Ranking-based Implicit Regularization for One-Class Collaborative Filtering

Defu Lian, Jin Chen, Kai Zheng, Enhong Chen, Senior Member, IEEE, and Xiaofang Zhou, Fellow, IEEE

Abstract—One-class collaborative filtering (OCCF) problems are ubiquitous in real-world recommendation systems, such as news recommendation, but suffer from data sparsity and lack of negative items. To address the challenge, the state-of-the-art algorithm assigns uninteracted items with smaller weights of being negative and performs low-rank approximation over the user-item interaction matrix. However, the prior ratings are usually suggested to be zero but may not be well-defined. To avert the direct utilization of prior ratings for uninteracted items, we propose a novel ranking-based implicit regularizer by hypothesizing that users’ preference scores for uninteracted items should not deviate a lot from each other. The regularizer is then used in a ranking-based OCCF framework to penalize large differences of preference scores between uninteracted items. To efficiently optimize model parameters in this framework, we develop the scalable alternating least square algorithm and coordinate descent algorithm, whose time complexity is linearly proportional to the data size. Finally, we extensively evaluate the proposed algorithms on six public real-world datasets. The results show that the proposed regularizer significantly improves the recommendation quality of ranking-based OCCF algorithms, such as BPRMF and RankALS. Moreover, the ranking-based framework with the proposed regularizer outperforms the state-of-the-art recommendation algorithms for implicit feedback.

Index Terms—One-class collaborative filtering, implicit regularization, item recommendation, coordinate descent, alternating least square.

1 INTRODUCTION

RECOMMENDER systems have become one of the most important ways to satisfy user interests and to increase the revenue of service providers. In recent years, plenty of research has concentrated on the recommendation for implicit feedback, since this data is more prevalent but full of noise. Implicit feedback is recorded whenever users interact with information systems. Recommendation for implicit feedback, also referred to as One-Class Collaborative Filtering (OCCF), aims to infer user interest from abundant yet noisy data and to provide each user a ranking list of items by matching user interest with items’ latent property.

However, OCCF is very challenging, since implicit feedback is usually very sparse and only includes users’ interacted items. Truly negative and potentially positive items are mixed together in uninteracted items [1], [2], [3]. Modeling uninteracted items plays a very important role in developing effective OCCF algorithms. Existing works treat interacted items as positive and uninteracted items as negative. These works can be categorized into two groups. The first regression-based methods optimize a pointwise weighted square loss, which sets the prior zero ratings to uninteracted items and assigns lower confidence to uninteracted items than interacted items [1], [2], [4], [5]. These works differ from each other in how many uninteracted items are considered for optimization. For example, all uninteracted items are considered negative in [1], [4] and partial samples from them are considered as negative in [2]. Due to the same confidence of uninteracted items being negative, the learning of parameters can be very efficient even when all uninteracted items are considered. The empirical comparison in [5] reveals that leveraging more uninteracted items leads to a higher quality of recommendation. According to the analysis in [6], [7], the explicit prior ratings suggest that prediction scores for uninteracted items should not deviate far from zero values. This constraint is formalized as an implicit regularizer, penalizing non-zero prediction scores for uninteracted items. However, the same and zero-valued prior ratings for all users may be questionable since the prior ratings may vary from user to user, but no other choices of prior ratings can be shown to work better.

Alternative ranking-based methods optimize a squared or non-linear pairwise loss, which also treats uninteracted items as negative. Instead of setting prior ratings to them, ranking-based methods assume that interacted items should be ranked higher than uninteracted items [3], [8], [9], [10]. In other words, the ranking-based methods avert the use of prior ratings. However, only when the square loss function is used, all uninteracted items can be efficiently incorporated [8]. When interacted items are ranked higher than uninteracted ones, we expect the ideal loss should be small, but the square loss may be still large. Therefore, they can not completely guarantee that interacted items are ranked higher than uninteracted items with respect to each user. When the non-linear loss functions are applied, we have to sample some items as negative from a large pool of uninteracted items, to achieve efficient optimization [3], [10]. Due to an extremely large number of candidate items,
only a small set of uninteracted items are used as negative for parameter learning. It is obvious that they also cannot guarantee the ranking between interacted items and other uninteracted items for each user.

To this end, in this paper, we propose a novel ranking-based implicit regularizer for ranking-based OCCF, by hypothesizing that prediction scores of uninteracted items for each user should not deviate a lot from each other [11]. The ranking-based implicit regularizer is formulated with squared differences of prediction scores between any two uninteracted items for each user. In this way, the proposed regularizer can implicitly determine personalized prior ratings for each user in an automatic way. Constrained by the regularizer, the ranking-based OCCF loss functions can guarantee that more uninteracted items are ranked lower than interacted items with respect to each user. This can satisfy the ranking-based requirement better and improve recommendation quality.

However, optimizing the ranking-based loss functions with the proposed regularizer is very challenging. For each user, every pair between uninteracted items is required to be considered. Since there are only a few interacted items, the total number of pairs between uninteracted items is extremely large. Therefore, it is time-consuming to directly calculate the regularization and its gradient with respect to latent factors. Without similar consideration to [7], gradient-based optimization approaches, which update parameters for users and items simultaneously, would be of low efficiency. To separate their update for simplification and improving efficiency, we develop alternating least square and coordinate descent algorithms to optimize pairwise loss functions with the proposed regularizer. For the squared pairwise loss function, parameters are updated in an exact closed-form. For the logit pairwise loss function, we alternate between deriving a local variational bound of the logit loss [12] and updating parameters in an approximately closed-form. Based on a novel caching strategy, the time complexity in each iteration of optimizing both loss functions is only linearly proportional to the data size.

To summarize, our main contributions are three-fold:

- We propose a novel ranking-based implicit regularizer and formulate it with squared differences of prediction scores between any two uninteracted items for each user. This regularizer plays an important role in better restricting ranking-based OCCF algorithms, such that more uninteracted items are ranked lower than interacted items with respect to each user. We apply the proposed regularizer for regularizing several ranking-based loss functions, and show that the regularizer significantly improves the quality of recommendation.
- We develop alternating least square and coordinate descent algorithms based on a strategy of first caching and continually updating for efficient optimization. The loss function can be either squared pairwise loss function or logit pairwise loss function. The time complexity of both algorithms is only in linear proportion to data size.
- We extensively evaluate the proposed algorithms on six real-world datasets, and show that the proposed algorithms outperform the competing baselines, including RankALS, BPR, WRMF and SQL_Rank, revealing the effectiveness of the proposed regularizer.

2 RELATED WORK

There are two main tasks in recommender systems, rating prediction and item recommendation. Rating prediction is investigated with explicit feedback datasets, where users rate some items with a specified set of ratings, while item recommendation is usually studied with implicit feedback datasets, where each user interacts with some items.

