## **Predicting Fuel Efficiency of Vehicles - Part 3**

### **Selecting and Training Models**

- 1. Select and Train a few Algorithms(Linear Regression, Decision Tree, RandomForest)
- 2. Evaluation using Mean Squared Error
- 3. Model Evaluation using Cross Validation
- 4. Hyperparameter Tuning using GridSearchCV
- 5. Check Feature Importance
- 6. Evaluate the Final System on test data
- 7. Saving the Model

```
In [1]: ##importing a few general use case libraries
import numpy as np
import pandas as pd
import matplotlib.pyplot as plt
import seaborn as sns

from sklearn.model_selection import StratifiedShuffleSplit
from sklearn.base import BaseEstimator, TransformerMixin
from sklearn.pipeline import Pipeline

from sklearn.impute import SimpleImputer
from sklearn.preprocessing import StandardScaler, OneHotEncoder
from sklearn.compose import ColumnTransformer

import warnings
warnings.filterwarnings('ignore')
```

# In [2]: pip install wget

Requirement already satisfied: wget in c:\users\lenovo\anaconda3\lib\site-packages (3.2) Note: you may need to restart the kernel to use updated packages.

Saved under auto-mpg (1).data

```
In [4]: ##segregate the feature and target variable
data = strat_train_set.drop("MPG", axis=1)
data_labels = strat_train_set["MPG"].copy()
data
```

#### Out[4]:

	Cylinders	Displacement	Horsepower	Weight	Acceleration	Model Year	Origin
145	4	83.0	61.0	2003.0	19.0	74	3
151	4	79.0	67.0	2000.0	16.0	74	2
388	4	156.0	92.0	2585.0	14.5	82	1
48	6	250.0	88.0	3139.0	14.5	71	1
114	4	98.0	90.0	2265.0	15.5	73	2
		•••	•••		•••		
147	4	90.0	75.0	2108.0	15.5	74	2
156	8	400.0	170.0	4668.0	11.5	75	1
395	4	135.0	84.0	2295.0	11.6	82	1
14	4	113.0	95.0	2372.0	15.0	70	3
362	6	146.0	120.0	2930.0	13.8	81	3

318 rows × 7 columns

```
In [5]: ##preprocess the Origin column in data
def preprocess_origin_cols(df):
    df["Origin"] = df["Origin"].map({1: "India", 2: "USA", 3: "Germany"})
    return df
```

```
In [6]: ##creating custom attribute adder class
acc_ix, hpower_ix, cyl_ix = 4,2, 0

class CustomAttrAdder(BaseEstimator, TransformerMixin):
    def __init__(self, acc_on_power=True): # no *args or **kargs
        self.acc_on_power = acc_on_power

def fit(self, X, y=None):
    return self # nothing else to do

def transform(self, X):
    acc_on_cyl = X[:, acc_ix] / X[:, cyl_ix]
    if self.acc_on_power:
        acc_on_power = X[:, acc_ix] / X[:, hpower_ix]
        return np.c_[X, acc_on_power, acc_on_cyl]

return np.c_[X, acc_on_cyl]
```

```
In [7]: def num_pipeline_transformer(data):
            Function to process numerical transformations
            Argument:
                data: original dataframe
            Returns:
                num attrs: numerical dataframe
                num pipeline: numerical pipeline object
            1.1.1
            numerics = ['float64', 'int64']
            num attrs = data.select dtypes(include=numerics)
            num pipeline = Pipeline([
                ('imputer', SimpleImputer(strategy="median")),
                ('attrs_adder', CustomAttrAdder()),
                ('std scaler', StandardScaler()),
                1)
            return num_attrs, num_pipeline
        def pipeline_transformer(data):
            Complete transformation pipeline for both
            nuerical and categorical data.
            Argument:
                data: original dataframe
            Returns:
                prepared_data: transformed data, ready to use
            cat attrs = ["Origin"]
            num attrs, num pipeline = num pipeline transformer(data)
            full pipeline = ColumnTransformer([
                ("num", num pipeline, list(num attrs)),
                ("cat", OneHotEncoder(), cat attrs),
            prepared_data = full_pipeline.fit_transform(data)
            return prepared data
```

#### From raw data to processed data in 2 steps

```
In [8]: ##from raw data to processed data in 2 steps
       preprocessed df = preprocess origin cols(data)
       prepared data = pipeline transformer(preprocessed df)
       prepared data
Out[8]: array([[-0.85657842, -1.07804475, -1.15192977, ..., 1.
               0. , 0.
             [-0.85657842, -1.1174582, -0.9900351, ..., 0.
               0. , 1. ],
             [-0.85657842, -0.3587492, -0.31547399, ..., 0.
                  , 0. ],
             [-0.85657842, -0.56566984, -0.53133355, \ldots, 0.
              1. , 0.
             [-0.85657842, -0.78244384, -0.23452666, ..., 1.
               0. , 0. ],
             [ 0.32260746, -0.45728283, 0.44003446, ..., 1.
               0. , 0. ]])
In [9]: prepared_data[0]
Out[9]: array([-0.85657842, -1.07804475, -1.15192977, -1.17220298, 1.21586943,
             -0.54436373, 1.70952741, 1.29565517, 1. , 0.
                      ])
              0.
```

