Credit Card Fraud Prediction Using XGBoost: An Ensemble Learning Approach

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ABSTRACT
With the development of technology, the internet and eCommerce online payment has become an essential mode of payment. Nowadays, credit card payment is a convenient mode of payment online as well as offline transactions. As online credit card payment increases, fraud transactions are likewise increasing day by day. Increasing fraud transactions in the online payment system became a more significant challenge for banks, companies, and researchers. Therefore, it is essential to have an efficient methodology to detect fraud transactions while payment has completed via credit card. Although many traditional approaches are already available for fraud transaction prediction, existing methods lack accuracy, and it can be increased by ensemble techniques such as XGBoost. In this paper, the authors use an ensemble approach that is XGBoost (eXtreme gradient boosting) for credit card fraud prediction. The results are compared with existing machine learning approaches.

KEYWORDS
Credit Card Fraud Prediction, Decision Tree, Ensemble Learning, Machine Learning, Random Forest, SVM, XGBoost

1. INTRODUCTION
With the exponential growth of technology, the internet, and the electronic market, credit card prediction has become an exciting research topic. Online transactions also play a vital role in the present scenario. Electronic commerce also employs Big data and artificial intelligence techniques to gain profits in their business (Maes et al., 2002; Niu et al., 2019). While most transactions and companies are running online, payment transactions are also online and generate massive data. Online payments are made using credit cards, debit cards, net banking, and UPI options. The online credit card payment system needs a predictive model to distinguish whether a transaction is a fraud or not (Maes et al., 2002). Although numerous modern prediction models are proposed to detect credit card fraud transactions, accuracy is still challenging (Niu et al., 2019; Odegua, 2020). The existing model lacks inefficiency because of the hugeness of transaction data and data imbalance problems. Imbalanced data refers to whether one class has more instances than another class category. It leads...
to the class imbalance problem (Divakar & Chitharanjan, 2019; Zhu et al., 2017). Another Challenge is a significant transaction because a large data set has increased heavy-tailed noise distribution and nonlinear patterns. Therefore, traditional approaches are not able to gain higher accuracy (Petropoulos et al., 2019). The preliminary study found that supervised and unsupervised learning techniques have been used to uncover credit card fraud and forecasting (Randhawa et al., 2018). Chan et al. recommended a cost model for fraud detection (Chan et al., 1999). They have combined multiple fraud detectors in the cost model and demonstrated that the loss had been reduced due to distributed data mining of fraud models. A fraud detection model is suggested by Bolton et al. using user behavior for a credit card transaction (Bolton & Hand, 2001). Zojaji et al. review various techniques, data set, and evaluation criteria for credit card fraud transactions (Zojaji et al., 2016).

Figure 1 shows the process of credit card fraud discovery. This process described those various essential conditions that are checked before completing the transaction. It includes enough balance, card number, expiry valid, PIN details, etc. According to the provided details prediction model, it classifies transactions as fraud or genuine. This paper investigated the analysis of credit card fraud prediction using Gradient Boosting algorithms XGBoost (Chen & Guestrin, 2016). In addition, we perform an extensive comparison of the prediction accuracy of the XGBoost approach with other machine learning approaches. We also presented evaluation matrices, such as accuracy, precision, recall, and F1 score.

Further, the paper is organized into four sections: Section 2 provides a relevant overview of credit card fraud discovery. Section 3 explains the selected models. In section 4, we have discussed our outcomes and discoveries, and section 5 concludes.

2. RELATED WORKS

Credit card fraud prediction is a huge problem for financial markets and banking systems. Various accurate and robust methods have been developed in recent years to identify credit card fraud transactions. Existing models use different statistical techniques, data mining algorithms, machine learning approaches, and deep learning to detect credit card fraud transactions (Randhawa et al., 2018). The linear regression model (Avery et al., 2004) and the probit model (Mizen & Tsoukas, 2012) were earlier used for credit card fraud detection. However, these models are not able to handle the no-linear dynamics of data. Another statistical model is known as the hazard rate model used for credit card rating. These models use the probability concept to detect defaulter customers from loan or portfolio data (Chava & Jarrow, 2004). Huang (2011) proposed a Gaussian process with a support vector machine (SVM) Vapnik, (1999) to detect credit card fraud transactions. This process requires high computation and has high complexity.
Shrivastava et al. (2008) used a Hidden Markov Model to detect credit card fraud transactions from the dataset. SVM and random forest were applied along with logistic regression to detect credit card fraud from international credit card transaction data Bhattacharyya et al., (2011). Decision tree-based methods are also employed by Sachin et al. (2013) to detect credit card fraud from real-life datasets. The problems of Class imbalance, concept drift, and verification latency were considered by Pozzalo et al. for credit card fraud detection Dal Pozzolo et al., (2017).

