

COMPREHENSIVE ANALYSIS OF CAR PRICES DATASET

FINAL TERM PROJECT

MACHINE LEARNING – I DR. REZA JAFARI

SUBMITTED BY
DIVAKARA RAO ANNEPU

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Abstract

This project undertakes a detailed exploration and analysis of a large dataset containing over 558,000 entries related to car sales, encompassing features such as make, model, year, odometer readings, condition, and selling prices. The objective of this study is to derive meaningful insights into factors influencing car prices and to develop predictive models that can accurately forecast car selling prices based on various attributes. The analysis involved extensive data preprocessing, including handling missing values, encoding categorical variables, and normalizing numerical data. Subsequent exploratory data analysis provided a foundational understanding of the distributions and relationships within the data, setting the stage for more complex analytical techniques.

The project progressed through several analytical phases, each addressing different aspects of data science and machine learning. In Phase I, exploratory data analysis and feature engineering were employed to prepare the dataset for modeling. Phase II focused on regression analysis, utilizing multiple linear regression to model the relationship between car features and their selling prices, achieving a high R-squared value indicative of a good model fit. Phase III explored various classification models to predict categorical outcomes, such as car condition, comparing the performance of models like Decision Trees, SVM, and Logistic Regression. The final phase, cluster analysis, applied K-means and DBSCAN to uncover natural groupings and patterns in the data. Each phase was supported by rigorous statistical testing and validation methods to ensure the robustness and accuracy of the findings. Recommendations were made based on the analysis results, emphasizing improvements in data collection, model tuning, and feature selection to enhance future predictions and analyses.

1 Introduction

The automotive industry stands as one of the most data-intensive sectors in the global economy, where every transaction generates a wealth of information that can be harnessed to predict market trends, optimize pricing strategies, and enhance consumer satisfaction. The advent of data science and machine learning offers profound opportunities to exploit this data, providing insights that can lead to more informed decisions. This project focuses on analyzing a substantial dataset of car sales, which includes over 558,000 records with 16 distinct attributes ranging from basic car details to sales information. The aim is to extract actionable insights from the dataset and develop predictive models to estimate car selling prices effectively.

The methodology of this project is structured into several distinct phases, each designed to address specific research questions and achieve the overarching goal of

understanding the determinants of car pricing in a detailed manner. Initially, the project begins with comprehensive data cleaning and preprocessing to handle missing data, encode categorical variables, and normalize distributions, ensuring the dataset is primed for robust analysis. Following this, exploratory data analysis (EDA) is conducted to identify trends, detect outliers, and understand the relationships between various car attributes and their selling prices. This foundational analysis sets the stage for more advanced statistical modeling.

In the subsequent phases, the project employs a variety of machine learning techniques. Phase II involves building and fine-tuning multiple linear regression models to predict car selling prices, assessing the model's performance through statistical metrics such as R-squared values and mean squared error (MSE). Phase III shifts focus to classification techniques, where models such as Decision Trees, Support Vector Machines (SVM), and Logistic Regression are developed to classify cars based on their condition. The performance of these models is meticulously compared using metrics like accuracy, precision, recall, and the Area Under the ROC Curve (AUC). The final analytical phase, cluster analysis, applies unsupervised learning methods to discover inherent groupings within the data, providing insights into market segmentation.

By integrating these diverse analytical techniques, the project not only enhances the understanding of factors that influence car prices but also demonstrates the application of data science methodologies in real-world industry settings. The insights gained from this study are intended to guide automotive companies in refining their pricing strategies and improving customer targeting. Through detailed data analysis and modeling, this project contributes to the evolving field of automotive data science, paving the way for further research and application in predictive analytics.

2 Description of the Dataset

The dataset under analysis comprises an extensive collection of over 558,837 car sales transactions, each detailed with 16 different attributes. This rich dataset provides a granular view of the various factors that influence car prices in the market. Each record in the dataset represents a unique car sale and includes both descriptive and quantitative information about the vehicle and the conditions of its sale.

2.1 Attributes Overview

• **Year:** The manufacture year of the car, crucial for pricing due to the age and technology of the vehicle.

- Make: The brand or manufacturer, such as Ford, Toyota, or BMW, which influences the car's market perception and price.
- **Model:** Specific model of the car affecting prices due to varying features, performance specifications, and market demand.
- **Trim:** Specific trim level of the car model, indicating variations in luxury and functional features that impact pricing.
- **Body:** The body style of the car, e.g., sedan, SUV, or hatchback, influencing consumer preference and thus pricing.
- **Transmission:** Type of transmission system (automatic, manual), a significant factor in driving experience and price.
- VIN: Vehicle Identification Number, a unique code that identifies individual vehicles.
- **State:** The U.S. state where the car was sold, influencing prices due to regional demand variations and economic conditions.
- **Condition:** A numeric rating of the car's overall physical and operational condition at the time of sale, influencing its market value.
- **Odometer:** Mileage on the car, where higher mileage usually decreases a vehicle's price due to wear and tear.
- Color: Exterior color of the car, affecting consumer preference and resale value.
- **Interior:** Color and type of the car's interior, contributing to aesthetic appeal and comfort.
- **Seller:** The entity or individual who sold the car, such as a dealer, an auction house, or a private seller.
- **MMR:** The Manheim Market Report value, an industry-standard wholesale price, serving as a benchmark for pricing used cars.
- **Selling Price:** The actual sale price of the car, the dependent variable in price prediction models.

• Sale Date: The date on which the transaction occurred, providing temporal context for the sale.

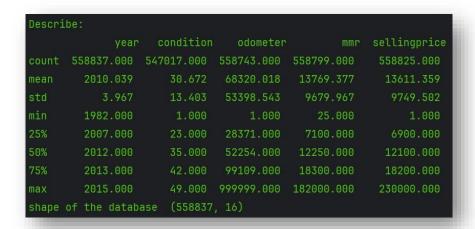


Figure 1

2.2 Data Utilization

The dataset serves multiple analytical purposes:

- **Price Prediction:** Utilizing attributes like year, make, model, and odometer readings to estimate the selling price.
- **Condition Assessment:** Classification models can predict the condition category of a vehicle based on its characteristics and historical data.
- Market Analysis: Analysis of sales trends over time and across different regions to inform market strategies for automotive firms.

2.3 Importance in Industry

In the automotive industry, a deep understanding of the factors influencing car prices and sales dynamics is crucial for inventory management, pricing strategies, and marketing. The dataset provides a comprehensive basis for such analyses, offering actionable insights applicable to car manufacturers, dealerships, and secondary market analysts.

3 Phase I: Feature Engineering & Exploratory Data Analysis

3.1 Data Cleaning

Data cleaning is a critical initial step in any data analysis project. For this dataset, addressing missing values is essential to ensure the integrity of subsequent analyses. Various methods such as imputation using the mean or median for numerical data, or mode for categorical data, can be applied. This dataset also required a check for duplicate entries, which were removed to prevent any bias or incorrect outcomes in the analysis.

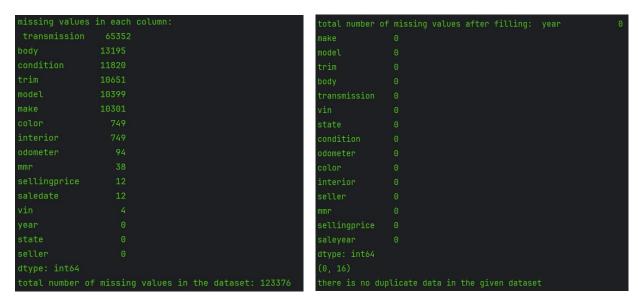


Figure 2: Missing Values before and after filling.

3.2 Aggregation

Aggregation was performed where applicable to summarize data features, enhancing the clarity and efficiency of the analysis. In the data preprocessing phase, it is often necessary to simplify or modify the data to enhance analytical accuracy and efficiency. One common transformation is extracting specific components from date fields. For our dataset, extracting the year from the 'saledate' column allows us to perform temporal analyses, such as year-on-year trends or seasonal patterns, more efficiently.

```
0 2014

1 2014

2 2015

3 2015

4 2014

Name: saleyear, dtype: int32
```

Figure 3: New extracted feature 'saleyear'

3.3 Down Sampling

Down sampling was considered to balance the dataset, particularly useful in scenarios where one class significantly outweighs another in a classification context. This ensures that the model trained on the data does not become biased towards the majority class. Here we have down sampled the data using the random sampling technique.

3.4 Dimensionality Reduction/Feature Selection

Dimensionality reduction and feature selection were conducted using several methods:

3.4.1 Random Forest Analysis

helped in identifying the most impactful features based on their importance scores.

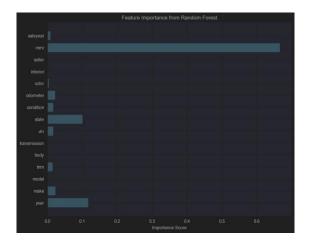


Figure 4: Feature importance using Random Forest.

3.4.2 Principal Component Analysis (PCA)

and condition number analysis were used to reduce the dimensionality while retaining the variance in the data.

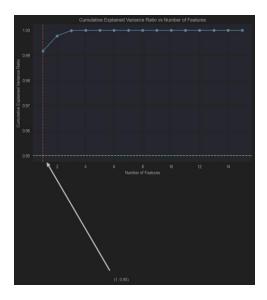


Figure 5: Principal Component Analysis.

3.4.3 Singular Value Decomposition (SVD)

offered insights into the data structure by decomposing it into singular vectors and singular values.

```
[1.13429410e+06 8.91372428e+04 5.29995997e+04 9.51280282e+03
1.90130776e+03 1.42460134e+03 8.99394118e+02 8.56783826e+02
7.70505241e+02 7.59132175e+02 7.50248791e+02 7.48140814e+02
7.15813121e+02 6.97626370e+02 6.44905276e+02] (558825, 15) (15, 15)
```

Figure 6: Single Value Decomposition Analysis

3.4.4 Variance Inflation Factor (VIF)

was calculated to check for multicollinearity among features, ensuring that the selected features are independent of each other.

