

Dataset

- https://huggingface.co/dat asets/scikit-learn/auto-mpg
- Auto Miles per Gallon (MPG) Dataset
- Contains data of fuel consumption of different cars from the 70's and 80's

Importing

```
displacement horsepower
                                                        acceleration model year
         cylinders
                                                weight
  18.0
                                                                 12.0
                            307.0
                                          130
                                                  3504
                                                                                70
  15.0
                            350.0
                                                  3693
                                                                 11.5
                                          165
                                                                                70
  18.0
                            318.0
                                          150
                                                  3436
                                                                 11.0
                                                                                70
   16.0
                                                                 12.0
                            304.0
                                          150
                                                  3433
                                                                                70
4 17.0
                            302.0
                                          140
                                                  3449
                                                                 10.5
```

car nai	origin	
chevrolet chevelle mali	1	0
buick skylark 3	1	1
plymouth satelli	1	2
amc rebel s	1	3
ford tori	1	4

```
# prompt: import pandas and this dataset from huggingface: https://huggingface.co/datasets/scikit-learn/auto-mpg
!pip install datasets
from datasets import load_dataset
import pandas as pd
```

```
dataset = load_dataset("scikit-learn/auto-mpg")
df = pd.DataFrame(dataset['train'])
print(df.head())
```

Data Cleaning

Getting rid of unnecesary columns, like:

- Car name
- Origin

We will work with:

- MPG
- Cylinders
- Displacement
- Horsepower
- Weight
- Acceleration
- Model Year

```
# prompt: delete car name and origin column

del df['car name']
 del df['origin']
 print(df.head())
 df.info()
```

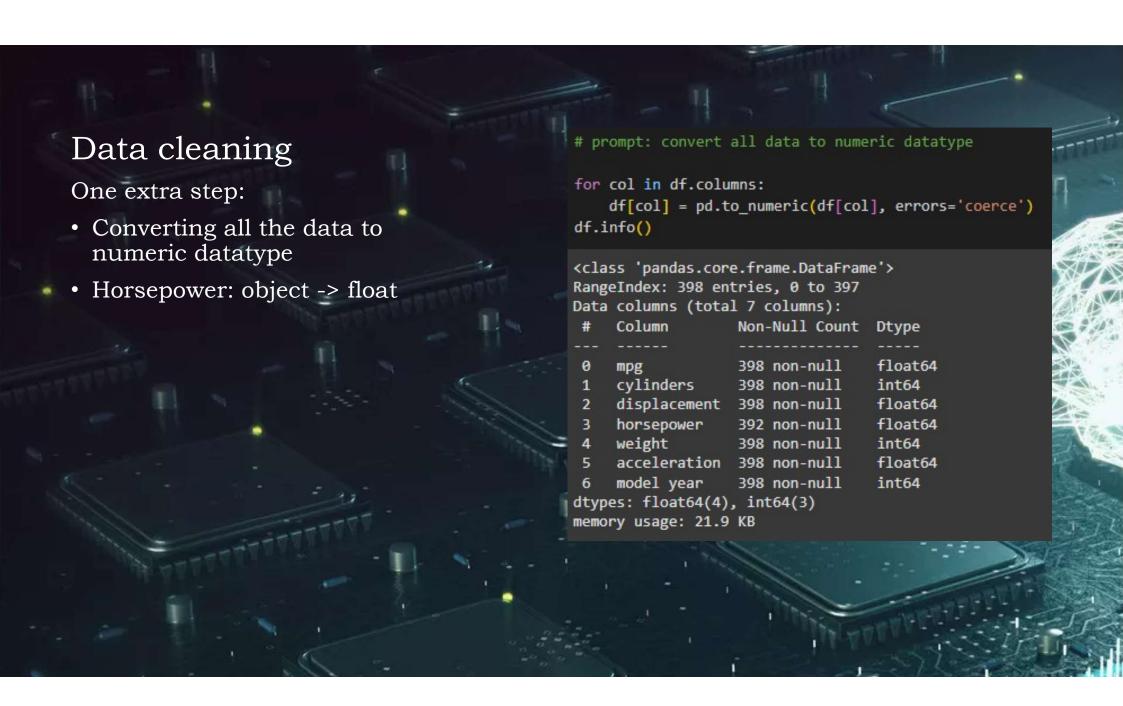
	mpg	cylinders	displacement	horsepower	weight	acceleration	model year
0	18.0	8	307.0	130	3504	12.0	70
1	15.0	8	350.0	165	3693	11.5	70
2	18.0	8	318.0	150	3436	11.0	70
3	16.0	8	304.0	150	3433	12.0	70
4	17.0	8	302.0	140	3449	10.5	70

