



## **Model Development Phase Template**

Date	July 2024
Team ID	Team-739859
Project Title	Auto Insurance Fraud Detection Using Machine Learning
Maximum Marks	10 Marks

## **Initial Model Training Code, Model Validation and Evaluation Report**

The initial model training code will be showcased in the future through a screenshot. The model validation and evaluation report will include a summary and training and validation performance metrics for multiple models, presented through respective screenshots.

## **Initial Model Training Code (5 marks):**

Paste the screenshot of the model training code

## **Model Validation and Evaluation Report (5 marks):**

Model	Summary	Training and Validation Performance Metrics
Model 1	Logistic regression model typically includes accuracy, precision, recall, F1 score to evaluate its predictive performance and generalization capability.	[ ] from sklearn.linear_model import LogisticRegression from sklearn.metrics import accuracy_score  # (your existing code)  def comparison(x_test,y_test):     # Create and fit a logistic regression model     lng = LogisticRegression()     lng.fit(X_train, y_train) # Assuming X_train and y_train are defined  # Now you can calculate predictions     lng.pred = lng.predict(x_test)     print("logistic Regression:",100*accuracy_score(y_test,lng_pred))     # (rest of your comparison function)  [ ] comparison(X_test, y_test) # Change 'x_test' to 'X_test' if it's defined that way





