Project 2_PuppalaSucharitha

February 5, 2023

0.0.1 Term Project
0.0.2 Project - 2
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0.0.5 DSC680-T301 Applied Data Science (2233-1)
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0.0.7 Date: 01/25/2023

```
[1]: # Importing all the necessary libraries.
     import numpy as np
     import pandas as pd
     import seaborn as sns
     import matplotlib as mpl
     import matplotlib.pyplot as plt
     from scipy.stats import skew
     from sklearn.model_selection import train_test_split
     from sklearn.preprocessing import StandardScaler
     from sklearn.linear_model import LogisticRegression
     from sklearn.neighbors import KNeighborsClassifier
     from sklearn.svm import SVC
     from sklearn import metrics
     from sklearn.ensemble import RandomForestClassifier, GradientBoostingClassifier
     from sklearn.metrics import confusion_matrix
     from sklearn.metrics import precision_score, recall_score, f1_score
     from sklearn.preprocessing import StandardScaler
     from sklearn.metrics import roc auc score
     from sklearn.metrics import confusion_matrix, accuracy_score,_
     →classification_report
     %matplotlib inline
     import warnings
     warnings.filterwarnings('ignore')
```

```
[2]: # Loading the dataset named water_potability.csv
waterdf = pd.read_csv('water_potability.csv')
```

```
[3]: # Getting the first 5 rows of the dataset.
waterdf.head()
```

```
Sulfate Conductivity \
[3]:
             ph
                   Hardness
                                   Solids Chloramines
            NaN 204.890455 20791.318981
                                              7.300212 368.516441
    0
                                                                     564.308654
    1 3.716080 129.422921 18630.057858
                                              6.635246
                                                                     592.885359
                                                              {\tt NaN}
    2 8.099124 224.236259 19909.541732
                                             9.275884
                                                              {\tt NaN}
                                                                     418.606213
    3 8.316766 214.373394 22018.417441
                                             8.059332 356.886136
                                                                     363.266516
    4 9.092223 181.101509 17978.986339
                                              6.546600 310.135738
                                                                     398.410813
```

```
Organic_carbon Trihalomethanes Turbidity Potability
0
       10.379783
                                    2.963135
                        86.990970
1
       15.180013
                        56.329076 4.500656
                                                      0
2
       16.868637
                        66.420093
                                    3.055934
                                                      0
                       100.341674
3
       18.436524
                                    4.628771
                                                      0
                                                      0
4
       11.558279
                        31.997993 4.075075
```

[4]: # Getting the shape of the dataset waterdf.shape

[4]: (3276, 10)

[5]: # Getting the size of the dataset waterdf.size

[5]: 32760

[6]: # Information about the dataset variables.
waterdf.info()

<class 'pandas.core.frame.DataFrame'>
RangeIndex: 3276 entries, 0 to 3275
Data columns (total 10 columns):

#	Column	Non-Null Count	Dtype
0	ph	2785 non-null	float64
1	Hardness	3276 non-null	float64
2	Solids	3276 non-null	float64
3	Chloramines	3276 non-null	float64
4	Sulfate	2495 non-null	float64
5	Conductivity	3276 non-null	float64
6	Organic_carbon	3276 non-null	float64
7	Trihalomethanes	3114 non-null	float64
8	Turbidity	3276 non-null	float64
9	Potability	3276 non-null	int64

dtypes: float64(9), int64(1)

memory usage: 256.1 KB

```
[7]: # Checking for null values waterdf.isna().sum()
```

```
[7]: ph
                         491
     Hardness
                           0
     Solids
                           0
     Chloramines
                           0
     Sulfate
                         781
     Conductivity
                           0
     Organic_carbon
                           0
     Trihalomethanes
                         162
     Turbidity
                           0
     Potability
                           0
     dtype: int64
```

Observations:

- We can see many null values present in the dataset.
- From the above we can see that the features ph, Sulfate and Trihalomethanes are having more null values. Let us fill the null values with the meaan values.

```
[8]: # filling the null values with the column mean values.
waterdf['ph']=waterdf['ph'].fillna(waterdf.groupby(['Potability'])['ph'].

→transform('mean'))
waterdf['Sulfate']=waterdf['Sulfate'].fillna(waterdf.

→groupby(['Potability'])['Sulfate'].transform('mean'))
waterdf['Trihalomethanes']=waterdf['Trihalomethanes'].fillna(waterdf.

→groupby(['Potability'])['Trihalomethanes'].transform('mean'))
```

```
[9]: # Checking for null values again waterdf.isna().sum()
```

```
[9]: ph
                         0
    Hardness
                         0
     Solids
                         0
     Chloramines
     Sulfate
     Conductivity
     Organic_carbon
     Trihalomethanes
                         0
     Turbidity
                         0
                         0
     Potability
     dtype: int64
```