Observations from these methods guided the selection of features that contribute meaningfully to the analysis.

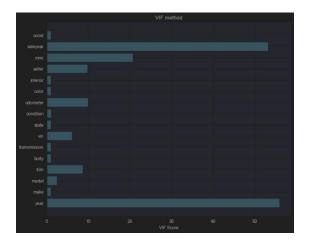


Figure 7: VIF Analysis.

After carefully analyzing all the methods mentioned above, below are final selected features for the Regression Analysis.

['year', 'make', 'model', 'trim', 'condition', 'odometer', 'sellingprice', 'mmr']

3.5 Discretization & Binarization

Label Encoding and One Hot Encoding were applied to convert categorical variables into numerical ones, necessary for modeling. Care was taken to avoid the dummy variable trap by dropping one category per feature, ensuring models do not encounter multicollinearity.

	year	make	model	trim	body	transmission	vin	state	condition	\
0	2015	24	658	998	36	1	403792	29	5.0	
1	2015	24	658	998	36	1	403784	29	5.0	
2	2014	3	9	283	37	1	506996	29	45.0	
3	2015	51	596	1399	37	1	546639	29	41.0	
4	2014	3	42	498	37	1	508029	29	43.0	
	odome	ter c	color	interio	r sel	ler mmr	sellingp	rice s	saleyear	
0	1663	9.0	43		1 7	201 20500.0	215	00.0	2014	
1	939	3.0	43		0 7	201 20800.0	215	00.0	2014	
2	133	1.0	33		1 4	978 31900.0	300	00.0	2015	
3	1428	2.0	43		1 13	794 27500.0	277	50.0	2015	
4	264	1.0	33		1 4	978 66000.0	670	00.0	2014	

Figure 8: Label Encoding and One Hot encoding.

3.6 Variable Transformation

Data normalization and standardization were used to bring all features into a similar scale, eliminating the domination of features with larger scales. Differencing was used to stabilize the mean of time series data by removing changes in the level of a time series, helping to stabilize the variance.

```
-----STANDARDIZE THE DATA------
                       trim
                              body transmission
                                                   vin state condition
          make model
   year
0 1.251 0.142 1.389 0.161 -0.235
                                           -0.18 0.809 -1.141
                                                                  -1.936
1 1.251 0.142 1.389 0.161 -0.235
                                           -0.18 0.809 -1.141
                                                                  -1.936
  0.999 -1.285 -1.566 -1.507 -0.179
                                           -0.18 1.459 -1.141
                                                                   1.080
  1.251 1.976 1.107 1.096 -0.179
                                           -0.18 1.708 -1.141
                                                                   0.779
  0.999 -1.285 -1.416 -1.005 -0.179
                                           -0.18 1.465 -1.141
                                                                   0.930
   odometer color interior seller
                                           sellingprice
                                      mmr
                                                       saleyear
0
    -0.968 1.145
                     -0.706
                             0.001 0.695
                                                  0.809
                                                          -0.059
1
    -1.104 1.145
                     -0.941
                             0.001 0.726
                                                  0.809
                                                          -0.059
                    -0.706 -0.545 1.873
    -1.255 -0.357
                                                  1.681
                                                           0.014
    -1.012 1.145
                             1.621 1.418
                     -0.706
                                                  1.450
                                                           0.014
    -1.230 -0.357
                     -0.706
                            -0.545 5.396
                                                  5.476
                                                          -0.059
```

Figure 9: Standardized Data

3.7 Anomaly Detection/Outlier Analysis and Removal

Anomaly detection was performed using distance-based and density-based methods to identify and remove outliers that could potentially skew the results. The impacts of these outliers were analyzed, and decisions were made based on their potential influence on the overall dataset.



Figure 10: Outlier Analysis Using Z-Score \geq 3.

3.8 Sample Covariance Matrix

The covariance matrix of the dataset was displayed using a heatmap to visualize the covariance between different variables. Observations from this analysis provided insights into the relationships and dependencies between variables.



Figure 11: Sample Covariance Matrix.

3.9 Sample Pearson Correlation Coefficients Matrix

Similarly, a heatmap of the Pearson correlation coefficients was generated to assess the linear relationship between pairs. This helped in understanding the bivariate relationships and the strength of the linear associations between variables. From the correlation matrix we can observe that 'mmr' and 'sellingprice' are highly correlated.



Figure 12: Pearson Correlation Matrix.

3.10 Balancing the Data

The dataset was evaluated for balance between different classes in the target variable. Given an imbalance, techniques such as SMOTE (Synthetic Minority Oversampling Technique) were employed to artificially balance the dataset, ensuring that the model's training process is not biased towards the majority class.

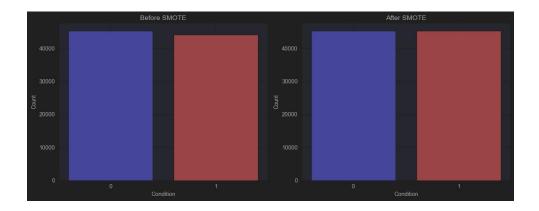


Figure 13: Data Balancing using SMOTE.

4 Phase II: Regression Analysis

This phase focuses on developing a multiple linear regression model to predict a continuous numerical feature of the dataset. The objective is to establish a reliable predictive model and evaluate its performance through various statistical measures and tests.

4.1 Regression Model Development

A multiple linear regression model was constructed using relevant predictors identified during the exploratory data analysis phase. This model aims to predict the selling price of cars based on their attributes such as year, make, model, and odometer readings.

4.2 T-test and F-test Analysis

- **T-test:** Each coefficient in the regression model was tested to determine if it significantly differed from zero. This helps in understanding which features have a statistically significant impact on the selling price.
- **F-test:** The overall significance of the regression model was evaluated using an F-test. This test checks the null hypothesis that no linear relationship exists between the dependent variable and the set of predictors.

4.3 Model Fitting and Diagnostics

The model was fitted using the ordinary least squares method. Key diagnostic plots were used to check for assumptions:

- Residual plots to assess homoscedasticity.
- Q-Q plots to check the normality of residuals.

4.4 Model Performance Evaluation

The model's performance was evaluated using the following criteria:

- **R-squared and Adjusted R-square:** Measures of how well the variations in selling prices are explained by the car attributes.
- AIC and BIC: Information criteria for model comparison, balancing the model fit and complexity.
- MSE (Mean Squared Error): A measure of the average of the squares of the errors—i.e., the average squared difference between the estimated values and the actual value.

4.5 Visualization of Model Predictions

A plot was created to show actual vs. predicted prices, including both training and testing data sets to visualize model effectiveness. The plot helps in visually assessing the accuracy of predictions made by the regression model.

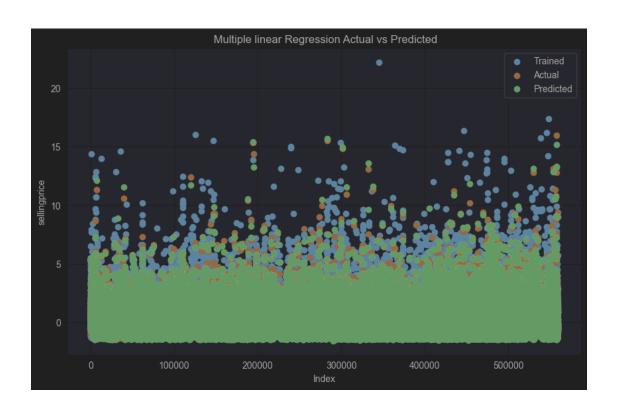


Figure 14: Actual vs Predicted vs Trained.

START OF MULTIPLE LINEAR REGRESSION							
OLS Regression Results							
=======================================							
Dep. Variable:	sellingprice	R-squared:	0.970				
Model:	OLS	Adj. R-squared:	0.970				
Method:	Least Squares	F-statistic:	2.041e+06				
Date:	Tue, 30 Apr 2024	Prob (F-statistic):	0.00				
Time:	00:49:57	Log-Likelihood:	1.4720e+05				
No. Observations:	447050	AIC:	-2.944e+05				
Df Residuals:	447042	BIC:	-2.943e+05				
Df Model:	7						
Covariance Type:	nonrobust						

Figure 15: OLS Summary

	coef	std err	t	P> t	[0.025	0.975]		
const	8.049e-05	0.000	0.309	0.757	-0.000	0.001		
year	-0.0176	0.000	-40.776	0.000	-0.018	-0.017		
make	-0.0016	0.000	-6.037	0.000	-0.002	-0.001		
model	0.0009	0.000	3.518	0.000	0.000	0.001		
trim	-0.0010	0.000	-3.799	0.000	-0.002	-0.000		
condition	0.0506	0.000	181.789	0.000	0.050	0.051		
odometer	-0.0059	0.000	-13.862	0.000	-0.007	-0.005		
mmr	0.9766	0.000	2887.629	0.000	0.976	0.977		
=======			:=======					
Omnibus:		479229.	340 Durbir	n-Watson:		1.998		
Prob(Omnibu	us):	0.	.000 Jarque	Jarque-Bera (JB): Prob(JB):		5846010473.180		
Skew:		3.	.899 Prob(J			0.00		
Kurtosis:		563.	164 Cond.	No.		3.33		
Notes:								
[1] Standar	rd Errors ass	ume that th	ne covariance	matrix of	the errors i	s correctly		
specified.								
MSE: 0.0295	5817785944886	3						

Figure 16: OLS Summary

4.6 Confidence Interval Analysis

Confidence intervals for regression coefficients were calculated, providing a range of values that are likely to contain the true value of the coefficients with a certain probability.