<class 'pandas.core.frame.DataFrame'>
RangeIndex: 398 entries, 0 to 397

Data columns (total 7 columns):

#	Column	Non-Null Count	Dtype			
0	mpg	398 non-null	float64			
1	cylinders	398 non-null	int64			
2	displacement	398 non-null	float64			
3	horsepower	398 non-null	object			
4	weight	398 non-null	int64			
5	acceleration	398 non-null	float64			
6	model year	398 non-null	int64			
dtypes: float64(3), int64(3), object(1)						

memory usage: 21.9+ KB



Getting familiar with the data

at.d	escri	be (

		mpg	cylinders	displacement	horsepower	weight	acceleration	model year
	count	398.000000	398.000000	398.000000	392.000000	398.000000	398.000000	398.000000
	mean	23.514573	5.454774	193.425879	104.469388	2970.424623	15.568090	76.010050
	std	7.815984	1.701004	104.269838	38.491160	846.841774	2.757689	3.697627
	min	9.000000	3.000000	68.000000	46.000000	1613.000000	8.000000	70.000000
	25%	17.500000	4.000000	104.250000	75.000000	2223.750000	13.825000	73.000000
	50%	23.000000	4.000000	148.500000	93.500000	2803.500000	15.500000	76.000000
	75%	29.000000	8.000000	262.000000	126.000000	3608.000000	17.175000	79.000000
	max	46.600000	8.000000	455.000000	230.000000	5140.000000	24.800000	82.000000

Correlation matrix

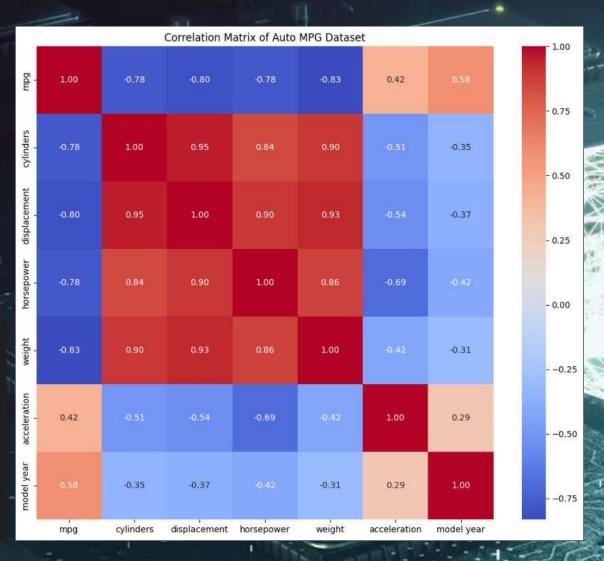
Shows the correlation between the variables

- 1 -> strong correlation
- -1-> strong inverse correlation
 - 0 -> no correlation

```
# prompt: create a correlation matrix of the dataset
import matplotlib.pyplot as plt
import seaborn as sns

# Calculate the correlation matrix
correlation_matrix = df.corr()

# Create a heatmap
plt.figure(figsize=(12, 10))
sns.heatmap(correlation_matrix, annot=True, cmap='coolwarm', fmt=".2f")
plt.title('Correlation Matrix of Auto MPG Dataset')
plt.show()
```