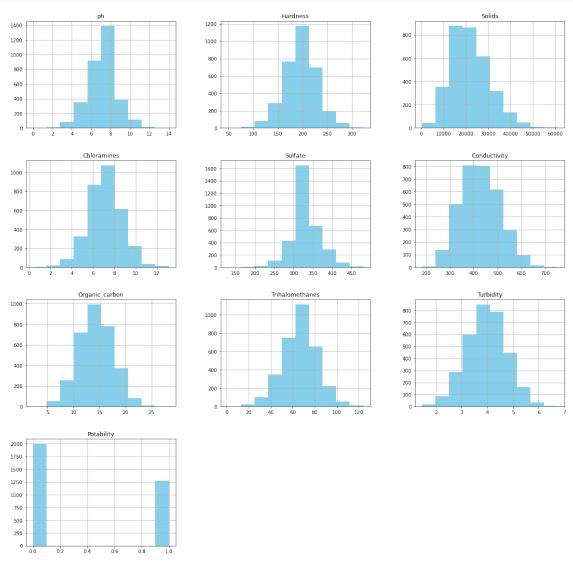
Observations: From the above we can see that the null values are being removed.

```
[10]: # Checking if any duplicated values present in the data set. waterdf.duplicated()
```

```
[10]: 0
              False
              False
      1
      2
              False
      3
              False
      4
              False
      3271
              False
      3272
              False
      3273
              False
      3274
              False
      3275
              False
      Length: 3276, dtype: bool
     Observations: Check for duplicate values is done and there are no duplicate values in the dataset.
[11]: # Getting the number of unique columns in the data set
      waterdf.nunique()
[11]: ph
                          2787
      Hardness
                          3276
      Solids
                          3276
      Chloramines
                          3276
      Sulfate
                          2497
      Conductivity
                          3276
      Organic_carbon
                          3276
      Trihalomethanes
                          3116
      Turbidity
                          3276
      Potability
                             2
      dtype: int64
[12]: # Initially lets's use the describe() to get an idea on the dataset.
      waterdf.describe()
[12]:
                              Hardness
                                               Solids
                                                        Chloramines
                                                                          Sulfate \
                       ph
                           3276.000000
                                          3276.000000
                                                        3276.000000
                                                                     3276.000000
      count
             3276.000000
      mean
                7.080855
                            196.369496
                                         22014.092526
                                                           7.122277
                                                                       333.785123
      std
                 1.469958
                             32.879761
                                          8768.570828
                                                           1.583085
                                                                        36.145701
                             47.432000
                                                                       129.000000
      min
                 0.000000
                                           320.942611
                                                           0.352000
      25%
                 6.277673
                            176.850538
                                         15666.690297
                                                           6.127421
                                                                       317.094638
      50%
                7.085378
                            196.967627
                                         20927.833607
                                                           7.130299
                                                                       334.564290
      75%
                7.870050
                            216.667456
                                         27332.762127
                                                           8.114887
                                                                       350.385756
               14.000000
                            323.124000
                                         61227.196008
                                                          13.127000
                                                                       481.030642
      max
                            Organic_carbon
             Conductivity
                                             Trihalomethanes
                                                                  Turbidity
                                                                              Potability
              3276.000000
                               3276.000000
                                                  3276.000000
                                                               3276.000000
                                                                             3276.000000
      count
      mean
               426.205111
                                  14.284970
                                                    66.395671
                                                                   3.966786
                                                                                0.390110
                 80.824064
                                   3.308162
                                                    15.769901
                                                                   0.780382
                                                                                0.487849
      std
```

```
2.200000
                                              0.738000
                                                            1.450000
                                                                         0.00000
min
         181.483754
25%
         365.734414
                           12.065801
                                             56.647656
                                                            3.439711
                                                                         0.00000
50%
         421.884968
                           14.218338
                                             66.303555
                                                            3.955028
                                                                         0.000000
75%
         481.792304
                           16.557652
                                             76.666609
                                                            4.500320
                                                                         1.000000
max
         753.342620
                           28.300000
                                            124.000000
                                                            6.739000
                                                                         1.000000
```

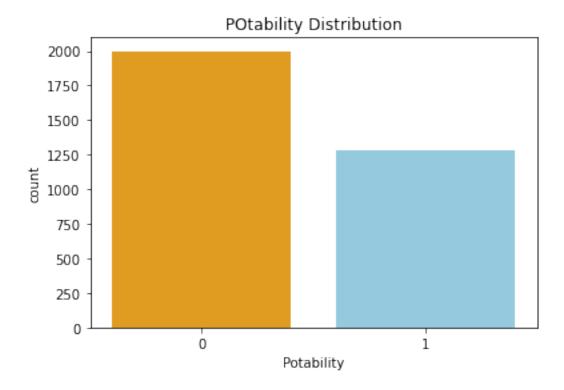
```
[13]: # Univariate analysis.
# plot histograms for each numerical variable
waterdf.hist(figsize = (20, 20),color='skyblue')
plt.show()
```



Observations: From the above histogram plot of all the variables we see that few features are slightly skewed.

```
[14]: #Let's cheeck for the distribution of the target variable Potability.
sns.countplot(data=waterdf,x='Potability' , palette=['orange','skyblue']);
#Title of the plot
plt.title('POtability Distribution')
```

[14]: Text(0.5, 1.0, 'POtability Distribution')