4.7 Stepwise Regression

Stepwise regression was employed to refine the model further:

- Features were selected based on their statistical significance and contribution to the model's explanatory power.
- Adjusted R-square values were monitored to ensure that the model complexity did not unduly increase without substantial gain in model performance.

```
Process and Justification for Backward Stepwise Regression:

Eliminated Feature AIC BIC Adjusted R-squared P-value \
0 model -294375.235 -294287.151 0.97 4.343e-04
1 trim -294364.856 -294287.783 0.97 1.080e-03

Selected Features Iteration
0 [const, year, make, trim, condition, odometer,... 1
1 [const, year, make, condition, odometer, mmr] 2

Final Selected Features: ['const', 'year', 'make', 'condition', 'odometer', 'mmr']
```

Figure 17: Stepwise Regression Iterations Summary

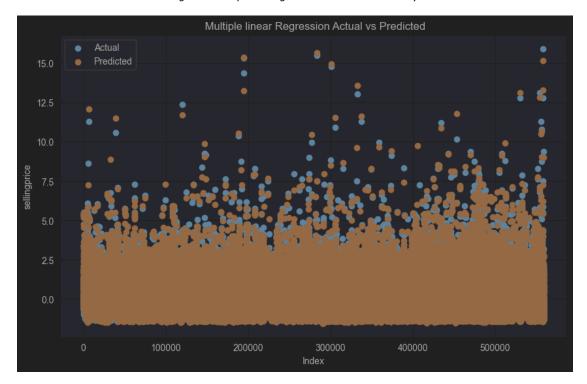


Figure 18: Stepwise Regression Actual vs Predicted.

4.8 Final Regression Model and Prediction

The final model was presented along with the regression equation. Predictions from this model were used to estimate the selling prices based on the car attributes. Observations and insights derived from the model predictions were discussed, providing actionable intelligence for stakeholders.

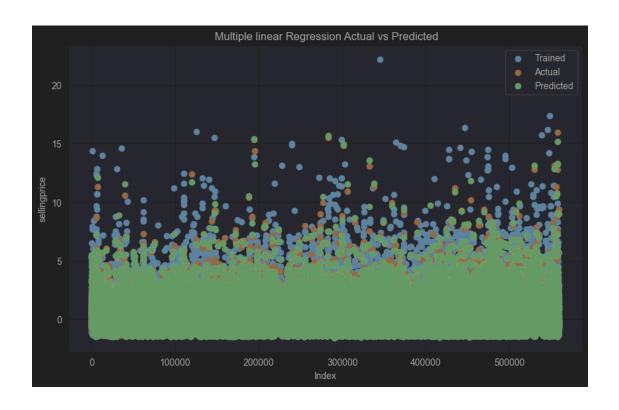


Figure 19: Final Model Actual vs Predicted vs Trained Data

Regression Equation:

$$y = 0.0001 + -0.0176*X1 + -0.0016*X2 + 0.0009*X3 + -0.0010*X4 + 0.0506*X5 + -0.0059*X6 + 0.9766*X7$$

4.9 Model Summary Table

A comprehensive table summarizing the regression analysis results was included. This table featured R-squared, adjusted R-square, AIC, BIC, and MSE, facilitating a clear understanding of the model's efficacy.

Metric	Value
R-squared	0.97
Adjusted R-squared	0.97

AIC	-0.02944
BIC	-0.02935
MSE	0.0295

Table 1: Summary of Regression Model Metrics

The linear regression model demonstrates a very high level of accuracy and reliability in predicting the target variable, as evidenced by the following key metrics:

- R-squared: 0.97 This value indicates that 97% of the variance in the dependent variable is explained by the independent variables included in the model. An R-squared value close to 1 suggests that the model has a very high level of predictive accuracy.
- Adjusted R-squared: 0.97 Adjusted R-squared also stands at 0.97, which is
 particularly noteworthy because it adjusts the R-squared value to account for
 the number of predictors in the model. This high value confirms that the model
 fits the data very well, even after the adjustment for the number of variables.
- MSE (Mean Squared Error): 0.0295 The MSE, which measures the average of
 the squares of the errors, is quite low. This indicates that the predictions made
 by the model are very close to the actual data points, leading to a minimal
 deviation in the predicted values from the real values.
- AIC (Akaike Information Criterion): -0.02944 AIC is a criterion used for model selection among a finite set of models. The lower the AIC, the better the model is at explaining the variation in the data while keeping the model simplicity. The negative value here suggests a model that balances model complexity and fit efficiently.
- **BIC (Bayesian Information Criterion): -0.02935** Similar to AIC, BIC is used for model selection, but it includes a penalty term for the number of parameters in the model. The negative BIC value indicates a model that is optimal for balancing between goodness of fit and model complexity, in a Bayesian context.

Overall, the combination of high R-squared and Adjusted R-squared values, along with a low MSE, and very favorable AIC and BIC values, strongly suggests that the linear regression model provides an excellent fit to the data. It effectively captures the

essential patterns and can be considered a reliable tool for prediction within this specific context. Future work might explore potential improvements or validations with additional data, but the current model stands as a robust analytical tool.

5 Phase III: Classification Analysis

In this phase, various machine learning classifiers were applied to predict a target variable, with an emphasis on evaluating and optimizing each model to achieve the best possible performance without overfitting or underfitting. Extensive hyperparameter tuning was conducted through grid search for each classifier.

5.1 Decision Tree Classifier

5.1.1 Model Optimization

Decision Trees were optimized using pre-pruning, post-pruning, and hyperparameter tuning via grid search. Parameters such as 'criterion', 'splitter', 'max depth', 'min samples _split', 'max features', and 'ccp alpha' were systematically tested.

```
Decision Tree Classifier Results:

| Confusion Matrix | Precision | Recall | F1-Score | ROC AUC |
| [[7789, 3539], | 0.675 | 0.667 | 0.671 | 0.677 |
| [3667, 7358]] | | | | |
```

Figure 20: Basic Decision Tree Modal Metrics

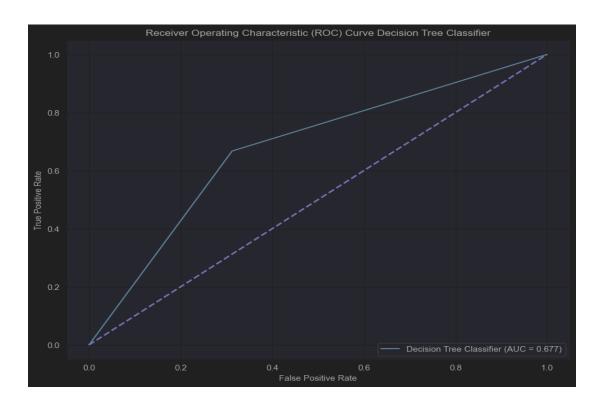


Figure 21: without tuning.

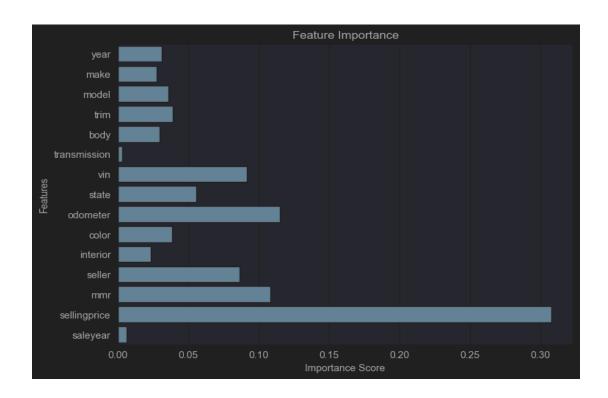


Figure 22: Feature Importances in Decision Tree

5.1.2 Cost Complexity Optimization

The Cost Complexity Pruning parameter, ccp alpha, was adjusted to find the optimal balance between tree depth and error reduction, minimizing overfitting.

Pre-Pruning

{'criterion': 'gini', 'max_depth': 2, 'max_features': None, 'min_samples_leaf': 2, 'min_samples_split': 2, 'splitter': 'best'}

Accuracy of pre tuned tree 0.707.

Figure 23: Pre pruning Decision Tree Metrics

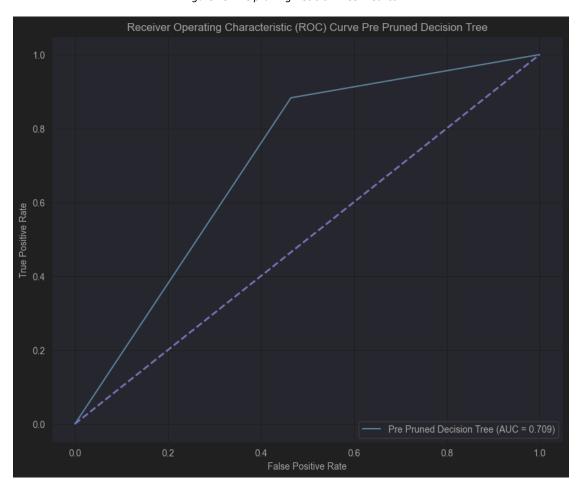


Figure 24: Pre-Pruning Decision Tree Roc Curve

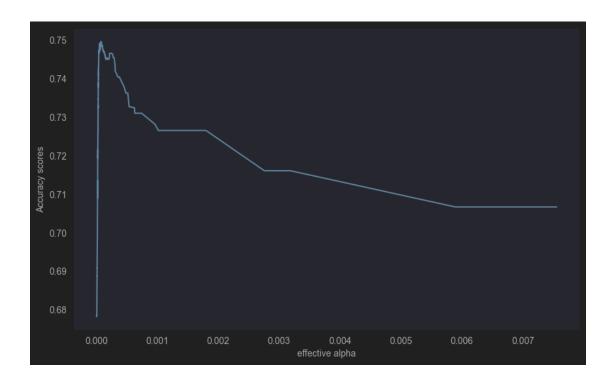


Figure 25: Effective Alpha values Vs Accuracy scores.