Rating prediction aims to predict missing ratings in a user-item rating matrix, based on intuitions that like-minded users share many rating patterns and that a user is likely to rate similar items to her rated ones. The KNN methods explicitly predict missing ratings with user-user similarity or item-item similarity [13], and have been improved by removing global effects from the data and globally learning interpolation weights for all nearest neighbors [14], [15]. The matrix factorization methods assume the user-item rating matrix is of low-rank, and directly apply SVD, stochastic gradient descent and alternating least square for learning user and item latent factors [16]. These methods may differ in whether incorporating missing values into loss functions. The matrix factorization methods were extended to model in a Bayesian way [17], such that hyperparameters are automatically learned from rating data. Following that, implicit feedback and side information were incorporated for promoting expressiveness of user latent factors as well as item latent factors [18], [19], [20]. However, these learning-based methods follow a Missing At Random (MAR) assumption, indicating the probability that a rating is missing does not depend on the value of that rating, or the value of any other missing ratings [21]. However, this assumption is usually violated in recommender systems, and the missing ratings are Missing At Not Random (MANR). Therefore, these learning-based methods were improved with error-imputation-based methods [22], [23], an inverse-propensity-scoring based method [24] or a doubly robust method of considering both of them [25].

Item recommendation aims to provide a ranking list of items for each user by learning from implicit feedback datasets. Since no ratings are in implicit feedback, we do not encounter the rating prediction problem, and the MANR assumption is not applicable. However, item recommendation faces the lack of negative items, since only each user’s interacted items are provided. The whole community almost agreed with treating interacted items as positive, but disputed much about how to use uninteracted items. As discussed in the introduction, uninteracted items are considered as negative with a lower confidence than interacted items as positive [1], [2], [4], [6], [26], [27], [28], or considered to be ranked in lower positions than interacted items [3], [8], [10], [29]. The ranking-based methods averted the usage of prior ratings for uninteracted items, but how to effectively exploit uninteracted items is very challenging. Though square loss can yield a closed-form updating equation in the alternating least square method when all uninteracted items are considered [8], accurately ranking pairs between interacted items and uninteracted ones may have large loss values. Though logit loss or hinge loss is almost zero for accurately ranking pairs, a small number of items should be sampled from uninteracted items for the sake of efficiency [3], [10]. The sampler should be efficient
for sampling and effective to pick informative samples. The common samplers include the uniform based sampler [3], the popularity based sampler [30], rejection sampling with the uniform proposal [10], dynamic negative sampler [31] and decomposable sampler [32], [33]. In this paper, we do not focus on negative sampling, but on a ranking-based implicit regularizer, which penalizes large differences of prediction scores between any two uninteracted items, to better distinguish interacted items from uninteracted items.

3 RANKING-BASED IMPLICIT REGULARIZER

3.1 Implicit Regularizer

In this paper, we investigate item recommendation for implicit feedback dataset. Assume user-item interaction matrix $R \in \{0, 1\}^{M \times N}$ available, where $M$ denotes the number of users and $N$ denotes the number of items. Each element $r_{ui}$ indicates whether the user $u$ interacts with the item $i$. $U = \{u_1, u_2, ..., u_M\}$ denotes all users and $E = \{e_1, e_2, ..., e_N\}$ denote all items. $E_u = \{i \in E | r_{ui} > 0\}$ denotes the items which the user $u$ interacts with and $U_i = \{u \in U | r_{ui} > 0\}$ denotes the users who have interacted with the item $i$.

The SOTA regression-based methods treat all uninteracted items as negative, but assign them a lower confidence being negative [1], [4], [6]. In particular, they optimize the objective function as follows:

$$\sum_{u,i} w_{ui} (r_{ui} - p_u^T q_i)^2 + \lambda \|p\|_F^2 + \|q\|_F^2, \quad (1)$$

where $p_u$ is the latent factor of user $u$, and $q_i$ is the latent factor of item $i$. $P$ and $Q$ respectively stack $p_u$ and $q_i$ by row. $w_{ui}$ equals $\alpha$ when user $u$ interacts with item $i$ and 1 otherwise. $\alpha$ is generally greater than 1.

According to [6], [34], Eq (1) is decomposed into a prediction error term and an implicit regularizer term, which is defined as:

$$\sum_{u} \sum_{i \notin E_u} (p_u^T q_i - 0)^2. \quad (2)$$

The recommendation for implicit feedback is interpreted as explicit problems with the additional implicit regularizer [6], which penalizes non-zero prediction scores of uninteracted items for each user, and plays an important role in improving deep matching network for recommendation [7].

3.2 Ranking-based Implicit Regularizer

It may be questionable to set the prior ratings of all uninteracted items for all users to zero, since prior ratings for uninteracted items should be as least different from user to user. However, it is difficult to set appropriate prior ratings for uninteracted items. Observing the implicit regularizer actually implies some “similarity” among uninteracted items (preference scores approaching the prior rating), we propose a ranking-based implicit regularizer, by hypothesizing each user’s preference scores for uninteracted items are close to each other. Formally, the ranking-based implicit regularizer is defined as follows:

$$\Omega = \frac{1}{2} \sum_{u} \sum_{i < j \notin E_u} (p_u^T q_i - p_u^T q_j)^2$$

$$= \frac{1}{4} \sum_{u} \sum_{i < j \notin E_u} (p_u^T q_i - p_u^T q_j)^2. \quad (3)$$

The expression can be written in a symmetric way due to no contribution of the case $i = j$ and symmetry between $i < j$ and $i > j$. Compared with Eq (2), the ranking-based regularizer eliminates the usage of prior ratings for uninteracted items through difference of preference scores between any two uninteracted items. When $\Omega$ is decreased, preference scores of uninteracted items are closer to each other. Therefore, the ranking-based implicit regularizer penalizes large differences of preference scores between any two of uninteracted items, playing a similar role to the implicit regularizer.

However, directly computing the regularizer and deriving its gradient is very time-consuming, since the regularizer involves $O(MN^2)$ differences of preference scores. Below we investigate how to reduce a great number of double computations. Before delving into details, we first rewrite the ranking-based implicit regularizer as follows:

$$\Omega = \frac{1}{4} (\Omega_0 - 2\Omega_1 + \Omega_2),$$

where

$$\Omega_0 = \sum_{u} \sum_{i,j \notin E_u} (p_u^T q_i - p_u^T q_j)^2$$

$$\Omega_1 = \sum_{u} \sum_{i \notin E_u} \sum_{j \in E_u} (p_u^T q_i - p_u^T q_j)^2$$

$$\Omega_2 = \sum_{u} \sum_{i \notin E_u} \sum_{j \notin E_u} (p_u^T q_i - p_u^T q_j)^2$$

The regularizer is then decomposed into summation over a full set of items and summation over each user’s interacted items. We will follow such a decomposition to derive gradient and derivative of the regularizer with respect to parameters. The gradient is used for alternating least square computed prior to updating user latent factors.

