Instance Selection and Optimization of Neural Networks

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performances of automatic credit rating are always hindered by imbalanced datasets, complex attributes and shortage of the

chosen data mining algorithms.

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Abstract— Credit scoring is an important tool in financial institutions, which can be used in credit granting decision. Credit applications are marked by credit scoring models and those with high marks will be treated as "good", while those with low marks will be regarded as "bad". As data mining technique develops, automatic credit scoring systems are warmly welcomed for their high efficiency and objective judgments. Many machine learning algorithms have been applied in training credit scoring models, and ANN is one of them with good performance. This paper presents a higher accuracy credit scoring model based on MLP neural networks trained with back propagation algorithm. Our work focuses on enhancing credit scoring models in three aspects: optimize data distribution in datasets using a new method called Average Random Choosing; compare effects of training-validation-test instances numbers; and find the most suitable number of hidden units. Another contribution of this paper is summarizing the tendency of scoring accuracy of models when the number of hidden units increases. The experiment results show that our methods can achieve high credit scoring accuracy with imbalanced datasets. Thus, credit granting decision can be made by data mining methods using MLP neural networks.

Keywords— Back propagation; credit scoring; multilayer perceptron; neural network

I. INTRODUCTION

Credit scoring, or credit rating, has been widely used in banks and other financial institutes. It is the set of decision models and their underlying techniques that help lenders judge whether an application of credit should be approved or rejected [1, 2]. Currently credit rating can be divided into two general categories: new credit application judgement and prediction of bankrupt after lending. The first kind uses personal information and financial status of a loan applicant as inputs to calculate a score. If the score is higher than a "safe level", the applicant has high possibility to preform good credit behaviour. On the contrary, a low score indicates high risk loan so the lender needs to take careful consideration of the application. The other kind of credit scoring focuses on the credit record of existing customers. This paper focuses on application scoring.

Compared with traditional credit scoring which is achieved by professional bank managers, automatic scoring has some obvious advantages: it saves costs and time for evaluating new credit applications; it is consistent and objective [3]. However, some current computational approaches are not as capable as experienced loan experts on judgement accuracy. The

In recent years, artificial intelligence has shown its advantages in credit scoring comparing with linear probability models, discriminant analysis and other statistical techniques [4]. Among those artificial intelligent models, Multilayer perception (MLP) models are widely utilized [2, 5, 6] and shown competitive prediction ability against other methods [7, 8]. In [9] back-propagation (BP) algorithm was developed and now has been mostly used in training MLP. Improvements of neural networks include altering the ratios of training and testing datasets, the number of hidden nodes, and the training iterations. A nine learning schemes with different training-tovalidation data ratios was investigated and got the implementation results with the German datasets [10]. They concluded that the learning scheme with 400 cases for training and 600 for validation performed best with an overall accuracy rate of 83.6%. Emotional neural network [11] is a modified BP learning algorithm. It has additional emotional weights that are updated using two additional emotional parameters: anxiety and confidence. When comparing emotional neural networks with conventional networks for credit risk evaluation, experimental results showed that both models were effective, but the emotional models outperformed the conventional ones in decision making speed and accuracy [12].

As reported in [3], RBF, LS-SVM and BP classifiers yield high performances on eight different credit scoring datasets. At the same time LDA and LOG also get good results, which indicate that the linear classifiers can also be available [4]. Several automatic scoring models also prove this result using the German, Australian and Japanese credit datasets [5]. In their experiments, C4.5 decision tree performed a little better than MLP on credit scoring but both of them are with high accuracies. On the contrary, Nearest Neighbor and Naïve Bayes classifiers appeared to be the worst in their tests.

Optimization on the ratios of training and testing datasets, the number of hidden nodes, and the training iterations can have positive effect on refining the performance of MLP. Using the German datasets, an experiment is designed to test the performances of MLPs with various ratios of training-validation instances [6]. In their tests, the learning scheme with 400 cases for training and 600 for validation performed best regarding to the overall accuracy rate. The number of hidden nodes is also considered in their tests, as they conduct their

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tests with MLPs that contains three different numbers of hidden nodes. However, their tests only consider some cases and could not lead to a general conclusion. Emotional neural network [7] has additional emotional weights that are updated using two additional emotional parameters: anxiety and confidence. It is a modified BP learning algorithm. The emotional models, as tested in experiments, outperformed the conventional ones in both speed and accuracy [8]. Artificial meta plasticity MLP is efficient when fewer patterns of a class are available or when information inherent to low probability events is crucial for a successful application. The accuracy of this model is 84.67% for the German dataset, and 92.75% for the Australian dataset [9].

