Features Selection Based on Uni-Variante ROC\_AUC for Classification and MSE for Regression. What is the ROC\_AUC? AUC :- Area Under the Curve. AUROC: - Area Under the Receiver Operating Characteristic curve. MSE: Mean square error ROC is a probability curve and AUC represents degree or measure of separability. Higher the AUC, better the model is at predicting 0s as 0s and 1s as 1s. By analogy, Higher the AUC, better the model is at distinguishing between patients with disease and no disease. An excellent model has AUC near to the 1 which means it has good measure of separability. A poor model has AUC near to the 0 which means it has worst measure of separability. In fact it means it is reciprocating the result. It is predicting 0s as 1s and 1s as 0s. And when AUC is 0.5, it means model has no class separation capacity whatsoever. • The receiver operating characteristics(ROC) curve is well known in evaluating classification performance. Owing to its superiority in dealing with imbalaced and cost-sensitive data, the ROC curve has been exploited as a popular metric to evaluate the ML models. ROC\_CURVE is drawn against the **sensitivity and 1-specificity** area under ROC curve shows the AUC which is a metric. Exsisting ROC based on feature selection apporoches are simple and effecive in individual feature selection. This is uni\_variate feature selection technique. That means while selecting the one feature it doesnt found out its effect any on the current features. That means we find the features important only based on AUC and ROC curve charectestics or metrics on UC metric. This is been widely used to determine the classification accuracy in supervised learning. • This kind of the ROC\_AUC classification can be used for the binary although it used for many classification (i.e.more than the binary) but this will be the difficult to use. Use of ROC\_AUC in Classification Problem. import numpy as np import matplotlib.pyplot as plt import pandas as pd import seaborn as sns %matplotlib inline from sklearn.feature selection import VarianceThreshold from sklearn.ensemble import RandomForestClassifier from sklearn.preprocessing import LabelEncoder from sklearn.model selection import train test split from sklearn.metrics import accuracy score,roc auc score data = pd.read csv('titanic.csv') data.head() survived pclass age sibsp parch fare embarked class who adult\_male deck embark\_town alive alone sex 0 22.0 7.2500 S Third NaN male 0 man True Southampton False no 1 38.0 71.2833 C 1 female First woman **False** Cherbourg False yes 2 1 3 female 26.0 0 7.9250 S Third woman **False** NaN Southampton True yes 3 35.0 53.1000 False C female First woman Southampton False yes 4 0 3 male 35.0 8.0500 S Third True NaN Southampton True man no In [9]: data.dropna(inplace=True) label = LabelEncoder() data['sex'] = label.fit transform(data['sex']) data['embarked'] = label.fit transform(data['embarked']) data['class'] = label.fit\_transform(data['embarked']) data['who']=label.fit transform(data['who']) data['adult\_male']=label.fit\_transform(data['adult\_male']) data['embark town']=label.fit transform(data['embark town']) data['alive']=label.fit transform(data['alive']) data['alone'] = label.fit transform(data['alone']) del data['deck'] x = data.drop(labels='alive',axis=1) y= data['alive'] x.shape, y.shape Out[16]: ((182, 13), (182,)) In [17]: x train, x test, y train, y test = train test split(x, y, train size = 0.8, random state =0) x train.shape,x test.shape Out[17]: ((145, 13), (37, 13)) In [ ]: #remove the constant and quasi constant In [18]: constant\_filter = VarianceThreshold(threshold=0.01) constant\_filter.fit(x\_train) constant filter.fit(x\_test) x\_train\_fiter = constant\_filter.transform(x\_train) x\_test\_filter = constant\_filter.transform(x\_test) x\_train\_fiter.shape,x\_test\_filter.shape Out[18]: ((145, 13), (37, 13)) In [ ]: #we want to remove duplicate data In [19]:  $x_{train_T} = x_{train_T}$ x test T = x test filter.T x train T = pd.DataFrame(x train T) x test T = pd.DataFrame(x test T) duplicate feature = x train T.duplicated() keep them = [not index for index in duplicate feature] x train unique = x train T[keep them].T x\_test\_unique = x\_test\_T[keep\_them].T x\_train\_unique.shape,x\_test\_unique.shape Out[20]: ((145, 11), (37, 11)) Now Calculate ROC\_AUC Score In [21]: #we need to create first upon the empty list like auc roc #It will read the one by one all the features from the columns of the X train unique.columns for feature in x train unique.columns: clf = RandomForestClassifier(n estimators=100, random state=0) #select the feature list from the x train unique and make it