BANGLADESH ARMY INTERNATIONAL UNIVERSITY OF SCIENCE & TECHNOLOGY (BAIUST)



Course Title: Artificial intelligence Sessional

Course Code: CSE-404

Project report on used car price prediction using machine learning

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Acknowledgement

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we look into matters that would have been overlooked otherwise.

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Abstract

In many branches of materials science it is now routine to generate data sets of such large size and dimensionality that conventional methods of analysis fail. Paradigms and tools from data science and machine learning can provide scalable approaches to identify and extract trends and patterns within voluminous data sets, To understand what a dataset is, we must first discuss the components of a dataset. A single row of data is called an instance. Datasets are a collection of instances that all share a common attribute. Machine learning models will generally contain a few different datasets, each used to fulfill various roles in the system.

For machine learning models to understand how to perform various actions, training datasets must first be fed into the machine learning algorithm, followed by validation datasets (or testing datasets) to ensure that the model is interpreting this data accurately.

Once you feed these training and validation sets into the system, subsequent datasets can then be used to sculpt your machine learning model going forward. The more data you provide to the ML system, the faster that model can learn and improve.

Work Summary

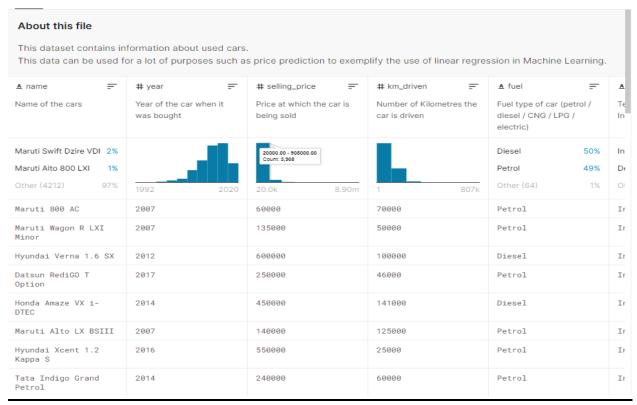
A data set is a collection of data. ... In Machine Learning projects, we need a training data set. It is the actual data set used to train the model for performing various actions

This dataset contains information about used cars.

This data can be used for a lot of purposes such as price prediction to exemplify the use of linear regression in Machine Learning.

The columns in the given dataset are as follows:

- 1. name
- 2. year
- 3. selling_price
- 4. km_driven
- 5. fuel
- 6. seller_type
- 7. transmission
- 8. Owner

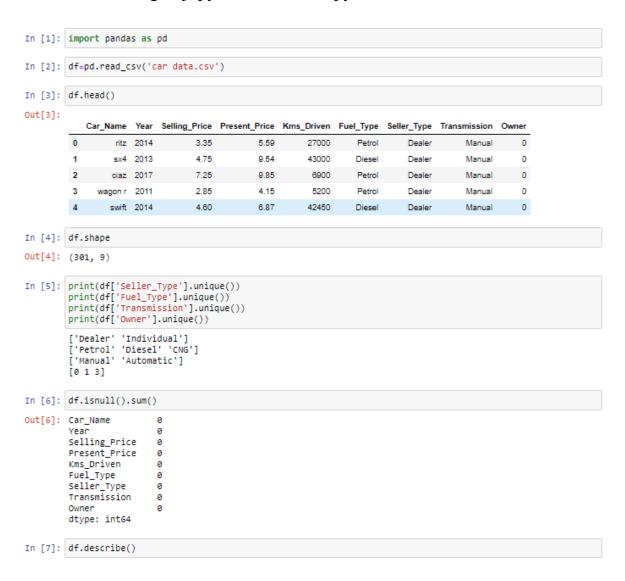


This data set from https://www.kaggle.com/nehalbirla/vehicle-dataset-from-cardekho contains information about used car by using this data set implement a prediction model to predict selling price for car.

Work description

We have created a new environment in anaconda prompt as every model has different dependencies.

Then we start working in jupyter notebook in python 3.7



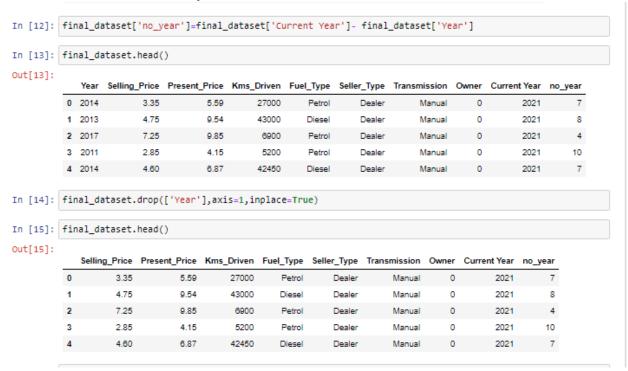
As there is no null value no need to determine mean value.