3.3 Regularizer Gradient

We take each user’s latent factor or each item’s latent factor as a whole and derive the gradient of the regularizer with respect to user latent factors and item latent factors. Firstly, we consider the gradient of $\Omega$ with respect to $p_u$, latent factor of user $u$, which is derived as follows:

$$\frac{\partial \Omega}{\partial p_u} = 2 \sum_{i \in E_u} \sum_{j} (q_i - q_j)(p_u^T q_i - p_u^T q_j)$$

$$= 2 \left( N Q_u^T q_u + N_u Q^T Q - \hat{q} q_u^T - \hat{q} q_u^T \right) p_u$$

where $\hat{q} = \sum_j q_j$ and $\hat{q}_u = \sum_{i \in E_u} q_i$. $Q_u \in \mathbb{R}^{N_u \times K}$, denoting $Q_{[E_u]}$, is a submatrix of $Q$ extracted by $E_u$. $N_u = |E_u|$ denotes the number of interacted items of user $u$ and $C_u = |N - N_u|$ denotes the number of uninteracted items. Deriving gradient of other two parts with respect to $p_u$ is similar. Aggregating them together, we obtain gradient of the ranking-based implicit regularizer with respect to $p_u$:

$$\frac{\partial \Omega}{\partial p_u} = C_u (Q^T Q - Q_u^T Q_u) p_u - (\hat{q} - \hat{q}_u)(\hat{q} - \hat{q}_u)^T p_u$$

$$= C_u (S - S_u) p_u - (\hat{q} - \hat{q}_u)(\hat{q} - \hat{q}_u)^T p_u \quad (4)$$

where $S = Q^T Q$ and $S_u = Q_u^T Q_u$. Due to independence of user latent vectors, the matrix $S$ and the vector $\hat{q}$ can be computed prior to updating user latent factors.
The gradient of the ranking-based implicit regularizer with respect to item latent factor is as follows:

\[
\frac{\partial \Omega}{\partial q_l} = \sum_u C_u p_u^T q_l - \sum_u p_u p_u^T (\tilde{q} - \tilde{q}_u) \\
- \sum_{u \in U_l} C_u p_u^T q_l + \sum_{u \in U_l} p_u p_u^T (\tilde{q} - \tilde{q}_u) \\
= (H - H^T) q_l - (T - T^T) \tilde{q} + P^T \tilde{r} - P_l^T \tilde{r}_l.
\]

where \( H = P^T \text{diag}(c) P, H^T = P^T \text{diag}(c^T) P, T = P^T P, \) and \( T^T = P_l^T P_l. \) The \( u \)-th element of \( c \) is \( C_u \).

4.1 Loss function

Ranking-based methods do not use explicit prior ratings for uninteracted items, and exactly align with the ranking-based implicit regularizer. Basically, ranking-based methods assume interacted items are ranked higher than uninteracted items. Therefore, ranking-based OCCF usually optimizes the following objective function:

\[
L_1 = \sum_u \sum_{i \in E_u, j \notin E_u} \ell(r_{ui} - r_{uj}, \hat{r}_{ui} - \hat{r}_{uj}) + \lambda R(P, Q)
\]

where \( r_{ui} \) denotes the true rating, \( \hat{r}_{ui} = p_u^T q_i \) denotes the predicted rating and \( R(P, Q) \) is the regularization term of \( \ell_1 \)-norm or \( \ell_2 \)-norm. Note that we do not use the absolute ratings but the relative ratings and aver explicit prior ratings for uninteracted items. In the RankALS \([8]\), \( r_{ui} - r_{uj} = 1 \) and \( \ell(y, \hat{y}) = (y - \hat{y})^2 \), where item popularity is incorporated for weighting each loss. In this case, all uninteracted items can be incorporated with efficiency guarantee, but correctly ranking pairs may also lead to large loss values. In the BPR \([3]\), \( \ell(y, \hat{y}) = -(y \log \sigma(\hat{y}) + (1 - y) \log(1 - \sigma(\hat{y}))) \) and \( r_{ui} - r_{uj} = 1 \). In this case, sampling from uninteracted items is a necessary condition of efficient training. In other words, whichever of these loss functions is used, they cannot guarantee that interacted items are ranked higher than uninteracted ones. Therefore, we will optimize the ranking-based loss function with the proposed ranking-based implicit regularizer as follows

\[
L = \sum_u \sum_{i \in E_u, j \notin E_u} \ell(1, \hat{r}_{ui} - \hat{r}_{uj}) + \lambda R(P, Q) + \alpha \Omega
\]

where we assume \( r_{ui} - r_{uj} = 1 \). \( \alpha \) controls effect of the ranking-based implicit regularizer.

However, we observe that the magnitude of implicit regularizer is much larger than the loss, particularly the BPR loss, in Eq (8). Therefore, the coefficient \( \alpha \) may be too sensitive to tune. According to our experience, the coefficient is usually very small to achieve best recommendation performance in the validation set. To reduce sensitivity of the coefficient, and intrinsically address the imbalance between interacted items and uninteracted items, we introduce a normalization strategy, which assign user-specific weight to each loss and each implicit regularizer. Particularly, we finally optimize the following objective function,

\[
L' = \sum_u \sum_{i \in E_u, j \notin E_u} z_u \ell(1, \hat{r}_{ui} - \hat{r}_{uj}) + \lambda R(P, Q) + \frac{1}{2 \alpha} \sum \sum_{i < j \notin E_u} y_u (\hat{r}_{ui} - \hat{r}_{uj})^2
\]

where \( z_u = f(C_u \times N_u) \) and \( y_u = f(C_u \times C_u) \) is a normalization coefficient for the loss and the regularizer, respectively. Note that \( N_u \) is the number of interacted items of user \( u \) and \( C_u \) is the number of uninteracted items. Here, we focus on
three strategies of the normalization coefficient, by setting \( f(x) = \frac{1}{x} \), \( f(x) = \frac{1}{\log x} \) and \( f(x) = \frac{1}{x^2} \). When \( f(x) = \frac{1}{x} \), it simply computes the average loss and regularizer for each user, but weakens contribution of users who have interacted with many items to the loss. The other two strategies can alleviate the problem to some extent.

Though we introduce a normalization coefficient into the regularizer, the gradient and derivative of regularizer with respect to parameters can be easily obtained by following the chain rule. Particularly, letting \( \tilde{p}_u = \sqrt{y_u p_u} \) and explicitly showing parameters of the regularizer, \( \Omega(P, Q) = \Omega(\hat{P}, \hat{Q}) \), and the gradient can be derived as follows:

\[
\frac{\partial \Omega(P, Q)}{\partial p_u} = \sqrt{y_u} \frac{\partial \Omega(\hat{P}, \hat{Q})}{\partial \tilde{p}_u} = y_u \frac{\partial \Omega(P, Q)}{\partial \tilde{p}_u}
\]

(10)

The derivation can be obtained similarly. However, such a trick is difficult to use in the loss functions, so we consider the loss functions one by one.