in the form of one diamensional array clf.fit(x train unique[feature].to frame(),y train) y pred = clf.predict(x test unique[feature].to frame()) #so, now we are appending value of roc auc in the empty list roc auc.append(roc auc score(y test, y pred)) In [22]: print(roc\_auc) [1.0, 0.5, 0.8888888888888888, 0.5518518518518519, 0.5, 0.5, 0.5259259259259259, 0.48148148148148145, 0.9074074 074074074, 0.9074074074074074, 0.51 In [24]: | #to put this array of roc\_auc in the pandas Dataframe Series roc values = pd.Series(roc auc) roc values.index = x train unique.columns roc values.sort values(ascending= False,inplace = True) • If we check then we find the series which is in sorted manner and also we will get the few values which are more than the 0.5 and these values are the most important than the other becasue if we use the value which having magnitude equal to or lesser than the 0.5 that time It will give us random prediction and not accurate so to get the better result/prediction we have to choose more than the o.5 so that the predict should not go wrong or random. • We are making here binary roc\_auc technique so that those who are giving the 0.5 probablity will not be counted because they will not previde the imformation which is required to accurate prediction. roc values 1.000000 0 0.907407 10 9 0.907407 2 0.888889 3 0.551852 6 0.525926 12 0.500000 5 0.500000 4 0.500000 0.500000 0.481481 dtype: float64 Those are the values are equal to the 0.5 and less than 0.5 are not required to classifiction. So for the we need to remove that kind of values. roc values.plot.bar(figsize = (15,6)) <AxesSubplot:> 1.0 0.8 0.6 0.4 0.2 0.0 10 sel= roc values[roc values>0.5] 1.000000 0 10 0.907407 9 0.907407 2 0.888889 0.551852 0.525926 dtype: float64 len(sel) Out[28]: 6 In [29]: x\_train\_roc = x\_train unique[sel.index] x test roc = x test unique[sel.index] Build the Model and compare the performance def run\_randomforest(x\_train, x\_test, y\_train, y\_test): clf = RandomForestClassifier(random state=0,n jobs=-1,n estimators=1000) clf.fit(x\_train,y\_train) y\_pred = clf.predict(x\_test) print('Accuracy : ',accuracy\_score(y\_test,y\_pred)) run randomforest(x train roc, x test roc, y train, y test) Accuracy: 1.0 Wall time: 2.01 s In [34]: x train roc.shape Out[34]: (145, 6) • If we compare with the previous x\_train,y\_train then see. • From that we can only say that there is no kind of impact on the accuracy but the time required to compute the things which is taken more than the x\_train\_roc,x\_test\_roc. run randomforest(x train, x test, y train, y test) Accuracy: 1.0 Wall time: 1.9 s • In the filtering method or any feature selection does not give us garantee of accuracy but it can give us time to compute the tthings which we want to calculate like accuracy. • Diffrenet algorithm can give us diffrent output. If we here that first randomfoorest with the data x\_train, x\_test with the roc and then x\_train,x\_test with the no any substitution. The first one does not gievn the accuracy as much given by the roc data but it given the less computational time voice versa. According our requirement we choose the either first one or last one that having better accuracy but less high computational time. Feature Selection using RMSE in Regression. RMSE = Root mean square error from sklearn import datasets from sklearn.linear model import LinearRegression from sklearn.metrics import mean\_absolute\_error,mean\_squared\_error,r2\_score boston = datasets.load boston() In [37]: print(boston.DESCR) .. boston dataset: Boston house prices dataset \*\*Data Set Characteristics:\*\* :Number of Instances: 506 :Number of Attributes: 13 numeric/categorical predictive. Median Value (attribute 14) is usually the targe t. :Attribute Information (in order): - CRIM per capita crime rate by town - ZN proportion of residential land zoned for lots over 25,000 sq.ft. proportion of non-retail business acres per town - INDUS - CHAS Charles River dummy variable (= 1 if tract bounds river; 0 otherwise) nitric oxides concentration (parts per 10 million) - NOX average number of rooms per dwelling proportion of owner-occupied units built prior to 1940 - AGE - DIS weighted distances to five Boston employment centres index of accessibility to radial highways - RAD - TAX full-value property-tax rate per \$10,000 - PTRATIO  $% \left( 1\right) =\left( 1\right) +\left( 1\right) +$ 1000(Bk - 0.63)^2 where Bk is the proportion of black people by town - LSTAT % lower status of the population - MEDV Median value of owner-occupied homes in \$1000's :Missing Attribute Values: None :Creator: Harrison, D. and Rubinfeld, D.L. This is a copy of UCI ML housing dataset. https://archive.ics.uci.edu/ml/machine-learning-databases/housing/ This dataset was taken from the StatLib library which is maintained at Carnegie Mellon University. The Boston house-price data of Harrison, D. and Rubinfeld, D.L. 