5.1.3 Performance Metrics

The performance of the optimized Decision Tree model was assessed using metrics such as precision, recall, F1-score, and the area under the ROC curve (AUC). Stratified K-fold cross-validation was used to ensure robustness against overfitting.

```
accuracy for post pruning : 0.75

pre metrics [[6061 5267]

[1290 9735]] 0.8829931972789116 0.7066612982597414 0.7090195506168571

Pruning Method Accuracy Confusion Matrix Recall AUC

0 Pre-Pruned 0.707 [[6061, 5267], [1290, 9735]] 0.883 0.709

1 Post-Pruned 0.750 [[7961, 3367], [2232, 8793]] 0.798 0.750

Logistic Regression Accuracy: 0.735
```

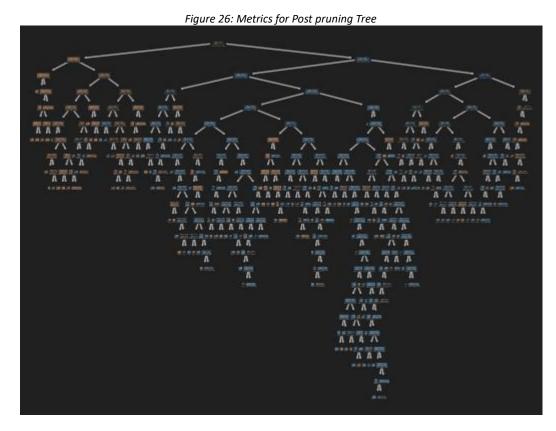


Figure 27: Post pruning Tree.

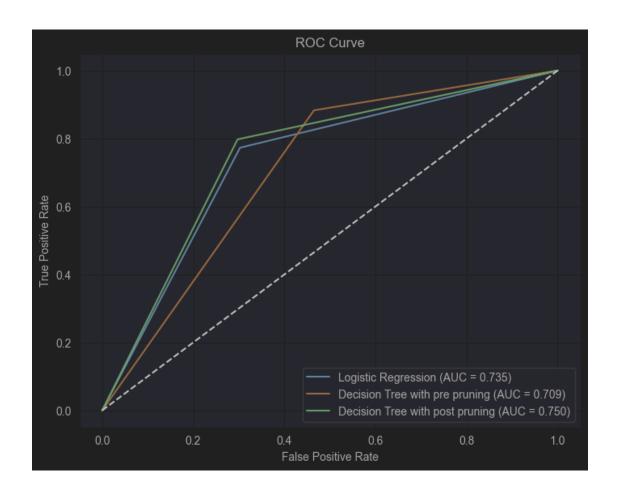


Figure 28: Comparing Post Pruning, Pre-Pruning, Logistic Regression

5.2 Logistic Regression

5.2.1 Hyperparameter Tuning

Grid search was employed to optimize the regularization strength and the choice of solver in Logistic Regression.

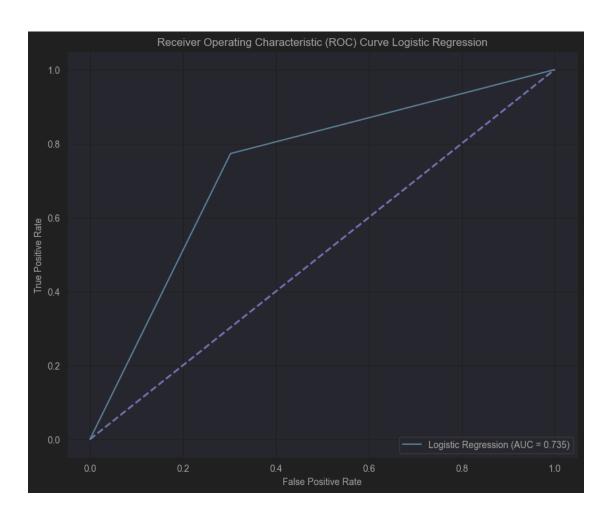


Figure 29: Logistic regression ROC Curve

5.2.2 Performance Metrics

The efficacy of the Logistic Regression model was evaluated using a confusion matrix, precision, recall, specificity, F1-score, and ROC AUC analysis.

Figure 30: Logistic Regression Metrics

5.3 K-Nearest Neighbors (KNN)

5.3.1 Optimum K Determination

The optimal number of neighbors, K, was determined using the elbow method, which involved plotting error rates for different values of K and choosing the K at the elbow point, where the decrease in error rate diminishes.

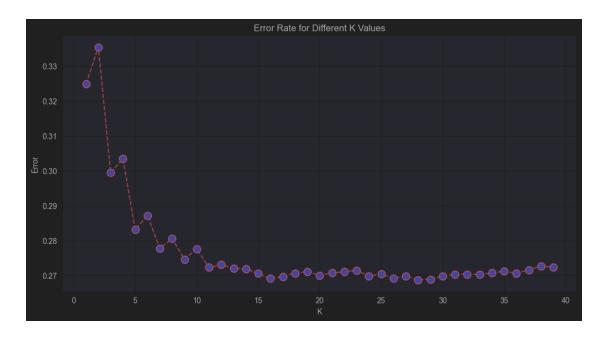


Figure 31: Error rate vs K

5.3.2 Performance Metrics

KNN performance was detailed through the same set of metrics, including sensitivity and specificity, supported by graphical representation of ROC curves.

```
Minimum error: 0.26868876660850893

Optimum K: 28

KNN Results:

| Confusion Matrix | Precision | Recall | F1-Score | ROC AUC |
| [7739, 3589], | 0.706 | 0.781 | 0.741 | 0.732 |
| [2417, 8608]] | | | | |
```

Figure 32: Performance Metrics of KNN

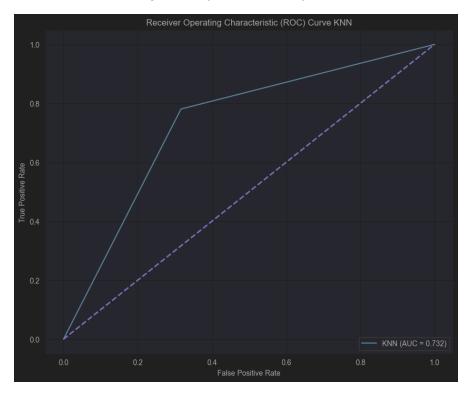


Figure 33: ROC Curve for KNN

5.4 Support Vector Machine (SVM)

5.4.1 Kernel Optimization

SVM was tested with different kernels: linear, polynomial, and radial base function (RBF) to determine which kernel provides the best separation between classes.

5.4.2 Performance Metrics

The performance was quantified and compared across different kernels using precision, recall, specificity, F1-score, and the area under the ROC curve.

Figure 34:Performance Metrics of SVM

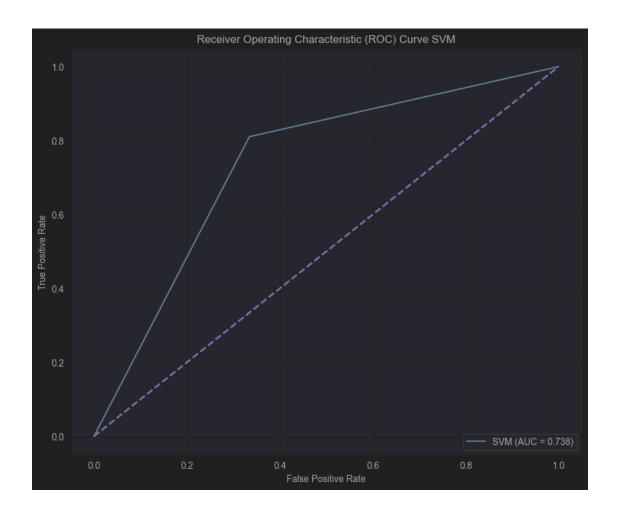


Figure 35: ROC Curve for SVM

5.5 Naïve Bayes

5.5.1 Model Application

The Naïve Bayes classifier was applied, given its effectiveness in high-dimensional datasets.

Figure 36: Naive Bayes Performance Metrics

5.5.2 Performance Metrics

Performance was evaluated using standard classification metrics and visualized through ROC curve analysis.

Stratified K-fold Cross-Validation F1-Score: 0.010

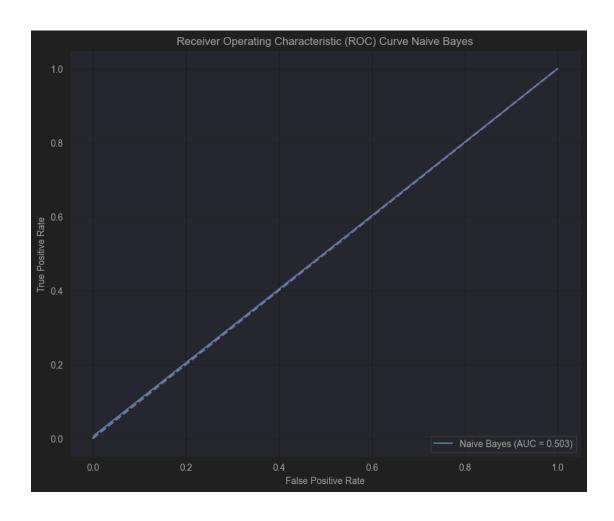


Figure 37: ROC Curve for Naive Bayes

5.6 Random Forest

5.6.1 Ensemble Techniques

Random Forest's performance was enhanced using bagging, stacking, and boosting to reduce variance and bias.

5.6.2 Performance Metrics

Metrics including MSE from bagging and boosting scenarios were analyzed, alongside traditional classification metrics.

Figure 38: Random Forest bagging metrics

Figure 39: Random Forest stacking metrics.

Figure 40: Random Forest Boosting metrics.

5.7 Neural Network

5.7.1 Network Configuration

A Multi-layer Perceptron (MLP) was configured with multiple layers and neurons to optimize learning capability.

5.7.2 Performance Metrics

Neural Network performance was critically assessed using confusion matrices, ROC curves, and cross-validation scores to ensure the model's generalization.

Figure 41: MLP metrics.

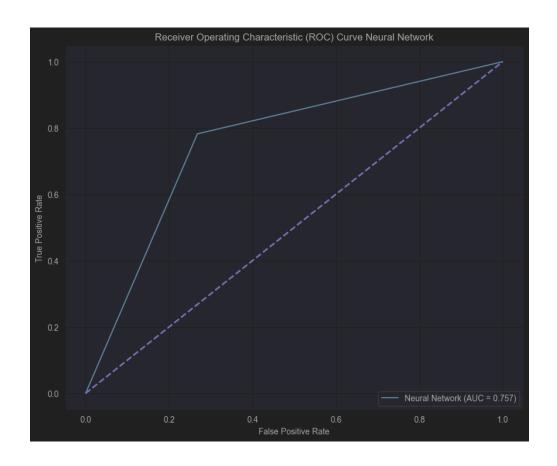


Figure 42: MLP ROC Curve

5.8 Comparative Analysis and Recommendations

A comprehensive comparative analysis was conducted to identify the classifier with the highest overall performance. Recommendations were made based on a combination of performance metrics, computational efficiency, and ease of model interpretation.

