# EECE 5698 Assignment 4: Matrix Factorization

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## **Problem 1**

1(a)

We can calculate the gradients as below

$$\nabla_{u_i} \operatorname{RSE}(U, V) = \nabla_{u_i} \left( \sum_{(i, j, r_{ij}) \in \mathcal{D}} \left( u_i^{\top} v_j - r_{ij} \right)^2 + \lambda \sum_{i=1}^n \|u_i\|_2^2 + \mu \sum_{j=1}^n \|v_j\|_2^2 \right)$$
$$= \sum_{j=1}^m 2 \left( u_i^{\top} v_j - r_{ij} \right) v_j + 2\lambda u_i$$

$$\nabla_{v_j} \operatorname{RSE}(U, V) = \nabla_{v_j} \left( \sum_{(i, j, r_{ij}) \in \mathcal{D}} \left( u_i^\top v_j - r_{ij} \right)^2 + \lambda \sum_{i=1}^n \|u_i\|_2^2 + \mu \sum_{j=1}^n \|v_j\|_2^2 \right)$$
$$= \sum_{i=1}^n 2 \left( u_i^\top v_j - r_{ij} \right) u_i + 2\mu v_j$$

1(b)

$$\nabla^{2}\ell(u,v) = \nabla^{2}(\left(u^{\top}v - r\right)^{2})$$

$$= \nabla^{2}(\left(uv - r\right)^{2})$$

$$= \begin{bmatrix} 2v^{2} & 2(2uv - r) \\ 2(2uv - r) & 2u^{2} \end{bmatrix}$$

$$\nabla^{2}\ell(0,0) = \begin{bmatrix} 0 & -2r \\ -2r & 0 \end{bmatrix}$$

It's eigenvalues are 2r, -2r, due to the noise,  $r \neq 0$ , so the eigenvalues could not be all non-negative. So  $\nabla^2 \ell(0,0)$  is not positive semi-definite. Thus  $\ell(u,v)$  is not convex.

1(c)

No. Because RSE is not convex. Gradient equals to zero doesn't mean global minimum.

1(d) Derive from results in 1(a)

$$\nabla_{u_i} \operatorname{RSE}(U, V) = \sum_{j=1}^{m} 2 \left( u_i^{\top} v_j - r_{ij} \right) v_j + 2\lambda u_i$$
$$= \sum_{j=1}^{m} 2 \left( 0^{\top} 0 - r_{ij} \right) 0 + 2\lambda 0$$
$$= 0$$

$$\nabla_{v_j} \operatorname{RSE}(U, V) = \sum_{i=1}^{n} 2 \left( u_i^{\top} v_j - r_{ij} \right) u_i + 2\mu v_j$$
$$= \sum_{i=1}^{n} 2 \left( 0^{\top} 0 - r_{ij} \right) 0 + 2\mu 0$$
$$= 0$$

# **Problem 2**

2(a)

$$RSE(U^{k}, V^{K}) = \sum_{(i,j,r_{ij})\in\mathcal{D}} \left(u_{i}^{k^{\top}} v_{j}^{k} - r_{ij}^{k}\right)^{2} + \lambda \sum_{i=1}^{n} \left\|u_{i}^{k}\right\|_{2}^{2} + \mu \sum_{j=1}^{n} \left\|v_{j}^{k}\right\|_{2}^{2}$$

$$= \sum_{i=1}^{n} \sum_{j=1}^{m} \delta_{ij}^{k^{2}} + \lambda \sum_{i=1}^{n} \left\|u_{i}^{k}\right\|_{2}^{2} + \mu \sum_{j=1}^{n} \left\|v_{j}^{k}\right\|_{2}^{2}$$

$$\nabla_{u_{i}} RSE(U^{k}, V^{k}) = \sum_{j=1}^{m} 2 \left(u_{i}^{k^{\top}} v_{j}^{k} - r_{ij}^{k}\right) v_{j}^{k} + 2\lambda u_{i}^{k}$$

$$= \sum_{j=1}^{m} 2\delta_{ij}^{k} v_{j}^{k} + 2\lambda u_{i}^{k}$$

$$\nabla_{v_{j}} RSE(U^{k}, V^{k}) = \sum_{i=1}^{n} 2 \left(u_{i}^{k^{\top}} v_{j}^{k} - r_{ij}^{k}\right) u_{i}^{k} + 2\mu v_{j}^{k}$$

$$= \sum_{i=1}^{n} 2\delta_{ij}^{k} u_{i}^{k} + 2\mu v_{j}^{k}$$

2(b) The implementation is as follow:

## 2(c)

We initialized the user and item profiles to random values instead of zero because we want to get rid of stuck at saddle point. If we initialize all profiles to be zero vectors, the RSE and its gradients will stay zero no matter how many iterations go, we will not get closer to our optimal solution.