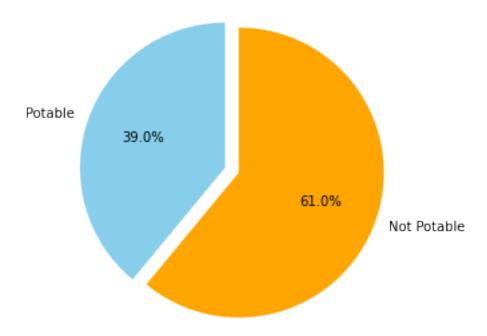


0.0.8 Observations:

From the above count plot we can see that the water that is potable i.e. safe for human consumption is less in count.

plt.title('Distribution of the target variable - Potability in %'); # $Title\ of_{\sqcup} \hookrightarrow the\ plot.$

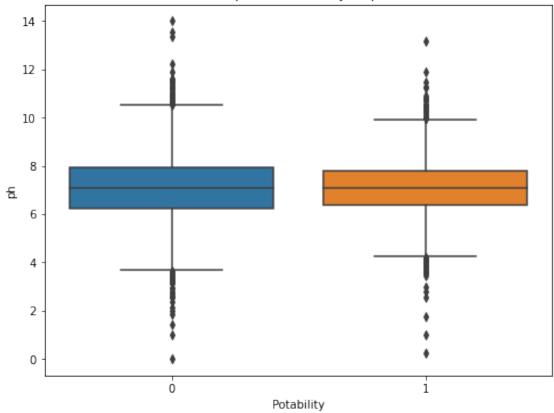
Distribution of the target variable - Potability in %



Observations: From the above plot we can say that about 39% of the water is potable i.e. safe for human consumption.

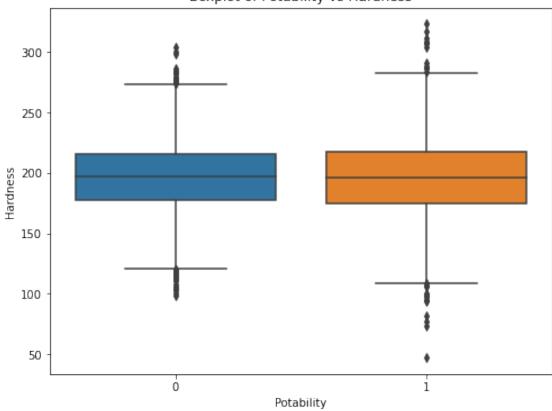
```
[16]: # bivariate analysis
# seaborn bar plot gives the variable average
# defining the plot size
plt.figure(figsize=(8,6))
sns.boxplot(x="Potability",y="ph",data=waterdf)
plt.title("Boxplot of Potability vs ph")
plt.show()
```





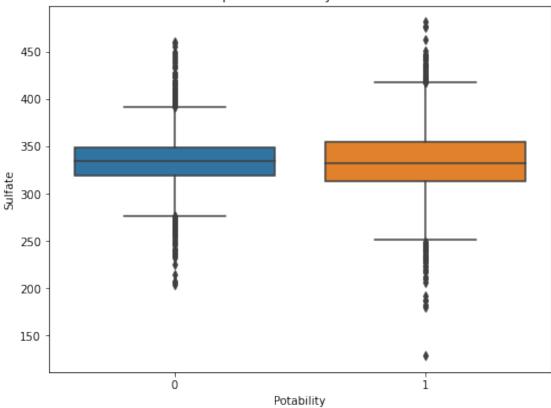
```
[17]: plt.figure(figsize=(8,6))
    sns.boxplot(x="Potability",y="Hardness",data=waterdf)
    plt.title("Boxplot of Potability vs Hardness")
    plt.show()
```

Boxplot of Potability vs Hardness

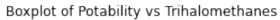


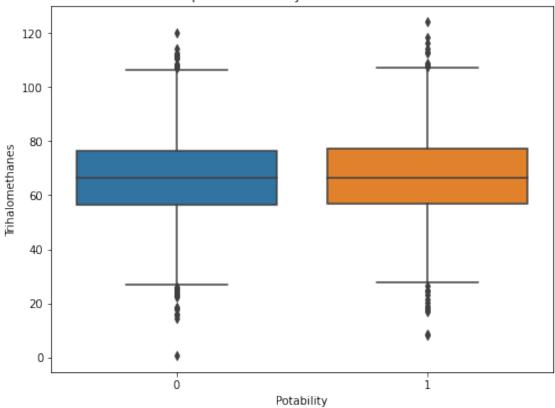
```
[18]: plt.figure(figsize=(8,6))
    sns.boxplot(x="Potability",y="Sulfate",data=waterdf)
    plt.title("Boxplot of Potability vs Sulfate")
    plt.show()
```