Next thing we have done, we used df.describe() to see some feature of data set such as standard deviation, count etc.

Preprocessing

Then we created our final data set we exclude car_name as it doesn't include mathematical importance.

After that we created no_year



And after we dropped Year column.

Convert categorical variable into dummy/indicator variables

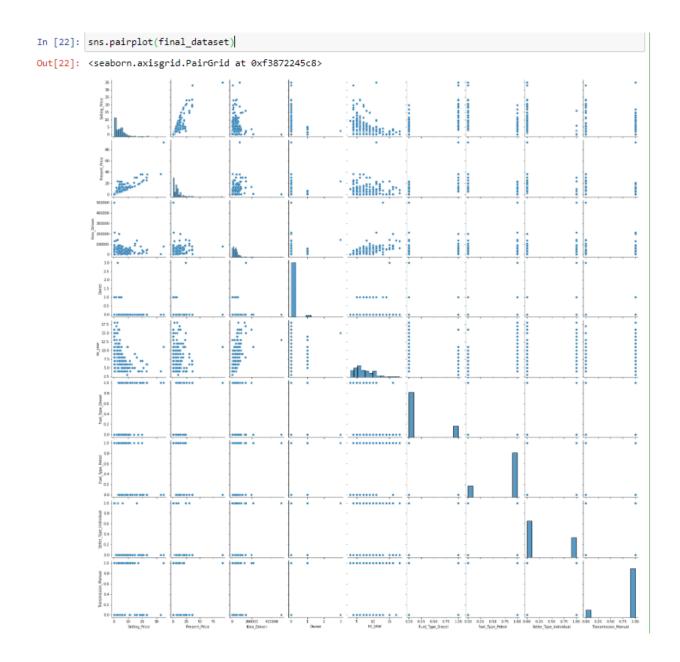
Here fuel_type is a categorical variable.

final_dataset=pd.get_dummies(final_dataset,drop_first=True) is used to convert it.

Next step is finding co-relation using .corr function. Co-relation say how one feature is connected to another.

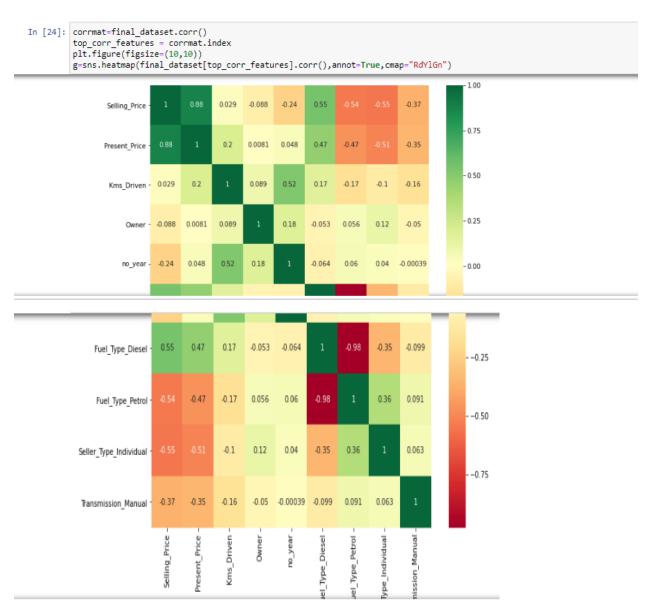
Using seaborn we can see pairplot

Which is shown bellow



Next step is plot it in heat map

A **heatmap** is a two-dimensional graphical representation of data where the individual values that are contained in a matrix are represented as colors. The seaborn **python** package allows the creation of annotated heatmaps which can be tweaked using Matplotlib tools as per the requirement.



