## ST1 Assignment 9 (Capstone Programming Project) By William Blucher and Kim Groves

```
from google.colab import drive
drive.mount('/content/drive')

   Drive already mounted at /content/drive; to attempt to forcibly remount, call drive.mount("/content/drive", force_remount=True).

%cd /content/drive/MyDrive/Colab_Project
   /content/drive/MyDrive/Colab_Project
!ls

DataForML.pkl final_LinearRegression.pkl NFLX.csv
```

This project is based on Netflix stock prices for the past five years. The data used for this project is on the Kaggle repository. (https://www.kaggle.com/datasets/jainilcoder/netflix-stock-price-prediction/data)

- · This data set contains (xyz) amount of data
- Our project is to develop a machine learning model that can predict the next stock price based on the previous days data.
- For solving this problem we (... fill wih process)

### Step 1: Reading the data Reading the data with python

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

import pandas as pd
import numpy as np

# This is simply used to access and shape the data so that it is easier for us to use later on.
NFLXData = pd.read_csv("/content/drive/MyDrive/Colab_Project/NFLX.csv", encoding='latin')
print("Shape before deleting duplicates:", NFLXData.shape)

# We remove any duplicates that may be within the data to ensure that no errors will occur later on.
NFLXData=NFLXData.drop_duplicates()
print("Shape after deleting duplicate views:", NFLXData.shape)

