RISK ANALYSIS SYSTEM FOR INVESTORS

Team 8 - ISTE 780: Data-Driven Knowledge Discovery by Dr. Qi Yu.

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Introduction to Checkpoint 4:

For this project, we worked on deciding the core algorithm to analyze the lending loan risk by classifying available loan data to categorize it as good loans and bad loans. We have considered 5 statistical machine learning models to get the analysis along with their accuracy and efficiency towards the available Lending Loan Club dataset. We worked on Decision trees Classification, KNN Classification, Logistic Regression, and Random Forest Classification. After observing the overall performance of these 4 models, for this instance, we have selected Random Forest Classification as the core algorithm for the system. Further, we worked on the 5th algorithm SVM, which is not covered in the class and tried to fine-tune it to get better outcome. We then compared all algorithms with each other and selected Random Forest Classifier as the core algorithm.

The Goal:

Investment in loan lending business is financially risky without a proper system to analyze the possibility of the existing loans being a good loan or bad loans. The investors should check the historic as well as current statistics of the borrower and deduce the result to invest more money towards improving bad loans or maintaining good loans. For this herculean task, we are proposing this model based on the historic and currently available data to find out the maximum possibility of existing loans becoming a good loan or a bad loan for investors.

Basic Steps of Data Preparation:

The different statistical machine learning models use a specific format of data and the absence of the proper format will lead to failure of the model. So, datasets generally require some amount of preparation before they can yield useful insights. Although our dataset has some missing values and irrelevant values that were taken care of in the previous checkpoint number 2, we have to select the particular set of features for each model. We have taken the measures not to miss out on relevant information for the final goal of the system as well as to not consider the most obvious and correlated attributes while features selection process.

Feature Selection:

After data pre-processing, we left with a cleaned dataset without any missing values or irrelevant values to perform classification models on it. The correlation between features can be found by performing linear regressions on some selective quantitative and qualitative attributes. After observing p-values from the lm() model, we have separated the set of suspiciously insignificant attributes. We will cross-check their importance in further processes of classification to find out if we are missing any important information or not.

On the observation of the remaining dataset, we needed a deciding feature (response) to perform the classification of loans to good or bad. From existing labels of loan status for

each case, we categorized the labels according to their description. This loan status attribute is now a response feature for our classification task. Now, we have to remove highly correlated features and direct dependant of loan status to avoid multicollinearity with loan status and overfitting our model with redundant features. Finally, we have the most important features to deploy classification models on them with loan status as the response.

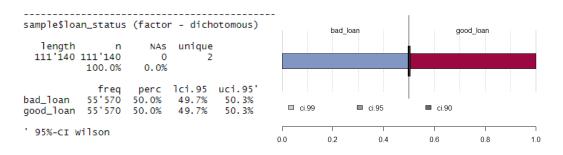
Class Balancing:

Imbalanced data means where the number of observations is not the same for all classes in the dataset. In our dataset, we have an 80:20 ratio of good_loans to bad_loans classes. However, this imbalanced condition is quite common in the real world and we must deal with it while performing a classification model on a dataset. Models like random forest fail to cope with such an imbalanced training dataset as they are sensitive to the proportions of the different classes. If performed on such a dataset, they always favor the majority class which may mislead the accuracies.

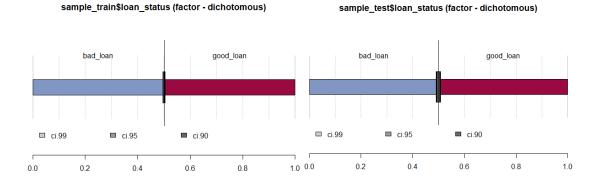
We have randomly selected the 50:50 portion of bad_loans and good_loans with an ample amount (around 100K) of observations in sample data and divide it to 80:20 portion for training and testing. This eliminates the imbalance of the class in the data so that the model will give actual accuracy.

