

# A novel hybrid data mining framework for credit evaluation

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**Abstract.** Internet loan business has received extensive attentions recently. How to provide lenders with accurate credit scoring profiles of borrowers becomes a challenge due to the tremendous amount of loan requests and the limited information of borrowers. However, existing approaches are not suitable to Internet loan business due to the unique features of individual credit data. In this paper, we propose a unified data mining framework consisting of feature transformation, feature selection and hybrid model to solve the above challenges. Extensive experiment results on realistic datasets show that our proposed framework is an effective solution.

## 1 Introduction

Internet finance has been growing rapidly in China recently. A number of online financial services, such as Wechat Payment and Yu'E Bao have receive extensive attentions. In addition to the payment services, Internet loan business has an explosive growth. On such platforms, borrowers request the loans online. The Internet loan service providers then help borrowers find proper loan agencies. However, it is critical for lenders to obtain the *credit worthiness* of borrowers so that they can minimize the loan risk (to avoid the loans to low credit users).

How to evaluate the *credit worthiness* of borrowers is one of challenges in Internet loan services. In conventional loan markets, banks (or other small firms) usually introduce credit scoring system [4] to obtain the credit worthiness of borrowers. During the credit evaluation procedure, the loan officer carefully checked the loan history of a borrower and evaluated the loan risk based on the officer's past experience (i.e., domain knowledge). However, the conventional credit evaluation procedure cannot be applied to the growing Internet loan markets due to the following reasons. First, the loan officers only have the limited information of borrowers through Internet loan service platform. Second, there are a tremendous amount of requests for Internet loan business every day, which demands the prompt approval (or disapproval) for customers. Thus, the tedious and complicated procedure of convention credit evaluations is no longer suitable for the fast growth of Internet loan business. Third, the conventional credit evaluation heavily depends on the judgment of loan officers. For example, the credit evalu-

ation is often affected by the knowledge, experience and the emotional state of the loan officer. As a result, there may exist misjudgments of loan officers. It is implied in [8] that computer-assisted credit evaluation approaches can help to solve the above concerns.

In fact, to distinguish the credit borrowers is equivalent to classifying all borrowers into two categories: the “good” borrowers who have good credits and are willing to pay their debts plus interest on time, and the “bad” users who may reject to pay their debts on time. Many researchers employ multiple supervised machine learning algorithms to solve the problem, such as Neural Network, Decision Tree and SVM. In particular, Huang et al. [6] utilize Support Vector Machine (SVM) and Neural Networks to conduct a market comparative analysis. Angelini et al. [1] address the credit risk evaluation based on two correlated Neural Network systems. Pang and Gong [9] also apply the C5.0 classification tree to evaluate the credit risk. Besides, Yap et al. [11] use data mining approach to improve assessment of credit worthiness. Moreover, several different methods have been proposed in [10, 12, 5].

Although previous studies exploit various models, there is no unified hybrid model that can integrate the benefits of various models. Besides, the existing models are not suitable for the growing Internet loan business due the following unique features of individual credit data: (i) *high dimension of features*, which can be as large as 1,000 ; (ii) *missing values*, which can significantly affect the classification performance; (iii) *imbalanced samples*, in which there are much more positive samples than negative samples. The above features result in the difficulties in analyzing credit data.

In light of the above challenges, we propose a unified analytical framework. The main contributions of this paper can be summarized as follows.

- We propose a novel hybrid data mining framework, which consists of three key phases: feature transformation, feature selection and hybrid model.
- We integrate various feature engineering methods, feature transformation procedures and supervised learning algorithms in our framework to maximize their advantages.
- We conduct extensive experiments on realistic data sets to evaluate the performance of our proposed model. The comparative results show that our proposed model has the better performance in terms of classification accuracy than other existing methods.

The remaining paper is organized as follows. We describe our proposed framework in Section 2. Section 3 shows experimental results. Finally, we conclude this paper in Section 4.

## 2 Our framework

In order to address the aforementioned concerns, we propose a hybrid data mining framework for credit scoring. As shown in Fig. 1, our framework consists

of three key phases: feature transformation, feature selection, hybrid model. We then describe the three phases in detail in the following sections.