# Next we check that the data is reflected the way we want.
NFLXData.head(10)

Shape before deleting duplicates: (1009, 7)
Shape after deleting duplicate views: (1009, 7)
```

Date 0pen High Low Close Adj Close Volume **0** 2018-02-05 262.000000 267.899994 250.029999 254.259995 254.259995 11896100 **1** 2018-02-06 247.699997 266.700012 245.000000 265.720001 12595800 265.720001 **2** 2018-02-07 266.579987 272.450012 264.329987 264.559998 264.559998 8981500 **3** 2018-02-08 267.079987 267.619995 250.000000 250.100006 250.100006 9306700 **4** 2018-02-09 253.850006 255.800003 236.110001 249.470001 249.470001 16906900 **5** 2018-02-12 252.139999 259.149994 249.000000 257.950012 257.950012 8534900 **6** 2018-02-13 257.290009 261.410004 254.699997 258.269989 258.269989 6855200 **7** 2018-02-14 260.470001 269.880005 260.329987 266.000000 266.000000 10972000 **8** 2018-02-15 270.029999 280.500000 267.630005 280.269989 10759700 280.269989 **9** 2018-02-16 278.730011 281.959991 275.690002 278.519989 278.519989 8312400

### A) Key Observations Regarding Data

There is a total of 1,009 rows, representing days and 7 categories of data for each day. The following are the descriptions of these categories:

- · Date Date stock open and closed (explain why each is continuous or not)
- · Open Stock price at market open

- · High high point of stock price
- · Low low point of stock price
- · Close stock price at market close
- · Adj Close adjusted close price
- · Volume amount of available stocks

### Step 2: Problem Statement Definition

- · Creating a prediction model to predict the price(Close) of the stocks.
- Target Variable: Close Predictors/Features: Open, High, Low, Adj Close and Volume.

### Step 3: Choosing the appropriate ML/AI Algorithm for Data Analysis.

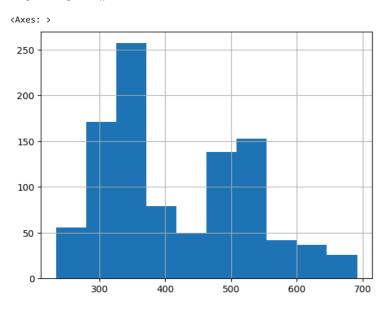
• Considering the problem statement, we need to create a supervised ML Regression model, as the target variable (and the other variables) are **continuous**. This hypothesis is further supported by the results further down.

## Step 4: Looking at the class distribution (Target variable distribution to check if the data is balanced or skewed.)

- This is used to check if the data is balanced or skewed (if the data is skewed too heavily, the accuracy of the ML algorithm may be negatively impacted.)
- · A bell curve would be the ideal result.
- The bell curve is ideally required to ensure that the Regression Model will work with the highest accuracy value, however slightly skewed data should still work effectively.
- If the data is too skewed the machine learning algorithm will struggle when trying to learn all possible scenarios for the data.

#### %matplotlib inline

- # This is used as the data we have chosen (close) is continuous.
- # The results of this, as stated above, will help us estimate the accuracy of our ML algorithm. NFLXData["Close"].hist()



#### Observations from Step 4

So based on this, the data distribution is fairly skewed. However, noting that this data set is based around stocks, and stocks are typically a volatile data set that is subject to influence from outside effectors (such as politics or standing of the company) this is to be expected and we can proceed regardless.

### Step 5: Basic Exploratory Data analysis

# From this, we look at a sample of the rows of data. NFLXData.head()

	Date	0pen	High	Low	Close	Adj Close	Volume
0	2018-02-05	262.000000	267.899994	250.029999	254.259995	254.259995	11896100
1	2018-02-06	247.699997	266.700012	245.000000	265.720001	265.720001	12595800
2	2018-02-07	266.579987	272.450012	264.329987	264.559998	264.559998	8981500
3	2018-02-08	267.079987	267.619995	250.000000	250.100006	250.100006	9306700
4	2018-02-09	253.850006	255.800003	236.110001	249.470001	249.470001	16906900

# Using tail and head (previous block) we are able to get a 'snapshot' of what the entire sample is like. NFLXData.tail()

	Date	Open	High	Low	Close	Adj Close	Volume
1004	2022-01- 31	401.970001	427.700012	398.200012	427.140015	427.140015	20047500
1005	2022-02- 01	432.959991	458.480011	425.540009	457.130005	457.130005	22542300
1006	2022-02- 02	448.250000	451.980011	426.480011	429.480011	429.480011	14346000
4007	2022-02-	404 440000	400 000040	40.4.070000	405 000000	405 000000	0005000

<sup>#</sup> This step allows us to inspect.

- # Using this we are able to identify the Data Type, the columns, and the number of empty cells.
- # In this instance there is no empty cells.
- # Additionally we can view what columns need to be removed (Qualitative data will not work for ML algorithms and thus needs to be remove NFLXData.info()

# Remove the data that was identified as qualitative in the previous step.
NFLXData = NFLXData.drop("Date", axis="columns")

### NFLXData.info()

# This step is another way to gain an understanding of the data within the set.
NFLXData.describe(include='all')

	0pen	High	Low	Close	Adj Close	Volume
count	1009.000000	1009.000000	1009.000000	1009.000000	1009.000000	1.009000e+03
mean	419.059673	425.320703	412.374044	419.000733	419.000733	7.570685e+06
std	108.537532	109.262960	107.555867	108.289999	108.289999	5.465535e+06
min	233.919998	250.649994	231.229996	233.880005	233.880005	1.144000e+06
25%	331.489990	336.299988	326.000000	331.619995	331.619995	4.091900e+06
50%	377.769989	383.010010	370.880005	378.670013	378.670013	5.934500e+06
75%	509.130005	515.630005	502.529999	509.079987	509.079987	9.322400e+06
max	692.349976	700.989990	686.090027	691.690002	691.690002	5.890430e+07

<sup>#</sup> This allows us to identify how many duplicates appear amongst each respective row.

Open 976 High 983 Low 989 Close 988 Adj Close 988 Volume 1005 dtype: int64

All the data values are > 20, therefore all the values are continuous and not categorical.

### Observations from Step 5 - Basic Data Exploratory Analysis

- · based on the exploration above we discovered the following
- · The Open column is continuos Selected for now
- · The Adj Close column is continuous Selected for now (Will be reviewd later as results are identical to close so far.
- · High Continuous and Selected
- · Low Continuous and Selected
- · Close Continuous, selected as target variable
- Volume Contiuous, selected for now.

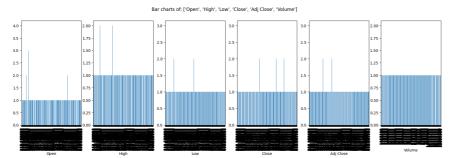
### Step 7: Removal of bad columns

There are no bad columns left to remove as of this stage, so none are removed.

### Step 8: Visual Exploratory Data Analysis

- · Visualising the data with a bar plot
- As previously mentioned, without categorical data, this step will not produce correct bar charts, however for the sake of exhaustion it will be done anyway.

<sup>#</sup> Any columns that are roughly 20 or less may be categories, not contiuous. NFLXData.nunique()



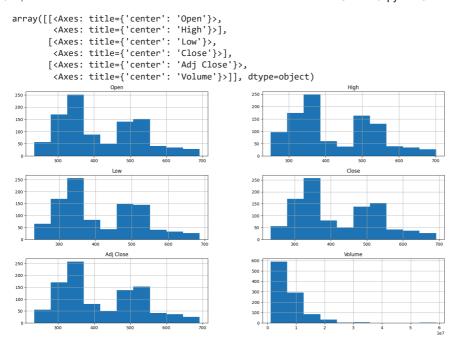
### Observations

As predicted, bar charts are not usable as only works with categorical data, of which there is none.

## Step 9: Visualising the distribution

Visualise the data with histograms.

# Displaying each of the data columns as histograms allows us to identify any outliers that may be present. NFLXData.hist(['Open', 'High', 'Low', 'Close', 'Adj Close', 'Volume'], figsize=(18,10))



From this step, it can be observed that within 'Volume" there is an outlier beyond roughly 3.5x10^7.

As this column will not be used as a predictor later on in the project, this fact can be ignored.

No other outliers are observed at this time.

## Step 10: Outlier Analysis

Outliers are as seen below for volume, as stated this is just to represent that we understand the process, however we will not be removing this data as it will not be used later in the models.

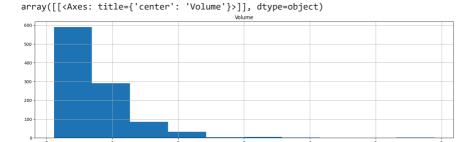
 $\mbox{\tt\#}$  This allows us to view the data, and identify where the outlier may appear.

NFLXData['Volume'][NFLXData['Volume']<(4\*(10\*\*7))].sort\_values(ascending=True)

