FUELWISE

Machine Learning Project



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TEAM MEMBERS

Team Leader

Deem Alrashidi

Group Member

Lama Alhujaili

Group Member

Sara Thaer

Group Member

Shahad Alsadah

Group Member

Sana Araj

Group Member

Shahad Adel

INTRODUCTION

Fuel consumption prediction utilizes advanced analytics and machine learning models to forecast the amount of fuel vehicles will use, enabling more efficient operations and reduced environmental impact.

Importance of Accurately Predicting Fuel Consumption:

Cost Efficiency

Operational Optimization

Reduced
Carbon
Footprints

THE PROBLEM

The problem of current fuel consumption prediction methods is:

- Lack of accuracy
- High variability in results
- Inefficiencies

And this can have a huge impact!

PROPOSED SOLUTION

We will use machine learning-based approach to predict fuel consumption more accurately.

Benefits of using machine learning model:

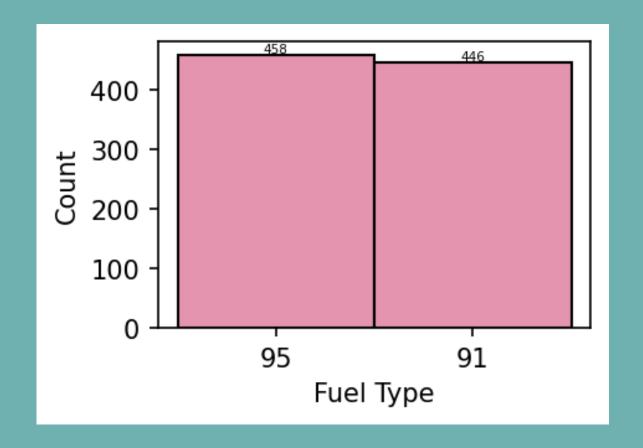
Adaptability to new data

Improvement in prediction accuracy

Automation of estimation processes.

DATA CLEANING

```
[ ] df = df.replace({'Fuel Type' : {'Z':'95', 'X': '91'}})
    df = df[~df['Fuel Type'].isin(['E','D'])]
    df
```