Classifier	Precision	Recall	F1-Score	AUC	Confusion Matrix	F1- Score (K-fold)
Decision Tree (Post pruning)	0.75	0.798	0.759	0.75	[[7961, 3367], [2232, 8793]]	0.758
Logistic Regression	0.713	0.773	0.742	0.735	[[7899, 3429],	0.753

KNN	0.706	0.781	0.741	0.732	[2503, 8522]] [[7739, 3589], [2417, 8608]]	0.723
SVM	0.702	0.811	0.752	0.738	[[7538, 3790], [2089, 8936]]	0.759
Naïve Bayes	0.744	0.008	0.016	0.503	[[11298, 30], [10938, 87]]	0.010
Random Forest	0.734	0.785	0.759	0.755	[[8196, 3132], [2365, 8660]]	0.767
Neural Network	0.740	0.782	0.760	0.757	[[8290, 3038], [2399, 8626]]	0.758

Table 2: Comparison of Classifier Performance

6 Phase IV: Clustering and Association [Independent Study]

This phase involves the application of clustering and association rule mining algorithms to analyze the dataset independently. The objective is to identify inherent groupings within the data and discover patterns or rules that could inform further analysis or operational strategies.

6.1 K-means or K-means++ Clustering

6.1.1 Algorithm Implementation

The K-means (or K-means++) algorithm was applied to segment the dataset into clusters. K-means++ was chosen for its efficiency in cluster center initialization, reducing the likelihood of poor cluster formation.

6.1.2 Silhouette Analysis

Silhouette analysis was conducted to determine the optimal number of clusters, *k*. This method assesses the consistency within clusters and is instrumental in validating the homogeneity of data points within the same cluster.

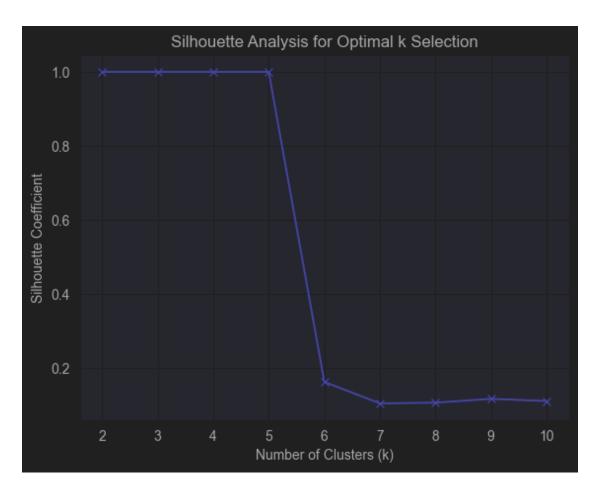


Figure 43:Silhouette Analysis for Optimal K Selection.

6.1.3 Within-Cluster Variation

The within-cluster sum of squares (WCSS) was plotted against the number of clusters to observe the point where the marginal gain in explained variance drops significantly (the elbow method), indicating the optimal k.

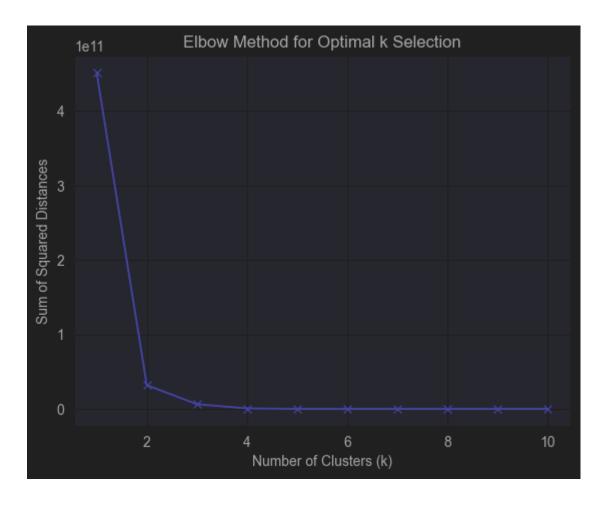


Figure 44: Elbow method for optimal K selection.

6.1.4 Observations

The clustering results were visualized using scatter plots, and clusters were analyzed based on the characteristics of data points within them. Observations about the clusters' composition and their practical implications were discussed.

6.2 DBSCAN Algorithm

6.2.1 Algorithm Implementation

The Density-Based Spatial Clustering of Applications with Noise (DBSCAN) was employed, which is excellent for identifying clusters of varying shapes and sizes in a dataset with noise.

6.2.2 Parameter Optimization

Parameters such as the minimum number of points per cluster and the epsilon distance were tuned to optimize clustering outcomes.

6.2.3 Observations

Clusters identified by DBSCAN were analyzed, and the algorithm's effectiveness in handling outliers was noted. Observations were detailed regarding the spatial distribution of clusters and their densities.

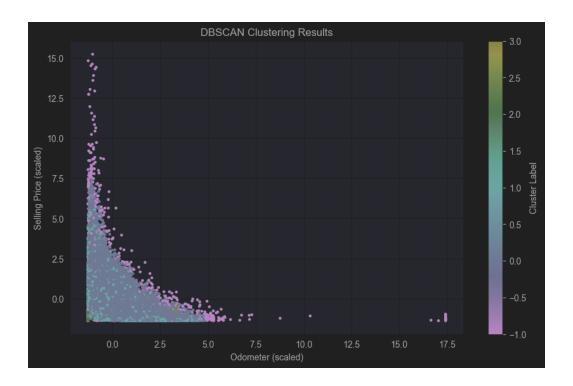


Figure 45: DBSCAN Clustering Results

6.3 Apriori Algorithm for Association Rule Mining

6.3.1 Algorithm Implementation

The Apriori algorithm was used to uncover relationships between features in the dataset. This analysis was aimed at finding frequent itemsets and deriving association rules that have significant support and confidence.

6.3.2 Rule Generation

Rules generated by the Apriori algorithm were evaluated based on their support, confidence, and lift. Only rules that exceeded predetermined thresholds were considered significant.

```
-----ASSOCIATION RULE MINING------
Processing 12 combinations | Sampling itemset size 2
               antecedents
                                        consequents antecedent support
               (make_Ford) (transmission_automatic)
0
          (make_Chevrolet) (transmission_automatic)
                                                                  0.108
               (trim_Base) (transmission_automatic)
                                                                  0.116
  (transmission_automatic)
                                        (make_Ford)
                                                                  0.970
  (transmission_automatic)
                                        (trim_Base)
                                                                  0.970
  (transmission_automatic)
                                    (make_Chevrolet)
                                                                  0.970
  consequent support confidence
                                           lift
                                                           conviction \
                                                  leverage
2
                                   0.976 1.006 1.101e-03
                                                                 1.248
               0.970
                        0.179
0
               0.970
                        0.105
                                   0.975 1.005 5.291e-04
                                                                 1.195
               0.970
                        0.111
                                   0.951 0.981 -2.184e-03
                                                                 0.616
               0.184
                        0.179
                                   0.185 1.006 1.101e-03
                                                                 1.001
               0.116
                        0.111
                                   0.114 0.981 -2.184e-03
                                                                 0.997
               0.108
                        0.105
                                   0.108 1.005 5.291e-04
                                                                 1.001
```

Figure 46: Association Rule Mining Using Apriori

6.3.3 Observations

The implications of the most relevant association rules were discussed, particularly in the context of their utility for cross-selling strategies or for segmenting the market based on feature associations.

6.3.4 Conclusion

This phase provided deep insights into the structural patterns of the dataset through clustering and explored potential relationships via association rule mining. The findings from this phase are intended to help in strategic decision-making and to highlight areas for further data-driven investigations.

7 Recommendations

This section synthesizes the insights gleaned from the project and offers strategic recommendations based on the analysis conducted in the previous phases. The recommendations are aimed at enhancing predictive modeling and providing actionable insights for operational or strategic implementations.

7.1 Learnings from the Project

- The project underscored the importance of rigorous data preprocessing and exploratory data analysis as foundational steps in predictive modeling.
- It demonstrated the effectiveness of various classifiers and their applicability depending on the nature and distribution of the data.
- The exercise of tuning hyperparameters and selecting features highlighted the delicate balance between model complexity and performance.

7.2 Best Performing Classifiers

The performance analysis of various classifiers on the dataset has produced insightful outcomes, especially in terms of precision, recall, F1-score, AUC (Area Under the Curve), and cross-validation results. Here's a summary of each classifier's performance:

 Decision Tree (Post pruning): This model shows a strong balance between recall and precision, resulting in an F1-score of 0.759 and an AUC of 0.75, indicating good model performance. The cross-validation F1-score of 0.758 confirms the model's stability across different subsets of data.

- Logistic Regression: The logistic regression model presents a slightly lower performance compared to the decision tree, with an F1-score of 0.742 and an AUC of 0.735. Its precision is slightly less but still competitive, and the F1-score in cross-validation is 0.753, suggesting consistency in performance.
- KNN (K-Nearest Neighbors): KNN shows a comparable recall to the logistic regression but has a lower precision which impacts its F1-score slightly, landing it at 0.741 with an AUC of 0.732. Its cross-validation score is the lowest among the competitive models at 0.723, indicating potential variability in performance across different data splits.
- **SVM (Support Vector Machine):** SVM performs notably well in terms of recall, which is the highest among the classifiers at 0.811. This high recall contributes to an F1-score of 0.752 and an AUC of 0.738. Its performance is robust as evidenced by a cross-validation F1-score of 0.759.
- Naïve Bayes: This model, while having a reasonable precision, shows a very low recall of 0.008, which drastically affects its F1-score, making it practically negligible at 0.016. The AUC is barely over 0.5, indicating performance close to a random classifier. The model is significantly underperforming in this scenario.
- Random Forest: Exhibiting strong performance, the Random Forest classifier
 has an F1-score and AUC closely mirroring those of the Decision Tree, both at
 0.759 and 0.755, respectively. Its cross-validation score is the highest at 0.767,
 highlighting its robustness and effectiveness across different data segments.
- **Neural Network:** The neural network shows balanced precision and recall, leading to an F1-score of 0.760 and an AUC of 0.757. Its performance is consistent across different folds of data, as shown by the cross-validation F1-score of 0.758.