```
1144000
982
        1287200
883
        1367800
        1595500
896
979
        1621100
999
       32346000
177
       32610900
744
       32637500
49
       33866500
428
       38258900
Name: Volume, Length: 1007, dtype: int64
```

# This allows us to single out the volume histogram and further view the outlier.

NFLXData.hist(['Volume'], figsize=(18,5))



# Step 12: (11 has been skipped as mentioned but number pattern has been followed for ease of reading) Missing Value Analysis

Check if any missing values have been introduced into the data.

# Next, once again we double check to ensure that there are no null values that could skew the results later on.

NFLXData.isnull().sum()

Open
High
Low
Close
Adj Close
Volume
dtype: int64

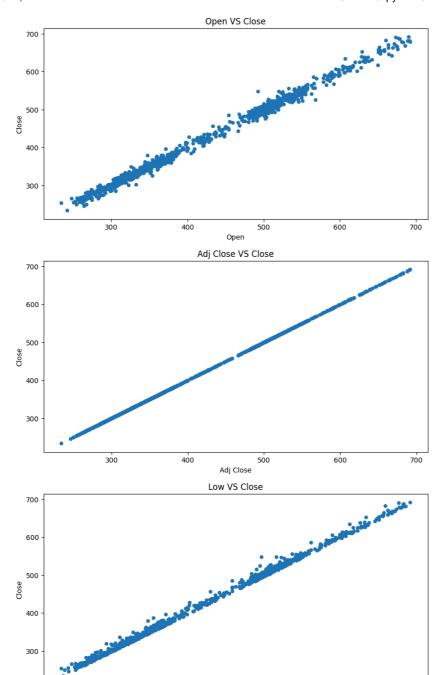
## Step 13: Feature selection - Scatter Plot and Correlation matrix.

Checking the correlation of the data to identify the columns of interest.

```
# Using scatter grams, we plot the current predictor columns against our selected 'Close' column
# This is used to visualise any positive or negative trends within the data.

ContinuousCols=['Open', 'Adj Close', 'Low', 'High', 'Volume']

for predictor in ContinuousCols:
    NFLXData.plot.scatter(x=predictor, y="Close", figsize=(10,5), title=predictor+" VS "+ 'Close')
```



## **Scatter Charts interpretation**

Upon reviewing the scatter plots, all data sets except volume have a strong positive correlation. Volume seems to have no identifiable correlation with our target variablel.

Finally, Adj Close ones again appears to be an exact copy of Close.

## Step 14: Feature selection with Correlation Value

Using the corr() function to identify the correlations of the data.

```
\mbox{\tt\#} Again, we use a corrolation function to check the correlations of the data.
```

# Greater than 0.5, or less than -0.5 ususally indicates a positive or negative correlation.

ContinuousCols=['Close', 'Open', 'Adj Close', 'Low', 'High', 'Volume']

CorrelationData=NFLXData[ContinuousCols].corr()
CorrelationData

	Close	Open	Adj Close	Low	High	Volume
Close	1.000000	0.996812	1.000000	0.998544	0.998551	-0.413362
Open	0.996812	1.000000	0.996812	0.998508	0.998605	-0.415838
Adj Close	1.000000	0.996812	1.000000	0.998544	0.998551	-0.413362
Low	0.998544	0.998508	0.998544	1.000000	0.998203	-0.432116
High	0.998551	0.998605	0.998551	0.998203	1.000000	-0.400699
Volume	-0.413362	-0.415838	-0.413362	-0.432116	-0.400699	1.000000

From the data above, we can see that open, low and high all have strong positive correlations with our focused 'Close'.

Additionally, we can identify that the 'ADJ Close' column has a correlation percentage of 100%, therefore meaning that it is, for this data set, functionally the same row. As such, it should be removed IOT not skew the accuracy of the ML model.

# This allows us to filter specifically for 'Close' and its correlations, specifically showing those that are a positive correlation. CorrelationData['Close'][abs(CorrelationData['Close']) > 0.5]

```
Close 1.000000
Open 0.996812
Adj Close 1.000000
Low 0.998544
High 0.998551
Name: Close, dtype: float64
```

Again this further shows that the 'Adj Close' column is an 100% match with our selected 'Close' column, and should therefore be discarded.

Additionally, from this we can further deduce that the volume column does not have a strong enough correlation, whether positive or negative, to be considered from this data set.

### Selecting Final Predictors/Features for building Machine Learning/Al model.

Step 15 and 16 are skipped as no categorical data.

# As we do not have categorical data, we are able to simply select the columns that we have good correlation.

```
SelectedColumns = ['Open', 'Low', 'High']
```

DataForML = NFLXData[SelectedColumns]
DataForML.head()

	0pen	Low	High
0	262.000000	250.029999	267.899994
1	247.699997	245.000000	266.700012
2	266.579987	264.329987	272.450012
3	267.079987	250.000000	267.619995
4	253.850006	236.110001	255.800003

# This saves the final data subset so that we can reference it later.

DataForML.to\_pickle('DataForML.pkl')

# Step 17: Data Pre-processing for Machine Learning Model Building or Model Development

Converting the nominal variable to numeric using get\_dummies()

```
# This treats all the nominal variables using dummy variables.
DataForML_Numeric = pd.get_dummies(DataForML)
```

# Include the target variable in the data.
DataForML\_Numeric['Close'] = NFLXData['Close']

# Print a sample of the first five rows
DataForML\_Numeric.head()

	0pen	Low	High	Close
0	262.000000	250.029999	267.899994	254.259995
1	247.699997	245.000000	266.700012	265.720001
2	266.579987	264.329987	272.450012	264.559998
3	267.079987	250.000000	267.619995	250.100006
4	253 850006	236 110001	255 800003	249 470001

### Step 18: Machine Learning Model Development

Splitting the data into training and testing sample

### Step 18: Standardisation/Normalisation of data