Boxplot of Potability vs Sulfate

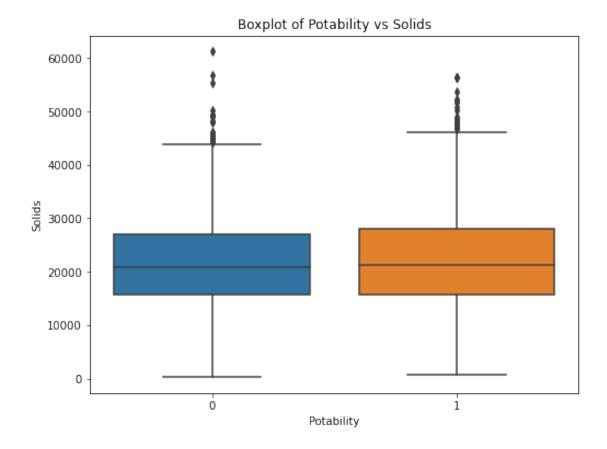


```
[19]: plt.figure(figsize=(8,6))
    sns.boxplot(x="Potability",y="Trihalomethanes",data=waterdf)
    plt.title("Boxplot of Potability vs Trihalomethanes")
    plt.show()
```



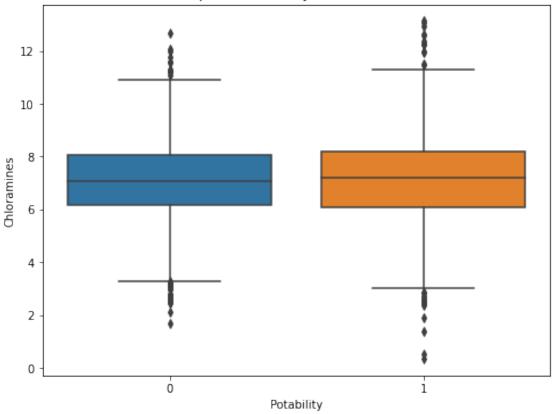


```
[20]: plt.figure(figsize=(8,6))
    sns.boxplot(x="Potability",y="Solids",data=waterdf)
    plt.title("Boxplot of Potability vs Solids")
    plt.show()
```

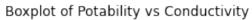


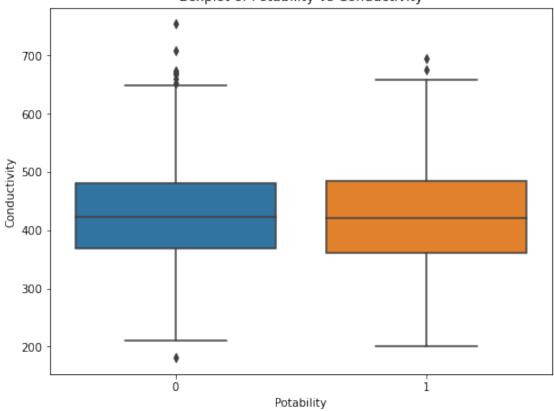
```
[21]: plt.figure(figsize=(8,6))
    sns.boxplot(x="Potability",y="Chloramines",data=waterdf)
    plt.title("Boxplot of Potability vs Chloramines")
    plt.show()
```

Boxplot of Potability vs Chloramines

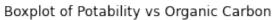


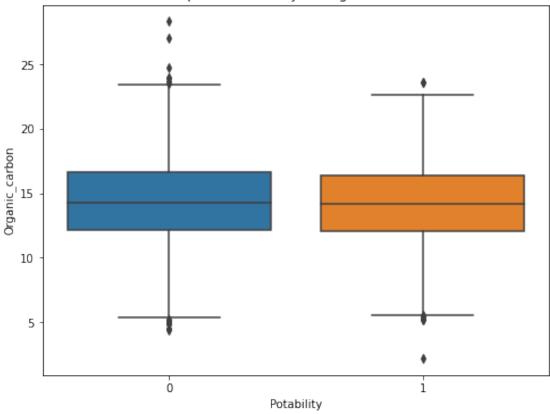
```
[22]: plt.figure(figsize=(8,6))
    sns.boxplot(x="Potability",y="Conductivity",data=waterdf)
    plt.title("Boxplot of Potability vs Conductivity")
    plt.show()
```



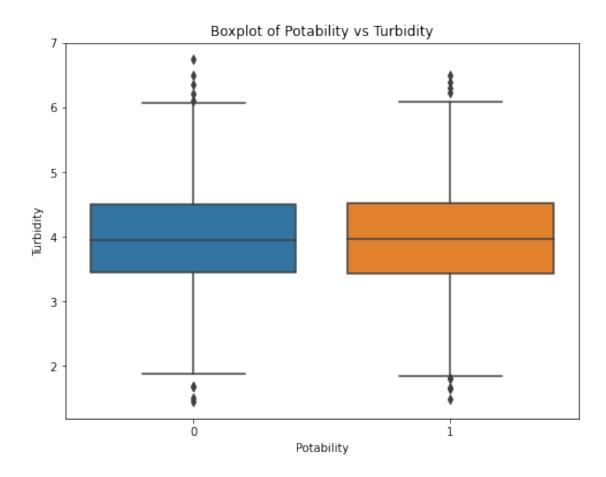


```
[23]: plt.figure(figsize=(8,6))
    sns.boxplot(x="Potability",y="Organic_carbon",data=waterdf)
    plt.title("Boxplot of Potability vs Organic Carbon")
    plt.show()
```





```
[24]: plt.figure(figsize=(8,6))
    sns.boxplot(x="Potability",y="Turbidity",data=waterdf)
    plt.title("Boxplot of Potability vs Turbidity")
    plt.show()
```

