Collaborative-Based Recommendation System

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Abstract—In this work, we study the problem of recommending items to users using a collaborative-based filtering approach. In particular, we are only given access to users and their ratings, and we would like to recommend new movies by predicting the missing ratings. To this end, we implement 16 different models. A modified version of Singular Value Decomposition (SVD++) performs the best among these models. It achieves a score of 0.97671 on Kaggle's validation set. In order to improve our predictions even further, we implement a ridge regression model that predicts ratings based on the predictions of 13 different models. We obtain a slight improvement over SVD++, increasing our accuracy score to 0.97368 on Kaggle's validation set. Finally, we explore and implement some further models proposed by Koren [4].

Keywords-Baseline, Collaborative-Based Filtering, K-Nearest Neighbors (KNN), Ridge Regression, Matrix Factorization, Netflix Data, Singular Value Decomposition (SVD, SVD++), Recommendation.

I. Introduction

Recommendation systems have became increasingly popular in recent years and have changed the way users interact with inanimate websites and applications. They are used in a variety of domains including (but not limited to) recommending movies, music, articles, search queries, and products/services in general. Recommendation systems can be broadly divided into three categories: collaborative-based systems, content-based systems, or a hybrid of both. Content-based systems examine the properties of the items and the user's preferences. On the other hand, collaborative-based filtering approaches are based on collecting and analyzing a large amount of information on users' behaviors, activities or preferences and predicting what users will like based on their similarity to other users.

In this work, we are given no access to any content information. Thus, we will stick to collaborative-based filtering systems. An important note to make is that in real life scenarios, one might be able to provide better recommendations if content data (name of movies, genre, actors/actress, etc..) or any additional features/statistics of user's activities (rating timestamps for example since users' preferences might vary over time) are available.

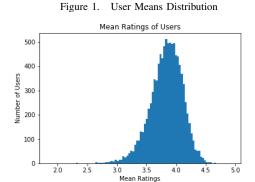
The remainder of this report is organized as follows: We first start by discussing some data statistics and exploration. We then discuss the various models we implemented, and the results we have obtained. We conclude the study with the model(s) we propose and some further thoughts.

II. DATA EXPLORATION

The training data consists of ratings provided by 10000 users on a subset of 1000 movies. The ratings are integer values that vary between 1 and 5, with 1 being the lowest possible rating and 5 being the highest. We extensively explored the data, but we will only discuss some of our findings.

Initially, we thought that among the many rating profiles, we might have two interesting extreme ones. In particular, we looked for users that strictly rated movies they hate, and those who strictly rated movies they loved. The reason behind this exploration was motivated by the idea that if we have many such users, then we might be able to extract additional features by labeling/learning whether a user might hate/like a movie. However, as shown in figure 1, the ratings are concentrated around 3.8. That is, on average users rate both movies they like or hate. In fact, there were no users who only gave ratings of 1. Similarly, there were no users who only gave ratings of 5.

Furthermore, we look at different groups of users which



we refer to as bloggers and new users. The bloggers are the top 5% and the new users are the lowest 5% of the users, where users are ordered by the number of their ratings. In particular, any user with more than 253 rating is considered a blogger, and any user with at most 19 ratings is considered a new user. An interesting behaviour is that there seem to be an almost increasing "linear" relationship between the ratings and their distribution, except for the following two regions: the region between the ratings 5 and 4 for the bloggers, and the region between the ratings 5 and 5 for the new users. This can base seen in figures 3 & 5, respectively. The data suggests that users tend to watch and thus rate

Figure 2. New Users Ratings Distribution

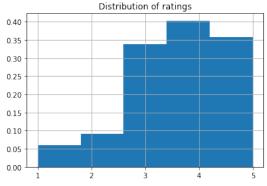
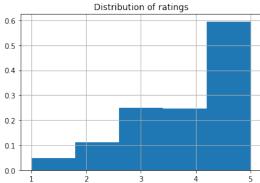


Figure 3. Bloggers Distribution



movies they believe they will like. There might be several explanations for the gaps between ratings 2 and 3 for the new user. One possible explanation we considered is that new users don't treat consecutive points as equidistant, and hence they regard rating 2 as too low for some movies. We thought of incorporating this into our models, by adjusting the ratings so that the distribution of ratings has an almost linear relationship with the ratings, however, we didn't have enough reason to justify such an assumption. Additionally, we were not able to explain the gap in the bloggers data, thus we decided to leave the data as is.