In [10], a two-stage hybrid modelling procedure with ANN and multivariate adaptive regression splines (MARS) is designed. After using MARS in building the credit scoring model, the obtained significant variables then served as the input nodes of ANN. However, ensemble system performed better only in one of the three datasets in the experiments of [11]. MLP is compared with multiple classifiers and classifier ensembles. Tests demonstrate that the ability of hybrid system is not better than usual methods.

As to other methods, support vector machine (SVM) and genetic algorithm (GA) are also used for credit rating with good performance. SVM model can be refined by reduction of features using F score and took a sample instead of a whole dataset to create the credit scoring model [12]. Test results showed that this method was competitive in the view of accuracy as well as computational time. In [13], they selected important variables by GA to combine bank's internal behavioral rating model and an external credit bureau model. This dual scoring model underwent more accurate risk judgment and segmentation to further discover the parts which were required to be enhanced in management or control from mortgage portfolio. Other than SVM and GA, Clustering-Launched Classification (CLC) is also available and may perform better than SVM [14]. A multi-criteria quadratic programming (MCQP) model was proposed based on the idea of maximizing external distance between groups and minimizing internal distances within a certain group [15]. It could solve linear equations to find a global optimal solution and obtained the classifier and at the same time used kernel functions to solve nonlinear problems. Comparing with SVM it seemed more accurate and scalable to massive problems. Decision tree (DT) is another good alternative method.

Recently, imbalanced datasets (where instances belonging to one class heavily outnumber instances in other classes) have attracted attention and some work has shown that appropriate ratios of different kinds of instances can augment classification accuracy. In [15], they used five real-world datasets to test the effect of good/bad credit instance ratio. Results showed that linear discriminant analysis (LDA) and logistic regression (LOG) performed acceptable rating accuracy with both slightly imbalanced datasets and highly imbalanced ones. To avoid the effect of imbalance data distribution, dynamic classifier and dynamic ensemble selection of features were added in the scoring model, which performed better than static classifiers [16].

Imbalance datasets can also bring bias to the evaluation of credit rating models. Traditionally, the performance of credit rating is evaluated by average accuracy. However, this value is highly susceptible by the imbalance number of instances, if one kind of instances is highly outnumbered the other. For example, if the credit rating model decided that every instance is "good", and the ratio of good and bad instances is 9:1, then the accuracy of this "silly" credit rating model will be 90%. This is obviously unacceptable. In this paper, both accuracy and another evaluation model called Area Under Curve (AUC) will be used in experiments and the details of AUC will be introduced.

In general, neural network with back propagation algorithm can score credit applications with high performance. However, there are still some problems for this model. Firstly, as the ratios of approved and rejected instances in the datasets are usually not balanced and sometimes highly imbalanced, the training process of a model may have to deal with more approved instances than rejected ones. This may cause bad performance when applying the learned model to the test datasets. Secondly, the numbers of instances used in test, validation and training sets are always limited. More data used in training means less used in validation and test, which may result in low accuracy. As all the three parts need as much instances as possible, the distribution of data is vital for the final test results. At last, the number of hidden units affects the accuracy and the time costs of a model. More hidden units may lead to high computing time but insufficient units cannot achieve high accuracy. The appropriate number can only be known by experiments with different models and choose the best according to the test results.

In order to solve these problems, three aspects of the MLP will be discussed in this paper: (1) Optimise the ratio of approved/rejected cases in input instances. (2) Test the effect of different ratios of training-validation-testing data. (3) Improve the structure of MLP network.

II. DATASET

There are many open credit dataset available on the website of UCI Machine Learning Repository [5]. Among those datasets, the German dataset is a real world dataset with 21 features including 20 attributes recording personal information and financial history of applicants. The last feature is labelled as approved (marked as 1) or rejected (marked as 2). This dataset contains 1000 instances, with 700 approved application instances and 300 rejected ones. These instances are presented randomly. Attribute details of this dataset are listed in Table 1. In this table, some attributes are numerical but others are qualitative and hard to be computed in training of neural networks. Thus, a numerical version of the dataset is used in this work. It transforms all qualitative variables to numeric and adds four more attributes.