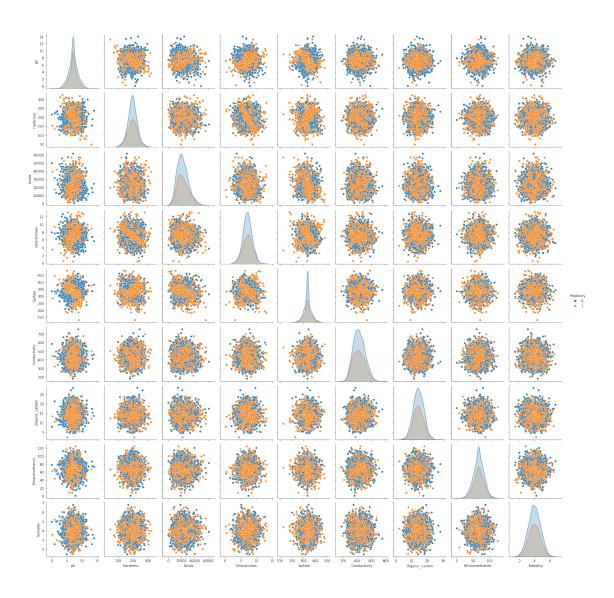


0.0.9 Observations:

From the above box plots for the distribution of the features of the dataset we see that the presence of all the features is more in the water that is not safe for drinking, i.e. Potability = 0.

```
[25]: # Let's use pairplot() function from seaborn to understand the relationship ⇒between all features.

sns.pairplot(waterdf,hue="Potability");
```



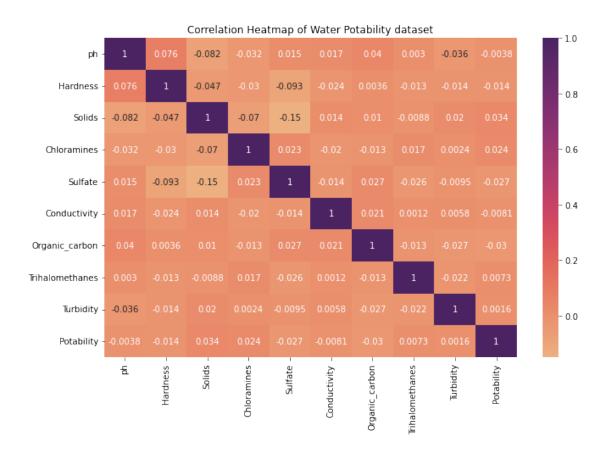
```
[26]: # Let's plot a correlation heatmap of the dataset.

# Calculate the correlation coefficient with corr().
corr_number = waterdf.corr()

# Create the heatmap for the correlation coefficients calculated above.
fig, ax = plt.subplots(1, 1, figsize=(10,7), tight_layout = True)
sns.heatmap(corr_number, annot = True, cmap = 'flare')

# Title of the plot
plt.title('Correlation Heatmap of Water Potability dataset')
```

[26]: Text(0.5, 1.0, 'Correlation Heatmap of Water Potability dataset')

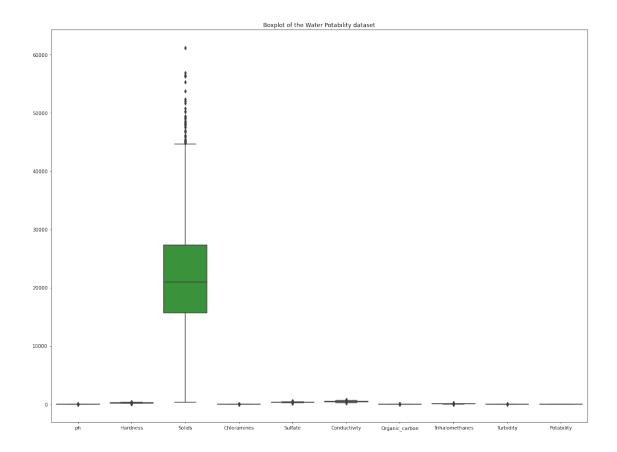


Observations:

- From the above correlation heat map we can see that there is no feature that is strongly correlated to the target variable "Potability".
- But from the above plot we can say that 'Hardness' has good correlation with the target variable 'Potability' but with very low correlation coefficient value.

```
[27]: # let's plot a boxplot for the diabetes dataset.
plt.figure(figsize=(20,15))
sns.boxplot(data = waterdf)
# Plot title
plt.title('Boxplot of the Water Potability dataset')
```

[27]: Text(0.5, 1.0, 'Boxplot of the Water Potability dataset')