#### **Selecting and Training Models**

- 1. Linear Regression
- 2. Decision Tree
- 3. Random Forest
- 4. SVM regressor

```
In [10]: from sklearn.linear model import LinearRegression
         lin reg = LinearRegression()
         lin reg.fit(prepared data, data labels)
Out[10]: LinearRegression(copy X=True, fit intercept=True, n jobs=None, normalize=False)
In [11]: ##testing the predictions with the
         sample data = data.iloc[:5]
         sample labels = data labels.iloc[:5]
         sample data prepared = pipeline transformer(sample data)
         print("Prediction of samples: ", lin_reg.predict(sample_data_prepared))
         Prediction of samples: [29.08069379 27.78336755 26.08031176 12.70419279 22.23454159]
In [12]: print("Actual Labels of samples: ", list(sample_labels))
         Actual Labels of samples: [32.0, 31.0, 26.0, 18.0, 26.0]
         Mean Squared Error
In [13]: from sklearn.metrics import mean_squared_error
         mpg predictions = lin reg.predict(prepared data)
         lin_mse = mean_squared_error(data_labels, mpg_predictions)
         lin_rmse = np.sqrt(lin_mse)
         lin rmse
Out[13]: 2.9590402225760863
```

#### **Decision Tree**

But no model is perfect, this means that our model has overfit the data to a great extent.

We won't be touching out test data until we finalize our model. So, how do we check for what's happening?

#### **Model Evaluation using Cross Validation**

Scikit-Learn's K-fold cross-validation feature randomly splits the training set into K distinct subsets called folds, then it trains and evaluates the model K times, picking a different fold for evaluation every time and training on the other K-1 folds.

The result is an array containing the K evaluation scores:

```
In [16]: from sklearn.model selection import cross val score
         scores = cross val score(tree reg,
                                  prepared data,
                                  data labels,
                                  scoring="neg mean squared error",
                                  cv = 10)
         tree reg rmse scores = np.sqrt(-scores)
In [17]: | tree_reg_rmse_scores
Out[17]: array([2.91510077, 2.77302858, 3.03320169, 3.37374977, 2.19003425,
                2.9557148 , 2.70918576, 5.2122452 , 4.20560302, 2.6146609 ])
In [18]: tree_reg_rmse_scores.mean()
Out[18]: 3.19825247393513
In [19]: | scores = cross_val_score(lin_reg, prepared_data, data_labels, scoring="neg_mean_squared_error", cv = 10)
         lin_reg_rmse_scores = np.sqrt(-scores)
         lin_reg_rmse_scores
Out[19]: array([3.43254597, 3.45157629, 3.6621715, 2.59652976, 2.48023405,
                2.74798115, 3.32524647, 2.42208917, 3.78133275, 2.8573747 ])
In [20]: lin_reg_rmse_scores.mean()
Out[20]: 3.0757081793709324
```

#### **Random Forest model**

Out[21]: 2.5685941515759576

#### **Support Vector Machine Regressor**

Out[22]: 3.08659162080283

**Hyperparameter Tuning using GridSearchCV** 

```
In [23]: from sklearn.model selection import GridSearchCV
         param grid = [
             {'n estimators': [3, 10, 30], 'max features': [2, 4, 6, 8]},
             {'bootstrap': [False], 'n estimators': [3, 10], 'max features': [2, 3, 4]},
         forest reg = RandomForestRegressor()
         grid_search = GridSearchCV(forest_reg, param_grid,
                                    scoring='neg mean squared error',
                                    return train score=True,
                                    cv=10,
         grid search.fit(prepared data, data labels)
Out[23]: GridSearchCV(cv=10, error score=nan,
                      estimator=RandomForestRegressor(bootstrap=True, ccp alpha=0.0,
                                                       criterion='mse', max_depth=None,
                                                      max features='auto',
                                                      max_leaf_nodes=None,
                                                      max samples=None,
                                                      min_impurity_decrease=0.0,
                                                      min_impurity_split=None,
                                                      min samples leaf=1,
                                                      min_samples_split=2,
```

min\_weight\_fraction\_leaf=0.0,
n\_estimators=100, n\_jobs=None,
oob\_score=False, random\_state=None,

verbose=0, warm\_start=False),

iid='deprecated', n jobs=None,

param grid=[{'max features': [2, 4, 6, 8],

scoring='neg mean squared error', verbose=0)

