

Computational Intelligence Laboratory Project: Collaborative Filtering

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Abstract—

I. INTRODUCTION

II. MODELS AND METHODS

In this section we describe what pre-existing algorithms we based our work on and how we used blending in the end to combine the predictions of several independent approaches. Furthermore we explain how we optimized our grade by finding a good quality / CPU time trade-off.

A. Preliminaries

Throughout this section we make frequent use of the following notation. We denote by \mathcal{K} the set of all existing ratings, i.e. $\mathcal{K} = \{(u, i) \mid \text{user } u \text{ rated item } i\}$. Additionally we define M and N to be the number of all users and items, respectively. In addition we define the set of ratings for user's and item's, $R(u) = \{i \mid (u, i) \in \mathcal{K}\}$ and $R(i) = \{u \mid (u, i) \in \mathcal{K}\}$. Furthermore we define the true rating of an item i by user u as r_{ui} and denote this rating's approximation by \hat{r}_{ui} . Moreover we make frequent use of Greek letters, where $\mu = |\mathcal{K}|^{-1} \sum_{(u,i) \in \mathcal{K}} r_{ui}$ is the global mean of all issued ratings, λ denotes regularization constants and γ is the learning rate in an Stochastic Gradient Descent algorithm.

B. Data Imputation and Baseline Estimators

Given that standard Collaborative Filtering algorithms such as K-Means and Singular Value Decomposition (SVD) require dense instead of sparse matrices different strategies exist to fill in the missing values. The most simple one is to replace all missing values with a constant zero value. Accuracy can be improved by considering other constants such as the global mean μ of all existing entries. More advanced techniques try to approximate a given missing value for a user u and item i by taking a combination the global mean and the specific user and item means b_u and b_i [1]–[4]:

$$b_{ui} = \mu + b_u + b_i. \quad (1)$$

b_u and b_i can be directly computed from the data, however it is advisable to introduce regularize terms to overcome over-fitting when only very few ratings are available. This leads to the following equations, where $R(u)$ and $R(i)$

denote the sets of all ratings by user u and item i respectively [1]–[4]:

$$b_i = \frac{\sum_{u \in R(i)} (r_{ui} - \mu)}{\lambda_i + |R(i)|} \quad b_u = \frac{\sum_{i \in R(u)} (r_{ui} - \mu - b_i)}{\lambda_u + |R(u)|} \quad (2)$$

λ_i and λ_u are regularize parameters that should be found via cross-validation. Finally it is also possible to estimate b_u and b_i by solving the regularized least squares problem [1]–[4]

$$\min_{b_*} \sum_{(u,i) \in \mathcal{K}} (r_{ui} - \mu - b_u - b_i)^2 + \lambda \left(\sum_u b_u^2 + \sum_i b_i^2 \right). \quad (3)$$

Possible solution approaches include Alternating Least Squares (ALS) or Stochastic Gradient Descent (SGD). The baseline estimators are able to capture a lot of signal present in the data so that it is advisable to make this preprocessing step part of every more involved algorithm. For example, the RMSE score for a simple SVD solution improved drastically when missing data was imputed with optimized baseline estimators.

C. Factorized Models

Motivated by the progress a simple SVD approach could make we investigated more sophisticated approaches.

1) *Regularized SVD*: Shortly after the Netflix Challenge started in 2006 Simon Funk proposed the idea of a “regularized SVD” [5]. In contrast to an ordinary Singular Value Decomposition this approach does not rely on the imputation of missing values, but only considers actual present ratings for training. The algorithm tries to find two matrices $P \in \mathbb{R}^{M \times K}$ and $Q \in \mathbb{R}^{N \times K}$ that accurately represent user-item interactions. Both users and items get transformed to the same latent factor space of a fixed dimension K where their dot product is taken to measure their compatibility. A rating r_{ui} is then approximated by $\hat{r}_{ui} = b_{ui} + q_i^T p_u$. In order to avoid over-fitting to the data a regularize term is added that penalizes large magnitudes of p_u and q_i . The associated least squares problem is the following:

$$\min_{q^*, p^*} \sum_{(u,i) \in \mathcal{K}} (r_{ui} - b_{ui} + q_i^T p_u)^2 + \lambda (\|q_i\| + \|p_u\|) \quad (4)$$

This problem again can be solved using either ALS or SGD. Here we assume that the biases have been estimated in a preprocessing step, however it is also possible to learn them at the same time as P and Q .

2) *SVD++*: Based on the work of Simon Funk, several improvements to regularized SVD were suggested, that also consider the item-item relations present in the data. These methods include “NSVD” [6] and “Asymmetric-SVD” [1] which both aim to model the user vector p_u based on the items that given user rated. For example in NSVD p_u is modeled via the sum $(\sum_{j \in R(u)} x_j) / \sqrt{|R(u)|}$ where $X \in \mathbb{R}^{N \times K}$ is a item dependent factor matrix learned from the data. The term $|R(u)|^{-1/2}$ serves as a user dependent normalization constant, which stabilizes the sum’s variance across the range of observed values of $|R(u)|$ [4].

Building on top of these ideas two very similar models were developed, that again include an explicit p_u vector for every user. These models are dubbed “MF-NSVD1” [7] and “SVD++” [1], [4] respectively. In SVD++ in addition to the user vector p_u the sum over all rated items is kept, which leads to the following approximation formula:

$$\hat{r}_{ui} = b_{ui} + q_i^T \left(p_u + |R(u)|^{-1/2} \sum_{j \in R(u)} x_j \right) \quad (5)$$

Observing that the modified user term is completely independent of the current item leads to a very quick learning and prediction algorithm that has a linear time complexity with regard to the input size. Similar to previous models we solved the regularized least squares problem using SGD.

D. Neighborhood Models

Historically neighborhood models enjoyed a huge popularity in Collaborative Filtering applications. These models include user-user and item-item based models. In user-user models one tries to find similar users to a given user u and use their existing ratings to recommend new items to u . Conversely, in item-item based models one tries to find similar items to the items a given user already rated, and recommend those. Usually the number of distinct items are less than the number of distinct users which results in item-item models being more efficient, because it is faster to compute similarities between all items than it is to compute them between all users. In addition, item-item models tend to provide a better user experience, because they offer an explanation for their recommendations and users are usually more familiar with their previous rated items than other users that are supposedly similar to them.

E. Blending of Results

III. RESULTS

A. Dataset

The dataset which was used throughout the development of our algorithm was made available on the project course webpage [8]. It contains the ratings of 10 000 users on 1000 movies on a scale between 1 and 5. Naturally, not every user rated every movie, in fact only 1 388 107 of the 10 000 000 possible ratings were present leading to a data sparsity of 86%.

B. Evaluation of Results

In order to train and test the developed algorithm the present ratings were randomly split in two disjoint sets of equal size. The first set then formed the training set, i.e. data that was trained on, and the second set was used for testing the algorithm. The chosen metric is “Root Mean Squared Error” (RMSE) which is defined in the following way [1], [3], [4]:

$$\sqrt{\sum_{(u,i) \in \text{TestSet}} \frac{(r_{ui} - \hat{r}_{ui})^2}{|\text{TestSet}|}} \quad (6)$$

Here r_{ui} denotes the real rating while \hat{r}_{ui} is the approximation through the algorithm. With ratings restricted to the interval $[1, 5]$ valid RMSE values range between 0 and 4, with lower scores being better. In addition, the raw CPU time was evaluated for the project grade. This favors solutions that are efficient with regard to time.

IV. DISCUSSION

V. SUMMARY

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