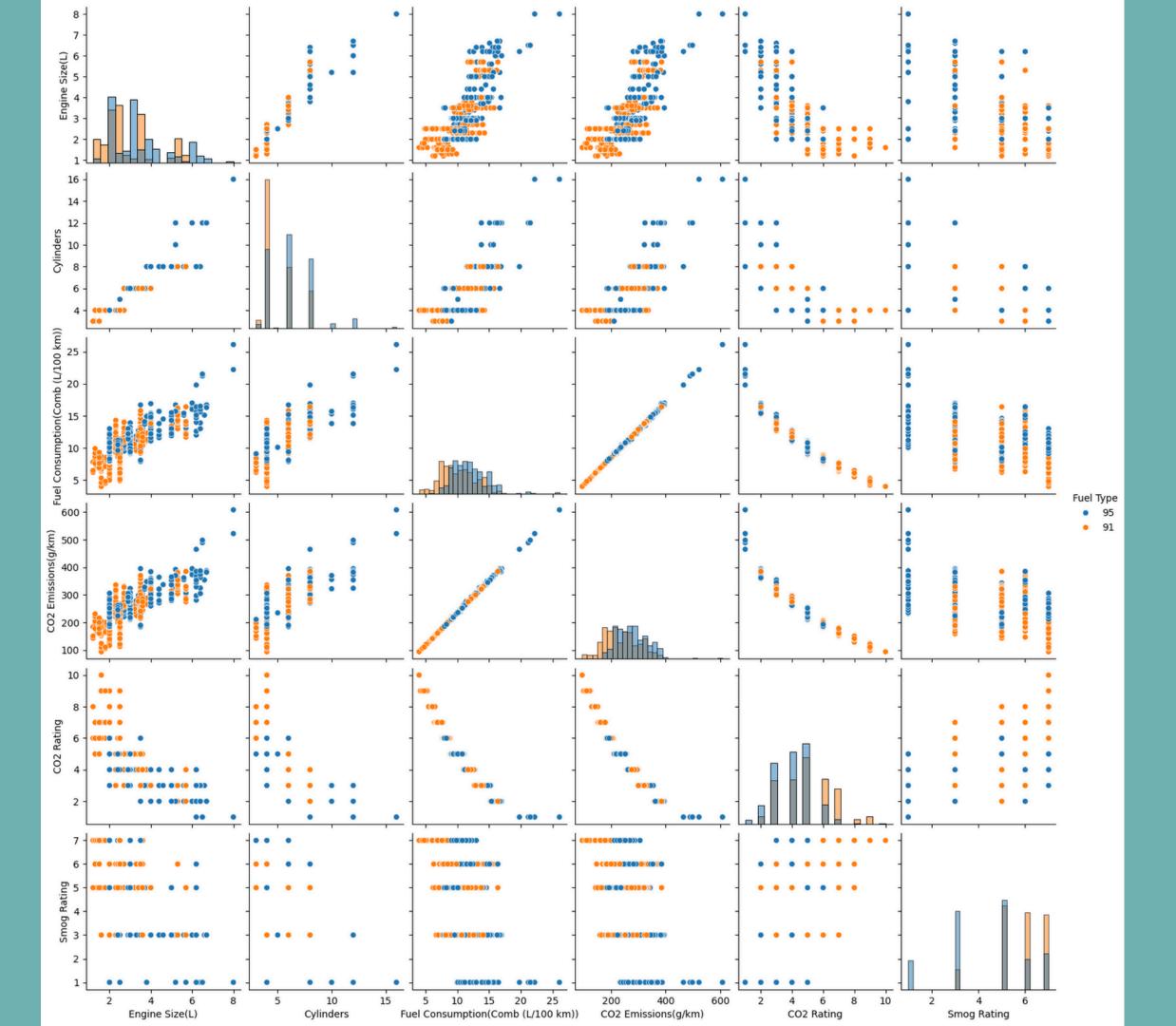
```
df.isna().sum()
Make
                                       0
Mode1
                                       0
Vehicle Class
                                      0
Engine Size(L)
                                       0
Cylinders
                                       0
Fuel Type
                                      0
Fuel Consumption(Comb (L/100 km))
                                      0
CO2 Emissions(g/km)
                                      0
CO2 Rating
                                      0
Smog Rating
                                       0
dtype: int64
```

Correlation= data[['Engine Size(Correlation	L)','CO2 Emissio	ons(g/km)', 'Fuel Cons	sumption(Comb (L/100 km))', 'Smog R	ating','Cylin	ders']].corr()
	Engine Size(L)	CO2 Emissions(g/km)	Fuel Consumption(Comb (L/100 km))	Smog Rating	Cylinders
Engine Size(L)	1.000000	0.824188	0.818694	-0.448239	0.920698
CO2 Emissions(g/km)	0.824188	1.000000	0.971671	-0.520437	0.833241
Fuel Consumption(Comb (L/100 km))	0.818694	0.971671	1.000000	-0.490473	0.821718
Smog Rating	-0.448239	-0.520437	-0.490473	1.000000	-0.502149
Cylinders	0.920698	0.833241	0.821718	-0.502149	1.000000