Bigdata technologies, Casandra, Spark, and Kafka, were used by Carcillo et al. (2018) in scalable real-time fraud finder (SCARFF) to detect fraud transactions. This model works in a real-time scenario and can handle a considerable volume of data. Different machine learning and deep learning methods were used by Addo et al. (2018) for credit risk scoring. They build a binary classification model to predict defaulters from massive data sets.

Six primary commercial bank data from January 2009 to December 2013 are used by Butaru et al. (2016) to detect credit risk. They have used consumer tradeline, credit bureau, and macroeconomic parameters to predict crime using classifiers such as decision tree, logistic regression, and random forest. This result suggested that a more customized approach is needed for financial institutions that include more features for prediction.

In recent years, deep learning approaches have encouraged outcomes in many fields, such as image processing, classification, and predictions Wang et al., (2015). These approaches are also appropriate to predict credit card fraud. Jurgovsky et al. (2018) formulated the credit card fraud prediction for a sequence classification problem. It is based on a supervised learning concept. They have used the LSTM (Long Short-Term Memory) approach to detect credit card fraud transactions. Roy et al. (2018) proposed a framework of deep learning with various tuning parameters to detect credit card fraud.

Seera et al. suggested using publicly accessible and real transaction records and employed 13 statistical and machine learning approaches for payment card fraud detection. The findings of the original and aggregated features are compared and examined Seera et al., (2021). The majority of data scientists regard credit risk assessment as a classification challenge, and they believe it helps uncover hidden patterns in credit scoring data analysis. It also aids specialists in improving their understanding of credit risk assessment. A wide range of machine learning algorithms has been used to demonstrate risk assessment systems in this field. Classification, ensemble classifiers, and hybrid classifiers are the three types of classification strategies Tripathi et al., (2021).

Although, there are multiple studies carried out for credit card fraud detection, like cardholder’s behavior prediction, enhancing the system’s processing time, etc. We are motivated by all the above works. We have used the XGBoost ensemble learning technique to improve the accuracy of the credit card fraud prediction system. It can create threads for parallel processing, have a cross-validation method, and handle missing values Mishra & Pandey, (2021). It reduces computation time and allocates memory resources efficiently. Because it was built and developed solely for the goal of model performance and computational speed, XGBoost has proved to push the limits of processing power for boosted trees algorithms. It was designed specifically for tree boosting techniques to use all available memory and hardware resources fully. Its implementation includes several additional features for model tuning, computing environments, and algorithm improvement. Both XGBoost and gradient boosting machines are ensemble tree approaches that employ the gradient descent architecture to boost weak learners. XGBoost, on the other hand, improves the base gradient boosting framework through system optimization and algorithm development Chen & Guestrin, (2016). We explore the XGBoost approach to gain the highest efficiency in large data sizes.

3. SELECTED MODELS

This section discusses various models for credit card fraud predictions. Our forecast and classification problem has considered Naïve Bayes, Support vector machine, Logistic regression, Random forest, Decision tree, and XGBoost Pandey et al., (2021).
3.1 Naïve Bayes Classifier

It is a famous machine learning model based on the Bayes theorem of conditional probability Rane & Kumar, (2018). It passes each transaction and calculates the posterior probability for available classes. According to the highest probability, the transaction is allocated to a class. Naïve Bayes classification is described using Equation (1):

$$P(C / x) = \frac{P(x / C) P(C)}{P(x)}$$  \hspace{2cm} (1)

Where,

- \(C\): specified class
- \(x\): transaction used for classification
- \(P(C)\) and \(P(x)\): prior probabilities
- \(P(C / x)\): posterior probability

Given the Naïve assumption, which says that a data point \(X=\{x_1, x_2, x_3, \ldots, x_i\}\), the possibility of each of its features taking place in a given class, the Equation can be rewritten as:

$$P(C / x) = P(C) . P(x_i / C)$$  \hspace{2cm} (2)

In this experiment, the following parameters (Table 1) are tuned for the Naïve Bayes classifier to get the results.

| Parameter       | Value   |
|-----------------|---------|
| priors          | None    |
| var_smoothing   | 1e-09   |

3.2 SVM Classifier

To solve regression and classification problems, Vapnik first introduced a support vector machine Cortes & Vapnik, (1995). This classifier is popular because of its high reliability, varied applications, and less vulnerability for the overfitted model Vapnik, (2006). It derives the finest hyperplane that maximizes the margin between two classes. It can be used for linear and nonlinear datasets. It can discover a nonlinear decision boundary by projecting the data with a nonlinear function to space with a higher dimension.