Observations: From the above boxplot we can see that there are few outliers present in the 'Solids' feature.

```
[28]: #Let's define a function for detecting outliers using IQR method.

def outliers_IQR(df):
    q1 = df.quantile(0.25)
    q3 = df.quantile(0.75)

    IQR = q3-q1
    outlier = df[((df<(q1-1.5*IQR)) | (df>(q3+1.5*IQR)))]
    return outlier
```

```
[29]: # Now let's check for the outliers in features : Glucose.
outlier = outliers_IQR(waterdf['Solids'])
print('Total number of outliers in feature Solids : '+ str(len(outlier)))
```

```
print('Max. outlier value : '+ str(outlier.max()))
print('Min. outlier value : '+ str(outlier.min()))
```

Total number of outliers in feature Solids : 47

Max. outlier value : 61227.19600771213 Min. outlier value : 44868.45836802399

Observations:

- We can see that the outliers are 47 present in the Solids feature.
- There are no zeros present when observed the outliers, let us proceed with the analysis without changing the values of the outliers present.

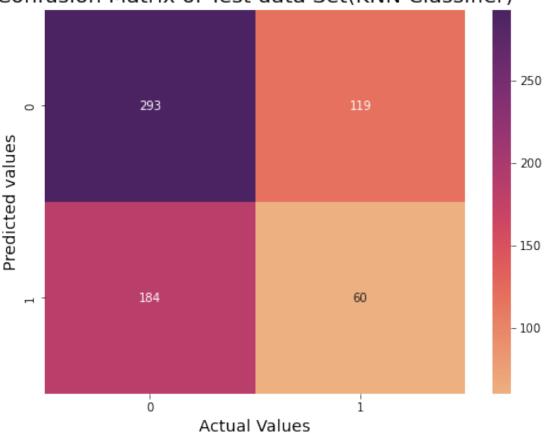
0.0.10 Modelling

```
[30]: # Let's define the features and target variables X and y respectively.
      X = waterdf.drop('Potability', axis = 1)
      y = waterdf['Potability']
[31]: # Let's split the dataset into 80% train and 20% test datasets using
      \rightarrow train\_test\_split().
      # Test size is 0.2
      X_train, X_test, y_train, y_test = train_test_split(X,y, test_size = 0.2, __
       \rightarrowrandom_state = 42)
[32]: # Let's get the size of the test and train datasets.
      print("X_train shape : {} rows and {} columns.".format(X_train.shape[0], X_train.
       \hookrightarrowshape [1]))
      print("y_train shape : {} rows.".format(y_train.shape[0]))
      print("X_test shape : {} rows and {} columns.".format(X_test.shape[0], X_test.
       \rightarrowshape[1]))
      print("y_test shape : {} rows.".format(y_test.shape[0]))
     X_train shape : 2620 rows and 9 columns.
     y_train shape : 2620 rows.
     X_test shape : 656 rows and 9 columns.
     y_test shape : 656 rows.
[33]: # Using the standard scaler on X_train and X_test datasets.
      # Fit the transform.
      standscaler = StandardScaler()
      X_train_std = standscaler.fit_transform(X_train)
      X_test_std = standscaler.transform(X_test)
```

0.0.11 KNN Classifier

```
[34]: # Create KNN classifier
      knnclass = KNeighborsClassifier()
      # Fit the model to train datasets.
      knnclass = knnclass.fit(X_train, y_train )
      # Create prediction of the model using the test data.
      knnclass_pred = knnclass.predict(X_test)
      # Create prediction of the model using the train data.
      knnclass_pred_train = knnclass.predict(X_train)
[35]: | # Let's create a Confusion Matrix for the test set predictions.
      knnconmatrix = confusion_matrix(y_test, knnclass_pred)
      # Print the Confusion matrix.
      print('Confusion Matrix of Test set predictions(KNN Classifier) : \n', \
       →knnconmatrix)
     Confusion Matrix of Test set predictions(KNN Classifier) :
      [[293 119]
      [184 60]]
[36]: # Plot the confusion matrix.
      # Define the size of the plot
      plt.figure(figsize=(8,6))
      # Confusion matrix heat map.
      sns.heatmap(knnconmatrix, annot=True,cmap = 'flare', fmt='d')
      # Plot Title
      plt.title('Confusion Matrix of Test data Set(KNN Classifier)', fontsize = 18)
      # x- Label
      plt.xlabel('Actual Values', fontsize = 14)
      # y-label
      plt.ylabel('Predicted values', fontsize = 14)
```





The accuracy score of the KNN Classifier on test dataset: 0.538109756097561
The accuracy score of the KNN Classifier on train dataset: 0.7145038167938931
Precision Score of the test set for the KNN Classifier: 0.335
Recall Score of the test set for the KNN Classifier: 0.246
F1 Score of the test set for the KNN Classifier: 0.284