### 4.2 Optimizing Square Loss

The square loss is widely used in many machine learning models, due to yielding closed-form solutions and being convenience for theoretical analysis. When it is applied for ranking-based methods in one-class collaborative filtering, it is formulated as follows:

\[
L' = \frac{1}{2} \sum_u \sum_{i \in E_u} \sum_{j \notin E_u} z_{ui} (1 - (\hat{r}_{ui} - \hat{r}_{uj}))^2
\]

(11)

Because of the large number of uninteracted items, straightforwardly computing gradient is very time-consuming. The use of square loss can make gradient computation efficient. Since we observe that both alternating least square and coordinate descent work better than gradient descent for optimizing WRMF [1], the well-known method for recommendation from implicit feedback, below we develop an alternating least square method and a coordinate descent method for efficient parameter learning.

#### 4.2.1 Alternating Least Square

The optimization procedure of the alternating least square alternates between updating user latent vector and updating item latent vector until convergence. The ALS method requires us to derive the gradient of the objective function with respect to latent vectors. The gradient of the regularizer is given in Section 3.3, so below we derive the gradient of the loss \( L' \) in Eq. (11). The gradient with respect to \( p_u \) is elaborated as follows:

\[
\frac{\partial L'}{\partial p_u} = \sum_{i \in E_u} \sum_{j \notin E_u} -z_u (q_i - q_j) \left( 1 - (p_u^T q_i - p_u^T q_j) \right)
\]

\[
= z_u (N_u q_i - q_i z_u) + z_u (N_u S + (N - 2 N_u) S^u) p_u - z_u (q_i - \hat{q}_i) + \frac{1}{\lambda} p_u.
\]

(12)

Setting \( \frac{\partial L}{\partial p_u} = \frac{\partial L'}{\partial p_u} + \alpha \frac{\partial \alpha}{\partial p_u} + \lambda \frac{\partial \lambda}{\partial p_u} \) to zero, we can obtain the closed-form solution for updating latent vector of user \( u \) in a form of \( A_u p_u = b_u \). Note that \( \hat{q} \) and \( Q^T Q \) should be precomputed for efficient computation. Regarding the gradient with respect to the item latent vector \( q_i \), we elaborate the similar process as follows:

\[
\frac{\partial L'}{\partial q_i} = \sum_{u \notin U_i} \sum_{j \in \bar{E}_u \setminus E_{u_i}} -z_u (q_i - q_j) \left( 1 - (p_u^T q_i - p_u^T q_j) \right)
\]

\[
= N (T - T^l) q_i - (H - 2 T^l) q_\hat{i} - P^T (z \circ n) - N P^T z_i.
\]

(13)

where the \( u \)-th element of \( n \) is \( N_u \). \( T \) and \( H \) is obtained by substituting \( p_u \) with \( \tilde{p}_u = \sqrt{y_u p_u} \) in \( T \) and \( H \), respectively. \( T^l \) and \( H^l \) respectively denote submatrix of \( T \) and \( H \) extracted by the user set \( U_i \). Setting \( \frac{\partial L}{\partial q_i} = \frac{\partial L'}{\partial q_i} + \alpha \frac{\partial \alpha}{\partial q_i} + \lambda \frac{\partial \lambda}{\partial q_i} \) to zero, we can obtain the closed-form solution for updating latent vector of item \( i \) in a form of \( A_i q_i = b_i \). Due to independence of item \( i \), \( T \) and \( H \) can be precomputed. However, it is infeasible to precompute \( \hat{q} \) and \( \tilde{p}_u \) for continuously changing data. We develop a strategy of first caching and continual updating to address this challenge. The details of procedure are referred to Algorithm 1, where each line is commented with time cost.

### Algorithm 1: ALS for Square Loss

**Input:** The rating matrix \( R \in \mathbb{R}^{M \times N} \)

**Output:** latent factors \( P \in \mathbb{R}^{M \times K} \) and \( Q \in \mathbb{R}^{N \times K} \)

1. Randomly initialize \( P, Q \),
2. for \( u \in \{1, \ldots, M\} \) do
3.   \( Q_{u[i]} = Q_{u[i]}^1 N_u \),
4.   \( \tilde{p}[u] = \langle Q_{u[i]}, p_u \rangle \),
5.   \( \hat{q} = Q^T 1_N \),
6. repeat
7.   \( S = Q^T Q \),
8.   for \( u \in \{1, \ldots, M\} \) do
9.     \( S^u = Q_{u[i]}^2 \),
10.    Compute \( A_u \) and \( b_u \),
11.   \( A_u p_u = b_u \),
12.   Compute \( H_i = T_i \),
13.   Compute \( T_i = T_i \),
14.   Compute \( T_i = T_i \),
15.   Compute \( T_i = T_i \),
16.   Compute \( T_i = T_i \),
17.   Compute \( T_i = T_i \),
18.   Compute \( T_i = T_i \),
19.   until Convergent;
20. Compute \( H_i = T_i \), \( H_i \) and \( T_i \),
21. Compute \( A_i q_i = b_i \),
22. Solve \( A_i q_i = b_i \),
23. Compute \( q_i = q_i + q_i \),
24. Compute \( Q_{i[i]} = Q_{i[i]} - q_i + q_i \),
25. Compute \( q_i = q_i - q_i + q_i \),
26. Compute \( q_i = q_i + q_i \),
27. until Convergent;
28. until Convergent;}
Complexity Since $\hat{q}$ and $Q^T Q$ are precomputed, updating user latent factor $p_u$ only costs $O(N_u K^2 + K^3)$, where we assume solving the system of linear equation costs $O(K^3)$. By caching the global quantities and dynamically updating them, updating $q_i$ only costs $O(N K^2 + K^3)$. To summarize, the time complexity of the proposed learning algorithm is $O(||R||K^2 + (M + N)K^3)$, where $||R||_0$ denotes the number of non-zero entries in the interaction matrix $R$. In other words, the optimization algorithm is only in linear proportion to the total number of users’ interacted items.

Regarding of the space complexity, the algorithm needs $O(MK + N(K + K^2))$ memory spaces for storing recommendation models. More specifically, the user and item latent matrices occupies $O(MK)$ and $O(NK)$, respectively. We also need $O(K^2)$ extra space for storing $Q^T Q$. Since we exploit the caching and updating strategy, we need $O(MK)$, $O(K)$, $O(M)$ extra memory spaces for these matrices $Q$, $q$, $\tilde{r}_i$, respectively.