We traverse linearly separable classes using two-class problems. For a given Dataset S as \((P1, Q1), (P2, Q2), \ldots, (P|S|, Q|S|)\), where \(Q_j\) is the class label whose value is from +1 to -1 \((Q_j \in (-1, +1))\). \(Q_j\) is associated with the \(P_j\) set of training tuples.

A hyperplane can be described as a \(P\) set of points satisfying

$$W.P - B = 0$$  \hspace{2cm} (3)
Where $W$ is a normal vector to the hyperplane and $\frac{B}{||W||}$ is the offset of the plane from the origin and average vector $W$. The best line, which has the least classification error, in general. Best hyperplane by the maximum distance of the hyperplane to the closest of the negative instance and positive instance.

**Figure 2. Support vector machine Classifier**

![Figure 2. Support vector machine Classifier](image)

Figure 2 displays SVM optimal hyperplane in training with sample transactions to classify credit card fraud transactions, and not fraud transactions are represented as disk-shaped Kumar et al., (2019). In this experiment, the following parameters (Table 2) are tuned for the SVM Classifier to get the results.

**Table 2. SVM classifier parameters**

| Parameter          | Value         |
|--------------------|---------------|
| alpha              | 0.0001        |
| average            | False         |
| class_weight       | None,         |
| early_stopping     | False         |
| epsilon            | 0.1           |
| eta0               | 0.0           |
| fit_intercept      | True,         |
| l1_ratio           | 0.15          |
| learning_rate      | 'optimal'     |
| loss               | 'hinge'       |
| max_iter           | 1000          |
| n_iter_no_change   | 5             |
| n_jobs             | None          |

*Table 2 continued on next page*
3.3 Logistic Regression

Logistic regression is the most popular classification model in machine learning. In this model probability of each class is calculated for prediction. This model predicts output by combining the input variable with weight. Suppose $x$ is an independent variable and $y$ is a dependent variable, then linear regression can be represented as

$$ y = a_0 + a_1 x $$

(4)

Where $a_0$ is a bias term and $a_1$ is the weight for input variable $x$. If a logistic function is used for logistic regression, it predicts each class’s probability and will use a sigmoid or logistic function Shakya, (2018). Figure 3 displays an example of the sigmoid function.

![Figure 3. Sigmoid function](image)

This function predicts values between 0 and 1, and it can be expressed as

$$ sigm(z) = \frac{1}{1 + e^{-z}} $$

(5)

If we have an independent variable $(x)$ and a dependent variable $(y)$, then logistic regression can be stated as

$$ P(y = 1) = sigm(a_0 + a_1 * x) $$

(6)

Where $a_0$ and $a_1$ are the parameters of logistic regression model and learned while the training process. The predicted value can be expressed as (if threshold =0.5)
\[ y = 1 \text{ if } P(y=1) \geq 0.5 \]
\[ y = 0 \text{ if } P(y=1) < 0.5 \]

In this experiment, the following parameters (Table 3) are tuned for the Logistic regression approach to get the results.

### Table 3. Parameters used in Logistic regression classifier

| Parameter Name       | Value          |
|----------------------|----------------|
| C                    | 1.0            |
| class_weight         | None           |
| dual                 | False          |
| fit_intercept        | True           |
| intercept_scaling    | 1              |
| l1_ratio             | None           |
| max_iter             | 100            |
| multi_class          | ‘auto’         |
| n_jobs               | None           |
| penalty              | ‘‘2’’         |
| random_state         | None           |
| solver               | ‘‘lbfgs’’      |
| tol                  | 0.0001         |
| verbose              | 0              |
| warm_start           | False          |

### 3.4 Random Forest Classifier

This classifier is mainly used for classification and regression problems. It uses ensemble learning methods for prediction by averaging the prediction of several independent base models. The average of forecasts combined with several random decision trees Breiman, (2001). It shows excellent performance while more features are selected. However, it can be used for significant scale problems also. While working on the classification, random forest classifier \( m \) can be gained via the majority of votes among \( K \) classification trees with input \( x \) as

\[
m(x : C_1, C_2, \ldots, C_K) = \begin{cases} 
1 & \text{if } \frac{1}{K} \sum_{i=1}^{K} m(x; C_i) > \frac{1}{2} \\
0 & \text{otherwise}
\end{cases}
\]