The German credit dataset is widely used as a benchmark and has many scoring models. In recent years different models have been utilized on this dataset to solve credit scoring problems. The accuracies of some representative models are listed in Table 2.

TABLE I. ATTRIBUTES OF THE GERMEN CREDIT DATASET

Number	Description	Class
attribute 1	Status of existing checking account	qualitative
attribute 2	Duration in month	numerical
attribute 3	Credit history	qualitative
attribute 4	Purpose	qualitative
attribute 5	Credit amount	numerical
attribute 6	Savings account/bonds	qualitative
attribute 7	Present employment since	qualitative
attribute 8	Instalment rate in percentage of	numerical
	disposable income	
attribute 9	Personal status and sex	qualitative
attribute 10	Other debtors / guarantors	qualitative
attribute 11	Present residence since	numerical
attribute 12	Property	qualitative
attribute 13	Age in years	numerical
attribute 14	Other instalment plans	qualitative
attribute 15	Housing	qualitative
attribute 16	Number of existing credits at this bank	numerical
attribute 17	Job	qualitative
attribute 18	Number of people being liable to	numerical
	provide maintenance for	
attribute 19	Telephone	qualitative
attribute 20	Foreign worker	qualitative

TABLE II. ACCURACY OF SOME REPRESENTATIVE MODELS

Article	Scoring Models	Accuracy (%)			
[9]	MLP	84.67			
[16]	Ensemble	82.03			
[17]	LS-SVM	81.9			
[18]	MLP	81.3			
[8]	MLP	81.03			
[12]	SVM	80.42			
[19]	DT	78.52			
[20]	Re-Rx	78.47			
[21]	SVM	78.46			
[22]	Case-based reasoning model	77.4			
[23]	SVM	76.6			
[24]	SVM	75.4			
[25]	SVM	71.8			

Some of the accuracies as listed are average rates in a group of models and others are the best one. The "scoring models" in table 1 are basic models used in experiments and many of them have been improved. From this table we can see that the highest accuracy is $84.67\pm1.5\%$ achieved by [18] using MLP model. This table only includes some results from journal papers published between 2011 and 2013. There were lots of experiments published in previous years but the accuracies were not better.

III. METHEDOLOGY

A. Average Random Choosing Method

In this paper, we designed an average random choosing (ARC) method to select instances from datasets, which is used to improve training effects and model accuracy.

The imbalance of data classes (where instances belonging to one class heavily outnumber instances in other classes) usually exist in credit datasets. Thus, in the training of neural networks there should be more instances of approved applications in order to get a better scoring model. From the point of test, as the original dataset is imbalanced, it is

reasonable to keep the same ratio (approved/failed) in the test dataset.

Another problem of data processing is the ratio of training-validation-test sets. All three sets should have proper amount of instances. Usually, more instances for training can lead to better chance to get a better model. However, as the amount of data is limited, more data used for training means less for validation and test, which will cause unpredictable test performance.

Usually when given a dataset containing both training and test data, the instances that used for training, validation and test are chosen randomly. For balanced dataset, this will cause imbalance distribution slightly and occasionally, which will hardly bring any problems. However, when the dataset is imbalanced, random choosing method will easily aggravate the imbalance problem. For example, if the training instance group is less imbalanced, which seems to be good for training, then the test group will be higher imbalanced and makes the evaluation of credit rating model biased.

To solve these problems, we propose a method called ARC to pre-process data and generate average number of different instanced in each data group such as training and test. Suppose the total number of instances is n, and the ratio of good applications in the dataset is p. Then the numbers of good and bad applications are

Good applications: p * n, Bad applications: (1 - p) * n

Then suppose the ratio of data used in training is t and in validation is v. Then we have

Training data: n * tValidation data: n * vTest data: n * (1 - t - v)

As we want the ratio of good to bad applications stays the same in training data (as in original data), the training, validation and test data can be divided into good cases and bad cases. The flow of processing data is listed in Figure 1.

