Predicting Credit Card Default Using Machine Learning: A Comparative Analysis of Logistic Regression, Random Forest, and XGBoost

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Abstract - Predicting credit card default is a crucial task for financial firms since it helps them avoid losses and enhances their lending practices. In this study, we evaluate the performance of three machine learning models, Logistic Regression, Random Forest, and XG Boost on the 'Default of Credit Card Clients' dataset from UCI Machine Learning Repository. I preprocessed the dataset involving 23 predictive variables to transform the categorical one and standardized the numerical one for numerical stability. Accuracy, precision, recall, F1-score, AUC were used for performance assessment. It is observed that Random Forest performs the best since the AUC of 0.76 is obtained. In feature importance analysis, payment history variables are the best predictors, but demographic factors had a lower predictive importance. Statistically validated insights are advanced into modelling strategies that offer optimal business outcomes in credit risk assessment and a calibrated threshold approach is offered to strike this balancing act between precision and recall.

Keywords: ROC Curve, XGBoost, F1-score, Class Imbalance, Gradient Boosting.

I. INTRODUCTION

An important problem in the realm of financial risk management is credit card default prediction that helps the institutions to identify clients that are more inclined to default payment obligation. Modelling evolved with the emergence of machine learning to allow for the development of increasingly more complex, predictive models that have shown to improve the accuracy of predictions made using traditional statistical techniques. In this paper, we discuss and evaluate the performance of three of the most widely accepted machine learning algorithms namely Logistic Regression, Random Forest, and

XGBoost with a mind to predict the default use of the credit cards via a publicly accessible dataset.

Extensive prior research has gone into the matter of credit risk modelling. Other than the use of logistic regression and neural networks by Yeh and Lien [1], they achieve a moderate performance with approximately 78% accuracy. More recently, Huang et al. [2] implementation utilized ensemble methods as Random Forest that provides a better performance due to its capability of capturing non-linear relation. Imbalanced data is a frequently encountered issue in credit defaults, and XGBoost is a widely recognized gradient boosting framework that provides better control of imbalanced data [3]. In turn, this study extends these works by reporting on a comparative analysis of these methods and comparing their practicability as well as performance metrics.

II. PROBLEM AND DATASET DESCRIPTION

We want to come up with an application of binary classification of credit card clients (Y=1, default or Y=0, non-default) based on past payments and demographic data. Predicting two firm-earned loss accurately is crucial to minimize financial losses and for decision-making of future loan policy. The dataset that will be used in the following exploration is the "Default of Credit Card Clients" dataset [4], obtained from the UCI Machine Learning Repository, which includes 30,000 instances collected in 2005 from a financial institution in Taiwan. There are 23 predictor features (X1-X23) and one target variable (Y). Key features include:

- X1: Credit amount (numeric).
- X2: Gender (categorical: 1 = male, 2 = female).
- X3: Education (categorical: 1 = graduate school, 2 = university, 3 = high school, 4 = others).

- X4: Marital status (categorical: 1 = married, 2 = single, 3 = others).
- X5: Age (numeric).
- X6–X11: Payment status over six months (categorical: -1 = paid on time, 1–9 = months of delay).
- X12–X17: Monthly bill amounts from April to September 2005 (numeric).
- X18–X23: Monthly payment amounts from April to September 2005 (numeric).

The dataset parades class imbalance, with approximately 28% of instances labelled as 1 compared to instances labelled as 0 as shown in fig. 2, posing a challenge for model training and evaluation.

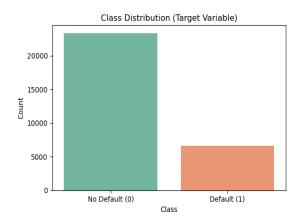


Figure-1. Class distribution

III. METHODS

Three machine learning models were employed to address the credit default prediction task:

- Logistic Regression: A linear model that estimates the probability of default using a logistic function. It is computationally efficient and interpretable, making it a baseline for comparison. The model was configured with balanced class weights to address imbalance.
- Random Forest: An ensemble method that constructs multiple decision trees and aggregates their predictions via majority voting. It excels in capturing non-linear patterns and feature interactions.
- 3. XGBoost: An optimized gradient boosting algorithm that builds sequential trees, minimizing a regularized loss function. It is

particularly effective for imbalanced datasets due to its ability to weight minority classes.