'n estimators': [3, 10, 30]},

pre dispatch='2\*n jobs', refit=True, return\_train\_score=True,

'n estimators': [3, 10]}],

{'bootstrap': [False], 'max features': [2, 3, 4],

```
In [24]: grid search.best params
Out[24]: {'max features': 6, 'n estimators': 10}
In [25]: cv scores = grid search.cv results
         ##printing all the parameters along with their scores
         for mean score, params in zip(cv scores['mean test score'], cv scores["params"]):
             print(np.sqrt(-mean score), params)
         3.5298036773871795 {'max features': 2, 'n estimators': 3}
         3.1543065463436575 {'max features': 2, 'n estimators': 10}
         2.9200301478012345 {'max_features': 2, 'n_estimators': 30}
         3.4367720636873687 {'max features': 4, 'n estimators': 3}
         2.86113945515678 {'max features': 4, 'n estimators': 10}
         2.7573866919109475 {'max_features': 4, 'n_estimators': 30}
         3.416158941425653 {'max_features': 6, 'n_estimators': 3}
         2.6743670980252503 {'max features': 6, 'n estimators': 10}
         2.7323235595460096 {'max features': 6, 'n estimators': 30}
         3.191919436072898 {'max_features': 8, 'n_estimators': 3}
         2.6796452995595543 {'max features': 8, 'n estimators': 10}
         2.68494337425188 {'max features': 8, 'n estimators': 30}
         3.2998308648412786 {'bootstrap': False, 'max features': 2, 'n estimators': 3}
         2.9058143534045002 {'bootstrap': False, 'max features': 2, 'n estimators': 10}
         3.1910324793160885 {'bootstrap': False, 'max features': 3, 'n estimators': 3}
         2.8726444715767814 {'bootstrap': False, 'max_features': 3, 'n_estimators': 10}
         3.4152733193520746 {'bootstrap': False, 'max_features': 4, 'n_estimators': 3}
         2.693486477642302 {'bootstrap': False, 'max features': 4, 'n estimators': 10}
```

#### **Checking Feature importance**

```
In [26]: # feature importances
         feature importances = grid search.best estimator .feature importances
         feature importances
Out[26]: array([0.28706325, 0.32659439, 0.09904163, 0.1015734 , 0.0182555 ,
                0.11371115, 0.03097858, 0.01759758, 0.00280047, 0.00114504,
                0.00123901])
In [27]: extra attrs = ["acc on power", "acc on cyl"]
         numerics = ['float64', 'int64']
         num attrs = list(data.select dtypes(include=numerics))
         attrs = num attrs + extra attrs
         sorted(zip(attrs, feature importances), reverse=True)
Out[27]: [('acc_on_power', 0.030978576378998753),
          ('acc on cyl', 0.01759757612432403),
          ('Weight', 0.10157340340308983),
          ('Model Year', 0.11371115169237556),
          ('Horsepower', 0.09904163488978267),
          ('Displacement', 0.3265943896835737),
          ('Cylinders', 0.28706324723924487),
          ('Acceleration', 0.018255500770649784)]
```

#### **Evaluating the entire system on Test Data**

```
In [28]: final_model = grid_search.best_estimator_

X_test = strat_test_set.drop("MPG", axis=1)
y_test = strat_test_set["MPG"].copy()

X_test_preprocessed = preprocess_origin_cols(X_test)
X_test_prepared = pipeline_transformer(X_test_preprocessed)

final_predictions = final_model.predict(X_test_prepared)
final_mse = mean_squared_error(y_test, final_predictions)
final_rmse = np.sqrt(final_mse)
```

```
In [29]: final_rmse
Out[29]: 3.137180461815992
```

#### Creating a function to cover this entire flow

```
In [30]: def predict mpg(config, model):
             if type(config) == dict:
                 df = pd.DataFrame(config)
             else:
                 df = config
             preproc df = preprocess origin cols(df)
             prepared_df = pipeline_transformer(preproc_df)
             y_pred = model.predict(prepared_df)
             return y pred
In [31]: ##checking it on a random sample
         vehicle_config = {
             'Cylinders': [4, 6, 8],
             'Displacement': [155.0, 160.0, 165.5],
             'Horsepower': [93.0, 130.0, 98.0],
             'Weight': [2500.0, 3150.0, 2600.0],
             'Acceleration': [15.0, 14.0, 16.0],
             'Model Year': [81, 80, 78],
             'Origin': [3, 2, 1]
         predict mpg(vehicle config, final model)
Out[31]: array([29.22, 18.3 , 19.24])
```

#### Save the Model

```
In [32]: import pickle

In [33]: ##saving the model
with open("model.bin", 'wb') as f_out:
    pickle.dump(final_model, f_out)
    f_out.close()

In [34]: ##loading the model from the saved file
with open('model.bin', 'rb') as f_in:
    model = pickle.load(f_in)
predict_mpg(vehicle_config, model)

Out[34]: array([29.22, 18.3 , 19.24])

In []:
```