Sample Data:

sample\$loan_status (factor - dichotomous)



Sample_Training Data and Sample_Testing Data



Structure of Sample Data

```
'data.frame': 111140 obs. of 40 variables:
                               : num 15000 6000 6000 31500 21000 ...
: num 15000 6000 6000 31500 21000 ...
 $ loan amnt
                                                      : num 15000 6000 6000 31500 21000 ...
: num 15000 6000 6000 31450 21000 ...
: Factor w/ 2 levels "36 months"," 60 months": 2 1 1 2 1 1 1 1 1 2 ...
: num 15.2 13.7 12.8 20.2 11 ...
: num 359 204 202 838 687 ...
: Factor w/ 7 levels "A","B","C","D",..: 3 3 2 5 2 4 1 5 3 5 ...
: Factor w/ 35 levels "A1","A2","A3",..: 13 14 9 23 7 17 2 24 12 22 ...
: Factor w/ 12 levels "A1","A2","A3",..: 13 14 9 23 7 17 2 24 12 22 ...
: Factor w/ 6 levels "A1","MORTGAGE",..: 2 6 6 6 2 5 2 2 6 5 ...
: num 45800 65000 38000 175000 160000 28000 90000 46000 38000 46000 ...
: Factor w/ 3 levels "Not Verified",..: 2 1 1 2 2 1 1 3 1 3 ...
: Factor w/ 2 levels "bad_loan","good_loan": 2 2 1 1 1 2 1 1 2 2 ...
: Factor w/ 2 levels "n","y": 1 1 1 1 1 1 1 1 1 ...
: Factor w/ 14 levels "car","credit_card",..: 3 8 3 5 3 10 3 3 3 3 ...
: num 11.45 18.08 15.6 6.86 16.85 ...
 $ funded_amnt
 $ funded_amnt_inv
 $ int_rate
 $ installment
 $ grade
 $ sub_grade
emp_length
shome_ownership
annual_inc
verification_status
loan_status
pymnt_nlan
 $ pymnt_plan
 $ purpose
                                                       : num 11.45 18.08 15.6 6.86 16.85 ...

: num 0 1 0 2 0 0 1 0 0 0 ...

: num 0 2 0 0 0 0 0 2 0 1 ...

: num 11 11 8 8 25 8 18 14 5 5 ...
 $ dti
 $ delinq_2yrs
 $ inq_last_6mths
$ open acc
 $ collection_recovery_fee : num 0 0 7.21 0 0 ...
                                                          : num 11604 204 202 838 687 ...
 $ last_pymnt_amnt
 $ collections_12_mths_ex_med: num 0 1 0 0 0 0 0 0 0 ...
$ application_type : Factor w/ 2 levels "INDIVIDUAL", "JOINT": 1 1 1 1 1 1 1 1 1 1 ...
$ acc now deline : num 0 0 0 0 0 0 0 0 0 ...
 $ tot_coll_amt
                                                         : num 0 875 0 0 0 0 0 86 0 0 .
                                                         : num 149643 23160 11712 23575 1088935 ...
 $ tot_cur_bal
 $ total_rev_hi_lim
                                                        : num 16400 33000 13800 10600 63300 19400 44000 20800 18900 18400 ...
```

Models Under Consideration:

1) Logistic Regression:

Logistic regression models the posterior probability that Y belongs to a category. Various models are trained using glm() to perform logistic regressions and p-values are taken into consideration in choosing the significant predictors which helps to reduce feature space and help in interpretability. The results of the best trained model is shown below.

Confusion Matrix:

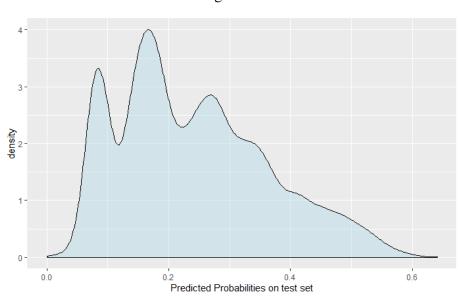
```
glm.probs1
0 1
0 48982 969
1 14924 1132
```

The accuracy obtained is 75.92%. As per confusion matrix, the Specificity of the model is 92.9% and Sensitivity is 98.06%, considering "0" as a positive class. (0 is Good Loan).

The coefficients of the following features are positive: Loan Amount, Interest Rate, Home Ownership - Other.

The coefficients of the following features are negative: Annual Income, Home Ownership - Own, Home Ownership - Rent,

There is no significant difference in the early years of employment. This means that the probability of defaulting is inversely proportional to the factors mentioned above.