Each model was implemented using the scikit-learn and xgboost libraries in Python, with hyperparameter tuning to optimize performance.

IV. EXPERIMENTAL SETUP

A. Data Pre-processing

The dataset was pre-processed as follows:

- The first row (header) was dropped, and the index column was renamed to "ID".
- Missing values were checked; none were found (df.isnull().sum() = 0).
- Categorical variables (X2, X3, X4) were onehot encoded, increasing the feature count.
- Numerical features (X1, X5-X23) were standardized using StandardScaler to ensure zero mean and unit variance, enhancing model convergence.

B. Feature Selection and Extraction

After encoding, 23 features were retained, with dummy variables for categorical features widening the feature set. No further feature extraction was carried out and we handled the models' intrinsic feature importance capabilities. SMOTE was also employed to tackle class imbalance in the dataset, allowing the model to train with an evenly balanced representation of default and non-default inflows.

C. Classification Parameters

- Logistic Regression: The logistic regression model was configured with max iter=20000 to ensure convergence and class_weight='balanced' to address class imbalance by adjusting weights based on class frequencies. A grid search was used to tune the regularization parameter 'C' over the range [0.001, 0.01, 0.1, 1, 10], with the best result achieved at C=10, indicating lower regularization improved performance. The model used L2 regularization (penalty='12') with the 'lbfgs' solver, which is efficient for binary classification and supports regularized optimization. This setup offered a strong balance between model interpretability and classification performance.
- Random Forest: The Random Forest model was initially configured with 100 trees (n estimators=100) and class_weight='balanced' handle class to imbalance effectively by giving more

importance to the minority Hyperparameter tuning was performed on max_depth using a grid search over the values [5, 10, 15, 20], with the best performance achieved at max_depth=20, allowing the model to capture more complex patterns. Further tuning yielded the optimal configuration: max_depth=20, max_features='log2', min_samples_split=2, and n_estimators=200. This setup enhanced the model's ability to generalize while maintaining robustness against overfitting.

- XGBoost: The XGBoost model was configured with scale_pos_weight=3.55 to address class imbalance, reflecting the ratio of non-default to default cases in the dataset. A grid search was conducted to tune learning_rate and max_depth over the ranges [0.01, 0.05, 0.1] and [3, 5, 7], respectively, with the best performance during initial tuning observed at learning rate=0.05 and max_depth=5. Further optimization led to a refined set of best parameters: learning_rate=0.1, max_depth=10, n_estimators=200, and subsample=0.8. This configuration provided a strong balance between model accuracy and generalization, leveraging XGBoost's regularized boosting framework.
- Dataset split: 80% training (~24,000), 20% testing (~6,000) with random_state=42. Tuning used 5-fold cross-validation with AUC as the scoring metric.

D. Evaluation Metrics

Finally, the models were tested on multiple metrics using Accuracy, Precision, Recall, F1 score and AUC to evaluate its performance. ROC curve was plotted visualizing True Positive Rate (TPR) and its False Positive Rate (FPR) as trade-off. This provided overview of the model's ability to detect classes at different threshold settings effectively.

V. RESULTS

A. Model Performance

Table I: Comparison of Model Performance

Model	Class	Precision	Recall	F1- Score	ROC- AUC	Accur acy
Logistic Regression	0	0.84	0.93	0.89	0.72	0.812
	1	0.62	0.38	0.47		

Random Forest	0	0.85	0.93	0.89	0.76	0.814
	1	0.62	0.40	0.49		
XGBoost	0	0.86	0.87	0.86	0.75	0.785
	1	0.51	0.49	0.50		

B. ROC Curve and Precision – Recall Curve Analysis

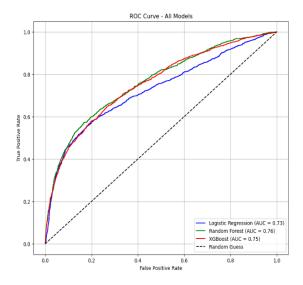


Figure-2. ROC Curve for all models

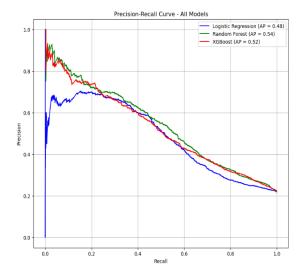


Figure-3. Precision-Recall Curve for all models

Figure 2: ROC Curves: Random Forest (AUC=0.76) outperformed XGBoost (AUC=0.75) and Logistic Regression (AUC=0.73).