### 4.2.2 Coordinate Descent

In spite of being linearly proportional to the number of interacted items, alternating least square is still cubic with the dimension $K$, due to the computation of matrix inversion. To improve the efficiency, we develop an coordinate descent algorithm, which updates an element of latent factor at a time given others fixed. Below we first obtain the derivative with respect to $p_{uf}$

$$\frac{\partial C}{\partial p_{uf}} = z_a N_a S_{ff} p_{uf} + z_u (N - 2 N_u) S_{uf} p_{uf} - 2 z_u \tilde{q}_u (\tilde{q}_f - \tilde{q}_uf) p_{uf} + z_u N_u \tilde{q}_uf - z_u N q_{uf} + z_u (N - 2 N_u) \sum_{k \neq f} p_{uk} S_{kf}^{u}$$

(14)

$$+ z_u N_u \sum_{k \neq f} p_{uk} S_{kf} - z_u \tilde{q}_uf \sum_{k \neq f} p_{uk} \tilde{q}_k - z_u \tilde{q}_uf \sum_{k \neq f} p_{uk} \tilde{q}_uf$$

We note that the time complexity of computing $S^{u}$ is $O(NuK^2)$. Since each time we only update one factor, the computation can be accelerated. In particular, we do not compute $S^{u}_{ff}$ in advance, instead we directly compute $S^{u}_{ff}$ and $\sum_{k \neq f} p_{uk} S^{u}_{kf}$ as follows:

$$S^{u}_{ff} = \sum_{i \in E_u} q_i^2,$$

$$\sum_{k \neq f} p_{uk} S^{u}_{kf} = \sum_{i \in E_u} p_{ui} q_i k_i f_i = \sum_{i \in E_u} \tilde{r}_{if} u_i q_i f_i.$$  

where $\tilde{r}_{if} u_i = \tilde{r}_{ui} - p_{uf} q_{if}$. As long as we first cache prediction scores for each user’s interacted items, these two quantities can be computed in $O(N_u)$ for each factor.

Next we obtain the derivatives with respect to $q_{if}$

$$\frac{\partial C}{\partial q_{if}} = N (\tilde{T}_{ff} - \tilde{T}_{ff}^T) q_{if} - (\tilde{H}_{ff} - 2 \tilde{H}_{ff}^T) q_{if} - \tilde{t}_{if} q_{if}$$

(15)

$$- \tilde{b}_f + P_i [:, f]^T \text{diag}(z_i) \tilde{r}_i + b_f - NP_i [:, f]^T z_i$$

We also do not compute $\tilde{H}$ and $\tilde{T}$ in advance, we directly compute the related quantities as follows.

$$\tilde{T}_{ff} = \sum_{u \in U} z_u p_{uf}^2,$$

$$\tilde{H}_{ff} = \sum_{u \in U} C_u z_u p_{uf},$$

$$\sum_{k \neq f} q_{ik} \tilde{T}_{kf} = \sum_{u \in U} z_u (\tilde{r}_{uf}) p_{uf},$$

$$\sum_{k \neq f} q_{ik} \tilde{H}_{kf} = \sum_{u \in U} C_u z_u (\tilde{r}_{uf}) p_{uf},$$

$$\tilde{t}_{if} q_{if} = \sum_{u \in U} z_u p_{uf} P_{i} q_{if}.$$  

As long as we additionally cache $\tilde{r}_u = p_{uf}^T \hat{q}$, these quantities are computed in $O(N_i)$. By observing only users in $U_i$ are
used for computation, we suggest updating \( \hat{r}_{ui} \) only for users in \( U_l \) instead of the whole users after updating each factor. This dramatically reduces the computational cost compared with the ALS method. The details of procedure are referred in Algorithm 2.

**Complexity** Compared with alternating least square method, matrix inversion is not required in the coordinate descent optimization method. By caching the global quantities and dynamically updating them, the time complexity of parameter learning is now reduced to \( O(|R|K + (M + N)K^2) \), \( K \) times faster than alternating least square. However, it may require more iterations for convergence. The space complexity of optimizing recommendation models is the same as ALS in Section 4.2.1, i.e., \( O(MK + NK + K^2) \).

### 4.3 Optimizing Logit Loss

The square loss can yield closed-form equations for updating parameters, but may lead to large loss values for accurately ranking pairs. In this section, we investigate how to optimize the ranking-based log loss with the proposed regularizer. Different from the square loss, it is impossible to incorporate all uninteracted items at a time. To address the issue, we sample the same size of uninteracted items as interacted items, and assume they should be ranked lower than interacted items in the ranking-based log loss, following previous work [3]. Then, the ranking-based log loss with the proposed regularizer is formulated as follows:

\[
\mathcal{L}' = \sum_u \sum_{i \in E_u} \sum_{j \notin E_u} z_u \log \left( 1 + \exp \left( -\left( \hat{r}_{ui} - \hat{r}_{uj} \right) \right) \right)
\]

where \( F_u \) is the set of sampled negative items for user \( u \). Different from the Bayesian Personalize Ranking (BPR) loss, which compares an interacted item with a sampled uninteracted item, the log loss function compares an interacted item with all items in \( F_u \).

Another issue of the ranking-based log loss lies in difficulty of directly deriving the closed-form updating equations. Here, we seek a upper variational quadratic bound of the log loss [12], [34] and then derive the closed-form updating equations for alternating least squares and coordinate descent methods. Particularly,

\[
\log \left( 1 + e^{-\hat{r}_{uij}} \right) = \log(1 + e^{\hat{r}_{uij}}) - \hat{r}_{uij}
\]

\[
\leq \lambda(u_{ij}) \left( \hat{r}_{uij} - \hat{r}_{uij}^2 - \frac{1}{2}(\hat{r}_{uij} + \hat{r}_{uij}) + \log(1 + e^{\hat{r}_{uij}}) \right)
\]

where \( \lambda(x) = \frac{1}{2} \tanh \left( \frac{x}{2} \right) = \frac{1}{2} \left( \sigma(x) - \frac{1}{2} \right) \) and \( \sigma(x) \) is the sigmoid function. The equality holds only if \( \hat{r}_{uij} = \hat{r}_{uij} \). By setting \( \hat{r}_{uij} = (\hat{r}_{ui} - \hat{r}_{uj}) \), the ranking-based log loss is bounded from the above by

\[
\hat{\mathcal{L}}' = \sum_u \sum_{i \in E_u} \sum_{j \notin E_u} z_u \left( \lambda(u_{ij}) (\hat{r}_{ui} - \hat{r}_{uj})^2 - \frac{1}{2}(\hat{r}_{ui} - \hat{r}_{uj}) \right)
\]

where \( \hat{r}_{uij} \) is the most recent estimation of \( \hat{r}_{uij} \). Next we elaborate the ALS and CD algorithms for optimizing the ranking-based log loss.