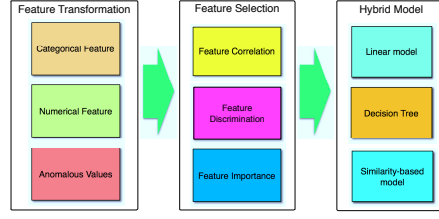


Fig. 1. Our proposed hybrid data mining framework consists of three phases

### 2.1 Feature Transformation

We categorize the features into two types: (i) *numerical* features are continuous real numbers, representing borrower’s age, height, deposit, income, *etc.*; (ii) *categorical* features are discrete integers, indicating borrower’s sex, educational background, race, *etc.* Since the two kinds of features cannot be treated as the same, we conduct a conversion so that they can be fit into an unified model.

**Categorical Feature Transformation** Regarding to categorical features, we exploit a simple *one-hot encoding*. For example, we use a four-bit one-hot binary to represent four seasons in a year. Specifically, ‘1000’, ‘0100’, ‘0010’ and ‘0000’ denote spring, summer, autumn and winter, respectively. The one-hot encoding conversion is intuitive and easy to be implemented. It converts a categorical feature with the unknown range into multiple binary features with value 0 or 1.

**Numerical Feature Transformation** The range of numerical features may differ vastly. For instance, the age is normally ranging from 1 to 100 while the deposit may vary from several hundred to several millions. We utilize the following mapping functions on original features and replace them with the mapped values so that we can reduce the differences between features.

$$Normalize(x_k) = \frac{x_k - mean}{std}, \tag{1}$$

$$Sigmoid(x_k) = \frac{1}{1 + e^{-x_k}}, \tag{2}$$

$$Tanh(x_k) = \frac{e^{x_k} - e^{-x_k}}{e^{x_k} + e^{-x_k}}, \tag{3}$$

$$Maxmin(x_k) = \frac{x_k - \min}{\max - \min}, \tag{4}$$

$$LogAll(x_k) = \log(x_k - \min + 1), \tag{5}$$

where  $x_k = \{x_k^{(1)}, x_k^{(2)}, \dots, x_k^{(n)}\}$  is a set of feature values indicating the  $k$ th dimension of the dataset,  $x_k^{(i)}$  indicates its value for the  $i$ th sample, *mean* denotes the mean value, *std* represents the standard deviation of  $x_k$ , *max* denotes the

maximum and min denotes the minimum value. Note that the above basic mapping functions can be nested. For example, a feature can be first transformed by *LogAll* function and can then be mapped into range (0, 1) by *Sigmoid* function.

**Anomalous Values Handling** Data sets may contain some values deviated from normal values (i.e., outliers) and some missing values. Specifically, we distinguish outliers by Eq. (6) according to the “3 sigma rules” in classical statistics:

$$Outlier(x_k^{(i)}) = \begin{cases} True, & \text{if } |x_k^{(i)} - mean| \geq 4 \times std \\ False, & \text{otherwise} \end{cases} . \quad (6)$$

Depending on the fraction of anomalous values in feature  $x_k$ , we first define the anomalous factor  $f = \frac{N_{missing} + N_{outlier}}{N_{sample}}$ , where  $N_{missing}$  represents the number of missing values,  $N_{outlier}$  denotes the number of outliers, and  $N_{sample}$  is the number of samples. We then propose three different methods to handle the outliers and the missing values: *replace*, *delete*, and *convert* based on different values of anomalous factor  $f$ ,

$$Anomalous(x_k) = \begin{cases} Replace, & \text{if } f \leq \alpha \\ Delete, & \text{if } f \geq \beta \\ Convert, & \text{otherwise} \end{cases} . \quad (7)$$

**Extra Feature Extraction** We also apply statistical methods to extract extra features. Specifically, we construct ranking features from numerical features and percentage features from categorical features. If the value of the  $k$ th numerical feature for the  $i$ th sample is  $x_k^{(i)}$ , the value of ranking feature for it is  $a_k^{(i)} = r_k^{(i)}$ , where  $r_k^{(i)}$  represents  $x_k^{(i)}$ 's ranking in  $x_k$ . However, this simple extension of numerical features significantly increases the dimension, which leads to the extra computational cost. To solve the problem, we use percentiles of the expanded features to represent them in a more concise way. If the extra features are  $A = \{a_1, a_2, \dots, a_n\}$ , we use 0th, 20th, 40th, 60th, 80th and 100th percentiles of  $A$  as final numerical extra features, which can be represented as  $e^{num} = \{a_{0\%}, a_{20\%}, a_{40\%}, a_{60\%}, a_{80\%}, a_{100\%}\}$ .