0.0.12 Random Forest Classifier

```
[38]: # Let's create Random Forest Model.
rdreg = RandomForestClassifier()

# Fit the Random Forest model to training datasets.
rdreg.fit(X_train, y_train)

# Creating the predictions on the test data for model validation.
rdreg_pred = rdreg.predict(X_test)

# Creating the predictions on the trianed data for model validation.
rdreg_pred_train = rdreg.predict(X_train)
```

```
[39]: # Let's create a Confusion Matrix for the test set predictions.

rdconmatrix = confusion_matrix(y_test, rdreg_pred)

# Print the Confusion matrix.

print('Confusion Matrix of Test set predictions(Support Vector Machine

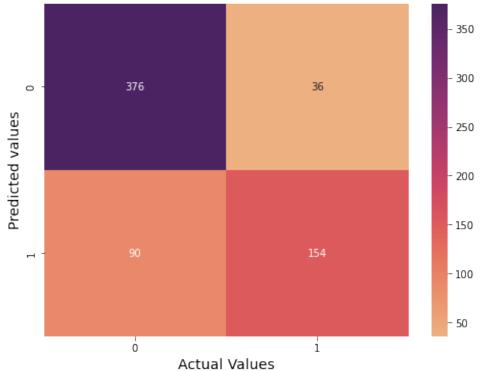
→Classifier): \n', rdconmatrix)
```

Confusion Matrix of Test set predictions(Support Vector Machine Classifier):
[[376 36]
[90 154]]

```
[40]: # Plot the confusion matrix.
    # Define the size of the plot
    plt.figure(figsize=(8,6))
    # Confusion matrix heat map.
    sns.heatmap(rdconmatrix, annot=True,cmap = 'flare', fmt='d')
```

[40]: Text(51.0, 0.5, 'Predicted values')

Confusion Matrix of Test data Set(Random Forest Classifier)



```
[41]: # Let's get the accuracy score, Precision, Recall and F1 Score of the Support

→ Vector Machine Classifier.

# Getting the accuracy score of the test dataset.

rdreg_accuracy = metrics.accuracy_score(y_test, rdreg_pred)

# Getting the accuracy score of the train dataset

rdreg_accu_train = metrics.accuracy_score(y_train, rdreg_pred_train)

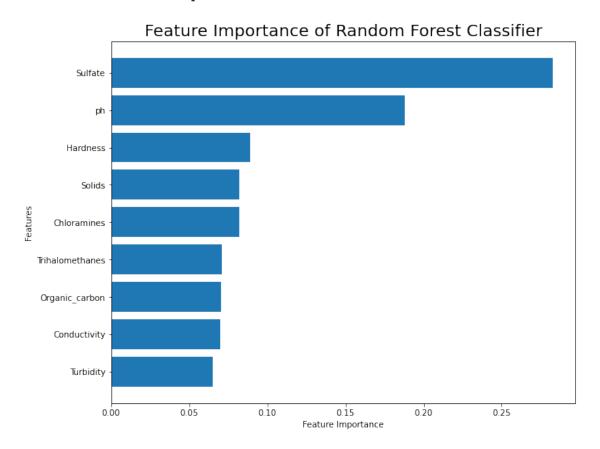
# Getting the precision score.

rdreg_precision = round(precision_score(y_test, rdreg_pred),3)
```

```
# Getting the Recall Score.
      rdreg_recall = round(recall_score(y_test, rdreg_pred),3)
      # Getting the F1 Score.
      rdreg_f1score = round(f1_score(y_test, rdreg_pred),3)
      # Printing the accuracy of the model.
      print('The accuracy score of the Random Forest Classifier on test dataset: {} '.
       →format(rdreg_accuracy))
      print('The accuracy score of the Random Forest Classifier on train dataset : {}_\|
       →'.format(rdreg_accu_train))
      print('Precision Score of the test set for the Random Forest Classifier : {}'.
      →format(rdreg_precision))
      print('Recall Score of the test set for the Random Forest Classifier : {}'.
       →format(rdreg recall))
      print('F1 Score of the test set for the Random Forest Classifier : {}'.
       →format(rdreg_f1score))
     The accuracy score of the Random Forest Classifier on test dataset:
     0.8079268292682927
     The accuracy score of the Random Forest Classifier on train dataset : 1.0
     Precision Score of the test set for the Random Forest Classifier: 0.811
     Recall Score of the test set for the Random Forest Classifier: 0.631
     F1 Score of the test set for the Random Forest Classifier: 0.71
[42]: # The importance of a feature is basically: how much this feature is used in
       →each tree of the forest.
      importantfeatures = rdreg.feature_importances_
      {\tt important features}
[42]: array([0.18781212, 0.08895214, 0.08228198, 0.08199801, 0.28297704,
             0.06966496, 0.07022538, 0.07091603, 0.06517234])
[43]: # initialize the idicies
      indices = np.argsort(importantfeatures)
      # Defining the size of the plot
      fig, ax = plt.subplots(figsize = (10,8))
      # Plotting the graph
      ax.barh(range(len(importantfeatures)), importantfeatures[indices])
      # Y-ticks
      ax.set_yticks(range(len(importantfeatures)))
      = ax.set_yticklabels(np.array(X_train.columns)[indices])
      # X-label
      plt.xlabel("Feature Importance", fontsize = 10)
      #Y - label
      plt.ylabel("Features", fontsize = 10)
```

```
# Title of the plot
plt.title("Feature Importance of Random Forest Classifier", fontsize = 20)
```

[43]: Text(0.5, 1.0, 'Feature Importance of Random Forest Classifier')