```
# Standardisation of data
# This is stated as optional, however for the sake of exhaustion it is needed for KNN and nueral networks.
from sklearn.preprocessing import StandardScaler, MinMaxScaler
# Chose MinMaxScaler over StandardScaler as the data doesnt have varying scales.
PredictorScaler = MinMaxScaler()
# Stored for use later.
PredictorScalerFit = PredictorScaler.fit(X)
# Generating the standardised values of X
X = PredictorScalerFit.transform(X)
# Once again splitting the data into test and train.
from sklearn.model_selection import train_test_split
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.3, random_state=42)
# This step shows us how much data is in each set.
print(X train.shape)
print(y_train.shape)
print(X_test.shape)
print(y_test.shape)
     (706, 3)
     (706,)
     (303, 3)
     (303,)
```

### Step 20: Multiple Linear Regression Algorithm

```
# Multiple Linear Regression
from sklearn.linear_model import LinearRegression
RegModel = LinearRegression()
# Print the parameters of the linear regression.
print(RegModel)
# Create the model using the training data.
LREG=RegModel.fit(X_train,y_train)
prediction=LREG.predict(X_test)
from sklearn import metrics
# Measurement of how well the data fits with the test.
print('R2 Value:',metrics.r2_score(y_train, LREG.predict(X_train)))
#Print the title of the output
print('\n######## Model Validation and Accuracy Calculations ########")
# Creating the sample data values that will be printed
TestingDataResults=pd.DataFrame(data=X test, columns=Predictors)
TestingDataResults[TargetVariable]=y_test
TestingDataResults[('Predicted'+TargetVariable)]=np.round(prediction)
# Printing five of the predicted values
print(TestingDataResults.head())
# Calculates the error for each of the rows
TestingDataResults['APE']=100 * ((abs(
  TestingDataResults['Close']-TestingDataResults['PredictedClose']))/TestingDataResults['Close'])
MAPE=np.mean(TestingDataResults['APE'])
MedianMAPE=np.median(TestingDataResults['APE'])
Accuracy =100 - MAPE
MedianAccuracy=100- MedianMAPE
print('Mean Accuracy on test data:', Accuracy) # If negative, indicative of an outlier present in the data.
print('Median Accuracy on test data:', MedianAccuracy)
# Custom function to ensure accuracy.
# Ensures there are no zeros in data when using MAPE
def Accuracy_Score(orig,pred):
    MAPE = np.mean(100 * (np.abs(orig-pred)/orig))
    return(100-MAPE)
# Using the make scorer of sklearn.metrics to have a custom MAPE scoring calculation.
from sklearn.metrics import make_scorer
custom_Scoring=make_scorer(Accuracy_Score, greater_is_better=True)
# Getting crossvalidation from sklearn
from sklearn.model_selection import cross_val_score
# Running a 10-fold cross validation to produce more accuracy values for comparison.
\# Finally passes the data of x and y as K-fold splits the data and authomatically chooses train/test.
# This step prints the main values we care about.
Accuracy_Values=cross_val_score(RegModel, X , y, cv=10, scoring=custom_Scoring)
print('\nAccuracy values for 10-fold Cross Validation:\n',Accuracy_Values)
print('\nFinal Average Accuracy of the model:', round(Accuracy_Values.mean(),2))
     LinearRegression()
     R2 Value: 0.9986731542285744
     ######## Model Validation and Accuracy Calculations #########
                               High
                                          Close PredictedClose
                      Low
     0 0.577471 0.587917 0.576875 509.640015
     1 0.592304 0.561601 0.572101 494.730011
     2 0.596449 0.589962 0.585891 500.859985
                                                           507.0
     3 0.330279 0.322451 0.303482 380.070007
                                                          381.0
     4 0.195188 0.182474 0.166985 315.100006
                                                           319.0
     Mean Accuracy on test data: 99.3029073385015
     Median Accuracy on test data: 99.47124720149543
     Accuracy values for 10-fold Cross Validation:
      [99.21860151 99.21910542 99.19755296 99.4083654 99.47040237 98.9092798
      99.20776329 99.40785386 99.62149833 99.22984505]
     Final Average Accuracy of the model: 99.29
```

### Decision Tree Regressor