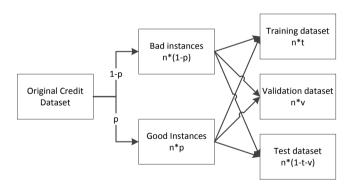


Fig. 1. Flow of data processing and the amount of instances in each group

From the original dataset, two different kinds of data, bad instances and good instances are divided into two groups. Then both of them are divided into training, validation and test datasets randomly. This step should be repeated for each new network training session, which can minimize the effect of

unordinary instances. This way, it is more likely to get a good data distribution which promotes a good scoring model. The algorithm can be described as follows:

Algorithm ARC

Original Dataset $D=\{x_i, t_i | i=0,1...n\}$

- 1. Select out all samples of class A A= $\{x_i, t_i | t_i=1\}$
- 2. Select out all samples of class B B= $\{x_i, t_i | t_i=2\}$
- 3. Select instances for test from class A $T_A=\{x_i, t_i \mid i = randperm(n*p*(1-t-v)), (x_i, t_i) \in A\}$
- 4. Select instances for test from class B $T_B=\{x_i, t_i \mid i=\text{randperm}(n*(1-p)*(1-t-v)), (x_i, t_i) \in B\}$
- 5. The test instances group $T=T_A \cup T_B$
- 6. Samples in class A that are used to train and validate $S_A=A-T_A$
- 7. Samples in class B that are used to train and validate $S_B = B T_B$
- 8. while(not reach the max epoch)
- 9. Choose out samples used for training in class A $S_{\Delta t} = \{x_i, t_i | i = randperm(n*p*t), (x_i, t_i) \in S_{\Delta}\}$
- 10. Choose out samples used for training in class B $S_{Bt}=\{x_i, t_i | i=randperm(n^*(1-p)^*t), (x_i, t_i) \in S_B\}$
- 11. The training group is $S_t = S_{At} + S_{Bt}$
- 12. The validation group is $S_v = (S_A S_{At}) \cup (S_B S_{Bt})$
- 13. Calculate output $y_i = \emptyset(\sum_{s=0}^h w_s \emptyset(\sum_{r=0}^n w_{sr} x_r)), x_r \in S_t$
- 14. Update MLP by BP. Error signal is $e(i)=t_i-y_i$ $t_i \in S_t$
- 15. Validate MLP $y_i = \emptyset(\sum_{s=0}^h w_s \emptyset(\sum_{r=0}^n w_{sr} x_r)), x_r \in S_v$
- 16. Performance of MLP $v = \frac{\sum_{i=0}^{n*v} e(i)}{n*v}$
- 17. $if(v>v_0)$ stop training
- 18. else $v_0=v$
- 19. end while

The function randperm(n) returns n different integer numbers. It is used to randomly select n different samples from an instance group.

By this method, all groups (training, validation and test) have the same ratio of good to bad instances. Traditional 10-fold validation method divides a dataset into 10 blocks of data randomly before training starts. It is not really random choosing and is not fit for imbalanced dataset. In some extreme cases, there could be one group with only one class of data. This is obviously unable to judge the performance of the model. Our average random method chooses instances randomly from both classes of data. It can choose data randomly which ensures more instance combinations can be used for training. Also the datasets of training, validation and test have the same ratio of instances.

B. NN and MLP Scoring Mode

A neural network is a massively parallel distributed processor made up of simple processing units, which has a natural propensity for storing experiential knowledge and making it available for use [26]. It was motivated by inspecting the human brain, which has high efficiency in computing and

recognizing [27]. According to the structure of networks, it can be divided as single layer feed forward networks, multilayer feed forward networks and recurrent networks. Feed forward networks do not have any feedback loop, which are different from recurrent networks.

ANNs are made of neurons, or called simple perceptron. The structure of neuron is shown in Fig 2. Each neuron is composed of two units. First unit adds products of weights coefficients and input signals. The second unit realizes nonlinear function, called neuron activation function. Signal e is adder output signal, and y = f(e) is output signal of nonlinear element. Signal y is also output signal of neuron.

A typical feed forward network is multilayer perceptron (MLP). Other kinds of feed forward networks include single layer networks, radial basis function networks, support vector machine and others. Among them, MLP has better performance when applying on some models, especially for complex models and nonlinear classification problems [28]. This research use MLP neural networks to construct a credit rating model.

A three layer MLP model will be developed in this research. The basic structure is demonstrated in Fig 3. The first layer is the input layer with 24 neurons since the dataset contains 24 input features (attributes). The last layer is the output layer with only one neuron, which stands for the score of an application. Only one hidden layer is used to reduce computing complexity.

The number of hidden neurons can have great impact on the performance of the network. Here we build 34 different MLP models, with the number of hidden units varying from 6 to 39. The model with highest classification accuracy will be chosen out. Our experiment results also prove that this scope is large enough since some of the models can achieve high accuracies.