Splitting the data into training and testing datasets

```
# prompt: separate the dataset into training and test data
from sklearn.model selection import train test split
# Assuming 'df' is your DataFrame
X = df.drop('mpg', axis=1) # Features (all columns except 'mpg')
y = df['mpg'] # Target variable ('mpg')
# Split the data into training and testing sets
X train, X test, y train, y test = train test split(X, y, test size=0.2, random state=42) # 80% training and 20% test
print("X_train shape:", X train.shape)
print("y train shape:", y train.shape)
print("X_test shape:", X test.shape)
print("y test shape:", y test.shape)
X train shape: (318, 6)
y train shape: (318,)
X test shape: (80, 6)
y test shape: (80,)
```

```
Another emerging problem: missing data
 # prompt: exclude missing data from the dataset then create a linear regression model
 from sklearn.linear model import LinearRegression
 from sklearn.metrics import mean squared error, r2 score
 # Drop rows with missing values
 df.dropna(inplace=True)
 # Assuming 'df' is your DataFrame
 X = df.drop('mpg', axis=1) # Features (all columns except 'mpg')
 y = df['mpg'] # Target variable ('mpg')
 # Split the data into training and testing sets
 X train, X test, y train, y test = train test split(X, y, test size=0.2, random state=42) # 80% training and 20% test
 # Create and train the linear regression model
 model = LinearRegression()
 model.fit(X train, y train)
 # Make predictions on the test set
 y pred = model.predict(X test)
 # Evaluate the model
 mse = mean squared error(y test, y pred)
 r2 = r2_score(y_test, y_pred)
 print(f"Mean Squared Error: {mse}")
 print(f"R-squared: {r2}")
Mean Squared Error: 10.502370329417325
```

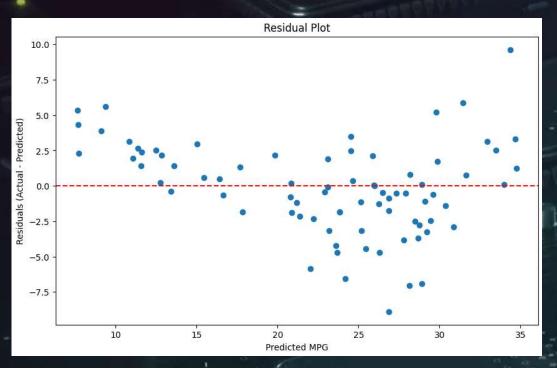
R-squared: 0.7942349075428586

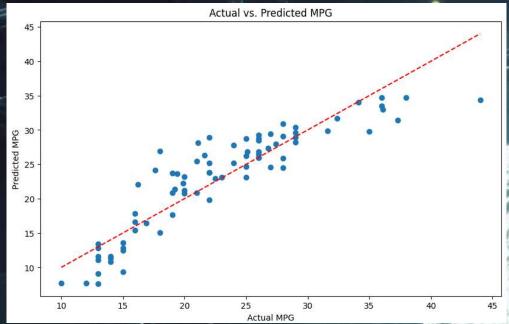
Showing the result of our Model

```
# prompt: create a simple example where you show the difference between the predicted and the real values
# Make predictions on the test set
y pred = model.predict(X test)
# Create a DataFrame to compare actual and predicted values
comparison df = pd.DataFrame({'Actual': y test, 'Predicted': y pred})
# Display the first few rows of the comparison DataFrame
print(comparison df.head(20))
# Calculate the difference between actual and predicted values
comparison df['Difference'] = comparison df['Actual'] - comparison df['Predicted']
# Display the updated DataFrame with the difference column
print(comparison df.head(20))
#Plot the difference
plt.figure(figsize=(10, 6))
plt.scatter(y test, y pred)
plt.plot([min(y test), max(y test)], [min(y test), max(y test)], color='red', linestyle='--') # Add a diagonal line
plt.xlabel('Actual MPG')
plt.ylabel('Predicted MPG')
plt.title('Actual vs. Predicted MPG')
plt.show()
# Plot the residuals (differences between actual and predicted)
plt.figure(figsize=(10, 6))
plt.scatter(y pred, comparison df['Difference'])
plt.axhline(y=0, color='red', linestyle='--') # Add a horizontal line at y=0
plt.xlabel('Predicted MPG')
plt.vlabel('Residuals (Actual - Predicted)')
plt.title('Residual Plot')
plt.show()
```

```
Predicted
                  Difference
Actual
 26.0
       25.932796
                     0.067204
 21.6 26.299279
                    -4.699279
 36.1 32.965489
                    3.134511
                    -0.854430
 26.0 26.854430
 27.0 29.453726
                    -2.453726
 28.0 29.086995
                    -1.086995
 13.0 7.638748
                    5.361252
 26.0 29.237044
                    -3.237044
                    -1.877360
 19.0 20.877360
      28.915829
                    0.084171
      12.485037
                    2.514963
  15.0
       23.698983
                    -4.698983
  19.0
  16.9
       16.418892
                    0.481108
  29.0
      28.196688
                    0.803312
 16.2 22.039679
                    -5.839679
 28.0 30.881524
                    -2.881524
 20.0 21.183111
                    -1.183111
 32.4 31.646623
                    0.753377
 27.4 27.916498
                    -0.516498
 35.0 29.781482
                     5.218518
```