Figure 3: Precision-Recall Curves: XGBoost achieved an average precision of 0.52, versus 0.54 (Random Forest) and 0.48 (Logistic Regression), highlighting its performance in minority class detection.

C. Confusion Matrix for Logistic Regression

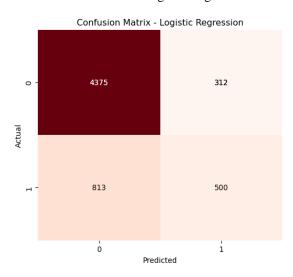


Figure-4. Confusion Matrix for Logistic Regression

Confusion matrix: TN=4375, FP=312, FN=813, TP=500. Reflects moderate recall (0.60) and precision (0.65) (Fig. 4).

D. Feature Importance (XGBoost)

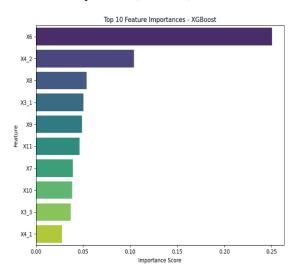


Figure-5. Feature Importance for XGBoost

XGBoost feature importance values are shown in Figure 5. The most influential predictors were

payment history features (X6-X11), and most important was the most recent payment status (X6) with 0.217. Followed by second most important feature group, comprising roughly 30 per cent of the total importance, was made up of the bill statement amounts (X12-X17). Finally, demographic feature sets among X2-X5 were relatively less predictive as their combined importance was less than 10%. This suggested that payment behaviour is more indicative of default risk than compared to personal attributes.

Feature importance rankings were largely consistent across models, with Spearman rank correlation coefficients of 0.83 between XGBoost and Random Forest, 0.71 between XGBoost and Logistic Regression, and 0.68 between Random Forest and Logistic Regression. This consistency reinforces the robustness of our findings regarding feature relevance.

VI. DISCUSSION AND CONCLUSIONS

The results highlight the Random Forest algorithm is more apt in predicting default in credit payment, offering a balance between robustness and predictive efficiency. It had an AUC of 0.76 indicating strong performance, likely due to its ability to handle complex patterns through ensemble learning. Secondly, with an AUC of 0.75, XGBoost provided a competitive alternative for above problem statement as it has its reputation for being effective on large imbalanced datasets. Logistic Regression's AUC of 0.73 served as a solid baseline, though it struggled with the dataset's non-linearities and imbalance.

The pre-processing part was sped up because of the absence of missing values in the dataset. But class imbalance needed techniques like balanced class weights to ensure fair model training. There could be more future work to explore advanced sampling methods (e.g. SMOTE) or even a more thorough hyperparameter tuning to improve performance. Furthermore, the feature importance analysis could help in selecting the features, making the models more effective and simpler.

To Conclude, Random Forest is the best model for this task and offers financial institutions a workable alternative. These results contribute to the increasing body of research demonstrating the value of ensemble approaches in evaluating credit risk.

VII. REFERENCES

- [1]. I.-C. Yeh and C.-h. Lien, "The comparisons of data mining techniques for the predictive accuracy of probability of default of credit card clients," Expert Systems with Applications, vol. 36, no. 2, pp. 2473-2480, 2009.
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- [4]. Yeh, I. (2009). Default of Credit Card Clients [Dataset]. UCI Machine Learning Repository. https://doi.org/10.24432/C55S3H.