Model 2	Decision tree classifier model commonly includes accuracy, precision, recall, F1 score which help assess the model's prediction accuracy and generalizability.	from sklaarm.tree import DecisionTreeClassifier from sklaarm.metrics import accuracy_score dtcoDecisionTreeClassifier()  # Replace Ellipsis with your actual training date X_train = [[1, 2, 1], [3, 4], [5, 6]] # Sample date, replace with your own y_train = [0, 1, 0] # Example labels, replace with your own  # Define X_test = replace with your actual test data X_test = [[7,0], [0,10], [1,1,2]] # Sample data, replace with your own # Define Y_test = replace with your actual test labels y_test = [1, 0, 2] # Sample data, replace with your own dtc.fft(X_train,y_train) y_pred-dtc.pred(ctX_test), replace with your own dtc.fft(X_train,y_train) y_test = [1, 0, 2] # Sample data, replace with your own dtc.fft(X_train,y_train) y_test = [1, 0, 2] # Sample data, replace with your own dtc.fft(X_train,y_train) y_test = [1, 0, 2] # Sample data, replace with your own dtc.fft(X_train,y_train) y_test = [1, 0, 2] # Sample data, replace with your own dtc.fft(X_train,y_train) y_test = [1, 0, 2] # Sample data, replace with your own dtc.fft(X_train,y_train) y_test = [1, 0, 2] # Sample data, replace with your own dtc.fft(X_train,y_train) y_test = [1, 0, 2] # Sample data, replace with your own dtc.fft(X_train,y_train) y_test = [1, 0, 2] # Sample data, replace with your own dtc.fft(X_train,y_train) y_test = [1, 0, 2] # Sample data, replace with your own dtc.fft(X_train,y_train) y_test = [1, 0, 2] # Sample data, replace with your own dtc.fft(X_train,y_train) y_test = [1, 0, 2] # Sample data, replace with your own dtc.fft(X_train,y_train) y_test = [1, 0, 2] # Sample data, replace with your own dtc.fft(X_train,y_train) y_test = [1, 0, 2] # Sample data, replace with your own dtc.fft(X_train,y_train) y_test = [1, 0, 2] # Sample data, replace with your own dtc.fft(X_train,y_train) y_test = [1, 0, 2] # Sample data, replace with your own dtc.fft(X_train,y_train) y_test = [1, 0, 2] # Sample data, replace with your own dtc.fft(X_train,y_train) y_test = [1, 0, 2] # Sample data, replace with your own dtc.fft(X_train,y_train) y_test = [1, 0, 2] # Sam
Model 3	Random forest classifier model often encompasses accuracy, precision, recall, F1 score to measure its prediction quality and robustness.	from sklearm.ensemble import RandomForestClassifier import pandas as pd import numpy as np from sklearm.metrics import accuracy_score # Import accuracy_score  # Define X_train and y_train here with your actual data X_train * [[1, 2], [3, 4], [5, 6]] # Example data, replace with your own y_train * [0, 1, 0] # Example data, replace with your own  # Define X_test - replace with your actual test data X_test - [[7,8], [9,10], [11], [2]] # Example data, replace with your own # Define Y_test - replace with your actual test labels y_test = [1, 0, 1] # Example data, replace with your own  # Initialize and fit the RandomForestClassifier (replace with your code)  # Initialize and fit the RandomForestClassifier (replace with your code)  # Fc. # Fallom # Fc. # Fallom # Fc.
Model 4	K-nearest neighbors' classifier model typically includes accuracy, precision, recall, F1 score to evaluate its prediction performance and generalization ability.	from sklearm.neighbors import KNeighborsClassifier from sklearm.neighbors import confusion_matrix, classification_report import numpy as np # import numpy for array manipulation  # Assuming you have a larger x_train and corresponding y_train defined elsewhere, # use a subset of it that matches the size of your y_train in this example x_train_subset = x_train[i3] # Select the first 3 samples from your larger x_train  # Now x_train_subset and y_train have the same number of samples if x_train_subset and y_train have the same number of samples if x_train_subset.shape[0] > 10 and y_train.shape[0] > 0:     km = XbelghborsLassifier(n_elshbors-10)     km.fit(x_train_subset, y_train) # fit the model with the subset y_prad = knn.predict(x_test)     print(confusion_matrix(y_test, y_pred))     print(classification_report(y_test, y_pred))  # If you don't have a larger x_train, you need to define it here with 800 samples # to match the size mentioned in the error message.
Model 5	Naive Bayes classifier model typically includes accuracy, precision, recall, F1 score to evaluate its prediction performance and generalization.	from sklearm.naive_bayes import CategoricalNB.GaussianNB inport numpy as 0p # Import numpy # Define X_train - replace with your actual training data X_train = np.array([[1, 2], [2, 4], [5, 6]]) # Convert X_train to a NumPy array  # Check if X_train has data before fitting the model if X_train.shape[0 > 0 # Use 'X_train' instead of 'x_train' gnb-GaussianNB() # Reshape y_train to be a 1D array if it's not empty if len(y_train) > 0 # Check if y_train has data using len() for lists y_train_reshaped = np.array(y_train).real() # # Reshape y_train to 1D using numpy model_lagn.fit(X_train_y_train_reshaped) # Use X_train # Assuming X_test and y_test are defined elsewhere as NumPy arrays predict_log-wood_la.predict(T_test) # Use X_train print('training accoracy','100*accoracy_score(model_t_predict(X_train),(y_train_reshaped))) # Use X_train print('testining accoracy','100*accoracy_score(y_test,predict_log))





def comparison(x\_test, y\_test):
 # Create and fit a logistic regression model
 lrg = LogisticRegression()
 lrg.fit(X\_train, y\_train) # Assuming X\_train and y\_train are defined # Now you can calculate predictions lrg\_pred = lrg.predict(x\_test) return lrg\_pred # Return the lrg\_pred variable lrg\_pred = comparison(X\_test, y\_test) Confusion matrix is used to evaluate # Now you can print the classification report
print(confusion\_matrix(y\_test, y\_pred))
print(classification\_report(y\_test, lrg\_pred)) the model's performance by showing Model 6 Show hidden output the actual versus predicted [ ] Start coding or generate with AI. classifications. print("confusion matrix\n",confusion\_matrix(y\_test,y\_pred),"\n")
print("classification\_report\n",classification\_report(y\_test,y\_pred)) Start coding or generate with AI. [ ] print(confusion\_matrix(y\_test,y\_pred))
 print(classification\_report(y\_test,y\_pred)) precision recall f1-score support