```
# Decision Trees (Bulk Else-If statements)
from sklearn.tree import DecisionTreeRegressor
RegModel = DecisionTreeRegressor(max_depth=5,criterion='friedman_mse')
# A good range of Max_depth is roughly between 2 - 20
# Printing all parameters of the decision tree.
print(RegModel)
# Create the mdodel using the training data.
DT=RegModel.fit(X_train,y_train)
prediction=DT.predict(X_test)
from sklearn import metrics
# Measuring how well the data fits.
print('R2 Value:',metrics.r2_score(y_train, DT.predict(X_train)))
# Using Matplotlib, plot the 10 largest importance pieces.
%matplotlib inline
feature_importances = pd.Series(DT.feature_importances_, index=Predictors)
feature_importances.nlargest(10).plot(kind='barh')
# Print statement model title.
print('\n##### Model Validation and Accuracy Calculations ########")
# Setting up the variables to be printed.
TestingDataResults=pd.DataFrame(data=X_test, columns=Predictors)
TestingDataResults[TargetVariable]=y_test
TestingDataResults[('Predicted'+TargetVariable)]=np.round(prediction)
# Printing the sample of predicted values.
print(TestingDataResults.head())
# Calculating the individual error for each row.
TestingDataResults['APE']=100 * ((abs(
  TestingDataResults['Close']-TestingDataResults['PredictedClose']))/TestingDataResults['Close'])
# Calculating mean and median MAPE
MAPE=np.mean(TestingDataResults['APE'])
MedianMAPE=np.median(TestingDataResults['APE'])
# Calculating accuracy as a percentage.
Accuracy =100 - MAPE
MedianAccuracy=100 - MedianMAPE
print('Mean Accuracy on test data:', Accuracy) # If negative, outlier present in data set.
print('Median Accuracy on test data:', MedianAccuracy)
# Custom function to ensure accuracy.
# Also ensure no zeros are present that may skew data.
def Accuracy_Score(orig,pred):
    MAPE = np.mean(100 * (np.abs(orig-pred)/orig))
    return(100-MAPE)
# Using sklearn.metrics's make_scorer to have a custom MAPE scoring calculation.
from sklearn.metrics import make scorer
custom_Scoring=make_scorer(Accuracy_Score, greater_is_better=True)
# Getting the crossvalidation function from sklearn
from sklearn.model_selection import cross_val_score
# Printing the 10-fold cross validation so to give further values to compare.
\# Finally, pass the data of the x and y as K-fold splits the data and authomatically decides test or train.
Accuracy_Values=cross_val_score(RegModel, X , y, cv=10, scoring=custom_Scoring)
print('\nAccuracy values for 10-fold Cross Validation:\n',Accuracy_Values)
print('\nFinal Average Accuracy of the model:', round(Accuracy_Values.mean(),2))
```

509.0

386.0

322.0

DecisionTreeRegressor(criterion='friedman\_mse', max\_depth=5) R2 Value: 0.9976613969339876

##### Model Validation and Accuracy Calculations ##########

Open Low High Close PredictedClose

0 0.577471 0.587917 0.576875 509.640015 509.0

1 0.592304 0.561601 0.572101 494.730011 496.0

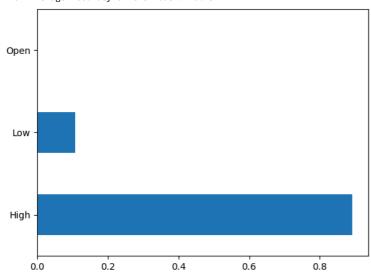
4 0.195188 0.182474 0.166985 315.100006 Mean Accuracy on test data: 98.56869716791199 Median Accuracy on test data: 98.8354057430968

2 0.596449 0.589962 0.585891 500.859985

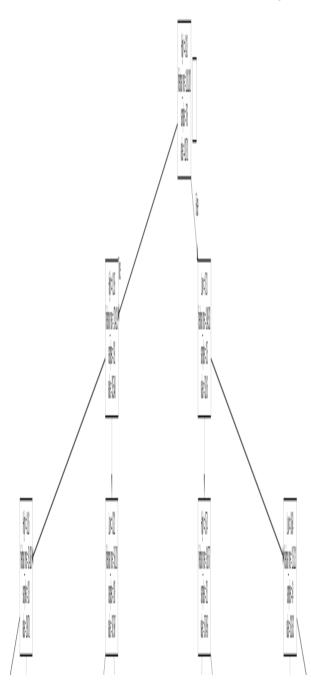
3 0.330279 0.322451 0.303482 380.070007

Accuracy values for 10-fold Cross Validation: [98.40322063 98.50561057 98.48597259 98.84648187 98.33407053 97.24234762 97.95457369 98.81150986 98.98729668 95.69681323]

Final Average Accuracy of the model: 98.13



### Visualising the Decision tree



Random Forest Regressor

```
# Random Forest
from sklearn.ensemble import RandomForestRegressor
RegModel = RandomForestRegressor(max_depth=4, n_estimators=400,criterion='friedman_mse')
# A good range for the max_depth is generatlly 2-10 and the should be between 100-1000
# Printing all the parameters
print(RegModel)
# Create the model using the training data
RF=RegModel.fit(X_train,y_train)
prediction=RF.predict(X test)
from sklearn import metrics
# Measurment of how well the data fits within the model.
print('R2 Value:',metrics.r2_score(y_train, RF.predict(X_train)))
# Plot of the 10 largest importance on a bar graph
%matplotlib inline
feature_importances = pd.Series(RF.feature_importances_, index=Predictors)
feature_importances.nlargest(10).plot(kind='barh')
print('\n######## Model Validation and Accuracy Calculations ########")
# Creating the prediction value variables
TestingDataResults=pd.DataFrame(data=X test, columns=Predictors)
TestingDataResults[TargetVariable]=y_test
TestingDataResults[('Predicted'+TargetVariable)]=np.round(prediction)
# Printing a sample of the prediction value variables
print(TestingDataResults.head())
# Calculating the APE percentage
TestingDataResults['APE']=100 * ((abs(
 Testing Data Results ['Close'] - Testing Data Results ['Predicted Close']))/Testing Data Results ['Close'])
#Create the mean and median APE
MAPE=np.mean(TestingDataResults['APE'])
MedianMAPE=np.median(TestingDataResults['APE'])
#Using the MAPE to calculate the accuracy percentage of the model
Accuracy = 100 - MAPE
MedianAccuracy= 100 - MedianMAPE
print('Mean Accuracy on test data:', Accuracy) # This value results as negative if outliers are present.
print('Median Accuracy on test data:', MedianAccuracy)
# Making a custom function to calculate the accuracy score using the original and predicted.
# Make sure no zeros are present within.
def Accuracy_Score(orig,pred):
   MAPE = np.mean(100 * (np.abs(orig-pred)/orig))
   return(100-MAPE)
# A custom score calculation using the make_scorer from sklearn
from sklearn.metrics import make_scorer
custom_Scoring=make_scorer(Accuracy_Score, greater_is_better=True)
# Getting the cross validation value from sklearn
from sklearn.model_selection import cross_val_score
# Printing the 10-fold cross validation so to give further values to compare.
# Finally, pass the data of the x and y as K-fold splits the data and authomatically decides test or train.
Accuracy\_Values = cross\_val\_score(RegModel, \ X \ , \ y, \ cv=10, \ scoring = custom\_Scoring)
print('\nAccuracy values for 10-fold Cross Validation:\n',Accuracy_Values)
print('\nFinal Average Accuracy of the model:', round(Accuracy_Values.mean(),2))
```

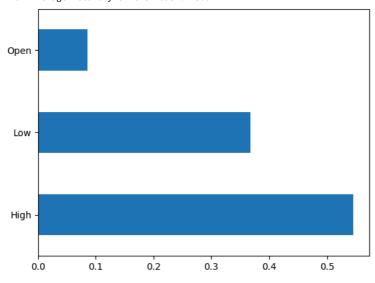
RandomForestRegressor(criterion='friedman\_mse', max\_depth=4, n\_estimators=400)
R2 Value: 0.9971395316156282