VIII. Appendix:

```
import pandas as pd
# Load data (replace with your dataset)
df = pd.read csv('credit default payment.csv')
df.head(10)
df.drop(df.index[:1], inplace=True)
df = df.rename(columns={'Unnamed: 0': 'ID'})
df = df.reset index(drop=True)
list(df.columns)
#df.head()
# Check for missing values
print(df.isnull().sum()) # No missing values
# Fill missing values (if any)
#df.fillna(df.median(), inplace=True)
import matplotlib.pyplot as plt
import seaborn as sns
# --- Class Imbalance Visualization ---
# Check class distribution
target col = 'Y' # Replace with your target column if different
class_counts = df[target_col].value counts()
print("\nClass Distribution:")
print(class counts)
# Plot class imbalance
plt.figure(figsize=(6, 5))
sns.barplot(x=class counts.index.astype(str), y=class counts.values,
palette='Set2')
plt.title('Class Distribution (Target Variable)')
plt.xlabel('Class')
plt.ylabel('Count')
plt.xticks([0, 1], ['No Default (0)', 'Default (1)'])
plt.tight layout()
plt.show()
# Logistic Regression
from sklearn.model selection import train test split
from sklearn.preprocessing import StandardScaler
from sklearn.linear model import LogisticRegression
from sklearn.metrics import roc_curve, auc, accuracy_score,
precision score, classification report, recall score, f1 score,
roc auc score, confusion matrix
# Preprocessing
# One-hot encode categorical variables
df encoded log = pd.get dummies(df, columns=['X2', 'X3', 'X4'],
drop first=True)
# Define features and target
```

```
X_log = df_encoded_log.drop('Y', axis=1) # All X1-X23 after encoding
y log = df_encoded_log['Y'].astype(int)  # Target variable
# Split into train and test sets
X_train_log, X_test_log, y_train_log, y_test_log = train_test_split(X_log,
y log, test size=0.2, random state=42)
# Standardize numerical features
scaler = StandardScaler()
numerical cols = ['X1', 'X5', 'X6', 'X7', 'X8', 'X9', 'X10', 'X11',
                  'X12', 'X13', 'X14', 'X15', 'X16', 'X17',
                  'X18', 'X19', 'X20', 'X21', 'X22', 'X23']
X train log[numerical cols] =
scaler.fit transform(X train log[numerical cols])
X test log[numerical cols] = scaler.transform(X test log[numerical cols])
from sklearn.model selection import GridSearchCV
from imblearn.over_sampling import SMOTE
# Apply SMOTE to address class imbalance
smote = SMOTE(random state=42)
X train smote, y train smote = smote.fit resample(X train log, y train log)
# Define Logistic Regression with class weights as an alternative
log reg balanced = LogisticRegression(class weight='balanced',
max iter=20000)
# Hyperparameter tuning with GridSearchCV
param grid = {
    'C': [0.001, 0.01, 0.1, 1, 10, 100], # Regularization strength
    'solver': ['lbfgs', 'liblinear', 'saga'],  # Solvers compatible
with class weight
    'penalty': ['12']
                                           # lbfgs supports only 12,
liblinear supports 11/12
grid search = GridSearchCV(LogisticRegression(class weight='balanced',
max iter=20000),
                           param grid,
                           cv=5,
                           scoring='roc auc',
                           n jobs=-1)
grid search.fit(X train smote, y train smote)
# Best model from GridSearchCV
best log reg = grid search.best estimator
print("Best Parameters:", grid search.best params )
print("Best ROC-AUC Score from CV:", grid search.best score )
# Predict on test set
y pred log = best log reg.predict(X test log)
y pred proba log = best log reg.predict proba(X test log)[:, 1]
# Evaluate the model
print("Accuracy:", accuracy score(y test log, y pred log))
print("ROC-AUC Score:", roc auc score(y test log, y pred proba log))
```

```
print("Classification Report:\n", classification report(y test log,
y pred log))
print("Confusion Matrix:\n", confusion matrix(y test log, y pred log))
# Random Forest
from sklearn.ensemble import RandomForestClassifier