#### 4.3.1 Alternating Least Square

The gradient of \( \hat{\mathcal{L}}' \) with respect to user latent vector \( p_u \) is derived as follows:

\[
\frac{\partial \hat{\mathcal{L}}'}{\partial p_u} = \sum_{i \in E_u} \sum_{j \notin E_u} 2z_u \lambda(u_{ij}) \left( q_i - q_j \right) \left( q_i - q_j \right)^T p_u
\]

\[-\sum_{i \in E_u} \sum_{j \notin E_u} \frac{1}{2} \left( q_i - q_j \right)
\]

\[= 2z_u \sum_{i \in E_u} \left( \sum_{j \notin E_u} \lambda(u_{ij}) \right) q_i q_i^T p_u
\]

\[+ 2z_u \sum_{j \in E_u} \left( \sum_{i \notin E_u} \lambda(u_{ij}) \right) q_j q_j^T p_u
\]

\[-2z_u \sum_{i \in E_u} \sum_{j \notin E_u} \lambda(u_{ij}) (\hat{r}_{uij} + q_i q_j^T p_u) + q_j q_j^T p_u
\]

\[-\frac{1}{2} z_u N_u \sum_{i \in E_u} q_i + \frac{1}{2} z_u N_u \sum_{j \notin E_u} q_j
\]

For efficient computation, we cache prediction scores for positive items in a sparse matrix \( \hat{\mathcal{R}}^+ \) and negative items in a sparse matrix \( \hat{\mathcal{R}}^- \), so that \( \lambda(u_{ij}) \) is computed in \( O(1) \). Setting \( \frac{\partial \mathcal{L}}{\partial p_u} = \frac{\partial \hat{\mathcal{L}}'}{\partial p_u} + \alpha \frac{\partial \alpha}{\partial p_u} + \lambda \hat{q}_u \) to zero, we can obtain the closed-form solution for updating latent vector of user \( u \) in a form of \( A_u p_u = b_u \). Computing a part of \( A_u \) in Eq (16) is achieved with a two-layer loop, which costs \( O(N u^2 K^2) \).

The gradient of \( \hat{\mathcal{L}}' \) with respect to item latent vector \( q_i \) is derived as follows:

\[
\frac{\partial \hat{\mathcal{L}}'}{\partial q_i} = \sum_{u \notin E_i} \sum_{j \notin E_i} 2z_u \lambda(u_{ij}) (p_u q_i - p_u q_j) p_u - \frac{1}{2} z_u p_u
\]

\[-\sum_{u \notin E_i} \sum_{j \notin E_i} 2z_u \lambda(u_{ij}) (p_u q_i - p_u q_j) p_u + \frac{1}{2} z_u p_u
\]

\[= 2z_u \sum_{u \notin E_i} \lambda(u_{ij}) \hat{q}_i p_u^T q_l
\]

\[+ 2z_u \sum_{u \notin E_i} \lambda(u_{ij}) \hat{q}_l p_u^T q_l
\]

\[-2z_u \sum_{u \notin E_i} \lambda(u_{ij}) \hat{r}_{uij}
\]

\[-2z_u \sum_{u \notin E_i} \lambda(u_{ij}) \hat{r}_{uij} + q_i q_j^T p_u
\]

\[-\frac{1}{2} z_u N_u p_u + \frac{1}{2} \sum_{u \notin E_i} z_u N_u p_u
\]

where \( V_i \) is a set of users without interacting item \( l \). We follow the similar first-caching-and-continual-updating strategy to efficiently compute high-cost quantities. Setting \( \frac{\partial \mathcal{L}}{\partial q_i} = \frac{\partial \hat{\mathcal{L}}'}{\partial q_i} + \alpha \frac{\partial \alpha}{\partial q_i} + \hat{q}_i \) to zero, we can obtain the closed-form solution for updating \( q_i \) in a form of \( A_q q_i = b_i \). The details of procedure are referred in Algorithm 3.

**Complexity** According to Algorithm 3, it is easy to understand that time complexity of updating user latent factor is \( O(N_u N^2 K^2 + MK^3) \). Regarding to updating latent factor for all items, the two-layer loop costs \( O \left( K^2 (N \sum_{l=1}^N \sum_{u=1}^M \delta_{ul} + \sum_{l=1}^N \delta_{ul} \sum_{j=1}^N \delta_{uj}) \right) \)
Algorithm 3: ALS for Logit-Loss

Input: The rating matrix $R \in \mathbb{R}^{M \times N}$
Output: Latent factors $P \in \mathbb{R}^{M \times K}$, and $Q \in \mathbb{R}^{N \times K}$
1 Randomly initialize $P$, $Q$;
2 for $u \in \{1, \ldots, M\}$ do
3 \[ Q[u, \_] \leftarrow Q^T u N_u, \]
4 \[ \hat{r}[u] \leftarrow \langle Q[u, \_], p_u \rangle; \]
5 \[ q \leftarrow Q^T 1_N; \]
6 repeat
7 \[ \text{for } u \in \{1, \ldots, M\} \text{ do} \]
8 \[ \text{Sample set item } E_u \text{ from } E \setminus E_u; \]
9 \[ S \leftarrow Q^T Q; \quad \text{ // } O(NK^2) \]
10 \[ \text{for } u \in \{1, \ldots, M\} \text{ do} \]
11 \[ \hat{r}_{E_u}, \hat{r}_{F_u} \leftarrow Q[E_u] ; p_u, Q[F_u] ; p_u; \]
12 \[ A_u \leftarrow \text{Hessian matrix of } (\alpha \Omega + \frac{1}{2} \lambda \|p_u\|_2^2); \]
13 \[ \text{for } i \in E_u \text{ do} \]
14 \[ \text{for } j \in F_u \text{ do} \]
15 \[ \lambda(\epsilon_{u,i,j}) \leftarrow \lambda(\hat{r}_{u,i}^+ - \hat{r}_{u,j}^-); \quad \text{ // O(1)} \]
16 \[ \text{Update } A_u; \quad \text{// } O(K^2) \]
17 \[ \text{Compute } b_u; \quad \text{// } O(N_u K) \]
18 \[ \text{Solve } A_u b_u = b_u; \quad \text{// } O(K^3) \]
19 \[ \hat{r}_{E_u}, \hat{r}_{F_u} \leftarrow Q[E_u] ; p_u, Q[F_u] ; p_u; \]
20 \[ H \leftarrow P^T \text{diag}(y \odot c) P; \quad \text{ // } O(K^2) \]
21 \[ \tilde{T} \leftarrow P^T \hat{y}(y) P; \quad \text{// } O(K^2) \]
22 \[ b \leftarrow P^T \hat{y}(y) \hat{r}; \quad \text{// } O(K) \]
23 \[ \text{for } l = 1 : N \text{ do} \]
24 \[ \hat{r}_{U_l}, \hat{r}_{V_l} \leftarrow P[U_l] ; q_l, P[V_l] ; q_l; \]
25 \[ A_l \leftarrow \text{Hessian matrix of } (\alpha \Omega + \frac{1}{2} \lambda \|q_l\|_2^2); \]
26 \[ b_l \leftarrow -\frac{1}{2} \sum_{u \in U_l} z_u N_u p_u + \frac{1}{2} \sum_{v \in V_l} z_v N_v p_v; \]
27 \[ \text{for } u \in U_l \text{ do} \]
28 \[ \text{for } j \in F_u \text{ do} \]
29 \[ \lambda(\epsilon_{u,i,j}) \leftarrow \lambda(\hat{r}_{u,i}^+ - \hat{r}_{u,j}^-); \quad \text{ // O(1)} \]
30 \[ \text{Update } A_l, b_l; \quad \text{// } O(K^2) \]
31 \[ \text{for } u \in V_l \text{ do} \]
32 \[ \text{for } i \in E_u \text{ do} \]
33 \[ \lambda(\epsilon_{u,i,j}) \leftarrow \lambda(\hat{r}_{u,i}^+ - \hat{r}_{u,j}^-); \quad \text{ // O(1)} \]
34 \[ \text{Update } A_l, b_l; \quad \text{// } O(K^2) \]
35 \[ \text{Solve } A_l q_l = b_l; \quad \text{// } O(K^3) \]
36 \[ \text{for } u \in U_l \text{ do} \]
37 \[ \text{Update score of item } l \text{ in } \hat{r}_{E_u}, \hat{r}_{F_u}; \]
38 \[ q \leftarrow q - q^{old} + q_l; \quad \text{// } O(K) \]
39 \[ Q_{U_l} \leftarrow Q_{U_l} - q^{old} + q_l; \quad \text{// } O(N_l K) \]
40 \[ \hat{r}_{U_l} \leftarrow \hat{r}_{U_l} + P_l^T (-q^{old} + q_l); \quad \text{// } O(N_l K) \]
41 \[ b \leftarrow b + \tilde{T}(-q^{old} + q_l); \quad \text{// } O(K^2) \]
42 until Convergent;