We use a similar method to obtain extra features from categorical features. Suppose  $x_k^{(i)}$  represent the  $k$ th categorical feature for the  $i$ th sample, the value of extra feature for it is  $b_k^{(i)} = p_k^{(i)}$ , where  $p_k^{(i)}$  represents the percentage of category  $b_k^{(i)}$  in  $x_k$ . If the extra categorical features are  $B = \{b_1, b_2, \dots, b_m\}$ , we use 0th, 20th, 40th, 60th, 80th and 100th percentiles of  $B$  as final categorical features as  $e^{cat} = \{b_{0\%}, b_{20\%}, b_{40\%}, b_{60\%}, b_{80\%}, b_{100\%}\}$ .

After feature conversion, each  $x_k^{(i)}$  is within the same range, we then use statistics to describe them to capture a high level information  $e^{sat} = \{mean, std, perc\}$ , where *mean*, *std* and *perc* represent the mean value, the standard deviation of  $x^{(i)}$  and the percentage of missing values in  $x^{(i)}$ , respectively.

## 2.2 Feature Selection

After the feature transformation, the dimension of features can be significantly increased (e.g., 3,000 in our testing datasets), which lead to the high compu-

tational complexity. Thus, it is crucial for us to select the most important and informative features to train a good model. In this paper, we combine three different feature selection techniques to extract the most useful features.

**Feature Correlation** If two features are correlated to each other, it implies that they convey the same information. Therefore, we can safely remove one of them. Consider an example that a person who has the higher income will pay the more tax. So, we can remove the tax feature and only keep the income feature during model training. There are many methods to measure the correlation (or similarity) between features. In this paper, we use the Pearson Correlation Coefficient (PCC), which is calculated by Eq. (8).

$$r_{xy} = \frac{\sum_{i=1}^n (x_i - \bar{x})(y_i - \bar{y})}{\sqrt{\sum_{i=1}^n (x_i - \bar{x})^2} \sqrt{\sum_{i=1}^n (y_i - \bar{y})^2}}, \quad (8)$$

where  $x = x_1, x_2, \dots, x_n$  and  $y = y_1, y_2, \dots, y_n$  represent two features,  $x_i$  and  $y_i$  denote the corresponding values for features  $x$  and  $y$  in the  $i$ th sample, and  $\bar{x}$  and  $\bar{y}$  denote the means for  $x$  and  $y$ , respectively. In practice, for the feature pairs whose  $r_{xy}$  is higher than 0.95, we arbitrarily remove one of them.

**Feature Discrimination** In model training, our goal is to discriminate different categories based on feature information. If a feature itself can distinguish positive and negative samples, implying that it has a strong correlation with the label, we shall include it in model training since it is an informative feature. For instance, F-score [3] is a simple technique to measure the discrimination of two sets of real numbers. Specifically, F-score is calculated by Eq. (9) as follows,

$$\frac{(\bar{x}^+ - \bar{x})^2 + (\bar{x}^- - \bar{x})^2}{\frac{1}{n^+ - 1} \sum_{k=1}^{n^+} (x_k^+ - \bar{x}^+)^2 + \frac{1}{n^- - 1} \sum_{k=1}^{n^-} (x_k^- - \bar{x}^-)^2}, \quad (9)$$

where  $\bar{x}$ ,  $\bar{x}^+$ ,  $\bar{x}^-$  are the average values of the whole sets, the positive and negative data sets, respectively,  $x_k^+$  is the  $k$ th positive instance and  $x_k^-$  is the  $k$ th negative instance. The larger F-score is, the more likely feature  $x$  is more discriminative.

**Feature Importance** Before applying hybrid model training in Section 2.3, we need to evaluate the importance of every feature in training set. Specifically, we choose the features that contribute the most to our model. After each training, we assign a certain importance value  $v_k$  to each feature  $x_k$ . Taking all information into consideration, we use Eq. (10) to calculate Feature Importance index (FI),

$$FI_k = 0.6 \times v_k^{gbd\text{t}} + 0.2 \times v_k^{r\text{f}} + 0.2 \times f_k, \quad (10)$$

where  $v_k^{gbd\text{t}}$  and  $v_k^{r\text{f}}$  represent importance values given by Gradient Boosting Decision Tree (GBDT) and Random Forest (RF), respectively and  $f_k$  denotes F-score of  $x_k$ . Since  $v_k^{gbd\text{t}}$ ,  $v_k^{r\text{f}}$  and  $f_k$  may not be within the same range, we use function *Maxmin* defined in Eq. (4) to normalize them first.