DATA ANALYSIS

Correlation



DATA ANALYSIS

Bivariate analysis

THE FEATURES

Before Preprocessing and Cleaning

~	 Investigate data (Missing values, Descriptions, Data Types) 															
-										+ Code	+ Code + Text					<u> </u>
C	data=pd.read_csv("dataset.csv") data.head()															
2		Model Year	Make	Model	Vehicle Class	Engine Size(L)	Cylinders	Transmission	Fuel Type	Fuel Consumption (City (L/100 km)	Fuel Consumption(Hwy (L/100 km))	Fuel Consumption(Comb (L/100 km))	Fuel Consumption(Comb (mpg))	CO2 Emissions(g/km)	CO2 Rating	Smog Rating
	0	2022	Acura	ILX	Compact	2.4	4	AM8	Z	9.9	7.0	8.6	33	200	6	3
	1	2022	Acura	MDX SH-AWD	SUV: Small	3.5	6	AS10	Z	12.6	9.4	11.2	25	263	4	5
	2	2022	Acura	RDX SH- AWD	SUV: Small	2.0	4	AS10	Z	11.0	8.6	9.9	29	232	5	6
П	3	2022	Acura	RDX SH- AWD A- SPEC	SUV: Small	2.0	4	AS10	Z	11.3	9.1	10.3	27	242	5	6
	4	2022	Acura	TLX SH- AWD	Compact	2.0	4	AS10	Z	11.2	8.0	9.8	29	230	5	7

THE FEATURES

After Preprocessing and Cleaning

0	x.head()											
∑	Engin	ne Size(L)	Cylinders	Fuel Consumption(Comb (L/100 km))	CO2 Emissions(g/km)	CO2 Rating	Smog Rating	Vehicle Class_X	Make_X	Model_Y	91	95
	1	3.5	6	11.2	263	4	5	6.0	25.0	576.0	0	1
	2	2.0	4	9.9	232	5	6	6.0	25.0	511.0	0	1
	3	2.0	4	10.3	242	5	6	6.0	25.0	512.0	0	1
	4	2.0	4	9.8	230	5	7	2.0	25.0	513.0	0	1
Γ.	5	2.0	4	9.8	231	5	7	2.0	25.0	514.0	0	1

SVM

- SVM is a robust algorithm for classification and regression tasks.
- Accuracy is 99.3%

SVM

```
from sklearn.svm import SVR

# Assuming you've defined your SVM regressor with appropriate parameters
svm_regressor = SVR(kernel='rbf')
# Fit the SVR model to your training data
svm_regressor.fit(xtrain, ytrain)
# Predict on the test set
y_pred = svm_regressor.predict(xtest)

# Evaluate the model
mse = mean_squared_error(ytest, y_pred)
print("Mean Squared Error:", mse)

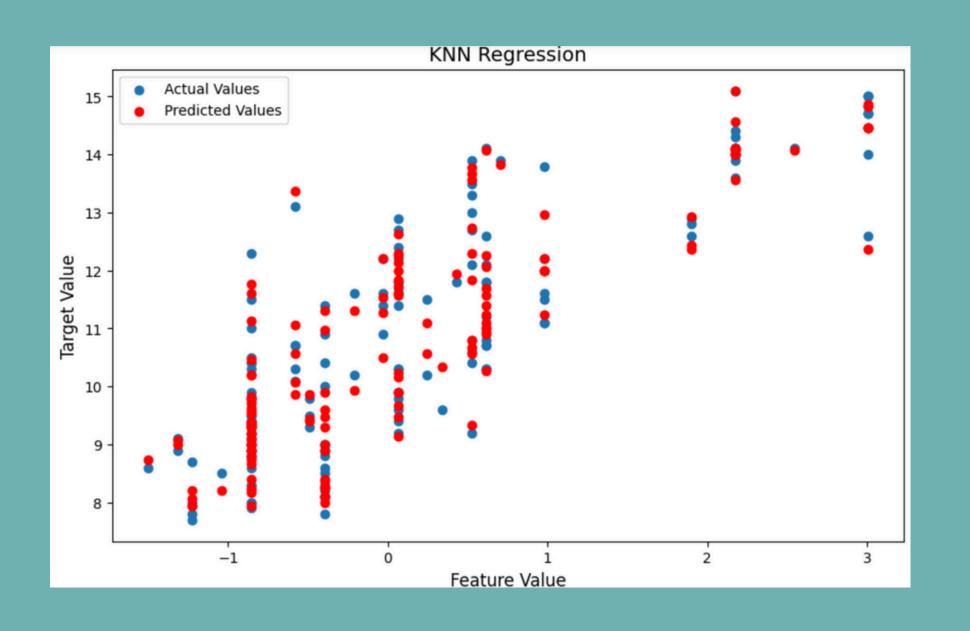
# Calculate accuracy (you may want to use a different metric for regression tasks)
accuracy = svm_regressor.score(xtest, ytest)
print("Accuracy:", accuracy*100)
```