#### **Problem 4**

4(a)

4(b)

```
def SE(joinedRDD):
    """ Receives as input a joined RDD as well as a λ and a μ and computes the MSE:
    SE(R,U,V) = Σ_{i,j} in data} (<ui,vj>-rij)^2
    The input is
    |-joinedRDD: an RDD with tuples of the form (i,j,δij,ui,vj), where δij = <ui,vj> - rij is the prediction difference.
    The output is the SE.
    SE = joinedRDD.map(lambda (i,j,sig,ui,vj): sig**2).reduce(add)
    return SE

def normSqRDD(profileRDD,param):
    """ Receives as input an RDD of profiles (e.g., U) as well as a parameter (e.g., λ) and computes the square of norms:
    λ Σ_i ||ui||_2^2
    The input is:
        -profileRDD: an RDD of the form (i,u), where i is an index and u is a numpy array
        | -param: a scalar λ>0
        The return value is:
        λ Σ_i ||ui||_2^2
    sum_u_norm = profileRDD.map(lambda (i, u): np.dot(u, u)).reduce(add)
    return param * sum_u_norm
```

In main, they are used as below to calculate the regularized objective function.

```
obj = SE(joinedRDD) + normSqRDD(U,args.lam) + normSqRDD(V,args.lam)
```

4(c)

#### **Problem 5**

5(a)

This part is for reading data for each fold if **args.folds** is specified, otherwise, read all data in a single fold.

This part constructs the training and testing set for each code.

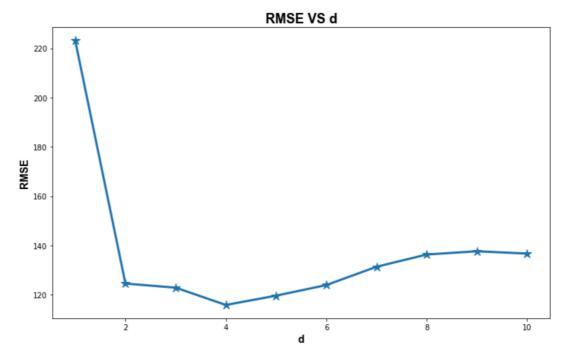
```
train_folds = [folds[j] for j in folds if j is not k ]

if len(train_folds)>0:
    train = train_folds[0]
    for fold in train_folds[1:]:
        it rain=train.union(fold)
        train.repartition(args.N).cache()
        test = folds[k].repartition(args.N).cache()
        Mtrain=train.count()
        Mtest=test.count()

print("Initiating fold %d with %d train samples and %d test samples" % (k,Mtrain,Mtest) )
else:
    train = folds[k].repartition(args.N).cache()
    test = train
        Mtrain=train.count()
        Mtest=test.count()
        print("Running single training over training set with %d train samples. Test RMSE computes RMSE on training set" % Mtrain )
```

```
Explaination of what each operation does is in the comments in red below:
# specify training folds as all folds except the test fold
train_folds = [folds[j] for j in folds if j is not k ]
if len(train folds)>0:
   # the next 3 lines merge the folds inside train folds into one
single list
   train = train_folds[0]
   for fold in train_folds[1:]:
       train=train.union(fold)
   # repartition the train data and cache it
   train.repartition(args.N).cache()
   # repartition the train data and cache it
   test = folds[k].repartition(args.N).cache()
   # calculate number of training samples
   Mtrain=train.count()
   # calculate number of test samples
   Mtest=test.count()
   # print info
   print("Initiating fold %d with %d train samples and %d test
samples" % (k,Mtrain,Mtest) )
else:
   # use all data as training set and repartition & cache it
   train = folds[k].repartition(args.N).cache()
   # use training set as test set
   test = train
   # calculate number of training samples
   Mtrain=train.count()
   # calculate number of test samples
   Mtest=test.count()
   # print info
   print("Running single training over training set with %d train
samples. Test RMSE computes RMSE on training set" % Mtrain )
5(b) step size is set as below, i is the number of iteration, args.gain and
args.power are user specified parameters:
gamma = args.gain / i**args.power
```

5(c) Plot the RMSE w.r.t. d as below:



We can see that b = 4 is the optimal value

5(d)

For --gain 0.1 --pow 0.0 --maxiter 5, we have the observation below

We can see that the step size is too big, so the RMSE is 'exploding'.