**Summary** We illustrate the whole feature selection procedure in Algorithm 1. In particular, after conducting feature transformation, we first remove features with large number of anomalous values. Then, we remove the highly correlated feature. Finally, we calculate Feature Importance Index and select the top  $K$  features based on the trained RF and GBDT values.

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**Algorithm 1** Feature Selection
 

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**Require:** a set of features  $X = \{x_1, x_2, \dots, x_n\}$ , selection threshold  $K$

**Ensure:** a subset of  $X$

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1: for each  $x_k$  in  $X$  do
2:   conduct feature transformation
3:   if  $Anomalous(x_k) == Delete$  then
4:     delete  $x_k$  from  $X$ 
5:   end if
6: end for
7: for each feature pairs  $(x_i, x_j)$  in  $X$  do
8:   calculate correlation  $r_{ij}$ 
9:   if  $r_{ij} \geq 0.95$  then
10:    delete  $x_i$  or  $x_j$  from  $x$ 
11:   end if
12: end for
13: train a RF and a GBDT with  $X$ 
14: for each  $x_k$  in  $X$  do
15:   obtain feature importance  $v_k^{rf}$  and  $v_k^{gbd}$ 
16:   calculate F-score  $f_k$ 
17: end for
18:  $v_k^{rf} \leftarrow Maxmin(v_k^{rf})$ 
19:  $v_k^{gbd} \leftarrow Maxmin(v_k^{gbd})$ 
20:  $f_k \leftarrow Maxmin(f_k)$ 
21:  $FI_k \leftarrow 0.6 \times v_k^{gbd} + 0.2 \times v_k^{rf} + 0.2 \times f_k$ 
22:  $X' \leftarrow K$  feature  $x_k$  in  $X$  with largest  $FI_k$ 
23: return  $X'$ 

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### 2.3 Hybrid Model

We first present the models that we use as follows:

- **Linear model.** To reduce the generalization error, we train 10 different Logistic Regression (LR) models with various parameters and blend their results.
- **Decision Tree model.** Gradient Boosting Decision Tree (XGBoost) [2] is a popular scalable Gradient Boosting approach. Like LR, we train 20 different XGBoost models and blend their prediction results.
- **Similarity-based model.** We use Pearson Correlation Coefficient (PCC) to evaluate the similarity between samples. Due the imbalance of samples, we identify the negative samples as many as possible. Therefore, we compare each sample in the test set with each negative sample in the training set and label those with high similarity as negative.

We then describe our proposed hybrid model. In particular, our model exploits one of ensemble classification algorithms - “bagging” [5, 7]. More specifically, we average the predictions from various models. With regard to a single model (e.g., LR), we average predictions of the same model with different parameters. We then average results from LR and XGBoost. The bagging method often

**Table 1.** Performance of AUROC on Different Methods during Different Phases

Factor		Classifier					
Data	Dimension	LR	RF	AdaBoost	GBDT	XGBoost	LR+XGBoost
Original	1138	0.6442 ± 0.0170	0.6549 ± 0.0313	0.6625 ± 0.0335	0.6574 ± 0.0350	0.6988 ± 0.0058	0.7048 ± 0.0065
Extended	1984	0.6566 ± 0.0235	0.6558 ± 0.0314	0.6624 ± 0.0319	0.6539 ± 0.0431	0.7000 ± 0.0060	0.7103 ± 0.0058
Refilled	1984	0.6755 ± 0.0198	0.6573 ± 0.0356	0.6649 ± 0.0329	0.6548 ± 0.0432	0.7025 ± 0.0059	0.7127 ± 0.0027
Selected	200	0.7025 ± 0.0257	0.6635 ± 0.0362	0.6772 ± 0.0291	0.6722 ± 0.0362	0.7200 ± 0.0049	0.7248 ± 0.0053

reduces overfit and smooths the separation board-line between classes. Besides, we also use PCC to identify the samples that are most likely to be negative.

Our hybrid model has a better performance than traditional single model due to the following reasons. Firstly, we exploit a diversity of models and the “bagging” method combines their results together so that their advantages are maximized and their generalization errors are minimized. Secondly, we utilize XGBoost library, which is an excellent implementation of Gradient Boosting algorithm, which is highly efficient and can prevent model from over-fitting.