Mean Squared Error: 0.028949171032108784
Accuracy: 99.28363455213638

KNN

- Straightforward and intuitive algorithm.
- Its performance depends on the choice of the hyperparameter k.
- The Accuracy is 97%

```
KNN
     from sklearn.neighbors import KNeighborsRegressor
     knn regressor = KNeighborsRegressor(n neighbors=3)
     knn_regressor.fit(xtrain, ytrain)
₹
             KNeighborsRegressor
      KNeighborsRegressor(n_neighbors=3)
     y_pred = knn_regressor.predict(xtest)
    mse = mean_squared_error(ytest, y_pred)
     print("Mean Squared Error:", mse)
     accuracy = knn_regressor.score(xtest, ytest)
     print("R^2 Score:", accuracy)
    Mean Squared Error: 0.12236781609195396
     R^2 Score: 0.9697193141449272
```

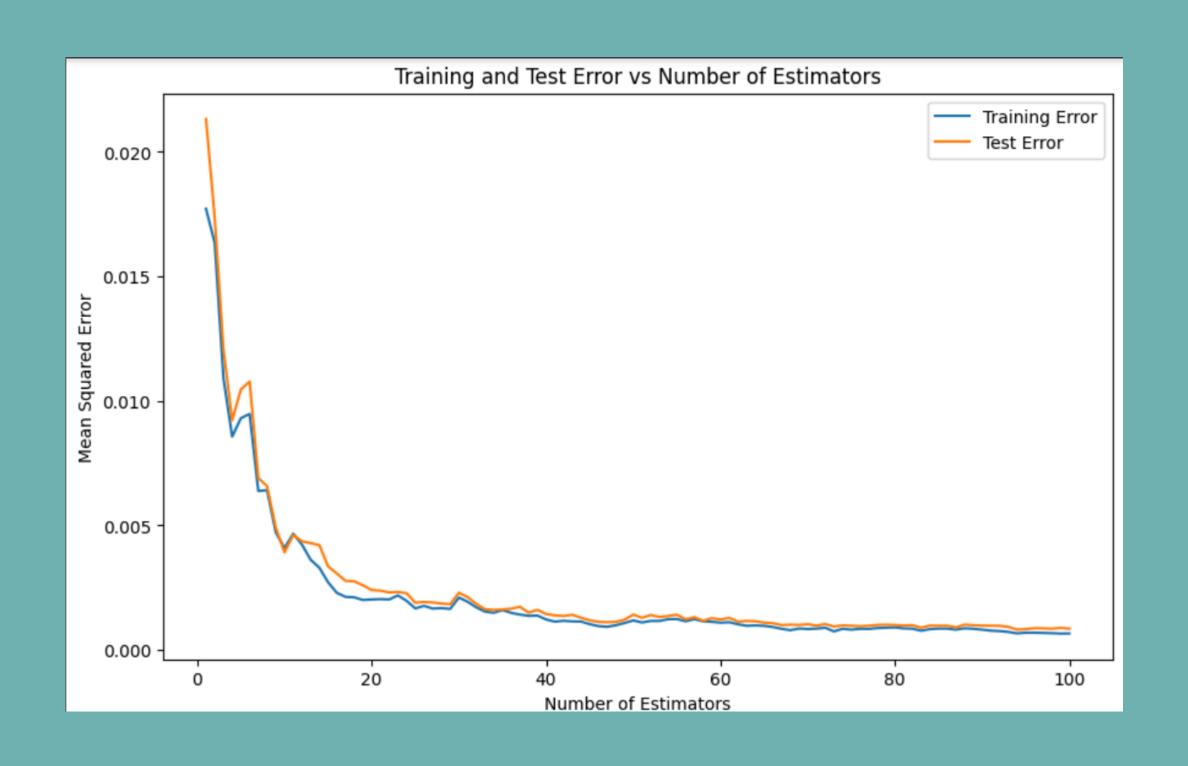


training score = 0.9996615319816969

AdaBoost Regressor

- Works on multiple weak models
- adjusts the weights
- deals with complex data and avoids overfitting
- The Accuracy is 99.9%

```
import numpy as np
     from sklearn.ensemble import AdaBoostRegressor
     from sklearn.tree import DecisionTreeRegressor
     from sklearn.metrics import mean_squared_error
     from sklearn.model_selection import train_test_split
     # Assuming xtrain, ytrain, xtest, ytest are already defined
     # Create a DecisionTreeRegressor as the base estimator
     base estimator = DecisionTreeRegressor(max depth=4)
     # Create the AdaBoost regressor
     ada_regressor = AdaBoostRegressor(base_estimator=base_estimator, n_estimators=50, learning_rate=1.0, random_state=42)
     # Fit the AdaBoost regressor to your training data
     ada regressor.fit(xtrain, ytrain)
     # Predict on the test set
     y pred = ada regressor.predict(xtest)
     # Evaluate the model
     mse = mean squared error(ytest, y pred)
     rmse = np.sqrt(mse)
     print("Root Mean Square Error:", rmse)
     # Calculate the R^2 score (coefficient of determination)
     accuracy = ada regressor.score(xtest, ytest)
     print("R^2 Score:", accuracy*100)
     print("training score = ",ada regressor.score(xtrain,ytrain))
→ Root Mean Square Error: 0.0377194013811472
     R^2 Score: 99.96479307606988
```