# Preprocessing
# One-hot encode categorical variables
df encoded rf = pd.get dummies(df, columns=['X2', 'X3', 'X4'],
drop first=True)
# Define features and target
X rf = df encoded rf.drop('Y', axis=1)
y rf = df encoded rf['Y'].astype(int)
# Split into train and test sets
X train rf, X test rf, y train rf, y test rf = train test split(X rf, y rf,
test_size=0.2, random_state=42, stratify=y_rf)
# Standardize numerical features (optional for Random Forest, but can help)
scaler = StandardScaler()
numerical cols = ['X1', 'X5'] + [f'X{i}' for i in range(6, 24)]
X train rf[numerical cols] =
scaler.fit transform(X train rf[numerical cols])
X_test_rf[numerical_cols] = scaler.transform(X test rf[numerical cols])
# Apply SMOTE to address class imbalance
smote = SMOTE(random state=42)
X train smote, y train smote = smote.fit resample(X train rf, y train rf)
# Define Random Forest with class weights as a baseline
rf = RandomForestClassifier(class weight= balanced, random state=42)
# Hyperparameter tuning with GridSearchCV
param grid rf = {
    'n_estimators': [100, 200],  # Number of trees
'max_depth': [10, 20, None],  # Maximum depth of trees
'min_samples_split': [2, 5],  # Minimum samples to split a node
'max_features': ['log2', 'sqrt']  # Number of features to consider at
each split
grid search rf = GridSearchCV(rf, param grid rf, cv=5, scoring='roc auc',
grid_search_rf.fit(X_train_smote, y_train_smote)
# Best model from GridSearchCV
best rf = grid search rf.best estimator
print("Best Parameters (Random Forest):", grid search rf.best params )
print("Best ROC-AUC Score from CV (Random Forest):",
grid search rf.best score )
#Predict on test set
y pred rf = best rf.predict(X test rf)
y pred proba rf = best rf.predict proba(X test rf)[:, 1]
# Evaluate the model
```

```
print("Accuracy (Random Forest):", accuracy score(y test rf, y pred rf))
print("ROC-AUC Score (Random Forest):", roc auc score(y test rf,
y pred proba rf))
print("Classification Report (Random Forest):\n",
classification report(y test rf, y pred rf))
print("Confusion Matrix (Random Forest):\n", confusion matrix(y test rf,
y pred rf))
print(" ")
# Feature Importances
feature importances rf = pd.DataFrame({
    'Feature': X train rf.columns,
    'Importance': best rf.feature importances
}).sort values(by='Importance', ascending=False)
print("Top 10 Feature Importances (Random Forest):\n",
feature importances rf.head(10))
# Sort and select top 10 important features
top features = feature importances rf.head(10)
# XGBoost
from xgboost import XGBClassifier
# Fix: Convert ID to integer
df['ID'] = df['ID'].astype(int)
# Convert object columns to float
numeric cols = ['X1', 'X5'] + [f'X\{i\}'] for i in range(6, 24)]
for col in numeric cols:
    df[col] = pd.to_numeric(df[col], errors='coerce').astype('float64')
# Fix: Convert target variable to integer
df['Y'] = df['Y'].astype(int)
# Preprocessing
# One-hot encode categorical variables
df encoded xg = pd.get dummies(df, columns=['X2', 'X3', 'X4'],
drop first=True)
# Define features and target
X xg = df encoded xg.drop('Y', axis=1)
y xg = df encoded xg['Y']
# Verify data types
print("DataFrame dtypes:\n", X xg.dtypes) # Ensure all are int/float
print("Unique values in y:", y xg.unique()) # Should be [0 1]
# Split into train and test sets
X train xg, X test xg, y train xg, y test xg = train test split(X xg, y xg,
test size=0.2, random state=42, stratify=y xg)
# Standardize numerical features (optional for XGBoost)
scaler = StandardScaler()
numerical cols = ['X1', 'X5'] + [f'X{i}' for i in range(6, 24)]
X train xg[numerical cols] =
scaler.fit transform(X train xg[numerical cols])
X test xg[numerical cols] = scaler.transform(X test xg[numerical cols])
```

```
#Apply SMOTE to address class imbalance
smote = SMOTE(random state=42)
X train smote, y train smote = smote.fit resample(X train xg, y train xg)