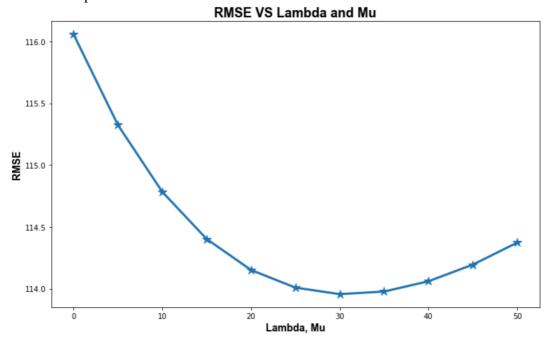
For **--gain 0.0001 --pow 1.0 --maxiter 20**, we have the observation below (for simplicity I just capture screen for 1 fold, the rest are similar):

```
Training set contains 10 users and 10 items
                Time: 9.031763 Objective: 1013500.042176
                                                             TestRMSE: 165.467171
Iteration: 1
                Time: 20.364278 Objective: 1010755.181194
                                                             TestRMSE: 165.464027
Iteration:
                Time: 31.279617 Objective: 1009408.319779
                                                              TestRMSE: 165.453452
Iteration:
                Time: 43.331643 Objective: 1008501.947583
                                                              TestRMSE: 165.443073
Iteration: 4
                Time: 53.420159 Objective: 1007813.094133
                                                              TestRMSE: 165.433514
Iteration:
                Time: 63.803859 Objective: 1007254.467580
                                                              TestRMSE: 165.424752
Iteration:
                Time: 76.152883 Objective: 1006782.808386
                                                              TestRMSE: 165.416680
Iteration:
Iteration: 8
                Time: 86.259500 Objective: 1006373.486731
                                                              TestRMSE: 165.409194
                Time: 97.272816 Objective: 1006011.123790
                                                              TestRMSE: 165.402207
Iteration: 9
                Time: 107.417636
Iteration: 10
                                     Objective: 1005685.461192
                                                                  TestRMSE: 165.395651
Iteration: 11
                Time: 117.996144
                                     Objective: 1005389.307289
                                                                 TestRMSE: 165.389468
Iteration: 12
                Time: 129.741703
                                     Objective: 1005117.420821
                                                                  TestRMSE: 165.383611
Iteration: 13
                Time: 139.845720
                                     Objective: 1004865.861072
                                                                  TestRMSE: 165.378044
Iteration: 14
                Time: 150.207647
                                     Objective: 1004631.588413
                                                                  TestRMSE: 165.372734
                                     Objective: 1004412.207617
                                                                  TestRMSE: 165.367655
Iteration: 15
                Time: 160.763873
                Time: 170.849454
                                     Objective: 1004205.796721
                                                                  TestRMSE: 165.362785
                Time: 181.421509
                                     Objective: 1004010.789316
                                                                  TestRMSE: 165.358104
Iteration: 18
                Time: 192.244272
                                     Objective: 1003825.891393
                                                                  TestRMSE: 165.353595
Iteration: 19
                Time: 202.592768
                                     Objective: 1003650.021234
                                                                  TestRMSE: 165.349245
Iteration: 20
                Time: 214.696185
                                     Objective: 1003482.265060
                                                                  TestRMSE: 165.345041
```

We can see that the step size is too small, so the optimization process is too slow.

#### 5(e)

Here as the searching space is continuous and so big, I just let  $\lambda = \mu = 0:5:50$  to search for the optimal value. However, a finer search may be conducted further around the optimal value I find.



From the figure, we can see that  $\lambda = \mu = 30$  is approximately the optimal solution.

## Bonus:

The method is similar as above.

- Set  $\lambda = \mu = 0$ , and set d = 1:10 to evaluate the least RMSE. If we can observe a first decreasing and then increasing curve, we take the minimum value of d, either do finer search or fix that value. Otherwise, give d a wider range to search.
- For investigating  $\lambda$  and  $\mu$ , I use a greedy search. I search  $\lambda$  in range [1,50] first, find the optimal value for  $\lambda$ . Then fix the optimal value for  $\lambda$ , search  $\mu$  in range [1,50]
- As it is too time consuming to do all this experiments, I just offer the general method here, more work needs to be done to find the exact optimal triplet.