Linear Regression

- A statistical method used for prediction
- It provides simple and interpretable way to understand the relationship between the variables.
- The Accuracy is 100%

```
Linear Regression

#To check if there is overfitting or under fitting
print("training score = ",lr.score(xtrain,ytrain))
print("testing score = ",lr.score(xtest,ytest))

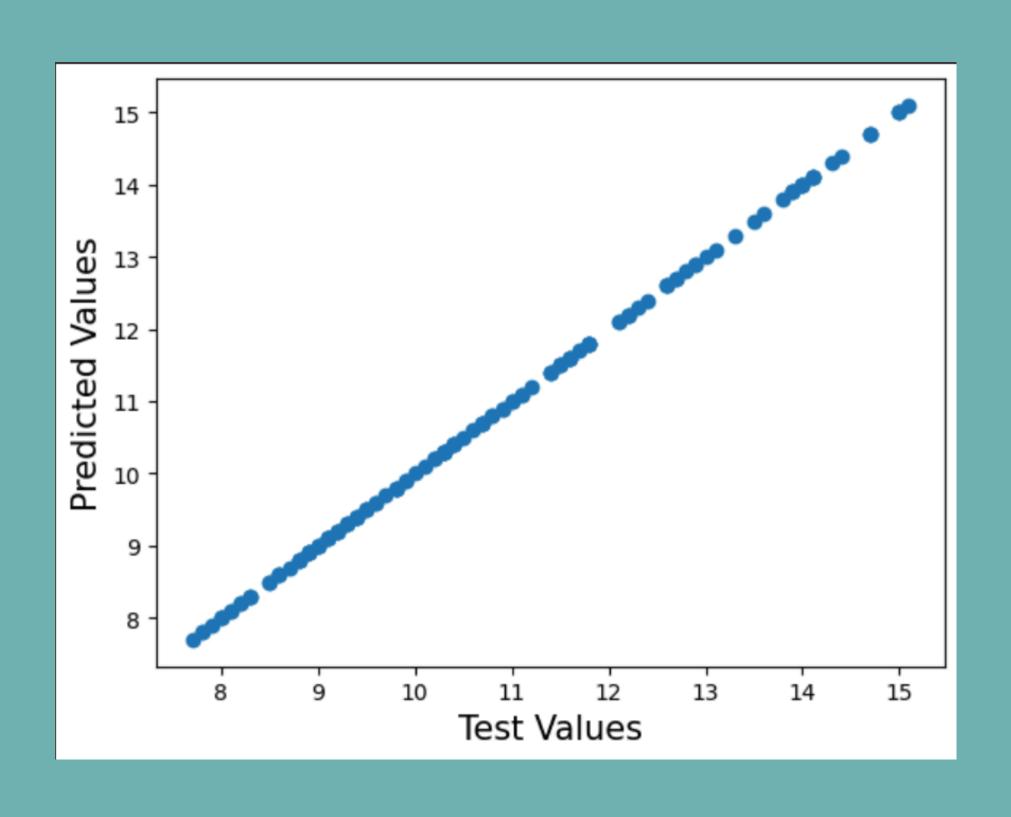
training score = 1.0

ypred = lr.predict(xtest)

r2_score(ytest, ypred)

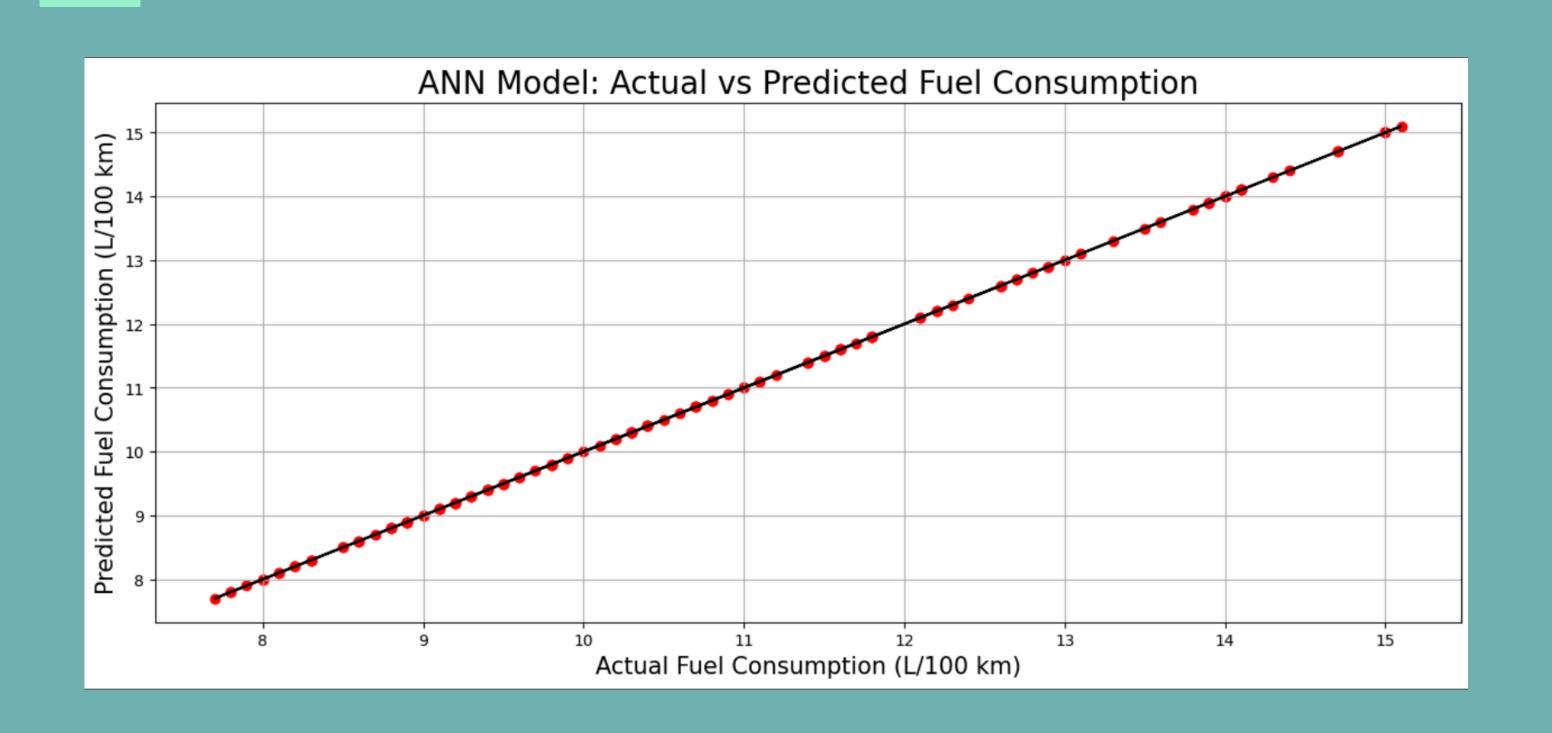
1.0

1.0
```