Showing the result of our Model





Could be better – let's try to make it better!

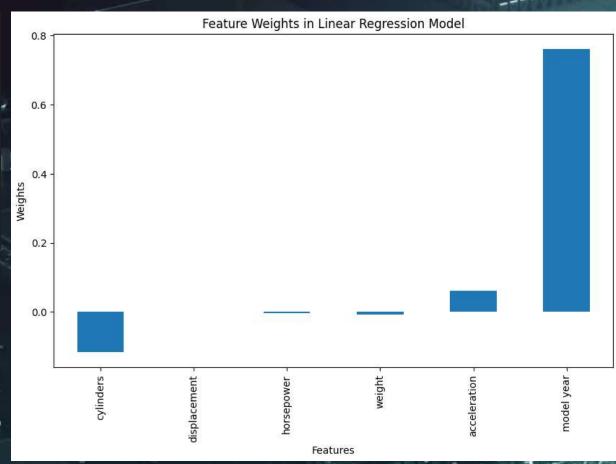
Showing the weights in the Linear Regression Model

```
# prompt: visualize the weights of the linear regression model
import matplotlib.pyplot as plt

# Get feature weights (coefficients) from the trained model
feature_weights = pd.Series(model.coef_, index=X.columns)

# Plot the feature weights
plt.figure(figsize=(10, 6))
feature_weights.plot(kind='bar')
plt.xlabel('Features')
plt.ylabel('Weights')
plt.title('Feature Weights in Linear Regression Model')
plt.show()
```

Model year with the biggest weight?
Interesting...



Polynomial Regression

Polynomial regression is a type of regression analysis used to model the relationship between a dependent variable and one or more independent variables when the relationship is nonlinear. It extends linear regression by fitting a polynomial equation to the data, which allows for modeling more complex patterns.

```
# prompt: try to make the prediction better
# ... (Your existing code)
from sklearn.preprocessing import PolynomialFeatures
from sklearn.pipeline import make pipeline
# Try polynomial regression
poly model = make pipeline(PolynomialFeatures(degree=2), LinearRegression()) #degree 2
poly model.fit(X train, y train)
y poly pred = poly model.predict(X test)
mse poly = mean squared error(y test, y poly pred)
r2 poly = r2 score(y test, y poly pred)
print(f"Polynomial Regression - Mean Squared Error: {mse poly}")
print(f"Polynomial Regression - R-squared: {r2 poly}")
# Compare with original Linear Regression
print(f"Linear Regression - Mean Squared Error: {mse}")
print(f"Linear Regression - R-squared: {r2}")
# ... (rest of your existing code, plotting etc.)
Polynomial Regression - Mean Squared Error: 7.0705240702947885
Polynomial Regression - R-squared: 0.8614725063570133
Linear Regression - Mean Squared Error: 10.502370329417325
Linear Regression - R-squared: 0.7942349075428586
```