III. MODELS

Table I summarizes the RMSE results we obtained for all the models we included in the final blending, the weights assigned when blending the different models using ridge regression (see section III-L), and the running times of the various models when run on an Intel Core i7-6700HQ, 8Gb RAM . In what follows, we give a brief overview of the algorithms we implemented.

¹This is only the overhead running time we get when blending the different models. That is, assuming we have predictions of the various models already computed.

²The RMSE of all the models is calculated using 6% of the training data set, except that of the blending model. The RMSE of the blending model is obtained from Kaggle's validation set.

# Method	RMSE	Computation Time (in minutes)	Weights
# Method	KWISE	Computation Time (in innutes)	weights
Baseline	0.9974	< 1	-0.2272
Global Mean	1.1143	< 1	0.1677
Item Mean	1.0285	< 1	-0.0914
Item Median	1.0992	< 1	-0.0154
User Mean	1.0915	< 1	-0.2250
User Median	1.1476	< 1	0.0239
KNN Item	0.9852	5	0.2175
KNN User	0.9893	50	0.3583
Matrix Factorization (ALS)	0.9847	13	-0.0066
Matrix Factorization (SGD)	0.9895	13	0.0111
Slope One	0.9982	< 1	-0.2586
SVD	0.9805	3	-0.8426
SVD++	0.9777	140	1.8864
Blending	0.97368	< 11	-

 $\label{eq:Table I} \mbox{Table I} \mbox{Summary on the Models Implemented} \ ^2$

A. Global Mean & Median

The simplest model is to compute the mean or the median value of the non-zero ratings in the training set, and return this value as the prediction. This gives a baseline score from which we can compare further models.

B. Users Mean & Median

Another simple model is to compute the mean or median of the ratings given by a user. The mean of the ratings of user n can be computed as follows.

$$\hat{x}_n := \frac{1}{|\Omega_{n:}|} \sum_{m \in \Omega_{n:}} x_{nm} \tag{1}$$

where Ω_n is the set of movies rated by the n^{th} user. The median can be computed in a similar fashion.

C. Items Mean & Median

Another simple model is to compute the mean or median of the ratings given to a movie. The mean of the ratings given to a movie m can be computed as follows.

$$\hat{x}_m := \frac{1}{|\Omega_{:m}|} \sum_{n \in \Omega_{:m}} x_{nm} \tag{2}$$

where Ω_m is the set of users who rated the m^{th} movie. The median can be computed in a similar fashion.

D. Baseline

The baseline predictions are the base on which most of the complex models are built on. It provides biases for each user and each item and combines them together with the global mean to make predictions. From this point onwards, another algorithm can add deviations from the baseline model using a different criteria, with the advantage of working with unbiased ratings. Normally, the biases would be extrapolated using ridge regression over the ratings. However, we use a model specified by Koren [4], based on ALS, which is lighter than ridge regression but gives similar results. This algorithm is implemented in the Python Surprise library [1].

E. KNN User Based

Before the Netflix Prize, it was the most common model in recommendation systems [4]. This model builds a similarity matrix over the users and utilizes it to find the k most similar users to a given one. For the similarity matrix, we use a baseline pearson similarity metric since this metric accounts for the biases of the ratings per user/movie. This criteria is computationally heavier than the other metrics (mean squared difference, pearson, cosine) because it makes some effort to deal with the biases. This algorithm is implemented in the Python Surprise library [1].

F. KNN Item Based

As the previous model, this builds a baseline pearson similarity matrix, but over the items and uses it to find the k most similar items to a given one. It is preferred to the user based KNN for the following reasons.

- Usually items are much less numerous than users, so the the algorithm is computationally less expensive.
- Adding a new user to the system doesn't involve the recomputation of all the weights, so it is easier to manage.
- Recommendations can be explained, improving the
 overall users' satisfaction. When the system suggests
 a movie, it can indicate the movie (or the movies)
 watched by the user in the past that brought the system
 to that particular recommendation. In a user based
 KNN, this kind of explanation is impossible due to
 privacy restrictions and is not useful anyway.

This algorithm implemented in the Python Surprise library [1].