```
######## Model Validation and Accuracy Calculations #########
                       High
                                  Close PredictedClose
      0pen
                Low
0 0.577471 0.587917 0.576875 509.640015
                                                  500.0
1 0.592304 0.561601 0.572101 494.730011
                                                  496.0
2 0.596449 0.589962 0.585891 500.859985
                                                  502.0
3 0.330279 0.322451 0.303482 380.070007
                                                  384.0
4 0.195188 0.182474 0.166985 315.100006
                                                  322.0
Mean Accuracy on test data: 98.74477854666928
Median Accuracy on test data: 98.99831266406315
```

Accuracy values for 10-fold Cross Validation: [98.4554302 98.42310631 98.37079947 98.82139029 98.60895763 97.43986628

Final Average Accuracy of the model: 98.01

97.72854863 98.56576993 99.02174248 94.64640591]



```
# Plotting the single Decision Tree from our Random Forest
```

# Load libraries used

 $\label{from:indep} \mbox{from IPython.display import Image}$ 

from sklearn import tree

import pydotplus

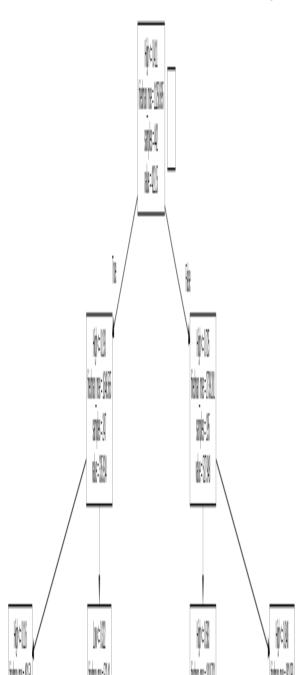
# Create DOT (graph description language) data for the our 6th Decision Tree in Random Forest dot\_data = tree.export\_graphviz(RegModel.estimators\_[5] , out\_file=None, feature\_names=Predictors, class\_names=TargetVariable)

# load the graph

graph = pydotplus.graph\_from\_dot\_data(dot\_data)

 $\mbox{\tt\#}$  Create the png of the graph

Image(graph.create\_png(), width=2000,height=2000)



Step 21: AdaBoost Algorithm

```
# Adaboost (Boost of multiple decision trees)
from sklearn.ensemble import AdaBoostRegressor
from sklearn.tree import DecisionTreeRegressor
# The 6 level decision tree is chosen as the "weak learner"
DTR=DecisionTreeRegressor(max depth=3)
RegModel = AdaBoostRegressor(n_estimators=100, base_estimator=DTR ,learning_rate=0.04)
# Printing the parameters of our Adaboost
print(RegModel)
# Creating the model using our training data
AB=RegModel.fit(X_train,y_train)
prediction=AB.predict(X_test)
from sklearn import metrics
# Measuring the quality of fit with training data.
print('R2 Value:',metrics.r2 score(y train, AB.predict(X train)))
# Using matplotlib to plot the 10 most highest importance columns.
%matplotlib inline
feature_importances = pd.Series(AB.feature_importances_, index=Predictors)
feature_importances.nlargest(10).plot(kind='barh')
print('\n####### Model Validation and Accuracy Calculations ########")
# Creating the variable to store our predicted values.
TestingDataResults=pd.DataFrame(data=X_test, columns=Predictors)
TestingDataResults[TargetVariable]=y_test
TestingDataResults[('Predicted'+TargetVariable)]=np.round(prediction)
# Printing a sample of the prediction values.
print(TestingDataResults.head())
# Calculation to calculate the error for each row.
TestingDataResults['APE']=100 * ((abs(
  TestingDataResults['Close']-TestingDataResults['PredictedClose']))/TestingDataResults['Close'])
MAPE=np.mean(TestingDataResults['APE'])
MedianMAPE=np.median(TestingDataResults['APE'])
Accuracy = 100 - MAPE
MedianAccuracy= 100- MedianMAPE
print('Mean Accuracy on test data:', Accuracy) # Negative value indicative of outlier in some cases
print('Median Accuracy on test data:', MedianAccuracy)
# Making a custom function to calculate the accuracy score using the original and predicted.
# Make sure no zeros are present within.
def Accuracy_Score(orig,pred):
    \texttt{MAPE = np.mean(100 * (np.abs(orig-pred)/orig))}
    return(100-MAPE)
# calculation using sklearn's make_scorer to make a custom scorer for accuracy.
from sklearn.metrics import make_scorer
custom_Scoring=make_scorer(Accuracy_Score, greater_is_better=True)
# Importing cross validation function from sklearn
from sklearn.model_selection import cross_val_score
# Running a 10 fold cross validation of the data.
# Passing full data of X and y because the K-fold will split the data into test or train for us.
Accuracy_Values=cross_val_score(RegModel, X , y, cv=10, scoring=custom_Scoring)
print('\nAccuracy values for 10-fold Cross Validation:\n',Accuracy Values)
print('\nFinal Average Accuracy of the model:', round(Accuracy_Values.mean(),2))
```

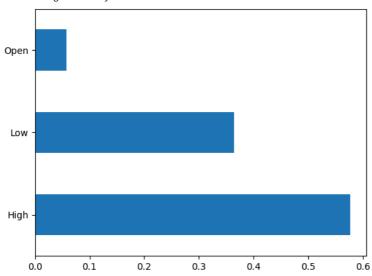
```
AdaBoostRegressor(base_estimator=DecisionTreeRegressor(max_depth=3), learning_rate=0.04, n_estimators=100)
```

R2 Value: 0.9894519230021294