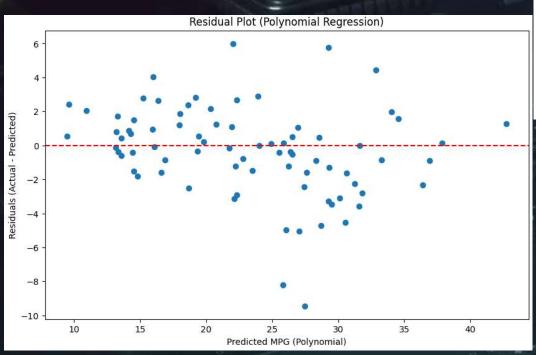
Showing the result of the Polynomial Model

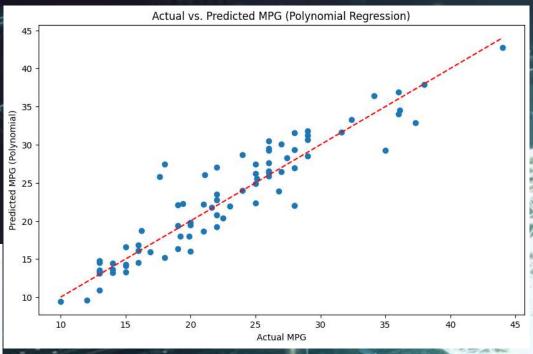
prompt: create a simple example where you show the difference between the predicted and the real values for the polynomial regression

```
# Create a DataFrame to compare actual and predicted values for polynomial regression
poly comparison df = pd.DataFrame({'Actual': y test, 'Predicted': y poly pred})
# Display the first few rows of the comparison DataFrame
print(poly comparison df.head(20))
# Calculate the difference between actual and predicted values
poly comparison df['Difference'] = poly comparison df['Actual'] - poly comparison df['Predicted']
# Display the updated DataFrame with the difference column
print(poly comparison df.head(20))
# Plot the difference for polynomial regression
plt.figure(figsize=(10, 6))
plt.scatter(y test, y poly pred)
plt.plot([min(y test), max(y test)], [min(y test), max(y test)], color='red', linestyle='--')
plt.xlabel('Actual MPG')
plt.ylabel('Predicted MPG (Polynomial)')
plt.title('Actual vs. Predicted MPG (Polynomial Regression)')
plt.show()
# Plot the residuals (differences between actual and predicted) for polynomial regression
plt.figure(figsize=(10, 6))
plt.scatter(y poly pred, poly comparison df['Difference'])
plt.axhline(y=0, color='red', linestyle='--')
plt.xlabel('Predicted MPG (Polynomial)')
plt.ylabel('Residuals (Actual - Predicted)')
plt.title('Residual Plot (Polynomial Regression)')
plt.show()
```

```
Actual Predicted
                   Difference
                    -0.520838
  26.0 26.520838
  21.6 21.751928
                    -0.151928
  36.1 34.544822
                    1.555178
  26.0 26.362115
                    -0.362115
  27.0 30.115101
                    -3.115101
                    -1.305105
       29.305105
  13.0 13.129743
                    -0.129743
                    -3.475287
  26.0 29.475287
                    -0.349260
  19.0 19.349260
  29.0 31.819244
                    -2.819244
                    -1.586001
  15.0 16.586001
                    -3.134139
  19.0 22.134139
  16.9 15.941869
                     0.958131
  29.0 28.530773
                     0.469227
  16.2 18.695331
                    -2.495331
  28.0 31.570395
                    -3.570395
                     0.208023
  20.0 19.791977
  32.4 33.265742
                    -0.865742
  27.4 28.307020
                    -0.907020
  35.0 29.240084
                     5.759916
```