Neural networks cannot work without "learning". The learning procedure is realized by learning algorithms, which calculate the weights of each neuron in the network. Based on the structure of NN, learning algorithms can be divided into supervised learning and unsupervised learning. In this paper, as the dataset has an outcome attribute, supervised learning will be used to obtain a MLP model for classification. The most popular learning algorithms for feed forward networks include back propagation (BP), radial basis function (RBF) and others [29].

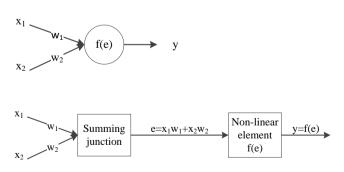


Fig. 2. Structure of neuron

return the network

In this paper, back propagation algorithm will be used in the training of MLP models. It is widely used in the training of MLP and proved to be effective for supervised training. The basic idea of BP is to update the weights of each connection by the difference of the outputs of the model and outputs in datasets. Initial weights will be given randomly in a range and the algorithm will stop if it meets the stopping criteria. It has been proved that the weights can converge in finite iterations. The algorithm of BP utilized on a MLP with one hidden layer can be described as follows:

```
Initialize network weights (often small random values)

Do

for each training example ex

prediction = neural-net-output(network, ex)

// forward pass

actual = teacher-output(ex)

compute error (prediction - actual) at the output units compute \Delta w_h for all weights from hidden layer to output layer // backward pass

compute \Delta w_i for all weights from input layer to hidden layer // backward pass continued update network weights

until all examples classified correctly or another stopping criterion satisfied
```

During the procedure of training, validation is used to control over-fitting. As the iteration goes on, the error rate of the network goes down until it reaches the lowest point. After that point, the error rate will not fall down any more and even rise. This is called over training or over fitting and the training procedure should be stopped when the error rate reaches the lowest point. Here we use some data to validate the network after each iteration. Validation data are not used to train the network, and they are more like the test data. After each iteration, we calculate the error rate with validation data and compare it with the result of the last iteration. If the new one is larger, then the iteration stops immediately. Else, the training and validation process goes on. In this way, we can avoid over fitting.

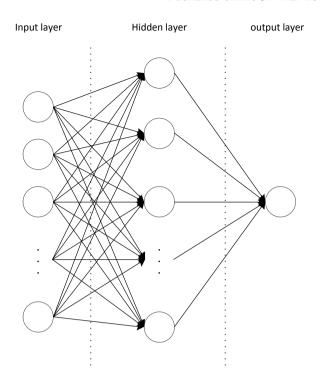


Fig. 3. Structure of a 3 layer MLP model.

C. Area Under Curve

Traditionally, classification accuracy is used to judge the "goodness" of feature group. Recently the Area Under Curve (AUC) have been proved to be a good alternative [30]. It was initially used as an evaluation function in Feature Selection to rank features in [31] and [32]. It was later extended to multiclass in [33] and test results showed that AUC has better performance than traditional overall accuracy (OA).

Define an experiment from P positive instances and N negative instances for some condition. The four outcomes can be formulated in a 2*2 contingency table or confusion matrix, as in table 3 [34].

		Cor			
		Positive	Negative		
				Precision=	
	Positive	True positive	False positive	$\sum Ture\ positve$	
Test		_	(Type 1 error)	Σ Test outcome positive	
outcome	Negative			Negative predictive value=	
		False negative (Type 2 error)	True negative	$\sum True\ negative$	
		(Type 2 ciror)		$\sum Test$ outcome negative	
		Sensitivity=	Specificity=		
		$\sum True\ positive$	$\sum Ture\ negative$	Accuracy	
		\sum Condition positive	\sum Condition negative	Accuracy	

TABLE III. FOUR OUTCOMES OF AN EXPERIMENT

A receiver operating characteristic (ROC) space is defined by false positive rate (FPR) and true positive rate (TPR) as x and y axes respectively, as shown in Fig 4. Write the probability for belonging in the class as a function of a decision/threshold parameter T as $P_1(T)$ and the probability of not belonging to the class as $P_0(T)$. The false positive rate FPR is given by FPR(T) = $\int_T^\infty P_0(T) \, dT$ and the true positive rate is TPR(T) = $\int_T^\infty P_1(T) \, dT$. The ROC curve plots parametrically TPR(T) versus FPR(T) with T as the varying parameter. The area under the curve (AUC) of ROC is equal to the probability that a classifier will rank a randomly chosen positive instance higher than a randomly chosen negative one (assuming 'positive' ranks higher than 'negative') [30].