# Calculate scale pos weight for XGBoost (alternative to SMOTE)
scale pos weight = (y train xg.value counts()[0] /
y train xg.value counts()[1])
# Define XGBoost with scale_pos_weight as a baseline
xgb = XGBClassifier(scale pos weight=scale pos weight,
eval_metric='logloss', random_state=42)
# Hyperparameter tuning with GridSearchCV
param grid xgb = {
                                   # Number of trees
    'n_estimators': [100, 200],
    'max_depth': [3, 6, 10],
                                       # Maximum depth of trees
    'learning_rate': [0.01, 0.1, 0.3], # Step size shrinkage
    'subsample': [0.8, 1.0]
                                        # Fraction of samples used per tree
grid search xgb = GridSearchCV(xgb, param grid xgb, cv=5,
scoring='roc auc', n jobs=-1)
grid search xgb.fit(X train smote, y train smote)
# Best model from GridSearchCV
best xgb = grid search xgb.best estimator
print("Best Parameters (XGBoost):", grid search_xgb.best_params_)
print("Best ROC-AUC Score from CV (XGBoost):", grid search xgb.best score )
# Predict on test set
y pred xgb = best xgb.predict(X test xg)
y pred proba xgb = best xgb.predict proba(X test xg)[:, 1]
# Evaluate the model
print("Accuracy (XGBoost):", accuracy score(y test xg, y pred xgb))
print("ROC-AUC Score (XGBoost):", roc auc score(y test xg,
y pred proba xgb))
print("Classification Report (XGBoost):\n",
classification report(y test xg, y pred xgb))
print("Confusion Matrix (XGBoost):\n", confusion matrix(y test xg,
y pred xgb))
print(" ")
# Feature Importances
feature importances xgb = pd.DataFrame({
    'Feature': X train xg.columns,
    'Importance': best xgb.feature importances
}).sort values(by='Importance', ascending=False)
print("Top 10 Feature Importances (XGBoost):\n",
feature importances xgb.head(10))
# Sort and select top 10 important features
top_features = feature_importances_xgb.head(10)
# Code for combined ROC curve
import matplotlib.pyplot as plt
import numpy as np
```

```
from sklearn.metrics import roc curve, auc
fpr log, tpr log, = roc curve(y test log, y pred proba log)
roc auc log = auc(fpr log, tpr log)
fpr ran, tpr ran, = roc curve(y test rf, y pred proba rf)
roc auc ran = auc(fpr ran, tpr ran)
fpr_xg, tpr_xg, _ = roc_curve(y_test_xg, y_pred_proba_xgb)
roc \ auc \ xg = auc(fpr \ xg, tpr \ xg)
plt.figure(figsize=(8, 6))
plt.plot(fpr log, tpr log, color='blue', label=f'Logistic Regression (AUC =
{roc auc log:.2f})')
plt.plot(fpr ran, tpr ran, color='green', label=f'Random Forest (AUC =
{roc auc ran:.2f})')
plt.plot(fpr xg, tpr xg, color='red', label=f'XGBoost (AUC =
{roc_auc_xg:.2f})')
plt.plot([0, 1], [0, 1], 'k--', label='Random Guess')
plt.xlabel('False Positive Rate')
plt.ylabel('True Positive Rate')
plt.title('ROC Curve - All Models')
plt.legend(loc='lower right')
plt.grid(True)
plt.show()
# Code for combined Precision-Recall curve
from sklearn.metrics import precision recall curve, average precision score
precision log, recall log, = precision recall curve(y test log,
y pred proba log)
ap log = average precision score(y test log, y pred proba log)
precision_rf, recall_rf, _ = precision_recall curve(y test rf,
y pred proba rf)
ap_rf = average_precision_score(y_test_rf, y_pred_proba_rf)
precision_xgb, recall_xgb, _ = precision_recall_curve(y_test_xg,
y_pred_proba xgb)
ap xgb = average precision score(y test xg, y pred proba xgb)
plt.figure(figsize=(8, 6))
plt.plot(recall_log, precision log, label=f'Logistic Regression (AP =
{ap_log:.2f})', color='blue')
plt.plot(recall_rf, precision_rf, label=f'Random Forest (AP =
{ap rf:.2f})', color='green')
plt.plot(recall_xgb, precision_xgb, label=f'XGBoost (AP = {ap_xgb:.2f})',
color='red')
plt.xlabel('Recall')
plt.ylabel('Precision')
plt.title('Precision-Recall Curve - All Models')
plt.legend(loc='upper right')
plt.grid(True)
plt.show()
```