(7)

In this experiment, \( C \) is a parameter set value. The following parameters (Table 4) are tuned for the Random Forest classifier approach to get the results in this experiment.
3.5 Decision Tree Classifier

It is a top-down method in which the root node produces a binary split up to a specific criterion. It is also used to solve classification problems in various domains. In the decision tree, internal nodes denote the test conditions, whereas leaf nodes denote class categories. Each produced terminal node indicates a predicted class. A decision tree creates a sequence of rules to identify a class of test transactions or instances according to classes. In this model, overfitting is handled by a post pruning strategy Nithyassik & Nandhini, (2010). The following parameters (Table 5) are tuned for the Decision Tree classifier approach to get the results in this experiment.

### Table 4. Parameters used in Random Forest classifier

| Parameter Name                  | Value |
|--------------------------------|-------|
| bootstrap                      | True  |
| ccp_alpha                      | 0.0   |
| class_weight                   | None  |
| criterion                      | ‘gini’|
| max_depth                      | None  |
| max_features                   | ‘auto’|
| max_leaf_nodes                 | None  |
| max_samples                    | None  |
| min_impurity_decrease          | 0.0   |
| min_impurity_split             | None  |
| min_samples_leaf               | 1     |
| min_samples_split              | 2     |
| min_weight_fraction_leaf       | 0.0   |
| n_estimators                   | 100   |
| n_jobs                         | None  |
| oob_score                      | False |
| random_state                   | None  |
| verbose                        | 0     |
| warm_start                     | False |
3.6 XGBoost Approach

XGBoost stands for eXtreme Gradient Boosting used for classification, regression, and problems related to ranking. This approach is an advanced version of the Gradient boosting approach Chen & Guestrin, (2016), which improves the results compared to the Gradient boosting approach. It uses ensemble techniques to improve results. Ensemble techniques are used to modify existing classification models to handle imbalanced class distributions. In ensemble learning, classification problems are solved by training multiple learners. Its central concept is to combine numerous weak learners into keen learners to boost the classifier’s performance. Figure 4 shows the idea of ensemble learning techniques.

Table 5. Parameters used in Decision tree classifier

| Parameter               | Value       |
|-------------------------|-------------|
| ccp_alpha               | 0.0         |
| class_weight            | None        |
| criterion               | ‘gini’      |
| max_depth               | None        |
| max_features            | None        |
| max_leaf_nodes          | None        |
| min_impurity_decrease   | 0.0         |
| min_impurity_split      | None        |
| min_samples_leaf        | 1           |
| min_samples_split       | 2           |
| min_weight_fraction_leaf| 0.0         |
| presort                 | ‘deprecated’|
| random_state            | None        |
| splitter                | ‘best’      |

Ensemble techniques can be categorized as bagging and boosting. It is shown in figure 5.
Bagging (Bootstrap Aggregation) is an ensemble technique. In this technique, various training samples are generated from original training sets known as a bootstrap sample, and this process is called bootstrapping. For prediction, each bootstrap sample is trained for an individual model. Final prediction calculates as aggregated by averaging in regression or voting in the classification of all bootstrap models.

Boosting is also an ensemble technique. It combines various weak learners to build a keen learner that provides improved results compared to individual learners. In each boosting step, weak learners are sequentially trained and correct their predecessor by adding weights to previously misclassified samples. In boosting, the technique bootstrapping is used to avoid variance and overfitting. XGBoost model uses boosting techniques for classification and predictions.

XGBoost adopts a more generalized method to control overfitting and contributes to improving the results. XGBoost works on parallel computing; hence it is fast as compared to the standard Gradient boosting approach. It can handle missing data and includes cross-validation features used to determine to boost round in each run. XGBoost needs a few parameters to tune for getting better results. In this experiment, the following parameters (Table 6) are tuned for the XGBoost approach to get the results.

### Table 6. Parameters used in XGBoost classifier

| Parameter Name       | Value            |
|----------------------|------------------|
| learning_rate        | 0.1              |
| n_estimators         | 1000             |
| max_depth            | 5                |
| min_child_weight     | 1                |
| gamma                | 0                |
| subsample            | 0.8              |
| colsample_bytree     | 0.8              |
| objective            | 'binary:logistic'|
| nthread              | 4                |
| scale_pos_weight     | 1                |
4. THE EMPIRICAL ANALYSIS OF CREDIT CARD PREDICTION

4.1 Data Source and Description

This paper uses a public credit card transaction dataset of European cardholders Dal Pozzolo et al., (2015), and it was also released in the Kaggle community. This dataset was made in September 2013. It comprises 284,807 transactions with only 492 transactions as a fraudulent class. It is a highly skewed dataset as the positive class contributes only 0.172% of the entire dataset. This unbalanced class distribution is presented in Figure 6.