G. Matrix Factorization SGD

Given D movies and N users, we define $X \in \mathbb{R}^{N \times D}$ to be the matrix containing all ratings. We aim to find two matrices $W \in \mathbb{R}^{K \times D}$ and $Z \in \mathbb{R}^{K \times N}$, such that $X \approx Z^T W$. In order to find these matrices we minimize the following quantity.

$$\frac{1}{2} \sum_{(d,n)\in|\Omega} \left[x_{nd} - (\mathbf{Z}^{\mathbf{T}}\mathbf{W})_{nd} \right]^2 + \frac{\lambda_w}{2} \|\mathbf{W}\|^2 + \frac{\lambda_z}{2} \|\mathbf{Z}\|^2$$
 (3)

where K is the number of latent features, and λ_z , λ_w are scalar constants that weigh the regularization terms.

The stochastic gradient descent is a stochastic approximation of the gradient descent optimization. The gradient of the objective function is computed by chosing a random element of the summation.

H. Matrix Factorization ALS

ALS is one of the main alternatives to the SGD to solve (3). It is a two-step iterative algorithm. In every iteration, it first fixes Z and solves for W, and then it fixes W and solves for Z. Fixing one matrix at a time, the problem is reduced to a linear regression and a simple least squares technique can be used.

I. SVD (Matrix Factorization)

Despite the name, which became common during the Netflix Prize, this model does not utilize a Singular Value Decomposition. Instead it is a matrix factorization with SGD as before, but it includes the learning of biases in the least square problem. This allows to take into account all those inclinations of users and items which are not easily accounted for by the matrix features (for example famous/unknown movies and enthusiastic/strict users). This algorithm is implemented in the Python Surprise library [1].

J. SVD++ (Matrix Factorization With Implicit Ratings)

This model is an extension of SVD, but it takes into account implicit ratings as proposed by Bell & Koren [2], [4]. In a real life scenario, especially in streaming services, lots of information can be considered as implicit ratings such as number of views, continuity and frequency of watching and so on. With our reduced dataset, we can only consider the fact that a user decided to rate a movie as an implicit rating, no matter what the rating is. We are not sure about what meaning to give to this factor, that's why the system automatically learns it for us. The SVD++ algorithm is very powerful, but it is also slow and difficult to tune. That is why we suggest not to use it unless one has more meaningful implicit factors. This algorithm is implemented in the Python Surprise library [1].

K. Slope One

The prediction \hat{r}_{ui} is set as follows.

$$\hat{r}_{ui} = \mu_u + \frac{1}{|R_i(u)|} \sum_{j \in R_i(u)} dev(i, j)$$
 (4)

where $R_i(u)$ is the set of relevant items. The set of relevant items is the set of items j rated by user u that have at least one common user with item i, and dev(i,j) is defined as the average difference between the ratings of i and those of j. That is,

$$dev(i,j) = \frac{1}{|U_{ij}|} \sum_{u \in U_{ij}} (r_{ui} - r_{uj})$$
 (5)

This algorithm is implemented in the Python Surprise library [1].

L. Blending

As explained in [3], the smallest RMSE can be achieved using a combination of different methods instead of an individual predictor. We decided to proceed in the same direction in order to obtain our final model. We approach blending as a linear regression problem. We use the Python module $sklearn.linear_model$ in order to find the weights for a linear combination of the models (referred to as linear blend in [3]). We choose the ridge regression as an estimator in order to avoid overfitting and assigning meaningless weights to the different models. Using a simple

linear regression estimator without regularization gave us huge weights (positive and negative) for the global mean and global median ³ models which however compensate each other, but did not give useful information on the real influence of these models.

For the blending, we split the data in 93.3% and 6.7% (equivalent to 15 folds in a cross validation). Then we use the 93.3% to train the models and the 6.7% to train the ridge regression. To compute the final ratings, we compute the ratings for each model, and then combine them using the weights found by ridge regression.

