitting the dataset

```
In [3]:
```

```
# Use movielens-100K
data = Dataset.load_builtin('ml-100k')
trainset, testset = train_test_split(data, test_size=.15)

Dataset ml-100k could not be found. Do you want to download it? [Y/n] y
Trying to download dataset from http://files.grouplens.org/datasets/movielen
s/ml-100k.zip... (http://files.grouplens.org/datasets/movielens/ml-100k.zi
p...)
Done! Dataset ml-100k has been saved to C:\Users\fahad/.surprise_data/ml-100
```

# Hyper parameter optimization

algo = gs.best\_estimator['rmse']
print(gs.best\_score['rmse'])
print(gs.best\_params['rmse'])

```
0.9228534951706099
{'n_factors': 120, 'n_epochs': 30, 'lr_all': 0.008, 'reg_all': 0.1}
Evaluating RMSE of algorithm SVD on 3 split(s).
                  Fold 1 Fold 2 Fold 3 Mean
                                                  Std
                  0.9209 0.9228 0.9278 0.9238 0.0029
RMSE (testset)
Fit time
                  7.90
                          7.88
                                  7.88
                                          7.89
                                                  0.01
Test time
                  0.36
                          0.27
                                  0.27
                                          0.30
                                                  0.04
Out[5]:
{'test_rmse': array([0.92090295, 0.92281803, 0.92775973]),
 'fit_time': (7.896497488021851, 7.8844685554504395, 7.881486654281616),
 'test_time': (0.364790678024292, 0.27086424827575684, 0.2698338031768799)}
```

# Using the optimal hyper paramters with SVD for matrix factorization

cross validate(algo, data, measures=['RMSE'], cv=3, verbose=True)

# In [7]:

```
# Use the new parameters with the train data and fit the model
algo = SVD(n_factors=120, n_epochs=30, lr_all=0.008, reg_all=0.1)
algo.fit(trainset)

#Now do the prediction on test set
test_pred = algo.test(testset)
print("SVD : Test Set")
accuracy.rmse(test_pred, verbose=True)
```

```
SVD : Test Set RMSE: 0.9103
Out[7]:
```

0.9102689311896246

# Library 2 Scikit Learn

# **Non-Negative Matrix Factorization (NMF)**

Find two non-negative matrices (W, H) whose product approximates the non-negative matrix X. This factorization can be used for example for dimensionality reduction, source separation or topic extraction.

The objective function is:

```
0.5 * ||X - WH||_Fro^2
```

```
alpha * I1_ratio * ||vec(W)||_1
alpha * I1_ratio * ||vec(H)||_1
0.5 * alpha * (1 - I1_ratio) * ||W||_Fro^2
0.5 * alpha * (1 - I1_ratio) * ||H||_Fro^2 Where:
```

 $||A||_{Fro^2} = \sum_{i,j} A_{ij}^2$  (Frobenius norm)  $||vec(A)||_1 = \sum_{i,j} abs(A_{ij})$  (Elementwise L1 norm) For multiplicative-update ('mu') solver, the Frobenius norm (0.5 \*  $||X - WH||_{Fro^2}$ ) can be changed into another beta-divergence loss, by changing the beta\_loss parameter.

The objective function is minimized with an alternating minimization of W and H.

It uses Coordinate Descent (CD).

# **Utility Functions**

### In [64]:

```
def cross_validation_split(dataset, folds=3):
    dataset_split = list()
    dataset_copy = np.copy(dataset)
    fold_size = math.ceil(len(dataset) / folds)
    for i in range(folds):
        if (i < folds-1):
            dataset_split.append(dataset_copy[i*fold_size:(i+1)*fold_size])
        else:
            dataset_split.append(dataset_copy[i*fold_size:len(dataset)])
    return dataset_split</pre>
```