Conclusion:

Overall, the **Random Forest** and **Neural Network** models demonstrate the most robust and consistent performance across all metrics, making them the preferable choices for classification tasks within this dataset. These models not only provide high F1-scores

and AUCs but also maintain performance stability across different subsets of data, as evidenced by their cross-validation scores. The SVM also presents itself as a strong contender, especially in scenarios where maximizing recall is crucial. Conversely, the **Naïve Bayes** classifier may require reconsideration or significant adjustments due to its inadequate performance in this specific context.

7.3 Improvements in Classification Performance

- Future work could explore more advanced ensemble techniques that might provide better generalization over the dataset.
- Increasing the dataset size or incorporating more diverse data sources could help in enhancing the model's learning and prediction accuracy.
- Continuous tuning of hyperparameters using automated techniques like grid search or random search could yield better results.

7.4 Features Associated with the Target Variable

- Analysis revealed specific features that are strongly associated with the target variable, aiding in understanding the drivers of the dependent variable.
- For instance, if year and odometer readings are identified as key predictors, these should be considered in pricing models or customer segmentation analyses.

7.5 Cluster Analysis

- The optimal number of clusters identified during the cluster analysis provides insights into market segmentation and customer behavior patterns.
- The characteristics of each cluster can be used to tailor marketing strategies or customize product offerings.