### 3 Experiment

We use the sample dataset from CashBus<sup>1</sup>, a micro credit company in China, which offers the online loaning service to individuals. The dataset contains 15,000 samples. Since all the samples are anonymous in order to protect user privacy, we cannot use any domain knowledge in problem analysis. The the dataset has the following features: (1) **High Dimension.** The dataset contains 1,138 features, including 1,045 numerical features and 93 categorical features. (2) **Missing values.** There are a total of 1,333,597 missing values in our dataset, making the missing rate 7.81%. The number of missing values for each feature is ranging from 19 to 14,517 and the number of missing values for each sample is ranging from 10 to 1,050. (3) **Imbalanced samples.** There are 13,458 positive samples while only 1,532 negative samples in the dataset.

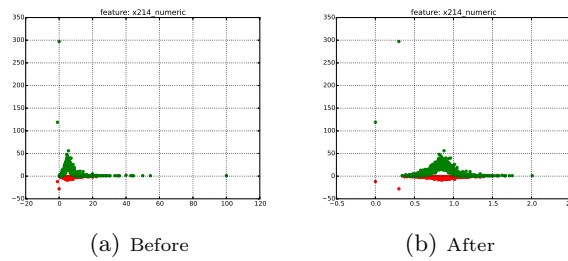
#### 3.1 Experimental Setup and Evaluation Metrics

We predict the probability that a user has a good credit and evaluate the prediction results by Area under the Receiver Operating Characteristic curve (AUROC), i.e.,  $AUROC = \frac{\sum_i S_i}{|P| \times |N|}$ , where  $P$  and  $N$  represent the positive samples and the negative samples in test set, respectively, and  $S_i$  is the score for the  $i$ th pairs between each positive sample and each negative sample, defined by

$$S_i = \begin{cases} 1, & score_{i-p} > score_{i-n} \\ 0.5, & score_{i-p} = score_{i-n} \\ 0, & score_{i-p} < score_{i-n}, \end{cases} \quad (11)$$

where  $score_{i-p}$  and  $score_{i-n}$  represent the scores for the positive and the negative sample, respectively. A higher value of AUROC means that the prediction result is more precise.

<sup>1</sup> <http://www.cashbus.com/>



**Fig. 2.** Impact of LogAll Transformation, where green points and red points represent positive and negative samples, respectively.

### 3.2 Performance Comparison

To investigate the prediction performance, we compare our proposed hybrid model (LR+XGBoost) with other five approaches (each with single model): Logistic Regression (LR), Random Forest (RF), AdaBoost, Gradient Boosting Decision Tree (GBDT) and XGBoost. Table 1 presents the comparative results of different models in different phases in terms of AUROC. Origin represents the raw data set. Extended represents feature transformation. Refilled represents the anomalous values handling process, where we set  $\alpha = 0.1$  to choose feature to refill them. Selected represents feature selection process, where we select the top  $K = 200$  features. We have the following observations: (1) in all four phases, our proposed hybrid model obtains a better AUROC score than any other methods; (2) our proposed model has a relatively small variation compared with other models, implying the stable performance; (3) LR + XGBoost outperform others, indicating that they are the right choices for constructing the hybrid model.

### 3.3 Impact of Feature Transformation

We then investigate the impact of feature transformation. Fig. 2 shows the impact of *LogAll* function on one numerical feature. After the feature transformation, the distribution of the features becomes more smooth and the extremely large values are minimized.

### 3.4 Impact of Refilling Anomalous Values

To deal with the large amount of missing values and outliers, we propose a method to refill the anomalous values based on anomalous value rate under  $\alpha$ . We set  $\alpha$  to be 0.02 to 0.6 to investigate the impact of  $\alpha$ . Table 2 presents the results, where Fill Features represent the number of features that are affected during this process. It is shown in Table 2 that AUROC values for both LR and XGBoost models first increase and then slowly decrease. This can be explained by the fact that filling the anomalous values can bring more information while too many extra filled values also cause noise. In fact, the best performance is obtained when  $\alpha = 0.1$ .