4.3.2 Coordinate Descent

Next we investigate how to optimize the objective function with coordinate descent. We first derive the derivative of the objective function with respect to $p_{uf}$ as follows:

\[
\frac{\partial \tilde{L}}{\partial p_{uf}} = \sum_{i \in E_u, j \in F_u} 2z_u \lambda(\epsilon_{uij}) (q_{if} - q_{jf})^2 p_{uf} + \sum_{u \in V_l} z_u \lambda(\epsilon_{uij}) (q_{if} - q_{jf}) \left( \hat{r}_{uf}^f - \hat{r}_{uf}^f \right) \tag{18}
\]

where $\hat{r}_{uf}^f = \hat{r}_{uf} - p_{uf} q_{if}$. The derivative with respect to $q_{if}$ is given as:

\[
\frac{\partial \tilde{L}}{\partial q_{if}} = \sum_{u \in U_l, j \in F_u} 2z_u \lambda(\epsilon_{uij}) \left( \hat{r}_{uf}^f + p_{uf} q_{if} - \hat{r}_{uf}^f \right) + \sum_{u \in V_l} z_u \lambda(\epsilon_{uij}) \left( \hat{r}_{uf}^f + p_{uf} q_{if} - \hat{r}_{uf}^f \right) - \sum_{u \in V_l} \frac{1}{2} \sum_{j \in F_u} z_u N_u p_{uf} + \sum_{u \in V_l} \frac{1}{2} \sum_{j \in F_u} z_u N_u p_{uf} \tag{19}
\]

Setting $\frac{\partial L}{\partial p_{uf}} = \frac{\partial \tilde{L}}{\partial p_{uf}} + \alpha \frac{\partial \eta}{\partial p_{uf}} + \lambda p_{uf} = 0$ and $\frac{\partial L}{\partial q_{if}} = \frac{\partial \tilde{L}}{\partial q_{if}} + \alpha \frac{\partial \eta}{\partial q_{if}} + \lambda q_{if} = 0$, we can take turn to update each factor of latent factors until convergence. Different from alternating learning square, the variational parameter $\lambda(\epsilon_{uij})$ should be updated whenever a factor is updated. This means $\lambda(\epsilon_{uij})$ is computed $K$ times as many as the alternating least square method. However, when only one factor is updated each time, the loss could be closer to the upper bound. According to Algorithm 2 and Algorithm 3, it is easy to figure out the detailed procedure of the coordinate descent algorithm, so that we omit it for space limitation. Based on the analysis of Algorithm 2, it is easy to deduce that the time complexity of the coordinate descent algorithm is $O(\sum u z_u^2 K + (M + N) K^3)$, which should be $K$ times faster than the alternating least square method. The space complexity of optimizing recommendation models is the same as ALS in Section 4.2.1, i.e., $O(MK + NK + K^2)$.

5 EXPERIMENTS

5.1 Experimental Settings

5.1.1 Datasets

We perform experiments on two types of datasets to verify the effectiveness of our methods. The first type of dataset is explicit feedback dataset, including MovieLens, Amazon and Yelp, where the ratings range from 1 to 5.

- **MovieLens**: The MovieLens dataset is from the famous MovieLens10M dataset, including 10,000,054 ratings from 71,567 users for 10,681 items.
- **Yelp**: The Yelp dataset includes 2,685,066 ratings from 409,117 users for 85,539 points of interest, such as restaurants, hotels, and shopping malls.
- **Netflix**: The Netflix dataset is another famous movie rating dataset. It contains 463,770 users and 17,764 items, which is larger than the MovieLens and Yelp dataset.
Datasets Ratings Users Items Sparsity

RANK baselines in a validation set, which consists of 5% training

5.2.1 Comparison with baselines

5.2 Results and Analysis

According to the metric of NDCG@200.

Gowalla

• 20%
during each iteration and treat them as negative samples.

• LastFM

as positive. The other type of dataset for experiments

datasets following [3], by considering user’s rated

data. Implicit feedback datasets are converted into implicit feedback datasets following [3], by considering user’s rated item as positive. The other type of dataset for experiments is implicit feedback dataset, including CiteUlike, Gowalla and LastFM, where user’s interacted items are considered as positive.

- CiteUlike: The CiteUlike dataset is from user’s individual library of articles [35].

- Gowalla: The Gowalla dataset includes users’ check-ins at locations in a location-based social network.

- LastFM: The dataset is a famous music recommendation dataset [36] with 360,000 users.

All the datasets are publicly available and we filter the users and items with fewer than 10 interactions. Table 1 summarizes the statistics of the datasets.

When optimizing the ranking-based logit loss, we randomly pick $N_i$ items from uninteracted ones for each user during each iteration and treat them as negative samples.

For each user, we randomly sample his 80% ratings as training set and the rest 20% as testing set. We fit a model to the training set and evaluate it in the test set. We repeat 5 random splits and report the averaged performances.

5.1.2 Baselines

- RankALS [8] optimizes the ranking-based square loss with an alternating least square algorithm. The parameter $s_j$ for each item is set to 1 so that they can be fairly compared with the proposed methods.

- WRMF [1], [2] optimizes regression-based square loss, by considering all uninteracted items as negative and assigning them lower confidence being negative. The optimization is based on alternating least square, which scales linearly with the number of interaction.

- BPR [3] optimizes the ranking-based logit loss with stochastic gradient descent. The negative samples are randomly picked from uninteracted items.

- SQL_RANK [37] exploits a listwise approach for recommendation which cast listwise collaborative ranking as maximum likelihood under a permutation model.

By presenting the top $k$ candidate items for each user, two standard metrics are used to assess the quality of ranking: NDCG and Recall. The parameters are fine-tuned according to the metric of NDCG@200.