8 Appendix

8.1 Python Code

The appendix includes all the Python code developed for this project. This includes scripts for data cleaning, model building, evaluation, and any additional analyses conducted. The code is well-commented to ensure that it is understandable and reusable.

```
import pandas as pd
import numpy as np
from mlxtend.preprocessing import TransactionEncoder
from sklearn.ensemble import RandomForestRegressor
from sklearn.preprocessing import LabelEncoder
import matplotlib.pyplot as plt
from sklearn.decomposition import PCA
from scipy.linalg import svd
from sklearn.preprocessing import StandardScaler
import seaborn as sns
from statsmodels.stats.outliers influence import
variance inflation factor
import statsmodels.api as sm
from sklearn.metrics import mean squared error, silhouette score
from sklearn.model selection import GridSearchCV, train test split,
StratifiedKFold
from sklearn.tree import DecisionTreeClassifier
from sklearn.linear model import LogisticRegression
from sklearn.svm import SVC
from sklearn.naive bayes import GaussianNB
from sklearn.neural network import MLPClassifier
from sklearn.metrics import confusion matrix, precision score,
recall score, f1 score, roc curve, auc
from sklearn.cluster import DBSCAN
from mlxtend.frequent patterns import apriori, association rules
from sklearn.tree import plot tree
from sklearn.metrics import accuracy_score, roc_auc_score
from sklearn.cluster import KMeans
from sklearn.ensemble import BaggingClassifier, StackingClassifier,
AdaBoostClassifier
from sklearn.neighbors import KNeighborsClassifier
from prettytable import PrettyTable
import warnings
from imblearn.over sampling import SMOTE
#%%
RED = "\033[91m"]
GREEN = "\033[92m"]
BLUE = "\033[94m""
RESET = " \setminus 033[0m"]
```

```
warnings.filterwarnings('ignore')
pd.set option('display.precision', 3)
print(GREEN + "\n ---------- Phase I: Feature
Engineering & EDA: -----\n" + GREEN)
car prices df = pd.read csv("car prices.csv")
print("info: \n", car_prices_df.info())
print("Describe: \n", car_prices_df.describe())
print("shape of the database ", car_prices_df.shape)
print("columns : ", car prices df.columns)
missing_values = car_prices_df.isna().sum()
print("\overline{m}issing values in each column:\n",
missing_values.sort_values(ascending=False))
print("total number of missing values in the dataset:",
car prices df.isna().sum().sum())
missing values per column = car prices df.isna().sum()
print("percentage of missing values per column:\n",
(missing values per column / car prices df.shape[0]) * 100)
car prices df = car prices df.dropna(subset=['sellingprice'])
print("missing values:", car prices df.isna().sum())
for column in
car prices df.select dtypes(include=['number']).columns:
    car prices df[column].fillna(car prices df[column].mean(),
inplace=True)
for column in
car prices df.select dtypes(exclude=['number']).columns:
    print("mode of the column", car_prices_df[column].mode()[0])
    car prices df[column].fillna(car prices df[column].mode()[0],
inplace=True)
car prices df['saledate'] = pd.to datetime(car prices df['saledate'],
errors='coerce', utc=True)
car prices df['saleyear'] = car prices df['saledate'].dt.year
car prices df['saleyear'] = car prices df['saleyear'].fillna(-
1).astype(int)
print(car prices df['saleyear'].head())
car prices df = car prices df.drop('saledate', axis=1)
print("total number of missing values after filling: ",
car prices df.isna().sum())
duplicate rows = car prices df[car prices df.duplicated()]
print(duplicate rows.shape)
print("there is no duplicate data in the given dataset")
```

```
print("\n----START OF FEATURE SELECTION & DIMENSIONALITY
REDUCTION----\n")
label_encoders = {}
data = car_prices_df.copy()
for column in data.select dtypes(include=['object']).columns:
    le = LabelEncoder()
    data[column] = le.fit_transform(data[column].astype(str))
label_encoders[column] = le
encoded data = data.copy()
print(data.head())
----\n")
scalar = StandardScaler()
data scaled = scalar.fit transform(data)
data scaled df = pd.DataFrame(data=data scaled, columns=data.columns)
data.update(data scaled df)
print(data.head())
print("\n-----Dimensionality reduction/feature selection-----
X = data.drop(['sellingprice'], axis=1)
y = data['sellingprice']
pca = PCA()
X_pca = pca.fit_transform(X)
print(X.columns)
cumulative variance ratio = np.cumsum(pca.explained variance ratio )
n features 95 = np.argmax(cumulative variance ratio >= 0.95) + 1
plt.figure(figsize=(10, 6))
plt.plot(np.arange(1, len(cumulative variance ratio) + 1),
cumulative variance ratio, marker='o', linestyle='-')
plt.xlabel('Number of Features')
plt.ylabel('Cumulative Explained Variance Ratio')
plt.title('Cumulative Explained Variance Ratio vs Number of
Features')
plt.grid(True)
plt.axvline(x=n features 95, color='r', linestyle='--')
plt.axhline(y=0.95, color='g', linestyle='--')
plt.annotate(f'({n features 95}, 0.95)', xy=(n features 95, 0.95),
xytext=(n features 95 + 5, 0.90),
            arrowprops=dict(facecolor='black', shrink=0.05))
plt.show()
```

```
print("Number of features needed for explaining more than 95% of the
dependent variance:", n features 95)
rf = RandomForestRegressor(n estimators=8, max depth=10,
random state=5805)
rf.fit(X, y)
importances = rf.feature importances
feature_names = X.columns
feature importance dict = dict(zip(feature names, importances))
sorted feature importance = sorted([[key,
feature_importance_dict[key]] for key in feature_importance_dict],
                                 key=lambda x: x[1], reverse=True)
print("feature importance dict: ", sorted_feature_importance)
features = list(feature_importance_dict.keys())
importances = list(feature importance dict.values())
plt.figure(figsize=(10, 8))
plt.barh(features, importances, color='skyblue')
plt.xlabel('Importance Score')
plt.title('Feature Importance from Random Forest')
plt.show()
print("\n----- START OF SINGLE VALUE DECOMPOSITION -----
----\n")
numerical cols = X.select dtypes(include=[np.number])
U, s, Vt = svd(numerical cols, full matrices=False)
print(s, U.shape, Vt.shape)
print("\n----- END OF SINGLE VALUE DECOMPOSITION -----
    ----\n")
print("\n-----\n")
X vif = X.select dtypes(include='number').assign(const=1)
vif data = pd.DataFrame({
    'Variable': X vif.columns,
    'VIF': [variance inflation factor(X vif.values, i) for i in
range(X vif.shape[1])]
vif sorted = vif data.sort values(by='VIF', ascending=False)
print(vif sorted)
plt.figure(figsize=(10, 8))
plt.barh(vif data['Variable'], vif data['VIF'], color='skyblue')
plt.xlabel('VIF Score')
```

```
plt.title('VIF method')
plt.show()
print("\n-----\n")
final_selected_features = ['year', 'make', 'model', _'trim',
'condition', 'odometer', 'sellingprice', 'mmr']
print("\n----- START OF ANOMALY REMOVAL/OUTLIER DETECTION
(Z-SCORE) -----\n")
cleaned carprices df = data[final selected features]
z scores = np.abs(
    (cleaned_carprices_df['sellingprice'] -
cleaned carprices df['sellingprice'].mean()) / cleaned carprices df[
        'sellingprice'].std())
data_cleaned_z = cleaned_carprices_df[z_scores < 3]</pre>
print("df shapes after anomaly deletion: ",
cleaned carprices df.shape, data cleaned z.shape)
z scores = np.abs((cleaned carprices df['sellingprice'] -
cleaned carprices df['sellingprice'].mean()) /
cleaned carprices df['sellingprice'].std())
outliers = z scores >= 3
plt.figure(figsize=(10, 6))
plt.scatter(cleaned carprices df.index,
cleaned_carprices_df['sellingprice'], c='blue', label='Non-outliers')
plt.scatter(cleaned carprices df[outliers].index,
cleaned carprices df[outliers]['sellingprice'], c='red',
label='Outliers')
plt.title('Outliers in Selling Price Using Z-score')
plt.xlabel('Index')
plt.ylabel('Selling Price')
plt.legend()
plt.show()
print("\n----- END OF ANOMALY REMOVAL/OUTLIER DETECTION
(Z-SCORE) -----\n")
print("\n----- START OF COVARIANCE MATRIX ------
---\n")
covariance matrix = encoded data[final selected features].cov()
plt.figure(figsize=(10, 10))
sns.heatmap(covariance matrix, annot=True, fmt=".2f",
cmap='coolwarm', cbar=True)
plt.title('Heatmap of the Sample Covariance Matrix')
plt.show()
print("\n----- END OF COVARIANCE MATRIX --------
--\n")
print("\n----- START OF CORRELATION MATRIX ------
```

```
----\n")
correlation matrix = data cleaned z.corr()
plt.figure(figsize=(10, 8))
sns.heatmap(correlation matrix, annot=True, fmt=".2f",
cmap='coolwarm', center=0)
plt.title('Heatmap of the Sample Pearson Correlation Coefficients
Matrix')
plt.show()
print("\n----- END OF CORRELATION MATRIX ------
---\n")
#응응
print(BLUE + "Phase II: Regression Analysis" + BLUE)
print("\n----- START OF MULTIPLE LINEAR REGRESSION -----
----\n")
X = data cleaned z.drop('sellingprice', axis=1)
y = data cleaned z['sellingprice']
X train, X test, y train, y test = train test split(X, y,
test size=0.2, random state=5805)
X train = sm.add constant(X train)
X test = sm.add constant(X test)
model = sm.OLS(y train, X train).fit()
model summary = model.summary()
print(model summary)
predictions = model.predict(X test)
mse = mean squared error(y test, predictions)
print(f"MS\overline{E}: {mse}\overline{"})
predicted values = model.predict(X test)
plt.figure(figsize=(10, 6))
plt.scatter(y_train.index, y_train, label='Trained')
plt.scatter(y_test.index, y_test, label='Actual')
plt.scatter(y test.index, predicted values, label='Predicted')
plt.title(' Multiple linear Regression Actual vs Predicted')
plt.xlabel('Index')
plt.ylabel('sellingprice')
plt.legend()
plt.show()
----\n")
```

```
#응응
def backward stepwise regression(X train, X test, y train, y test,
threshold=0.00001):
    selected features = list(X train.columns)
    all results = []
    iteration = 1
    while len(selected features) > 0:
        X_train_selected = X_train[selected_features]
        X train selected = sm.add constant(X train selected)
        X_test_selected = X_test[selected_features]
        X test selected = sm.add constant(X test selected)
        model = sm.OLS(y_train, X_train_selected).fit()
        p values = model.pvalues[1:]
        eliminated feature = None
        aic = model.aic
        bic = model.bic
        adj r2 = model.rsquared adj
        p value = None
        if p values.max() > threshold:
            eliminated feature = p values.idxmax()
            selected features.remove(eliminated feature)
            p value = p values.max()
        else:
            break
        results = {'Eliminated Feature': eliminated feature,
                   'AIC': aic, 'BIC': bic, 'Adjusted R-squared':
adj r2,
                   'P-value': p_value,
                   'Selected Features': selected features.copy()}
        all results.append(results)
        print("Model summary at iteration {} ".format(iteration))
        print(model.summary())
        iteration += 1
    final selected features = selected features
    return all results, final selected features
all results, final selected features =
backward stepwise regression(X train, X test, y train, y test)
results dfs = []
for i, results in enumerate(all results):
    results df = pd.DataFrame([results])
```

```
results df['Iteration'] = i + 1
    results dfs.append(results df)
results df = pd.concat(results dfs, ignore index=True)
print("Process and Justification for Backward Stepwise Regression:")
print(results_df)
print("\nFinal Selected Features:", final_selected_features)
X train, X test, y train, y test = train test split(X, y,
test size=0.2, random state=5805)
final model = sm.OLS(y train, X train).fit()
predicted values = final model.predict(X test)
plt.figure(figsize=(10, 6))
plt.scatter(y_test.index, y_test, label='Actual')
plt.scatter(y test.index, predicted values, label='Predicted')
plt.title(' Multiple linear Regression Actual vs Predicted')
plt.xlabel('Index')
plt.ylabel('sellingprice')
plt.legend()
plt.show()
mse = mean squared error(y test, predicted values)
print("Mean Squared Error (MSE) BACKWARD STEPWISE REGRESSION :", mse)
print("\n---- END OF MULTIPLE LINEAR REGRESSION
(BACKWARD STEPWISE REGRESSION) -----\n")
#%%
print("\n-----Phase III: Classification Analysis:-----
  ----\n")
print(car prices df.info())
downsampled df = data.sample(frac=0.2, replace=False, random state=1)
X = downsampled df.drop(['condition'], axis=1)
y = downsampled df['condition']
threshold value = y.median()
y = pd.cut(y, bins=[-float('inf'), threshold value, float('inf')],
labels=[0, 1])
X train, X test, y train, y test = train test split(
    X, y, test size=0.2, stratify=y, random state=5805
print(y.value counts())
```

```
fig, axes = plt.subplots(nrows=1, ncols=2, figsize=(12, 5))
# Plotting class distribution before SMOTE
axes[0].bar([0, 1], y_train.value_counts(), color=['blue', 'red'])
axes[0].set_title('Before SMOTE')