IV. EXPERIMENTS AND RESULTS

Experiments are designed to test the effective of our new instance selection method and find out a better MLP model. There are three aspects of the experiments. The first aspect is to find out the best amount of data used in training, validation and test, respectively. Then based on the results, the second aspect is to focus on our new instance choosing method. Thirdly, we discuss about the number of hidden neurons by training each model 20 times with different initial weights for each kind of model in all experiments. All instances for training, validation and test are randomly chosen. The best and average error rates are listed and discussed.

A. Ratio of Training-Validation-Test

In the experiments, we use 3 different ratios of data, 800:100:100, 900:50:50 and 600:200:200 to find out the most suitable one. To be more accurate, all groups of data are chosen randomly from the German dataset. The number of hidden neurons varies from 6 to 39, which ensures that every group can get their best model. Also each kind of model is trained 20 times. We record the lowest error rate and average rate of each kind of model. Test results are listed in Table 5.

The result shows that a ratio of 800:100:100 performs better in nearly all aspects in regards to accuracy rate. The lowest error rate, 0.17, is achieved with 15 hidden units which is also the best model of all. The average of error rate can indicate an overall performance of some model groups. The model with 10 hidden units seems more stable, with an average error rate of 0.2295 which is lower than the others. For all models, the average lowest error rate is around 0.208, indicating that it is more likely to get a very low error rate with this training-validation-test ratio.

The ratio of 600:200:200 gives more data to testing, which leads to insufficient data for training. Thus it gets high error

rates in regards to both the lowest value and the average value. Although the ratio 900:50:50 has more instances for training and the lowest error rate is very close to the best one, the shortness of testing data leads to a high average value, which means there is a low possibility to get a good model.

B. Average Random Choosing method

In this section, we compare models trained by data from our Average Random Choosing (ARC) method. Nothing has changed except that we control the percentage of approved/rejected instances in each dataset. The overall percentage is 70% for approved instances and 30% for rejected instances. So this ratio stays the same in training, validation and test data groups. The ratio of training-validation-test is 800:200:200, which performs best in our previous tests. The results are listed on Table 4.

Comparing with the traditional instances divide methods, which randomly divide different kinds of instances in test, validation and training, our new method obviously outperforms the others. As to the AUC, the ARC method performs much better than the other methods, especially the 900:50:50 group, because class imbalance problem is very serious in that group.

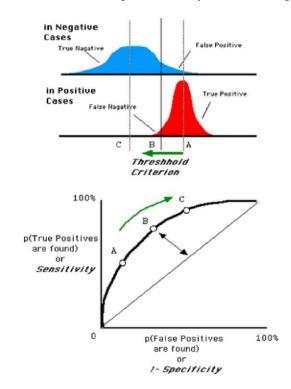


Fig. 4. Area Under Curve (AUC)

TABLE IV. AUCS AND ACCURACIES OF ARC METHOD COMPARING WITH OTHER METHODS

Divide rate	800:100:100)	900:50:50		600:200:200		ARC method	
Evaluation method	AUC	Average Accuracy	AUC	Average Accuracy	AUC	Average Accuracy	AUC	Average Accuracy
9 hidden neurons	0.71	0.762	0.58	0.769	0.69	0.724	0.76	0.80
Best of all	0.73	0.771	0.62	0.769	0.71	0.738	0.76	0.80

As we can see in Table 4, the accuracy may be higher than AUC value. That is because AUC is more objective than accuracy when the dataset is imbalanced. For example, if the MLP classifier output 1 in all case, no matter what the input is. Then the accuracy will be 70% since this class occupies 70% instances in the test dataset. In extreme cases, if there are 90% of instances in one class and 10% in the other, this "dull" classifier will have 90% classification accuracy, while the AUC is only 50%. Thus, if the dataset is imbalanced, AUC should be used as the evaluation method of classifiers.

Test results show that better organized data can enhance model performance especially when the dataset is imbalanced between its classes. As the German dataset is a real world application, our Average Random Choosing method can be easily generalized to handling other datasets. Also, the test results proved that AUC is a more objective evaluation method when the training data set has class imbalance problem. The average accuracy may be biased and cannot reflect the real performance of MLP models.