This dataset has a total of 30 features. 28 features represented as V1 to V28 obtained after principal component analysis transformation and two features time and amount without PCA transformation. Table 7 shows the description of the dataset features.

4.2 Performance Assessment

Precision, recall, accuracy, AUC, and f-measure are utilized to evaluate the performance of the models. Table 8 shows the confusion matrix for assessment.
4.3 Result Analysis and Discussion

This section describes the experimental results of the different models simulated on the credit card transaction dataset. After cleaning the transactional dataset, we have applied the normalization and data standardization process. The complete dataset is partitioned into a training dataset (70%) and testing datasets (30%) for the simulation study. The training dataset builds the models, and the testing dataset evaluates the performance of these models. In this paper, all experimentations are conducted using the Python programming language.

Firstly, the credit card transaction dataset is trained using different models such as Naïve Bayes, SVM, random forest, decision tree, logistic regression, and XGBoost; performances are evaluated. Table 10 shows the experimental results of the classification task.
Different models’ performance is assessed using different measures such as accuracy, precision, recall, and F2-score. The performance of various machine learning methods is compared with the ensemble learning-based XGBoost model. From the analysis of table 4, the accuracy of the XGBoost model is higher than in other traditional models. It indicates that the ensemble learning-based approach, such as XGBoost, performs better prediction over other traditional approaches. Figure 7 represents the comparative graphical results of these models.

![Figure 7. Comparative results graph](image)

Another comparison made by the AUC (Area Under Curve) values, which are generated from precision and recall, is also higher in the XGBoost model. This comparison is depicted in table 11.

### Table 10. Performance comparison

| Model       | Accuracy | Precision | Recall | F1-Score |
|-------------|----------|-----------|--------|----------|
| Naive Bayes | 0.8934   | 0.90      | 0.90   | 0.89     |
| SVM         | 0.9390   | 0.94      | 0.94   | 0.94     |
| Logistic Regression | 0.9443   | 0.94      | 0.95   | 0.94     |
| RandomForest | 0.9341   | 0.93      | 0.94   | 0.94     |
| DecisionTree | 0.9340   | 0.92      | 0.92   | 0.92     |
| Xgboost     | 0.9644   | 0.96      | 0.97   | 0.96     |

### Table 11. AUC Results comparison

| Model       | AUC  |
|-------------|------|
| Naive Bayes | 0.9021|
| SVM         | 0.9382|
| Logistic Regression | 0.9578|
| RandomForest | 0.9430|
| DecisionTree | 0.9349|
| Xgboost     | 0.9669|
According to the true positive and false positives, ROC (Receiver operating characteristic) graph is also prepared for result evaluations. This ROC graph is depicted in figure 8.

Figure 8. ROC comparison

In conclusion, the ensemble techniques-based model, such as XGBoost used in this paper, has a better performance than state-of-the-art machine learning models. Therefore, the XGBoost model is a valuable model for credit card fraud prediction.

5. CONCLUSION

In this study, a comparative study is conducted among the ensemble technique-based XGBoost model and other traditional machine learning models to classify a credit card transaction as a fraud or not. Our study significantly reduces the risks of financial losses and the uncertainty that institutions encounter in their day-to-day operations. Our research includes a benchmark credit card transaction dataset to measure performances. Performance evaluation was conducted using machine learning and ensemble methodologies. For this, a publicly available credit cardholder’s dataset is used. XGBoost produces better performance in the ensemble model’s prediction accuracy, precision, recall, and AUC values. Considering the poor performance of traditional models in dealing with unbalanced data, the XGBoost model uses ensemble techniques to improve performance, gain prediction accuracy up to 96%, and outperform other traditional models. This research is significant in light of the tremendous growth of e-commerce operations, resulting in the exponential growth of online transactions in this digital era. Consumers are turning to internet purchasing as a result of the Covid-19 epidemic. As a result, a financial institution’s ability to limit the risk of fraud transactions depends on the efficient system for the fraud detection system. Various models have dealt with credit card fraud predictions recently, including machine learning, AI, and deep learning approaches. However, there is still room to develop more accurate and efficient models. The current study can be improved from several perspectives. Hybrid models can be built by combining two or more models to improve the research. Furthermore, the detection algorithms’ online implementation will be examined.

FUNDING AGENCY

Publisher has waived the Open Access publishing fee.
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