Then we allocate all our independent feature

X = final_dataset.iloc[:,1:] these are final_dataset=df[[Present_Price, Kms_Driven, Fuel_Type, Seller_Type, Transmission, Owner, Current, Year no_year]]

{1: is representing column starting form Present_Price}
our dependent feature is selling price
y = final_dataset.iloc[:,0]

Feature Importance and Doing a train test split

As it is regression problem we are using ExtraTreesRegressor to find important feature

ExtraTreesRegressor()

```
print(model.feature_importances_)
[4.11946112e-01 4.15172773e-02 3.99537471e-04 7.78993520e-02
 2.04928051e-01 1.89998368e-02 1.21892158e-01 1.22417674e-01]
feat_importances = pd.Series(model.feature_importances_, index=X.columns)
feat_importances.nlargest(5).plot(kind='barh')
plt.show()
           no_year
 Seller_Type_Individual
 Transmission_Manual
    Fuel_Type_Diesel
       Present Price
                0.00
                     0.05
                           0.10
                                0.15
                                     0.20
                                           0.25
                                                 0.30
                                                       0.35
```

We are taking all feature but only top 5 is shown in chart.

Algorithm

Random forest algorithm

Random forest is a flexible, easy to use machine learning algorithm that produces, even without hyper-parameter tuning, a great result most of the time. It is also one of the most used algorithms, because of its simplicity and diversity (it can be used for both classification and regression tasks). In this post we'll learn how the random forest algorithm works, how it differs from other algorithms and how to use it.

HOW RANDOM FOREST WORKS

Random forest is a supervised learning algorithm. The "forest" it builds, is an ensemble of decision trees, usually trained with the "bagging" method. The general idea of the bagging method is that a combination of learning models increases the overall result.

Put simply: random forest builds multiple decision trees and merges them together to get a more accurate and stable prediction.

One big advantage of random forest is that it can be used for both classification and regression problems, which form the majority of current machine learning systems. Let's look at random forest in classification, since classification is sometimes considered the building block of machine learning

```
from sklearn.model_selection import RandomizedSearchCV
random_grid = {'n_estimators': n_estimators,
                'max_features': max_features,
               'max_depth': max_depth,
               'min_samples_split': min_samples_split,
               'min_samples_leaf': min_samples_leaf}
print(random grid)
{'n_estimators': [100, 200, 300, 400, 500, 600, 700, 800, 900, 1000, 1100, 1200], 'max_features': ['auto', 'sqrt'], 'max_dept
h': [5, 10, 15, 20, 25, 30], 'min_samples_split': [2, 5, 10, 15, 100], 'min_samples_leaf': [1, 2, 5, 10]}
rf = RandomForestRegressor()
rf_random = RandomizedSearchCV(estimator = rf, param_distributions = random_grid,scoring='neg_mean_squared_error', n_iter = 10, or representations = respectively.
rf_random.fit(X_train,y_train)
Fitting 5 folds for each of 10 candidates, totalling 50 fits
[CV] END max_depth=10, max_features=sqrt, min_samples_leaf=5, min_samples_split=5, n_estimators=900; total time=
[CV] END max_depth=10, max_features=sqrt, min_samples_leaf=5, min_samples_split=5, n_estimators=900; total time=
[CV] END max_depth=10, max_features=sqrt, min_samples_leaf=5, min_samples_split=5, n_estimators=900; total time=
[CV] END max_depth=10, max_features=sqrt, min_samples_leaf=5, min_samples_split=5, n_estimators=900; total time=
[CV] END max_depth=10, max_features=sqrt, min_samples_leaf=5, min_samples_split=5, n_estimators=900; total time=
```

Performing hyper parameter tuning using RandomizedSearchCV RandomizedSearchCV helps to find best parameter form n_estimators Next step if fitting

Prediction

predictions=rf_random.predict(X_test)

```
In [45]: plt.scatter(y_test,predictions)
Out[45]: <matplotlib.collections.PathCollection at 0xf38ea20288>
```

Plotting is linearly available which represent our prediction is efficient and better.