```
######## Model Validation and Accuracy Calculations #########
                         High
                Low
                                    Close PredictedClose
      0pen
0 0.577471 0.587917 0.576875 509.640015
  0.592304 0.561601 0.572101
                               494.730011
                                                    493.0
                               500.859985
                                                   505.0
2 0.596449 0.589962 0.585891
3 0.330279 0.322451 0.303482
                               380.070007
                                                    388.0
4 0.195188 0.182474 0.166985
                               315.100006
                                                   317.0
Mean Accuracy on test data: 97.74835457799438
Median Accuracy on test data: 98.04831337136687
```

Accuracy values for 10-fold Cross Validation:
[97.26972335 97.39495947 96.75205006 97.17342265 96.75582601 94.63849664 98.54092564 97.09591207 98.16570955 93.57097683]

Final Average Accuracy of the model: 96.74



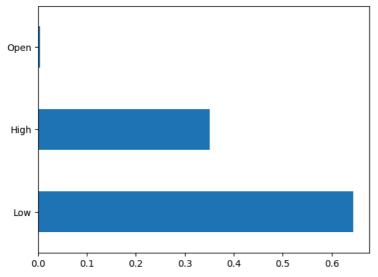
### XGBoost Regressor

```
# Xtreme Gradient Boosting (XGBoost)
from xgboost import XGBRegressor
RegModel=XGBRegressor(max_depth=2,
                      learning_rate=0.1,
                      n estimators=1000,
                      objective='reg:linear',
                      booster='gbtree')
# Print all XGBoost parameters
print(RegModel)
# Creating the model using out training data.
XGB=RegModel.fit(X_train,y_train)
prediction=XGB.predict(X_test)
from sklearn import metrics
# Measuring how well our training data works within the model.
print('R2\ Value:',metrics.r2\_score(y\_train,\ XGB.predict(X\_train)))
# Plotting the ten columns with the highest feature importance value.
%matplotlib inline
feature_importances = pd.Series(XGB.feature_importances_, index=Predictors)
feature_importances.nlargest(10).plot(kind='barh')
print('\n######## Model Validation and Accuracy Calculations ########")
# Creating a variable to store the values of prediction
TestingDataResults=pd.DataFrame(data=X_test, columns=Predictors)
TestingDataResults[TargetVariable]=y_test
TestingDataResults[('Predicted'+TargetVariable)]=np.round(prediction)
# Printing a sample of the prediction values
print(TestingDataResults.head())
# Making a calulation of the error.
TestingDataResults['APE']=100 * ((abs(
  TestingDataResults['Close']-TestingDataResults['PredictedClose']))/TestingDataResults['Close'])
```

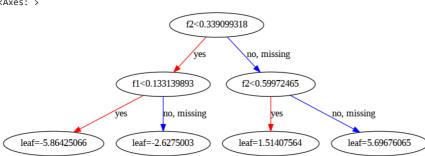
```
# Calculating the mean and median APE using the previous result.
MAPE=np.mean(TestingDataResults['APE'])
MedianMAPE=np.median(TestingDataResults['APE'])
#Calculating the accuracy as a percentage using the MAPE
Accuracy =100 - MAPE
MedianAccuracy=100- MedianMAPE
print('Mean Accuracy on test data:', Accuracy) # This can be negative if an outlier is present.
print('Median Accuracy on test data:', MedianAccuracy)
# Defining a custom function to calculate accuracy
# Make sure there are no zeros in the Target variable if you are using MAPE
def Accuracy_Score(orig,pred):
    MAPE = np.mean(100 * (np.abs(orig-pred)/orig))
    return(100-MAPE)
# Custom Scoring MAPE calculation
from sklearn.metrics import make scorer
custom_Scoring=make_scorer(Accuracy_Score, greater_is_better=True)
# Importing cross validation function from sklearn
from sklearn.model_selection import cross_val_score
# Running 10-Fold Cross validation on a given algorithm
# Passing full data X and y because the K-fold will split the data and automatically choose train/test
Accuracy_Values=cross_val_score(RegModel, X , y, cv=10, scoring=custom_Scoring)
print('\nAccuracy values for 10-fold Cross Validation:\n',Accuracy_Values)
print('\nFinal Average Accuracy of the model:', round(Accuracy_Values.mean(),2))
     XGBRegressor(base_score=None, booster='gbtree', callbacks=None,
                  colsample_bylevel=None, colsample_bynode=None,
                  colsample_bytree=None, device=None, early_stopping_rounds=None,
                  enable_categorical=False, eval_metric=None, feature_types=None,
                  gamma=None, grow_policy=None, importance_type=None,
                  interaction_constraints=None, learning_rate=0.1, max_bin=None,
                  max_cat_threshold=None, max_cat_to_onehot=None,
                  max_delta_step=None, max_depth=2, max_leaves=None,
                  \verb|min_child_weight=None, missing=nan, monotone_constraints=None, \\
                  multi_strategy=None, n_estimators=1000, n_jobs=None,
                  num_parallel_tree=None, objective='reg:linear', ...)
     R2 Value: 0.9995434001336165
     ######## Model Validation and Accuracy Calculations #########
                               High
                                           Close PredictedClose
                      Low
           0pen
       0.577471 0.587917 0.576875 509.640015
                                                           508.0
     1 0.592304 0.561601 0.572101 494.730011
                                                           493.0
     2 0.596449 0.589962 0.585891 500.859985
                                                           510.0
     3 0.330279 0.322451 0.303482 380.070007
                                                           386.0
     4 0.195188 0.182474 0.166985 315.100006
                                                           314.0
     Mean Accuracy on test data: 99.07659283944815
     Median Accuracy on test data: 99.34123847167325
     Accuracy values for 10-fold Cross Validation:
```

[98.92753273 98.92258095 98.9963492 99.31086522 99.14092335 98.04559908 98.86405587 99.12300896 99.47586872 95.58969741]

Final Average Accuracy of the model: 98.64



#Plotting a single Decision tree out of XGBoost from xgboost import plot tree import matplotlib.pyplot as plt



## K-Nearest Neighbour