M. Surprise Library Customization

The Surprise library was customized for performance enhancements. In particular we introduced the following functionalities:

- Shuffling of datapoints during the SGD of SVD and SVD++: In the original code, the order of the examined datapoints was the same at each iteration of the algorithms, depriving the Gradient Descent from the "stochastic" property.
- Scaling of learning rates at each iteration of the gradient descent in SVD and SVD++: This can help finding the minimum, especially in SVD++ where the high number of parameters makes the minimization very sensitive. We chose to scale by a factor of 0.9, as proposed by Koren et al [3].
- Computational speed improvement in SVD++: We compute user fixed information (number of ratings, items rated) once at the beginning of the algorithm, instead of calculating it inside the for loop over all the ratings in the train set (as done by the library). This modification allows to almost double the speed of the algorithm, while maintaining the same results.
- Initialization of baseline in SVD++: We initialize the
 baseline parameters with the ones computed by the
 BaselineOnly model instead of using an array of zeros.
 This initialization can bring some improvement in this
 complex model, where wrong biases can easily make
 the SGD inefficient.
- Optimization of baseline pearson similarity computation: Before estimating the baseline for the similarity matrix, we check if they were already computed. This fix can help especially when doing a grid search over the parameters

Two other completely custom algorithms have been implemented in the Surprise framework: FactorizedNeighborhood and WeightedNeighborhood. They were proposed by Koren et al. during the Netflix Prize [2], [4]. We didn't include these models in the blending because unfortunately we

couldn't obtain a satisfying RMSE when using them. This may be caused by an error in the implementation or by the fact that it is very difficult to find the right parameters to feed them with. Thus, more tests should be performed before proposing a push into the original library. In the next two sections, we provide a brief explanation of the models and some details on their implementations.

N. Weighted Neighborhood

This model was proposed by Bell and Koren during the Netflix Prize [2], [4]. It avoids using a specific similarity measure which could affect the ratings in an unnatural way. On the contrary, it lets the algorithm learn the weights with SGD and introduces new biases, learnt during the iterations, which respond to the new scale of the weights. The function that estimates the prediction is the following.

$$\hat{r}_{ui} = \mu + b_u + b_i + |R^k(i;u)|^{-\frac{1}{2}} \sum_{j \in R^k(i;u)} (r_{uj} - b_{uj}) w_{ij} + c_{ij}$$
(6)

where μ is the global mean, b_{uj} are the baseline estimations computed with the BaselineOnly algorithm, b_u , b_i , w_{ij} , c_{ij} are learnt with simple SGD over the regularized quadratic error (see [2], [4] for more details), w_{ij} corresponds to the weights assigned to the neighbors ratings and c_{ij} takes care of the implicit ratings.

O. Factorized Neighborhood

This model is an extension of the Weighted Neighborhood which also includes the matrix factorization estimation [2]. It somehow corresponds to a blending, but it is done directly during the learning, so it should be more powerful. The parameters are learnt using the linear SGD proposed in the paper, which tries to minimize the regularized quadratic error. The estimations are given by

$$\hat{r}_{ui} = \mu + b_u + b_i + q_i^T (p_u + |R(u)|^{-\frac{1}{2}} \sum_{j \in R(u)} y_j) + |R^k(i; u)|^{-\frac{1}{2}} \sum_{j \in R^k(i; u)} (r_{uj} - b_{uj}) w_{ij} + c_{ij}$$
(7)

Since this algorithm is computationally very heavy, efforts were made in order to optimize the code. The most important step in this direction was achieved using the Cython compiler. In order to make the compiler produce an efficient C code, we tried to observe (as much as possible) some rules while writing the functions critic for the computation time:

³ After the introduction of the regularizer, the weight assigned to the global median was zero, so we decided to exclude it from the models. That's why it does not appear in any table or in the code.

³Cython is an optimizing static compiler for both the Python programming language and the extended Cython programming language (based on Pyrex). The Cython language is very close to the Python language, but Cython additionally supports calling C functions and declaring C types on variables and class attributes. This allows the compiler to generate very efficient C code from Cython code(http://cython.org/).

- For the variables we used Cython fixed types instead of Python dynamic types.
- Indexing of numpy arrays were done without slicing, inside for loops over the single elements.
- We use memoization when possible, avoiding to compute the same quantities more than once.
- Bounds check and wrap around of arrays were disabled.

IV. CONCLUSIONS

In this paper, we showed that combining different methods can improve RMSE of the final prediction. While individual models score an RMSE of at least 0.977, blending achieves an RMSE of around 0.973. The blending improves the score since models can compensate for each other.

We achieved 0.97349 as final score on Kaggle's validation set using the blending technique we described. Further improvements could be possibly attained by either combining more methods, using a different blending technique, or even using the same models but with different parameters. The models are very sensitive to the variation of the parameters, so running a grid search with a very dense refinement for each parameter can improve the model. However the drawback is the high computational time of such a process.

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