5.2 Results and Analysis

5.2.1 Comparison with baselines

We tune hyperparameters of the proposed methods and baselines in a validation set, which consists of 5% training data. Note that the proposed method with the square loss is trained for 20 iterations and the proposed method with logit loss is trained for 50 iterations. Results are shown in Table 2, where we set the dimension of latent space in all methods to 32 and report NDCG and Recall at cutoffs 10, 50, and 200.

On all the datasets, our proposed algorithm outperforms all the baselines. In particular, the square loss with the ranking-based implicit regularizer (Square-RIR) is much better than RanKALS. The relative improvements are 120.1% on average with respect to NDCG@10. This shows the effectiveness of the proposed regularizer. The logit loss with the ranking-based implicit regularizer (Logit-RIR) is also much better than BPR. The relative improvements are 51.7% on average with respect to NDCG@10. The superior performance of BPR to RankALS may lie in logit loss, which can yield small loss values for accurately ranking pairs. WRMF is the best among the baselines, since it imposes an implicit regularizer $\sum_{i \in N} \sum_{j \in E_{ij}} (p_i q_j)^2$ to penalize non-zero preference prediction. However, the proposed algorithms still outperform WRMF, and the average relative improvements are 5.5% in the explicit feedback datasets, and up to 8.3% in the implicit feedback datasets in terms of NDCG@10. This indicates that the ranking-based implicit regularizer is more suitable for sparser implicit feedback datasets. The logit loss can not consistently beat the square loss, but BPR is better than RanKALS. This indicates that the ranking-based implicit regularizer benefits the square loss more than the logit loss, to reduce large loss values for accurately ranking pairs.

Furthermore, we investigate how the performance of the proposed algorithm varies with the increasing dimension of latent space. The results of NDCG@200 on MovieLens and Gowalla datasets are reported in Figure 1, where $K$ is varied in the set $\{16, 32, 48, 64, 128\}$. As the dimension of latent space increases, the performance of the proposed method with the logit loss gets close to the square loss in terms of NDCG@200 in both datasets. The proposed method with the logit loss outperforms BPR, and the relative improvements increase as the dimension grows. Compared to WRMF, the proposed methods are better even with the increase of dimension, especially in the MovieLens dataset. In the sparser Gowalla dataset, as the dimension increases, the relative improvements to WRMF algorithms gradually decrease. This shows the superiority of the proposed ranking-based implicit regularizer to the implicit regularizer in WRMF.

5.2.2 Effect of loss functions

As we exploit two kinds of loss functions and four normalization strategies in our framework, we conduct an ablation study to understand the effect of normalization strategies and loss functions in the four datasets. Alternating least square (ALS) is used for optimizing all the algorithms. “none”, “avg”, “sqrt” and “log” denotes $f(x) = 1$, $f(x) = \frac{1}{x}$, $f(x) = \frac{1}{\sqrt{x}}$ and $f(x) = \frac{1}{\log(x)}$, respectively.

Figure 2 summarizes the performance of NDCG@50 when the coefficient of the regularizer varies. The best parameter for different normalization strategies is different from each other. For the “avg” and “sqrt” strategy, $\alpha$ is much larger than the other two strategies and thus much easier to tune. Moreover, these two strategies show better recommendation performance. This indicates these two

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Fig. 1. NDCG@200 and Recall@200 on the MovieLens and Gowalla.

5.2.3 Effect of optimization methods

Figure 3 shows how loss values reduce in the CiteULike dataset when the loss functions are optimized with the alternating least square method and the coordinate descent method, respectively. We choose the normalization strategy \( f(x) = 1 \) and fine-tune the hyperparameters. In the figure, loss values are shown on the y-axis with a logarithmic scale. We observe that all methods converge within 30 iterations. According to this figure, we can see that the alternating least square method converges in fewer iterations than the coordinate descent method, particularly optimizing the logit loss. This is because the coordinate descent method updates one factor at a time while different factors are correlated with each other. Therefore, the coordinate descent method is much more efficient but requires more iterations for convergence. As analyzed before, the time cost of the coordinate descent method is K times less than the alternating least
square method. This is also evidenced in Figure 4. In addition, an interesting observation is that using the coordinate descent method to optimize the logit loss can converge to a smaller loss than the alternating least square method. A potential explanation is that in the coordinate descent method, the variational parameter $\lambda(\hat{e}_{uij})$ is recalculated in each iteration, so that the loss function is closer to the upper bound. This is also confirmed by using the coordinate descent method to optimize the square loss function converge to the approximately same loss to the alternating least square method, where there is no variational parameter.

In addition to loss values, we also report the recommendation performance in Table 3, where the square loss and the logit loss are optimized by these two optimization methods. Here, we utilize the "sqrt" normalization strategy and set the dimension to $K = 16$. From this table, we can observe that in the CiteULike and Gowalla datasets, optimizing both loss functions by the coordinate descent method is slightly better than the alternating least square method. In the MovieLens and Yelp datasets, optimization based on the coordinate descent method only leads to improvements for the logit loss but not for the square loss. This is also explained in the coordinate descent method, the variational parameter is recalculated each time, so the loss function is better approximated by its upper bound.

6 CONCLUSION AND FUTURE WORK

In this paper, we propose a novel ranking-based implicit regularizer and leverage them to improve the ranking-based one-class collaborative filtering algorithms, whose loss functions include the square loss and the logit loss. We develop an alternating least square method and a coordinate descent method to efficiently optimize the proposed objective functions. We then elaborate on the analysis of training time complexity, showing both of them scale linearly with data size, and the coordinate descent method is $K$ times faster than the alternating least square method in each iteration. Finally, we extensively evaluate the proposed method with 6 real-world datasets and show that the proposed regularizer significantly improves the recommendation performance of the ranking-based one-class collaborative filtering algorithms, and outperforms the competing baselines. We also show that the coordinate descent method is more efficient and even yields slightly better recommendation performance than the alternating least square method, particularly for the logit loss.

In the future, we would like to investigate how to optimize the proposed objective functions with stochastic gradient descent and apply them for advanced neural network recommenders [7]. Moreover, we will also try to relax the hypothesis that prediction scores of uninteracted items should not deviate a lot from each other. For example, we may consider only restricting score proximity among uninteracted items with similar popularity.

ACKNOWLEDGMENTS

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Enhong Chen (SM’07) received the PhD degree from USTC. He is a professor and vice dean of the School of Computer Science, USTC. His general area of research includes data mining and machine learning, social network analysis, and recommender systems. He was on program committees of numerous conferences including KDD, ICDM, and SDM. His research is supported by the National Science Foundation for Distinguished Young Scholars of China.

Xiaofang Zhou is a Chair Professor at the Department of Computer Science and Engineering (CSE), Hong Kong University of Science and Technology (HKUST). He received his BSc and MSc degrees from Nanjing University in 1984 and 1987 respectively, and PhD in Computer Science from University of Queensland (UQ) in 1994. His research focus is to find effective and efficient solutions for managing, integrating, and analysing large-scale complex data for business, scientific and personal applications.