```
#kNN
# K-Nearest Neighbor(KNN)
from sklearn.neighbors import KNeighborsRegressor
RegModel = KNeighborsRegressor(n_neighbors=3)
# Printing all the parameters of KNN
print(RegModel)
# Creating the model on Training Data
KNN=RegModel.fit(X_train,y_train)
prediction=KNN.predict(X_test)
from sklearn import metrics
# Measuring Goodness of fit in Training data
print('R2 Value:',metrics.r2_score(y_train, KNN.predict(X_train)))
# Plotting the feature importance for Top 10 most important columns
# The variable importance chart is not available for KNN
print('######## Model Validation and Accuracy Calculations ########")
# Printing some sample values of prediction
TestingDataResults=pd.DataFrame(data=X_test, columns=Predictors)
TestingDataResults[TargetVariable]=y test
TestingDataResults[('Predicted'+TargetVariable)]=np.round(prediction)
# Printing sample prediction values
print(TestingDataResults.head())
# Calculating the error for each row
TestingDataResults['APE']=100 * ((abs(
 Testing Data Results ['Close'] - Testing Data Results ['Predicted Close']))/Testing Data Results ['Close'])
MAPE=np.mean(TestingDataResults['APE'])
MedianMAPE=np.median(TestingDataResults['APE'])
Accuracy =100 - MAPE
MedianAccuracy=100- MedianMAPE
print('Mean Accuracy on test data:', Accuracy) # Can be negative sometimes due to outlier
print('Median Accuracy on test data:', MedianAccuracy)
# Defining a custom function to calculate accuracy
# Make sure there are no zeros in the Target variable if you are using MAPE
def Accuracy_Score(orig,pred):
    MAPE = np.mean(100 * (np.abs(orig-pred)/orig))
    #print('#'*70,'Accuracy:', 100-MAPE)
    return(100-MAPE)
# Custom Scoring MAPE calculation
from sklearn.metrics import make_scorer
custom_Scoring=make_scorer(Accuracy_Score, greater_is_better=True)
# Importing cross validation function from sklearn
from sklearn.model_selection import cross_val_score
# Running 10-Fold Cross validation on a given algorithm
# Passing full data X and y because the K-fold will split the data and automatically choose train/test
\label{lem:condition} Accuracy\_Values=cross\_val\_score(RegModel,\ X\ ,\ y,\ cv=10,\ scoring=custom\_Scoring)
print('\nAccuracy values for 10-fold Cross Validation:\n',Accuracy_Values)
print('\nFinal Average Accuracy of the model:', round(Accuracy_Values.mean(),2))
     KNeighborsRegressor(n_neighbors=3)
     R2 Value: 0.9989790942951331
     ######## Model Validation and Accuracy Calculations #########
                      Low
            0pen
                               High
                                           Close PredictedClose
     0 0.577471 0.587917 0.576875 509.640015
     1 0.592304 0.561601 0.572101 494.730011
                                                           488.0
     2 0.596449 0.589962 0.585891 500.859985
                                                           507.0
     3 0.330279 0.322451 0.303482 380.070007
                                                           382.0
     4 0.195188 0.182474 0.166985 315.100006
                                                           321.0
     Mean Accuracy on test data: 99.05814068358532
     Median Accuracy on test data: 99.31442221288626
     Accuracy values for 10-fold Cross Validation:
      [98.95644363 99.0577121 98.9613391 99.27343742 99.14276876 98.36592285
      98.83590812 99.26087998 99.45536694 95.95525821]
     Final Average Accuracy of the model: 98.73
```

## Support Vector Machine (SVM) Regressor