axes[0].set_xlabel('Condition')
axes[0].set_ylabel('Count')
axes[0].set xticks([0, 1])
# Applying SMOTE to balance the dataset
smote = SMOTE(random state=5805)
X train, y train = smote.fit resample(X train, y train)
# Printing class distribution after applying SMOTE
print(y train.value counts())
# Plotting class distribution after SMOTE
axes[1].bar([0, 1], y train.value counts(), color=['blue', 'red'])
axes[1].set title('After SMOTE')
axes[1].set xlabel('Condition')
axes[1].set_ylabel('Count')
axes[1].set xticks([0, 1])
plt.tight layout()
plt.show()
def evaluate model(preds, y test, model name, dt best):
    print(f'{model name} Results:')
    table = PrettyTable()
    # Adding headers for all metrics
    table.field names = ["Confusion Matrix", "Precision", "Recall",
"F1-Score", "ROC AUC"]
    # Compute metrics
    precision = precision score(y test, preds)
    recall = recall score(y test, preds)
    f1 = f1 score(y test, preds)
    # Compute ROC AUC
    fpr, tpr, _ = roc curve(y test, preds)
    roc auc = auc(fpr, tpr)
    # Format confusion matrix
    cm = confusion_matrix(y_test, preds)
    cm_str = f''[[\{cm[0, 0]\}, \{cm[0, 1]\}], \n [\{cm[1, 0]\}, \{cm[1, 0]\}, \]
1]}]"
    # Add rows for metrics
    table.add row([cm str, f"{precision:.3f}", f"{recall:.3f}",
f"{f1:.3f}", \overline{f}"{roc auc:.3f}"])
    print(table)
    plt.figure(figsize=(10, 8))
```

```
plt.plot(fpr, tpr, label=f'{model_name} (AUC = {roc_auc:.3f})')
plt.plot([0, 1], [0, 1], color="navy", lw=2, linestyle="--")
    plt.xlabel("False Positive Rate")
    plt.ylabel("True Positive Rate")
    plt.title("Receiver Operating Characteristic (ROC) Curve
{}".format(model_name))
    plt.legend(loc="lower right")
    plt.show()
    skf = StratifiedKFold(n splits=5, random state=42, shuffle=True)
    scores = []
    for train idx, test idx in skf.split(X, y):
        X train fold, X test fold = X.iloc[train idx],
X.iloc[test idx]
        y_train_fold, y_test_fold = y.iloc[train idx],
y.iloc[test idx]
        model = dt_best.fit(X_train_fold, y_train_fold)
        preds = model.predict(X test fold)
        scores.append(f1 score(y test fold, preds))
    print(f'Stratified K-fold Cross-Validation F1-Score:
{np.mean(scores):.3f}')
    print('----')
#%%
dt model = DecisionTreeClassifier(random state=5805)
dt model.fit(X train, y train)
y pred = dt model.predict(X test)
evaluate model(y pred, y test, "Decision Tree Classifier", dt model)
accuracy = accuracy_score(y_test, y_pred)
print("Accuracy:", accuracy)
importance = pd.Series(dt model.feature importances ,
index=X.columns)
print("Feature Importance:", importance)
print("feature to be removed :", importance.idxmin())
plt.figure(figsize=(8, 6))
sns.barplot(x=importance, y=importance.index)
plt.title("Feature Importance")
plt.xlabel("Importance Score")
plt.ylabel("Features")
plt.show()
selected dt features = importance.sort values()[::-1][:6]
print(selected dt features.index)
X2 = downsampled df[selected dt features.index]
X train, X test, y train, y test = train test split(
```

```
X2, y, test size=0.2, stratify=y, random state=5805
dt model2 = DecisionTreeClassifier(random state=5805)
dt model2.fit(X train, y train)
y pred dt2 = dt model2.predict(X test)
evaluate_model(y_pred_dt2, y_test, "Decision Tree Classifier_2",
dt model2)
#%%
tuned parameters = [{"max depth": range(2, 10, 1),
                     "min_samples_leaf": range(2, 10, 1),
                     "min_samples_split": range(2, 10, 1),
                     'max_features': [None, 'sqrt', 'log2'],
                     'splitter': ['best', 'random'],
                     'criterion': ['gini', 'entropy']}]
dt model pruned = DecisionTreeClassifier(random state=5805,
ccp alpha=0.09)
grid search = GridSearchCV(estimator=dt model pruned,
param grid=tuned parameters, cv=5, n jobs=-1)
grid search.fit(X train, y train)
print(grid search.best params )
dt model pre pruned = grid search.best estimator
dt model pre pruned.fit(X train, y train)
y pre predicted = dt model pre pruned.predict(X test)
print(f'Accuracy of pre tuned tree {accuracy score(y test,
y_pre_predicted):.3f}')
plt.figure(figsize=(20, 12))
plot tree(dt model pre pruned, rounded=True, filled=True)
plt.show()
path = dt model.cost complexity pruning path(X train, y train)
print(path)
ccp alphas, impurities = path.ccp alphas, path.impurities
clfs = []
for ccp alpha in ccp alphas:
    clf = DecisionTreeClassifier(random state=5806,
ccp alpha=ccp alpha)
    clf.fit(X train, y train)
    clfs.append(clf)
acc scores = [accuracy score(y test, clf.predict(X test)) for clf in
clfsl
max acc score index = acc scores.index(max(acc scores))
plt.figure(figsize=(10, 6))
plt.grid()
plt.plot(ccp alphas[:-1], acc scores[:-1])
plt.xlabel("effective alpha")
```

```
plt.ylabel("Accuracy scores")
plt.show()
post pruning clf = clfs[max acc score index]
post pruning clf.fit(X train, y train)
y_post_predicted = post_pruning_clf.predict(X_test)
post pruning acc = accuracy score(y test, y post predicted)
print("accuracy for post pruning : ", round(post pruning acc, 3))
plot tree(decision tree=post pruning clf, rounded=True, filled=True)
pre confusion matrix = confusion matrix(y test, y pre predicted)
pre_accuracy = accuracy_score(y_test, y_pre_predicted)
pre_recall = recall_score(y_test, y_pre_predicted)
pre_roc_auc = roc_auc_score(y_test, y_pre_predicted)
print(" pre metrics ", pre_confusion_matrix, pre_recall,
pre_accuracy, pre_roc_auc)
post confusion matrix = confusion matrix(y test, y post predicted)
post accuracy = accuracy score(y test, y post predicted)
post recall = recall score(y test, y post predicted)
post roc auc = roc auc score(y test, y post predicted)
metrics data = {
    'Pruning Method': ['Pre-Pruned', 'Post-Pruned'],
    'Accuracy': [pre accuracy, post accuracy],
    'Confusion Matrix': [pre confusion matrix,
post confusion matrix],
    'Recall': [pre recall, post recall],
    'AUC': [pre roc auc, post_roc_auc]
metrics df = pd.DataFrame(metrics data)
print(metrics df.head())
logreg model = LogisticRegression(random state=5805)
logreg model.fit(X train, y train)
y logistic pred = logreg model.predict(X test)
accuracy = accuracy score(y test, y logistic pred)
print("Logistic Regression Accuracy:", round(accuracy, 3))
logistic_auc = roc_auc_score(y_test, y_logistic_pred)
dt pre auc = roc auc score(y test, y pre predicted)
dt_post_auc = roc_auc_score(y_test, y_post_predicted)
fpr_dt_post, tpr_dt_post, _ = roc_curve(y_test, y_post_predicted)
plt.figure(figsize=(8, 6))
plt.plot(fpr lr, tpr lr, label=f'Logistic Regression (AUC =
{logistic auc:.3f})'
```

```
plt.plot(fpr_dt_pre, tpr_dt_pre, label=f'Decision Tree with pre
pruning (AUC = {dt_pre_auc:.3f})')
plt.plot(fpr_dt_post, tpr_dt_post, label=f'Decision Tree with post
pruning (AUC = {dt post auc: .3f})')
plt.plot([0, 1], [0, 1], 'k--')
plt.xlabel('False Positive Rate')
plt.ylabel('True Positive Rate')
plt.title('ROC Curve')
plt.legend()
plt.grid(True)
plt.show()
logistic_confusion_matrix = confusion_matrix(y_test, y_logistic_pred)
logistic_accuracy = accuracy_score(y_test, y_logistic_pred)
logistic_recall = recall_score(y_test, y_logistic_pred)
logistic auc = roc auc score(y test, y logistic pred)
metrics data = {
    'Pruning Method': ['Pre-Pruned', 'Post-Pruned', 'Logistic'],
    'Accuracy': [pre accuracy, post accuracy, logistic accuracy],
    'Confusion Matrix': [pre confusion matrix, post confusion matrix,
logistic confusion matrix],
    'Recall': [pre recall, post recall, logistic recall],
    'AUC': [pre roc auc, post roc auc, logistic auc]
metrics df = pd.DataFrame(metrics data)
evaluate model(y logistic pred, y test, "Logistic Regression",
logreg model)
evaluate model(y pre predicted, y test, "Pre Pruned Decision Tree",
dt model pre pruned)
evaluate model(y post predicted, y test, "Post Pruned Decision Tree",
post pruning clf)
print(metrics df)
#%%
print("\n----\n")
error = []
for i in range(1, 40):
    knn = KNeighborsClassifier(n_neighbors=i)
    knn.fit(X_train, y_train)
    pred i = knn.predict(X test)
    error.append(np.mean(pred i != y test))
plt.figure(figsize=(12, 6))
plt.plot(range(1, 40), error, color='red', linestyle='dashed',
marker='o', markerfacecolor='blue', markersize=10)
plt.title('Error Rate for Different K Values')
plt.xlabel('K')
```

```
plt.ylabel('Error')
plt.show()
knn = KNeighborsClassifier(n neighbors=error.index(min(error)) + 1)
knn.fit(X train, y train)
pred = knn.predict(X test)
print('Minimum error:', min(error))
print('Optimum K:', error.index(min(error)) + 1)
evaluate_model(pred, y_test, 'KNN', knn)
#응응
print("\n-----
                    ----SVM-----
-\n")
parameters = {
   'kernel': ['linear', 'rbf', 'poly'],
stratified kfold = StratifiedKFold(n splits=5, shuffle=True,
random_state=5805)
# Perform grid search with cross-validation
svm classifier = SVC(probability=True)
grid search = GridSearchCV(svm classifier, parameters,
cv=stratified kfold, n jobs=-1)
grid search.fit(X train, y train)
print("Best Hyperparameters:", grid search.best params )
svm = grid search.best estimator
svm.fit(X train, y train)
y svm pred = svm.predict(X test)
evaluate model(y_svm_pred, y_test, 'SVM', svm)
#응응
nb model = GaussianNB()
nb_model.fit(X_train, y_train)
y nb pred = nb model.predict(X test)
evaluate_model(y_nb_pred, y_test, 'Naive Bayes', nb_model)
#%%
# Bagging
bagging clf = BaggingClassifier(DecisionTreeClassifier(),
n_estimators=100, random_state=5805)
bagging clf.fit(X train, y train)
y_bagging_pred = bagging_clf.predict(X_test)
evaluate_model(y_bagging_pred, y_test, 'Bagging', bagging_clf)
#%%
```

```
# Stacking
estimators = [('dt', DecisionTreeClassifier()), ('lr',
LogisticRegression())]
stacking_clf = StackingClassifier(estimators=estimators)
stacking_clf.fit(X_train, y_train)
y_stacking_pred = stacking_clf.predict(X_test)
evaluate_model(y_stacking_pred, y_test, 'Stacking', stacking clf)
#응응
# Boosting
boosting clf = AdaBoostClassifier(DecisionTreeClassifier(),
n_estimators=100, random_state=5805)
boosting_clf.fit(X_train, y_train)
y_boosting_pred = boosting_clf.predict(X_test)
evaluate_model(y_boosting_pred, y_test, 'Boosting', boosting clf)
#응응
nn model = MLPClassifier(hidden layer sizes=(100, 50),
activation='relu', solver='adam', random state=5805)
nn model.fit(X train, y train)
y nn pred = nn model.predict(X test)
evaluate model(y nn pred, y test, 'Neural Network', nn model)
#응응
sum of squared distances = []
for k in range(1, 11):
    kmeans = KMeans(n clusters=k, init='k-means++',
random state=5805)
    kmeans.fit(X)
    sum of squared distances.append(kmeans.inertia )
plt.plot(range(1, 11), sum of squared distances, 'bx-')
plt.xlabel('Number of Clusters (k)')
plt.ylabel('Sum of Squared Distances')
plt.title('Elbow Method for Optimal k Selection')
plt.show()
#응응
silhouette coefficients = []
for k in range(2, 11):
    kmeans = KMeans(n clusters=k, init='k-means++',
random state=5805)
    kmeans.fit(X)
    silhouette coefficients.append(silhouette score(X,
kmeans.labels ))
plt.plot(range(2, 11), silhouette coefficients, 'bx-')
plt.xlabel('Number of Clusters (k)')
plt.ylabel('Silhouette Coefficient')
```

```
plt.title('Silhouette Analysis for Optimal k Selection')
plt.show()
#%%
print("\n----\n")
data = car prices df.copy().sample(frac=0.2)
data clean = data.dropna(subset=['condition', 'odometer', 'mmr',
'sellingprice', 'year'])
features = data clean[['year', 'condition', 'odometer', 'mmr',
'sellingprice']]
features_scaled = StandardScaler().fit_transform(features)
dbscan = DBSCAN(eps=0.5, min samples=10)
clusters = dbscan.fit predict(features scaled)
print(clusters)
plt.figure(figsize=(10, 6))
plt.scatter(features scaled[:, 2], features scaled[:, 4], c=clusters,
cmap='viridis', s=5)
plt.xlabel('Odometer (scaled)')
plt.ylabel('Selling Price (scaled)')
plt.title('DBSCAN Clustering Results')
plt.colorbar(label='Cluster Label')
plt.show()
print("\n----\n")
print("\n-----------ASSOCIATION RULE MINING------
--\n")
df apriori filtered = car prices df[['make', 'model', 'trim',
'transmission']].sample(frac=0.05)
df apriori filtered = pd.get dummies(df apriori filtered,
drop first=True)
frequent itemsets = apriori(df apriori filtered, min support=0.1,
use colnames=True, verbose=1)
rules = association rules(frequent itemsets, metric="lift",
min threshold=0.2)
rules = rules.sort values(['confidence'], ascending=False)
print(rules.head(10))
print("\n-----------ASSOCIATION RULE MINING------
--\n")
```

To maintain the clarity and continuity of the report, the full codebase is provided in a separate file linked within this section.

9 References

• https://ieeexplore.ieee.org/abstract/document/9740817

- https://scikit-learn.org/stable/
- https://pandas.pydata.org/docs/
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