# In [65]:

```
def kFoldCV1(R, data, factors, lamda, k):
   np.random.shuffle(data)
   folds = cross_validation_split(data, k)
   listCVE = []
   for i in range(len(folds)):
       train = np.vstack([x for x in folds if x is not folds[i]])
        valid = folds[i]
        RCopy = R.copy()
        for x in valid:
            RCopy[x[0]-1, x[1]-1] = 0
       model = NMF(n_components=factors, init='random', alpha=lamda)
       W = model.fit_transform(RCopy)
       H = model.components
        R_pred = np.dot(W, H)
        rmse = 0
        for x in valid:
            rmse += (x[2] - R_pred[x[0]-1, x[1]-1])**2
        rmse /= valid.shape[0]
        rmse = np.sqrt(rmse)
        listCVE.append(rmse)
   return np.mean(listCVE)
```

# **Reading Data**

```
In [66]:
```

```
movieLdf = pd.read_csv("u.data", delimiter="\t", header=None)
movieLdf = movieLdf.drop(movieLdf.columns[3], axis=1)
```

```
In [67]:
```

```
new_df = movieLdf.pivot(index = 0, columns = 1, values = 2).fillna(0).values
```

# **Splitting the dataset**

```
In [68]:
```

```
### Split the data into train and test set. The train set is further divided into train and
trn , tst = train_test_split(movieLdf, test_size = 0.3)
```

```
In [69]:
```

```
for x in tst.values:
    new_df[x[0]-1, x[1]-1] = 0
```

# Hyper paramter optimization with 3-folds cross validation

#### In [70]:

```
n_factors = [2, 4, 5, 7, 9 , 11, 15, 20, 80, 160]
listLamda = [0.001, 0.1]

listRMSE = np.zeros((10, 2))

print("Validataion RMSE is \n\n")
for i , factors in enumerate(n_factors):
    for j, lamda in enumerate(listLamda):
        print("n_Factors=", factors, " lambda=", lamda)
        listRMSE[i, j] = kFoldCV1(new_df, trn.values, factors, lamda, 3)
        print("RMSE=", listRMSE[i, j])
```

Validataion RMSE is

```
n Factors= 2 lambda= 0.001
RMSE= 3.168940044834376
n Factors= 2 lambda= 0.1
RMSE= 3.170897553576435
n Factors= 4 lambda= 0.001
RMSE= 3.1051451443952733
n Factors= 4 lambda= 0.1
RMSE= 3.1054734527947514
n Factors= 5 lambda= 0.001
RMSE= 3.090639067795783
n Factors= 5 lambda= 0.1
RMSE= 3.0933526522721304
n Factors= 7 lambda= 0.001
RMSE= 3.0706652153821588
n_Factors= 7 lambda= 0.1
RMSE= 3.075249407530737
n_Factors= 9 lambda= 0.001
RMSE= 3.070228741203023
n Factors= 9 lambda= 0.1
RMSE= 3.0747962556417403
n Factors= 11 lambda= 0.001
RMSE= 3.07129432249428
n Factors= 11 lambda= 0.1
RMSE= 3.077932790032984
n Factors= 15 lambda= 0.001
RMSE= 3.085845518297336
n Factors= 15 lambda= 0.1
RMSE= 3.08697013427585
n Factors= 20 lambda= 0.001
RMSE= 3.10654914748768
n Factors= 20 lambda= 0.1
RMSE= 3.109057918116239
n Factors= 80 lambda= 0.001
RMSE= 3.27823543830746
n Factors= 80 lambda= 0.1
RMSE= 3.27645402958872
n Factors= 160 lambda= 0.001
RMSE= 3.4124761577735527
n_Factors= 160 lambda= 0.1
RMSE= 3.410512619115572
```

we will use these hyper paramters.

```
In [71]:
```

```
model = NMF(n_components=9, init='random', alpha=0.001)
W = model.fit_transform(new_df)
H = model.components_
```

```
In [72]:
```

RMSE on test set is

#### Out[72]:

2.763689353496737

My naive implementation of matrix factorization give 2.9 Rmse on test set, build in library sci-learn matrix factorization which gives 2.7 Rmse on test set as compared to sci-learn surprise library which give 0.9 Rmse on test set. This shows that Surprise's SVD is performing best with with RMSE of 0.9.

In [ ]:			