Testing Error Plot

2) Decision Tree:

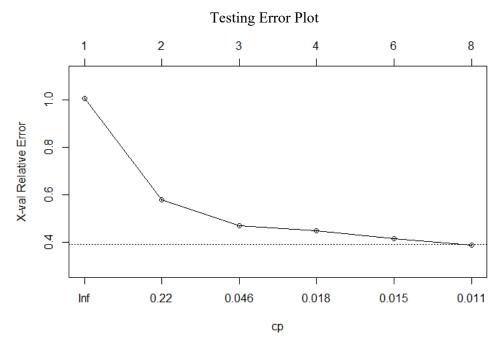
A classification tree is used to predict a qualitative response by assigning each observation to the most commonly occurring class of training observations in the region to which it belongs. The complexity parameter (cp) is used to control the size of the decision tree and to select the optimal tree size acts as tuning parameter for the bias-variance trade-off. If the cost of adding another variable to the decision tree from the current node > cp, then tree building does not continue. We prune the tree to avoid any overfitting of the data.

The x-error is the cross-validation error (R has built-in cross validation). We use rel_error(relative error i.e. training error rate), xerror and xstd together are used to help choose where to prune the tree. The one with least cross-validated error (xerror) is the optimal value of CP.

Confusion Matrix:

tree_pred		
loan_status	bad_loan	good_loan
bad_loan	7007	4095
good_loan	198	10928

The accuracy obtained is 80.68%. As per confusion matrix, the Specificity of the model is 36.8% and Sensitivity is 98.2%, considering "bad_loans" as a positive class.



As can be observed from the model the optimal complexity parameter value cp = 0.011 is chosen as this leads to low test error rate and an accuracy of 80.68%.

3) KNN Classifier:

In theory we would always like to predict qualitative responses using the Bayes classifier. But for real data, we do not know the conditional distribution of Y given X, and so computing the Bayes classifier is impossible. An attempt to estimate the conditional distribution of Y given X, and then classify a given observation to the class with highest estimated probability is the K-nearest neighbours (KNN) classifier.

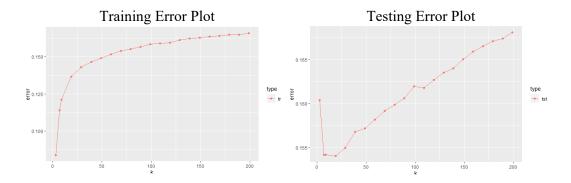
Given a positive in-K and a test observation x0, the KNN classifier first identifies the neighbours K points in the training data that are closest to x0, represented by N0. It then estimates the conditional probability for class j as the fraction of points in N0 whose response values equal j. For K=1 the training error rate is always equal to zero. For low values of K the bias of the model is low and the variance is high and for high values of K the bias of the model is high and variance low. Hence, we need to find a Bias-Variance trade-off by choosing optimal value of K.

Confusion Matrix for KNN with K=19:

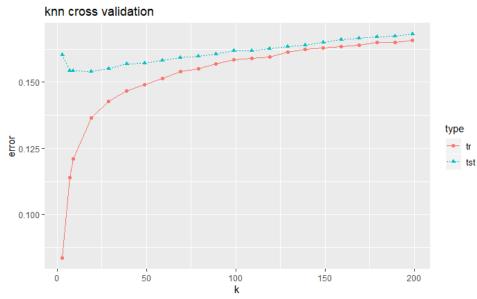
	pred		
sample_test\$status	0	1	Total
0	7493 37.5%	2481 12.4%	9974
1	4526 22.6%	5500 27.5%	10026

The accuracy obtained is 65%. As per confusion matrix, the Specificity of the model is 24.85% and Sensitivity is 54.8%, considering "0" as a positive class. (0 is Good Loan).

Results of K-Fold Cross Validation: 5 folds for KNN with K up to 200.



Comparing Training and Testing Plots:



As can be observed from the model the optimal value of K where the testing error rate is minimum is at K=19 and the accuracy obtained is 65%.

The Core Algorithm: Random Forest

The decision trees suffer from high variance. This means that if we split the training data into two parts at random, and fit a decision tree to both halves, the results that we get could be quite different. Bootstrap aggregation, or bagging, is a general-purpose procedure for reducing the bagging variance of a statistical learning method by averaging a set of observations which reduces variance.

A very straightforward way to estimate the test error of a bagged model, without the need to perform cross-validation or the validation set approach is Out-of-Bag Error Estimation (OOB). Bagging typically results in improved accuracy over prediction using a single tree. Bagging improves prediction accuracy at the expense of interpretability.

Although the collection of bagged trees is much more difficult to interpret than a single tree, we can obtain an overall summary of the importance of each predictor in the context of bagging classification trees, by adding up the total amount the Gini index is decreased by splits over a given predictor, averaged over all B trees. Random forests provide an improvement over bagged trees by decorrelating the trees by choosing a random sample of m predictors as split candidates from the full set of p predictors. The predictions from the bagged trees will be highly correlated. Unfortunately, averaging many highly correlated quantities does not lead to as large of a reduction in variance as averaging many uncorrelated quantities. This means that bagging will not lead to a substantial reduction in variance over a single tree in this setting. Random forests overcome this problem by forcing each split to consider only a subset of the predictors.