'Hedonic prices and the demand for clean air', J. Environ. Economics & Management, vol.5, 81-102, 1978. Used in Belsley, Kuh & Welsch, 'Regression diagnostics ...', Wiley, 1980. N.B. Various transformations are used in the table on pages 244-261 of the latter. The Boston house-price data has been used in many machine learning papers that address regression problems. .. topic:: References - Belsley, Kuh & Welsch, 'Regression diagnostics: Identifying Influential Data and Sources of Collinearity', Wiley, 1980. 244-261. - Quinlan, R. (1993). Combining Instance-Based and Model-Based Learning. In Proceedings on the Tenth Internat ional Conference of Machine Learning, 236-243, University of Massachusetts, Amherst. Morgan Kaufmann. In [38]: x = pd.DataFrame(data=boston.data,columns=boston.feature names) x.head() ZN INDUS CHAS NOX RM AGE DIS RAD TAX PTRATIO **B** LSTAT **0** 0.00632 18.0 0.0 0.538 6.575 65.2 4.0900 1.0 296.0 15.3 396.90 4.98 2.31 17.8 396.90 **1** 0.02731 0.0 7.07 0.0 0.469 6.421 78.9 4.9671 2.0 242.0 9.14 0.0 0.469 7.185 61.1 4.9671 **2** 0.02729 0.0 7.07 2.0 242.0 17.8 392.83 4.03 **3** 0.03237 0.0 2.18 0.0 0.458 6.998 45.8 6.0622 3.0 222.0 18.7 394.63 2.94 3.0 222.0 **4** 0.06905 0.0 2.18 0.0 0.458 7.147 54.2 6.0622 18.7 396.90 5.33 In [39]: y = boston.targetIn [40]: x\_train,x\_test,y\_train,y\_test = train\_test\_split(x,y,train\_size= 0.8,random\_state =0) x\_train.shape,x\_test.shape Out[40]: ((404, 13), (102, 13)) In [41]: #its method we have follow for univariant feature selection. mse = []for feature in x\_train.columns: clf = LinearRegression() clf.fit(x\_train[feature].to\_frame(),y\_train) y\_pred = clf.predict(x\_test[feature].to\_frame()) mse.append(mean squared error(y test,y pred)) In [42]: print(mse)  $[76.38674157646072,\ 84.66034377707906,\ 77.02905244667242,\ 79.36120219345942,\ 76.95375968209432,\ 46.907351627395]$ 718, 46.33630536002592] In [43]: mse = pd.Series(mse) mse.index = x train.columns mse.sort\_values(ascending = False,inplace = True) Out[43]: ZN 84.660344 DIS 82.618741 82.465000 RAD 81.794971 PTRATIO AGE 80.391548 CHAS 79.361202 TAX 78.308314 В 77.752856 INDUS 77.029052 NOX 76.953760 CRIM 76.386742 RM 46.907352 LSTAT 46.336305 dtype: float64 In [44]: len(mse) 13 Out[44]: Higher the mse means more the error and lower the mse means the low error if we can that the some features having the low mse those features are very important than the others.so we taking the two features to get the accuracy of linear regression. In [45]: mse.plot.bar(figsize = (15,5))<AxesSubplot:> Out[45]: 80 70 60 50 40 30 20 10 ĪΑΧ NDUS CRIIM N DIS RAD AGE 쮼 õ In [46]: x train 2 = x train[['RM','LSTAT']]  $x_{test_2} = x_{test[['RM', 'LSTAT']]}$ %%time In [47]: model = LinearRegression() model.fit(x\_train\_2,y\_train) y\_pred = model.predict(x\_test\_2) print('r2 score :',r2\_score(y\_test,y\_pred)) print('rmse :',np.sqrt(mean\_squared\_error(y\_test,y\_pred))) print('sd of house price :',np.std(y)) r2 score : 0.5409084827186417 rmse : 6.114172522817781 sd of house price : 9.188011545278203 Wall time: 96.9 ms From the above we can understand that the root\_mean\_square\_error is lesser than the standard deviation by selecting the two features. So, we can say that the present model is not bad model. If we do the prediction by using the original dataset then lets see what happen. %%time model = LinearRegression() model.fit(x\_train,y\_train) y\_pred = model.predict(x\_test) print('r2 score :',r2\_score(y\_test,y\_pred)) print('rmse :',np.sqrt(mean\_squared\_error(y\_test,y\_pred))) print('sd of house price :',np.std(y)) r2 score : 0.5892223849182507 rmse : 5.783509315085136 sd of house price : 9.188011545278203 Wall time: 98.2 ms Morever we can say that the feature selection by using the model it will give us of gurantee that we will get the accuracy or less error but will reduce our computational time. From the above we had tried first with the feature those who having the lower mean squared error it was the bad model afterwards • we had taken the prediction on the original data in that we get the r2 score maximum than the previous but rmsq we lower than the previous one but computational time was quite high. So,we just say that we doesnt have gurantee to get the accuracy but it can be our suitability what we will prefered either time or accuracy.

Sometimes, It depends on the model which we are using and data type and requirement.