C. Number Of Hidden Neurons

In Table 5 we also list the result of validation for each kind of model, including the lowest and average rates among the 20 different models. As to the error rates in tests, the lowest ones seem to have been achieved when the number of hidden neuron is 9, or 10, or 12, respectively. However, with these numbers, many other models also get competitive results. When the number of hidden units is 6, 14, 16, 19, 24, 26, 35, 36 and 39, the lowest rate is 0.14 which is only a little higher than 0.13 so it can still be regarded as good models. As to the average rate, the model with 9 hidden neurons gets the lowest one which is 0.2085. But the average of all models is only 0.223, which means that there is little difference between models when only accuracy is considered. So there is no ubiquitous principle for the relationship between the number of hidden neurons and the accuracy of a model.

However, computation time is also very important to neural network research. More hidden neurons can lead to more computation time. Thus, if some models produce the same accuracy, the one with less hidden neurons is preferred. In our experiments, the network with 9 hidden neurons wins out in most cases (Shown in Table 5). So this model is chosen as the most suitable for the German credit dataset in our experiments.

D. Summary

In the first round of experiments, the 800: 100: 100 combinations get the highest accuracy so this ratio is most suitable for building an acceptable model. After that, we use our Average Random Choosing method to optimize the dataset. The ratio of approved/rejected instances in German dataset, 7:3, is precisely implemented in the datasets for training, validation and testing. Results show that our new method can remarkably enhance the accuracy and AUC value, compared with the traditional methods. Finally, models with 9 hidden units perform best among all models, in consideration of high accuracy and low computational time. Also we find an interesting relationship between number of hidden neurons and validation accuracy: the more hidden units a model has, the higher accuracy it may get when validated.

Compared with the results in other relevant articles with the same benchmark dataset, our model achieves a high accuracy of 87%, which is almost higher by 5% than the best result reported in the relevant literature so far. Our best model contains 9 hidden neurons, using our Average Random Choosing method. The ratio of training-validation-test data is 800:100:100. If we take validation results into consideration, the highest accuracy reaches 92%, which is almost 10% higher than the best result from existing models.

It is more interesting when checking the validation error rates. As validation data is also a kind of test data (but with a different purpose), the error rate of validation is proper to reflect the prediction ability. It is always higher than accuracy of test data because a training process will not stop until a validation gets high accuracy. In our experiments, the accuracy of validation can reach almost 0.92 with 38 hidden units in the model. It is 0.05 higher than the best result with test data. Although this value is not as objective as the accuracy from pure test data, it indicates that MLP model has the ability of rating credit application more precisely. Also, there is an interesting tendency found in the validation dataset. As the number of hidden units gets larger, this error rates seems to be lower. It is shown in Fig. 5.

V. CONCLUSIONS

In this paper, we designed a new instance selection method called Average Random Choosing method. It is used to solve highly imbalanced dataset problems. Experiments demonstrate that this method can improve MLP model accuracy in some extent. It could also be used in other dataset and real world problems.

We also designed a MLP model for credit rating with high accuracy compared with other works that utilized the same dataset. Our experiment test 36 models with different number of hidden nodes, and 3 kinds of training with distinctive ratio of training-validation-test instances.

From the results of our experiments, the accuracy of classification has a trend of declining with raising number of hidden nodes. The most valuable contribution of this paper is presenting a MLP model with high credit rating accuracy.