Confusion Matrix: Using 100 ntrees.

The accuracy obtained is 83.85%. As per confusion matrix, the Specificity of the model is 5.5% and Sensitivity is 72.96%, considering "bad loan" as a positive class.

Experimentation with number of trees:

Accuracy	Number of trees	Mtry
0.6452672	2	1
0.7857207	20	1
0.8303041	50	2
0.8350279	80	2
0.838582	100	Default
0.8388069	200	Default
0.8393918	300	Default

Fine Tuning of Random Forest:

Number of selected variables is denoted by **mtry** in randomForest() function. We tried to fine tune model by selecting the mtry value with minimum OOB (out of bag) error. We used tuneRF() function from randomForest library of R. After getting best mtry value, mtry = 5, to run random forest with 500 number of trees which resulted in the following confusion matrix.

As can be observed from the model best accuracy is obtained by the random forest where at each split in the tree 5(mtry) variables are chosen from the set of predictors available.

Confusion Matrix for tuned model:

```
Type of random forest: classification Number of trees: 500
No. of variables tried at each split: 5

OOB estimate of error rate: 14.92%

Confusion matrix:

0 1 class.error
0 23798 1228 0.04906897
1 6231 18743 0.24949948
```

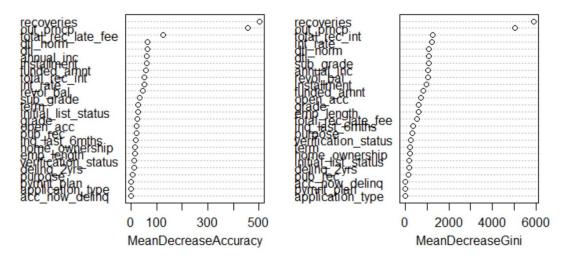
The accuracy obtained is 85.85%. As per confusion matrix, the Specificity of the model is 4.9% and Sensitivity is 75%, considering "0" as a positive class. (0 is Good Loan).

This shows 2% increase in the accuracy while maintaining the balance between specificity and sensitivity.

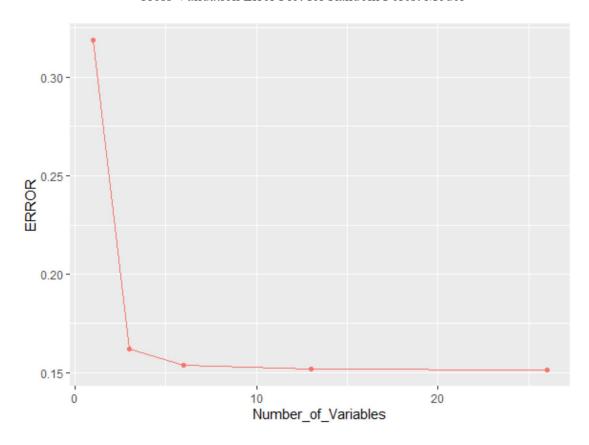
Important Variables:

0 1 MeanDecreaseAccuracy MeanDecreaseGi				MeanDecreaseGini
funded_amnt	59.062974	-9.99808369	59.4375095	8.225606e+02
term	25.220009	17.00610938	28.6329166	2.294903e+02
int_rate	41.187831	15.46472174	51.7064580	1.221226e+03
installment	64.381398	-10.52195514	61.9224445	9.871404e+02
grade	21.399746	10.04322715	23.7746139	6.105185e+02
sub_grade	30.695043	12.34153590	34.8491808	1.065119e+03
emp_length	12.222538	6.82629993	14.3095800	6.056729e+02
home_ownership	16.831934	1.44727322	14.5689830	2.197509e+02
annual_inc	58.399194	13.62109857	62.1901323	1.061248e+03
verification_status	13.661837	0.01781608	11.3422200	2.520511e+02
pymnt_plan	0.000000	0.00000000	0.000000	3.047619e-03
purpose	9.079724	-0.44879015	7.3088611	3.392163e+02
dti	63.128932	-18.66457785	62.9609215	1.072762e+03
delinq_2yrs	14.582744	-3.01136226	10.6742237	1.868868e+02
inq_last_6mths	15.848687	5.43737509	16.6559429	3.697262e+02
open_acc	25.347984	-2.94476260	20.2905139	7.204413e+02
pub_rec	24.207630	-4.83185467	20.1494673	1.527397e+02
revol_bal	55.311118	-16.94684166	44.9572881	1.039432e+03
initial_list_status	28.139929	7.71095230	27.1622863	1.873734e+02
out_prncp	433.859963	411.61494044	454.9763048	5.011962e+03
total_rec_int	31.905949	39.68577183	53.1169468	1.245195e+03
total_rec_late_fee	122.077035	42.61824331	126.4837268	5.466317e+02
recoveries	458.686587	468.48089102	501.3150562	5.880251e+03
application_type	0.000000	0.00000000	0.000000	0.000000e+00
acc_now_delinq	-1.158723	0.70013984	-0.6821246	6.858953e+00
dti_norm	65.187336	-18.12528534	64.3720870	1.078666e+03