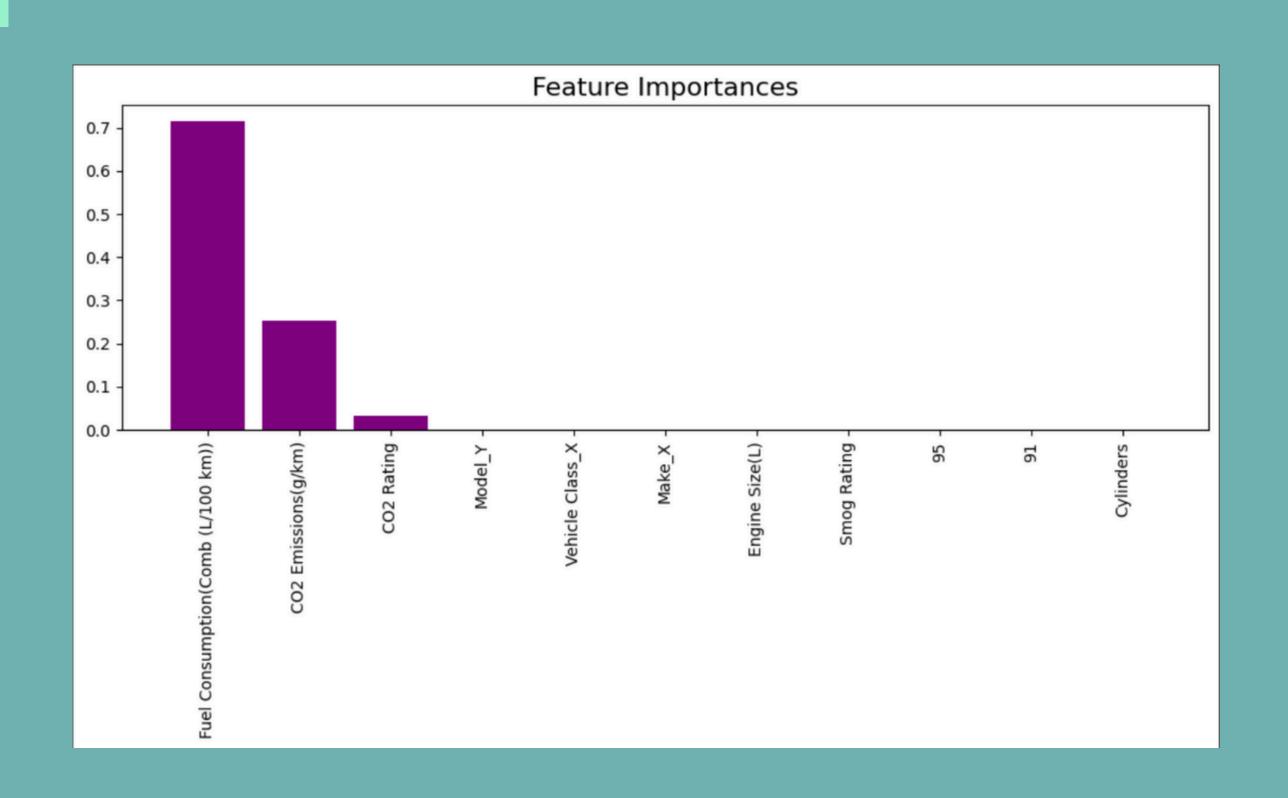
Artificial Neural Network

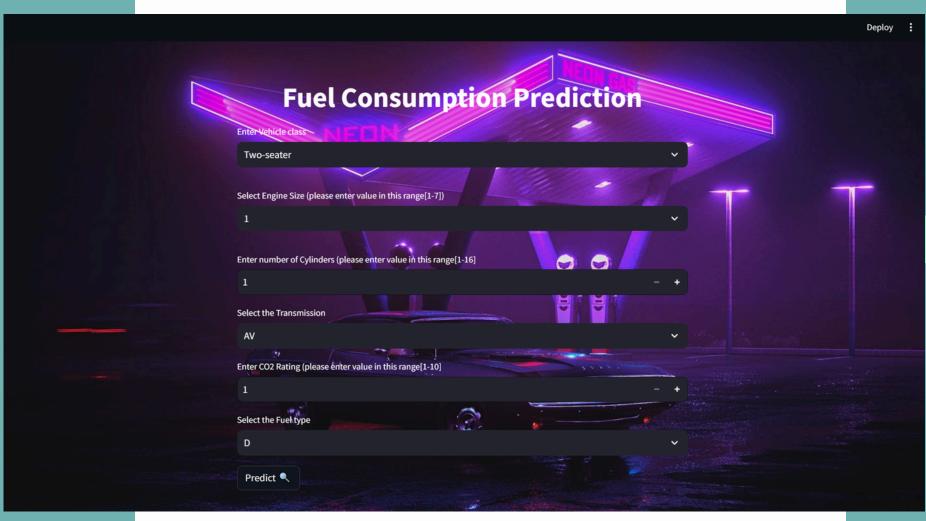
- particularly useful because of their ability to model complex relationships between input variables and the target output.
 The Accuracy is 99%
- Neural Network from sklearn.neural network import MLPRegressor nn = MLPRegressor(hidden layer sizes=(900,), activation='relu', solver='adam', max iter=1000) nn.fit(xtrain, ytrain) **₹** MLPRegressor MLPRegressor(hidden_layer_sizes=(900,), max_iter=1000) from sklearn.metrics import mean_squared error, r2_score, mean absolute error accuracy nn = nn.score(xtest, ytest) * 100 print("Accuracy:", accuracy nn) rmse nn = np.sqrt(mean squared error(ytest, ypred)) print("RMSE:", rmse nn) mae nn = mean absolute error(ytest, ypred) print("MAE:", mae nn) r2 nn = r2 score(ytest, ypred) print("R2:", r2 nn) → Accuracy: 99.39959865306272 RMSE: 1.8513413036726634e-15 MAE: 1.5435928397547003e-15 R2: 1.0



Random Forest

- The random forest method is an ensemble learning technique used primarily for classification and regression tasks.
 The Accuracy is 99%
- Random Forest from sklearn.ensemble import RandomForestRegressor rf = RandomForestRegressor() no of decision tree = [10,20,30,40,50,60,70,80,90,100] max no of features = ['sqrt','log2'] $\max_{depth} = [6,7,8,9,10,11,12,13,14,15]$ criterion of decision tree = ["squared error", "poisson"] min sample split=[2,3,4,5,6] [] random_grid = { 'n_estimators': no_of_decision_tree, 'max features': max no of features, 'max depth': max_depth, 'criterion': criterion_of_decision_tree, 'min samples split': min sample split [] from sklearn.model selection import RandomizedSearchCV rscv = RandomizedSearchCV(estimator = rf , param_distributions = random_grid , n_iter = 25 , cv = 5 ,n_jobs=-1) rscv.fit(xtrain, ytrain) RandomizedSearchCV ▶ estimator: RandomForestRegressor ▶ RandomForestRegressor





PROJECT OUTCOME

NEXT STEPS

DATA PROCESSING

Include more sophisticated data processing techniques.

LARGER DATESETS

Implement oour models on larger datasets

OUR WEBSITE

Improve our website to include an Al chatbot

EXPERIMENT

Experiment with various other factors and models.

CONCLUSION

STRIVE FOR SUSTAINABALITY