Deployment

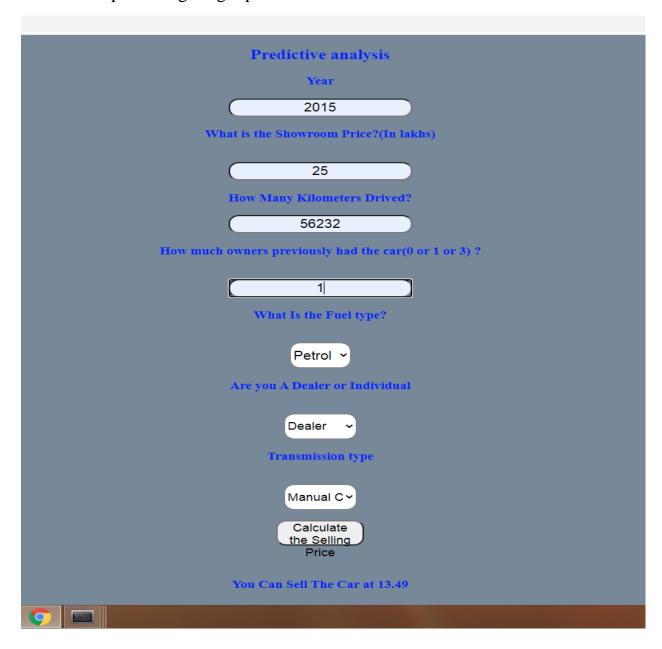
For frontend app.py is used. Using flask micro web framework we rendered index.html our model random_forest_regression_model.pkl we loaded the pickle file to get prediction value

```
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C:\Users\shubha\Desktop\machine learning\app.py
app.py × Billing-System.py ×
              from flask import Flask, render_template, request
              import jsonify
              import numpy as np
import sklearn
              from sklearn.preprocessing import StandardScaler
              model = pickle.load(open('random_forest_regression_model.pkl', 'rb'))
             @app.route('/',methods=['GET'])
def Home():
                  return render_template('index.html')
            standard_to = StandardScaler()
@app.route("/predict", methods=['POST'])
def predict():
    Fuel_Type_Diesel=0
    if request.method == 'POST':
        Year = int(request.form['Year'])
        Present_Price=float(request.form['Present_Price'])
        Kms_Driven=int(request.form['Kms_Driven'])
        Kms_Driven2=np.log(Kms_Driven)
        Owner=int(request.form['Owner'])
        Fuel_Type_Petrol='Petrol'):
        if(Fuel_Type_Petrol='Petrol'):
        Fuel_Type_Petrol=1
              standard_to = StandardScaler()
                                    Fuel_Type_Petrol=1
                                   Fuel_Type_Diesel=0
                             Fuel_Type_Petrol=0
Fuel_Type_Diesel=1
                        Seller_Type_Individual=request.form['Seller_Type_Individual']
if(Seller_Type_Individual=='Individual'):
    Seller_Type_Individual=1
                            Seller_Type_Individual=0
                        Transmission_Mannual=request.form['Transmission_Mannual']
if(Transmission_Mannual=='Mannual'):
                               Transmission_Mannual=1
                         prediction=model_predict([[Present_Price,Kms_Driven2,Owner,Year,Fuel_Type_Diesel,Fuel_Type_Petrol,Seller_Type_Individual,Transmission_Mannual]])
                         output=round(prediction[0],2)
                              return render_template('index.html',prediction_texts="Sorry you cannot sell this car")
                               return render_template('index.html',prediction_text="You Can Sell | the Car at {}".format(output))
                         return render_template('index.html')
              if __name__=="__main__":
    app.run(debug=True)
```

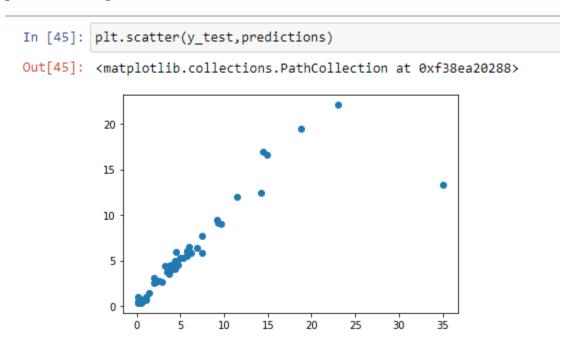
Output

As we used flask we are getting web output we run app.py in http://127.0.0.1:5000/
Show in output after giving input.



Performance

Plotting Cross-Validated Predictions for an un-aggregated model, an un-aggregated (i.e. the usual) cross validation can be used as approximation for predictive performance/generalization error estimate



In the x axis we have measured data and y axis we have prediction Line is liner indicate good accuracy of model.

Discussion: there are many external factor in measuring the value of car such as engine hp, torque produce, condition etc. can play an important role for determining the price of a used car and can play an important role in feature importance if they were included in csv.

Reference:

1. Vehicle dataset. Used Cars data form websites (https://www.kaggle.com/vehicle-dataset-from-cardekho)