Showing the result of the Polynomial Model

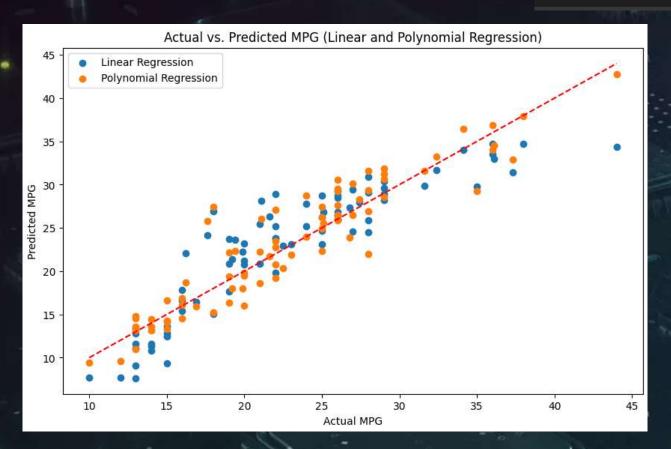




Linear VS Polynomial

```
plt.figure(figsize=(10, 6))
plt.scatter(y_test, y_pred, label='Linear Regression')
plt.scatter(y_test, y_poly_pred, label='Polynomial Regression')
plt.plot([min(y_test), max(y_test)], [min(y_test), max(y_test)], color='red', linestyle='--')
plt.xlabel('Actual MPG')
plt.ylabel('Predicted MPG')
plt.title('Actual vs. Predicted MPG (Linear and Polynomial Regression)')
plt.legend()
plt.show()
```

show the linear regression plot and the polynomial regression plot on one plot



Still could be better?

Maybe – Let's try!

Random Forest Regression

Random Forest Regression is an ensemble learning method that uses a collection of decision trees to perform regression tasks. It is an extension of the Random Forest algorithm, which is primarily known for classification but can also handle regression tasks effectively.

```
# prompt: make the prediction better
```

```
# ... (Your existing code)
```

```
from sklearn.preprocessing import StandardScaler
from sklearn.model_selection import GridSearchCV
from sklearn.ensemble import RandomForestRegressor
```

```
# Feature scaling
scaler = StandardScaler()
X_train_scaled = scaler.fit_transform(X_train)
X_test_scaled = scaler.transform(X_test)
```

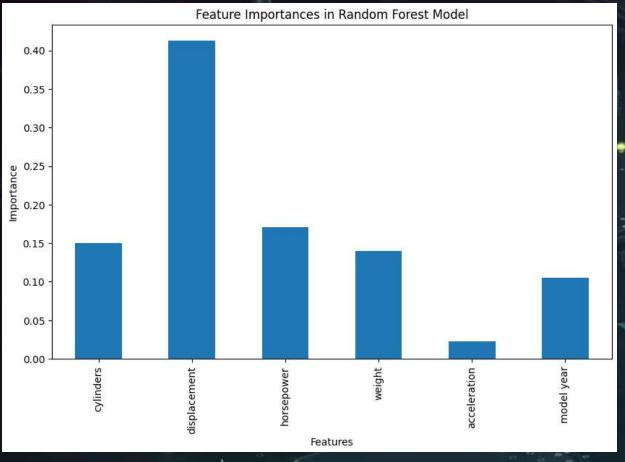
Try Random Forest Regressor with hyperparameter tuning
rf_model = RandomForestRegressor(random_state=42)

We got our best score:

Random Forest Regression - Mean Squared Error: 6.037425564634708

Random Forest Regression - R-squared: 0.8817132332469282

Random Forest weights



```
# prompt: show the weights used in the random forest regression
import pandas as pd

# Get feature importances from the best Random Forest model
feature_importances = pd.Series(best_rf_model.feature_importances_, index=X.columns).

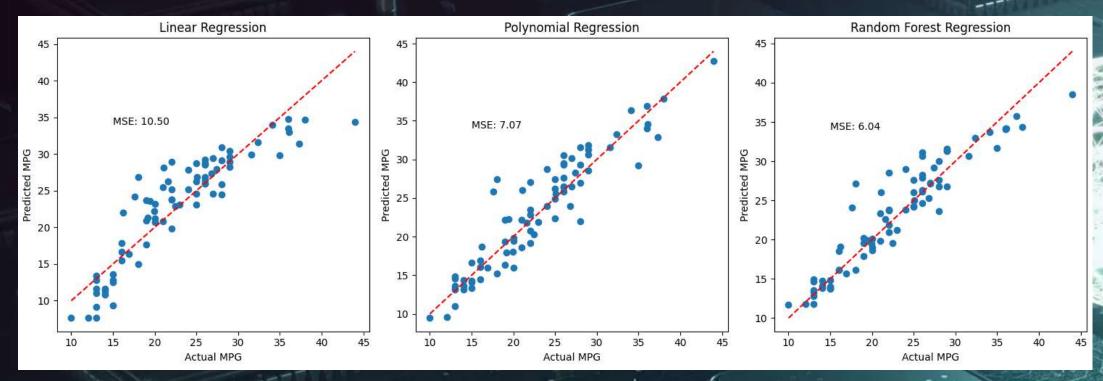
# Plot the feature importances
plt.figure(figsize=(10, 6))
feature importances.plot(kind='bar')
plt.xlabel('Features')
plt.ylabel('Importance')
plt.title('Feature Importances in Random Forest Model')
plt.show()
```

Now this one makes a lot more sense than the weights in the linear regression model!

Plotting all of our results

```
# prompt: plot all three regressions next to each other and their mse
plt.figure(figsize=(15, 5))
plt.subplot(1, 3, 1)
plt.scatter(y test, y pred)
plt.plot([min(y test), max(y test)], [min(y test), max(y test)], color='red', linestyle='--')
plt.title('Linear Regression')
plt.xlabel('Actual MPG')
plt.ylabel('Predicted MPG')
plt.text(min(y_test) + 5, max(y_test) -10, f"MSE: {mse:.2f}")
plt.subplot(1, 3, 2)
plt.scatter(y test, y poly pred)
plt.plot([min(y test), max(y test)], [min(y test), max(y test)], color='red', linestyle='--')
plt.title('Polynomial Regression')
plt.xlabel('Actual MPG')
plt.ylabel('Predicted MPG')
plt.text(min(y_test) + 5, max(y_test) - 10, f"MSE: {mse_poly:.2f}")
plt.subplot(1, 3, 3)
plt.scatter(y test, y rf pred)
plt.plot([min(y_test), max(y_test)], [min(y_test), max(y_test)], color='red', linestyle='--')
plt.title('Random Forest Regression')
plt.xlabel('Actual MPG')
plt.ylabel('Predicted MPG')
plt.text(min(y test) + 5, max(y test) - 10, f"MSE: {mse rf:.2f}")
plt.tight layout()
plt.show()
```

Plotting all of our results



Conclusion

prompt: give me a conclusion about this project

This project investigated the prediction of automobile fuel efficiency (MPG) using various regression models. The analysis began with data preprocessing, including handling missing values and removing irrelevant columns from the "auto-mpg" dataset. Exploratory data analysis, visualized through a correlation heatmap, provided insights into feature relationships.

Three regression models were trained and evaluated: Linear Regression, Polynomial Regression (degree 2), and Random Forest Regression. Model performance was assessed using Mean Squared Error (MSE) and R-squared. Visualizations, including actual vs. predicted MPG plots and residual plots, helped analyze the models accuracy and identify potential areas of improvement. Furthermore, feature importance was analyzed for the Random Forest model. A comparative plot was created to easily compare the performances of the three models. The results show that the Random Forest Regression model, after hyperparameter tuning, achieved the best performance based on the lowest MSE.