**Table 2.** Performance of AUROC on LR and XGBoost under Different Fill Criterion

Factor		Classifier	
Criterion $\alpha$	Fill Features	LR	XGBoost
0.02	661	0.6560 $\pm$ 0.0243	0.6991 $\pm$ 0.0063
0.05	883	0.6563 $\pm$ 0.0240	0.7001 $\pm$ 0.0062
0.1	881	0.6755 $\pm$ 0.0198	0.7025 $\pm$ 0.0059
0.2	885	0.6567 $\pm$ 0.0234	0.7021 $\pm$ 0.0053
0.3	905	0.6567 $\pm$ 0.0227	0.7008 $\pm$ 0.0054
0.4	1018	0.6561 $\pm$ 0.0231	0.7016 $\pm$ 0.0052
0.5	1024	0.6562 $\pm$ 0.0238	0.7015 $\pm$ 0.0053
0.6	1026	0.6558 $\pm$ 0.0239	0.7013 $\pm$ 0.0072

### 3.5 Impact of Feature Selection

After feature transformation, the dimension of features is significantly increased due to the introduction of extra features. We then exploit the feature selection algorithm to reduce the dimension of features. Specifically, we investigate the impact of the feature importance values given by different models and we set the feature selection threshold to be Top  $K = 200$ . It is shown in Table 3 that our proposed LR+XGBoost achieve the best performance.

**Table 3.** Performance of Different Models under Different Feature Selection Methods

Factor		Classifier					
Importance Calculation	Dimensions	LR	RF	AdaBoost	GBDT	XGBoost	LR+XGBoost
RF	200	0.6818 $\pm$ 0.0168	0.6626 $\pm$ 0.0372	0.6656 $\pm$ 0.0359	0.6602 $\pm$ 0.0350	0.6888 $\pm$ 0.0073	0.6927 $\pm$ 0.0423
XGBoost	200	0.6995 $\pm$ 0.0263	0.6732 $\pm$ 0.0267	0.6695 $\pm$ 0.0367	0.6641 $\pm$ 0.0333	0.7195 $\pm$ 0.0051	0.7214 $\pm$ 0.0031
FSore	200	0.6716 $\pm$ 0.0193	0.6586 $\pm$ 0.0373	0.6608 $\pm$ 0.0367	0.6572 $\pm$ 0.0454	0.6701 $\pm$ 0.0133	0.6843 $\pm$ 0.0245
Ensemble	200	0.7025 $\pm$ 0.0257	0.6635 $\pm$ 0.0362	0.6772 $\pm$ 0.0291	0.6722 $\pm$ 0.0362	0.7200 $\pm$ 0.0049	0.7248 $\pm$ 0.0053

### 3.6 Impact of Selection Threshold

In addition to the feature importance, the threshold top  $K$  also contributes to the final quality of the selected features. To investigate the impact of  $K$ , we set  $K$  to be 100 to 1500 and conduct experiments based on LR and XGBoost models. It is shown in Table 4 that AUROC values first increase and then slowly decrease as  $K$  increases. The best performance is obtained when  $K = 200$ .

## 4 Conclusion and Future Work

In this paper, we propose a novel hybrid data mining framework for individual credit evaluation. To address the challenging issues in individual credit data, such as the high dimension, the outliers and imbalanced samples, we exploit various feature engineering methods and supervised learning models to establish a unified framework. The extensive experimental results show that our proposed framework has a better classification accuracy than other existing methods. There are several future directions in this promising area. For example, we can apply the unsupervised algorithms to utilize the unlabeled data. Besides, we shall use the domain knowledge in finance to further improve the feature transformation and the feature selection procedure.

**Table 4.** Performance of LR and XGBoost under Different Thresholds

Factor Threshold	Classifier	
	LR	XGBoost
100	0.7019 ± 0.0235	0.7156 ± 0.0035
200	0.7025 ± 0.0257	0.7200 ± 0.0049
300	0.6996 ± 0.0184	0.7176 ± 0.0048
400	0.6912 ± 0.0217	0.7129 ± 0.0031
500	0.6853 ± 0.0186	0.7092 ± 0.0038
600	0.6810 ± 0.0157	0.7043 ± 0.0062
800	0.6775 ± 0.0174	0.7032 ± 0.0061
1200	0.6748 ± 0.0237	0.7003 ± 0.0051
1500	0.6721 ± 0.0246	0.7020 ± 0.0060

## Acknowledgment

The work described in this paper was supported by the National Key Research and Development Program (2016YFB1000101), the National Natural Science Foundation of China under (61472338) and the Fundamental Research Funds for the Central Universities.

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