Important Variables with sorted order: topmost are most important



Cross Validation Error Plot for Random Forest Model



As we observe from the Cross-Validation Error Plot for Random Forest Model the optimal number of predictors used for building the random forests are p = 26.

Model not covered in the class: Support Vector Machine (SVM)

The support vector machine (SVM) is an extension of the support vector classifier that results from enlarging the feature space in a specific way, using kernels. A kernel is a function that quantifies the similarity of two observations. Support vector machines are intended for the binary classification setting in which there are two classes. But there are the extensions of support vector machines to the case of more than two classes. The tuning parameter C controls the bias-variance trade-off of the statistical learning techniques. It is generally chosen via cross-validation. C bounds the sum of the epsilons and hence determines the severity of the violation of the margin that it will tolerate. The smallest distance from the observations to the separating hyperplane is known as the margin. The separating hyperplane that is farthest from the training observations after computing perpendicular distance from each training observations to the given separating hyperplane.

Observations that lie directly on the margin, or on the wrong side of the margin for their class, are known as support vectors. These observations affect the support vector classifier. When the tuning parameter C is large, then the margin is wide, many observations violate the margin, and so there are many support vectors. The advantage of using SVM is computational cost because using kernels, one need only compute the kernel function for all distinct pairs of inputs. This can be done without explicitly working in the enlarged feature space. For some kernels, such as the radial kernel, the feature space is implicit and infinite-dimensional, hence we can not do all the computations using our limited resources.

Training of SVM:

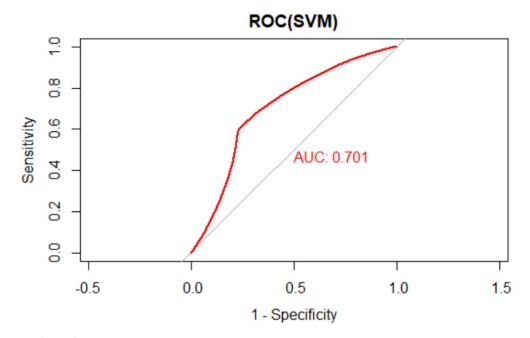
Initially we trained the SVM model using radial kernel with gamma value as 1 and cost as 10.

For SVM, we use Radial Basis as a kernel function. Because of limited computational power, we decided to use only a small subset of the data for training the model. We have sampled out 5% (3500 records) data out of 70000 equally balanced samples (equal number of positive and negative classes) for training the model.

Confusion Matrix:

```
predict_label 0 1
0 19226 7030
1 13975 26269
```

Above model when tested on a test data sample containing about 66000 records, it produced around 64% accuracy with 82 % specificity and 45% sensitivity.



Fine Tuning of SVM:

Later we tried to improvise the accuracy of SVM model by fine-tuning the gamma and cost values. In this step we tried giving cost values as 0.1, 1, 10, 100 and 1000, and gamma values as 0.5, 1, 2, 3 and 4.

```
# Fine-Tune the SVM to improve the accuracy.

tune.out=tune(svm , status~., data=train_data, kernel ="radial",
ranges =list(cost=c(0.1 ,1 ,10 ,100 ,1000),
gamma=c(0.5,1,2,3,4)))

summary(tune.out)
```

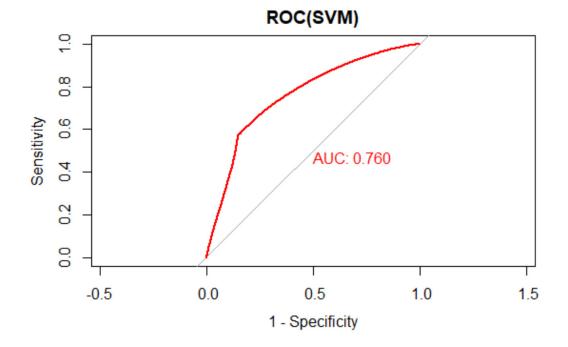
After this training SVM produced the best result at Cost=1 and gamma = 0.5.