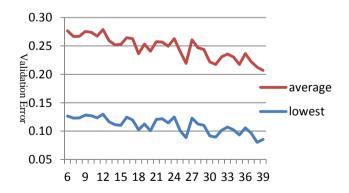


Fig. 5. Tendency of model's error rates when numbers of hidden units increase

Number of hidden neurons

TABLE V. TEST RESULTS OF GERMEN CREDIT DATASET

-	800:1	00:100	900:5	50:50	600:20	00:200	Av	erage Randor	n Choosing metl	Choosing method		
Number of	Lowest								lowest	average		
hidden	error	Average	Lowest	Average	Lowest	Average	Lowest test	Average	validate	validate		
neurons	rate	error rate	error rate	error rate	error rate	error rate	error rate	error rate	error rate	error rate		
6	0.21	0.2415	0.21	0.2925	0.245	0.273	0.14	0.2145	0.1265	0.1502		
7	0.21	0.251	0.2	0.251	0.225	0.273	0.15	0.211	0.1228	0.1441		
8	0.19	0.239	0.21	0.2765	0.22	0.2615	0.17	0.2215	0.1231	0.1438		
9	0.18	0.2385	0.18	0.2315	0.245	0.27675	0.13	0.2085	0.1284	0.1472		
10	0.18	0.2295	0.2	0.243	0.24	0.27075	0.13	0.221	0.1270	0.1471		
11	0.2	0.242	0.18	0.2635	0.225	0.26275	0.16	0.223	0.1234	0.1439		
12	0.21	0.246	0.19	0.2675	0.23	0.26575	0.13	0.225	0.1297	0.1491		
13	0.2	0.256	0.21	0.2705	0.2	0.26625	0.16	0.2225	0.1163	0.1431		
14	0.21	0.258	0.19	0.258	0.205	0.26525	0.14	0.2105	0.1113	0.1407		
15	0.17	0.2595	0.21	0.26	0.24	0.2715	0.16	0.219	0.1103	0.1422		
16	0.22	0.246	0.23	0.26	0.23	0.266	0.14	0.2085	0.1245	0.1399		
17	0.19	0.248	0.21	0.245	0.235	0.2745	0.19	0.2415	0.1196	0.1433		
18	0.21	0.2595	0.21	0.293	0.21	0.26725	0.17	0.221	0.1022	0.1342		
19	0.2	0.255	0.24	0.277	0.235	0.28	0.14	0.2135	0.1124	0.1407		
20	0.21	0.2585	0.21	0.262	0.22	0.2655	0.16	0.2215	0.1006	0.1402		
21	0.22	0.256	0.21	0.2835	0.24	0.2665	0.16	0.215	0.1207	0.1368		
22	0.2	0.2545	0.22	0.2985	0.23	0.26475	0.15	0.2195	0.1217	0.1351		
23	0.19	0.259	0.21	0.2675	0.24	0.27075	0.18	0.2265	0.1143	0.1353		
24	0.19	0.26	0.2	0.26	0.245	0.27375	0.14	0.215	0.1249	0.1377		
25	0.22	0.257	0.21	0.2845	0.24	0.27175	0.15	0.223	0.1007	0.1394		
26	0.22	0.2725	0.19	0.276	0.235	0.28575	0.14	0.224	0.0884	0.1310		
27	0.2	0.261	0.22	0.287	0.24	0.2735	0.17	0.2175	0.1227	0.1380		
28	0.21	0.2625	0.23	0.2755	0.235	0.2755	0.15	0.2235	0.1123	0.1345		
29	0.21	0.262	0.19	0.266	0.225	0.273	0.15	0.2265	0.1100	0.1339		
30	0.23	0.2765	0.23	0.278	0.21	0.2665	0.15	0.2275	0.0916	0.1304		
31	0.24	0.2715	0.23	0.2775	0.25	0.28575	0.18	0.228	0.0892	0.1283		
32	0.22	0.265	0.2	0.2775	0.245	0.272	0.18	0.234	0.1013	0.1295		
33	0.23	0.2755	0.22	0.273	0.245	0.27725	0.17	0.2345	0.1070	0.1287		
34	0.22	0.272	0.22	0.281	0.24	0.28575	0.17	0.228	0.1023	0.1285		
35	0.22	0.2675	0.22	0.2885	0.235	0.28225	0.14	0.229	0.0932	0.1243		
36	0.24	0.2835	0.19	0.285	0.245	0.2895	0.14	0.2275	0.1055	0.1311		
37	0.22	0.278	0.23	0.3175	0.225	0.29125	0.16	0.2265	0.0963	0.1258		
38	0.17	0.28	0.22	0.301	0.21	0.27825	0.19	0.2385	0.0801	0.1324		
39	0.24	0.294	0.2	0.284	0.235	0.284	0.14	0.2375	0.0854	0.1215		
Best	0.17	0.2295	0.18	0.2315	0.2	0.2615	0.13	0.2085	0.0801	0.1215		
Average	0.2082	0.25988	0.20941	0.27389	0.23161	0.27375	0.1553	0.2231	0.1102	0.1368		

Further works may focus on feature selection for neural networks, as attributes in dataset are not always relevant to classification feature. The elimination of redundant features is an important task in data mining. Also it should be noticed that the performance of credit rating model is always evaluated by accuracies. Other statistic methods, such as F-score, standard deviation (SD) and others should be used to help choose out better credit rating models.

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