Confusion Matrix:

Accuracy	Sensitivity	Specificity
<dbl></dbl>	<dbl></dbl>	<dbl></dbl>
68.367	57.849	78.888

Upon testing on a test data sample containing about 66000 records, it produced around 68.36% accuracy with 78.88% specificity and 57.85% sensitivity.

Hence, we can see about 4% improvement in the accuracy after fine-tuning the model.



CONCLUSION:

After working on fine tuning of all models, we can compare them with each other using outcomes and confusion matrices of them on testing dataset.

Confusion Matrices of Testing:

Logistic Regression	glm. probs1 0 1 0 48982 969 1 14924 1132 0 - good loan and 1 - bad loan	
Decision Tree Classifier	tree_pred loan_status bad_loan good_loan bad_loan 7007 4095 good_loan 198 10928	
KNN Classifier	- 0 1	
	0 7493 2481 37.5% 12.4%	
	1 4526 5500 22.6% 27.5% 0-good loan and 1- bad loan	
Random Forest Classifier	Confusion matrix: 0 1 class.error 0 23798 1228 0.04906897 1 6231 18743 0.24949948 0-good loan and 1- bad loan	
Support Vector Machine	31 0 1 0 19237 7019 1 14017 26227 0- good loan and 1- bad loan	

Accuracies of all models:

Logistic Regression	75.92 %
Decision Tree Classifier	80.68 %
KNN Classifier	65 %
Random Forest Classifier	85.85 %
Support Vector Machine	68.36 %

The specificity of the models suggests that the fraction of times that the model has wrongly predicted bad loans as good loans and having a large specificity for the model may have financial ramifications to the banking company like credit card companies. Logistic regression classifier has the highest specificity of the 5 models analysed and the value is 92.9%. Hence even though SVM classifier has lower accuracy companies would still prefer to use this model due to its low specificity as compared to Logistic regression. KNN has lower specificity than SVM but also low accuracy compared to it. The lowest specificity is obtained by random forest classifier and value is 4.9%.

The sensitivity of the models suggests the fraction of times that the model has correctly predicted bad loans from the total number of bad loans in the dataset. This indicator would still have some economic impact on the financial company but not as bad as wrongly predicting bad loans as good loans.

Ranking of the models based on their accuracies, specificity and sensitivity and assuming the economic impact of wrongly predicting bad loans as good loans is much higher.

1	Random Forest Classifier	85.85 %
2	Decision Tree Classifier	80.68 %
3	KNN Classifier	65 %
4	Support Vector Machine	68.36 %
5	Logistic Regression	75.92%

The KNN Classifier gives lowest accuracy of 65 % even after training and testing on different K values. SVM gives slightly better accuracy than KNN Classifier which is 68.36%. Hence, both SVM and KNN ruled out from consideration of the core algorithm.

Ultimately, models with admissible accuracies to be considered for the candidate of core algorithms are Logistic Regression, Decision Tree and Random Forest. Logistic regression gives 75.92 % accuracy which is obtained after experimentation of interactions and transformations of features to make them significant towards the response. However, Decision tree gives better accuracy than Logistic regression and hence Logistic regression ruled out from the candidates for core algorithm. Decision tree gives 80.68 % accuracy on the testing data after experimentation of pruning the tree and considering cp values for it. The decision trees have limited predictive power due to high variance and even after bagging, it may not be effective on test data due to highly correlated predictors and therefore, in general, the Random Forest model is preferred on the top of Decision trees.

In our case, Decision Tree ruled out all other models but Random Forest. At first Random Forest gave only 83.85 % accuracy which is just 3 % improvisation as compared to the tuned Decision tree. However, after fine tuning the Random Forest model and applying the best mtry value of lowest OOB validation error to it, the accuracy boosted up to 85.85 % on testing data which is very good improvisation over the Decision Tree. Also, the random forest provides default regularization by considering number of decision trees and the significant difference of their accuracies, we can safely say that Random Forest is the stronger and our core algorithm to analyze the loan lending risk to investors by classifying available loan data to categorize it as good loans and bad loans.