0.0.13 XGB Classifier

[44]: #from xgboost import XGBClassifier

```
from xgboost import XGBClassifier

[45]: # Let's create XGB classifier.
    xg = XGBClassifier()

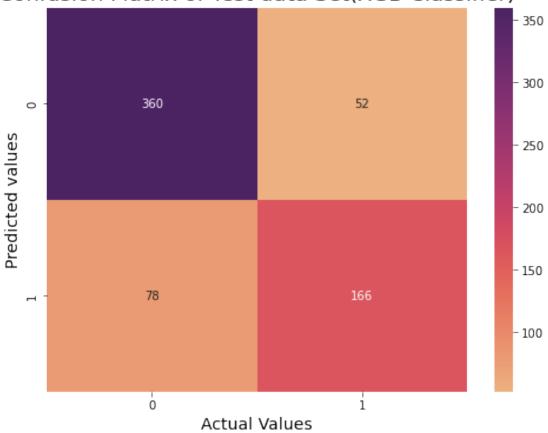
# Fit the Random Forest model to training datasets.
    xg.fit(X_train, y_train)

# Creating the predictions on the test data for model validation.
    xg_pred = xg.predict(X_test)

# Creating the predictions on the trianed data for model validation.
    xg_pred_train = xg.predict(X_train)
```

```
[46]: # Let's create a Confusion Matrix for the test set predictions.
      xgconmatrix = confusion_matrix(y_test, xg_pred)
      # Print the Confusion matrix.
      print('Confusion Matrix of Test set predictions(XGB Classifier): \n', __
       →xgconmatrix)
     Confusion Matrix of Test set predictions(XGB Classifier):
      [[360 52]
      [ 78 166]]
[47]: # Plot the confusion matrix.
      # Define the size of the plot
      plt.figure(figsize=(8,6))
      # Confusion matrix heat map.
      sns.heatmap(xgconmatrix, annot=True,cmap = 'flare', fmt='d')
      # Plot Title
      plt.title('Confusion Matrix of Test data Set(XGB Classifier)', fontsize = 18)
      # x- Label
      plt.xlabel('Actual Values', fontsize = 14)
      # y-label
      plt.ylabel('Predicted values', fontsize = 14)
[47]: Text(51.0, 0.5, 'Predicted values')
```





```
# Let's get the accuracy score, Precision, Recall and F1 Score of the Support
Vector Machine Classifier.
# Getting the accuracy score of the test dataset.
xg_accuracy = metrics.accuracy_score(y_test, xg_pred)

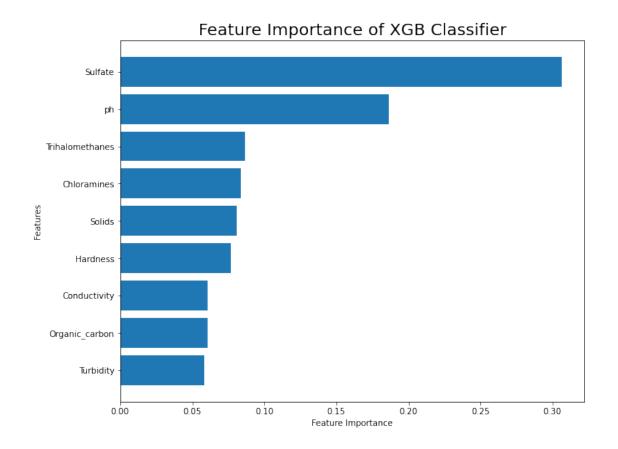
# Getting the accuracy score of the train dataset
xg_accu_train = metrics.accuracy_score(y_train, xg_pred_train)

# Getting the precision score.
xg_precision = round(precision_score(y_test, xg_pred),3)

# Getting the Recall Score.
xg_recall = round(recall_score(y_test, xg_pred),3)

# Getting the F1 Score.
xg_flscore = round(f1_score(y_test, xg_pred),3)
```

```
# Printing the accuracy of the model.
      print('The accuracy score of the XGB Classifier on test dataset: {} '.
      →format(xg_accuracy))
      print('The accuracy score of the XGB Classifier on train dataset : {} '.
       →format(xg_accu_train))
      print('Precision Score of the test set for the XGB Classifier : {}'.
      →format(xg_precision))
      print('Recall Score of the test set for the XGB Classifier : {}'.
      →format(xg_recall))
      print('F1 Score of the test set for the XGB Classifier : {}'.format(xg_f1score))
     The accuracy score of the XGB Classifier on test dataset: 0.801829268292683
     The accuracy score of the XGB Classifier on train dataset : 1.0
     Precision Score of the test set for the XGB Classifier: 0.761
     Recall Score of the test set for the XGB Classifier: 0.68
     F1 Score of the test set for the XGB Classifier: 0.719
[49]: # The importance of a feature is basically: how much this feature is used in
      \rightarrow each tree of the forest.
      xgimportantfeatures = xg.feature_importances_
      xgimportantfeatures
[49]: array([0.18623367, 0.07650602, 0.0807505, 0.08349272, 0.30664834,
             0.06081047, 0.06049719, 0.08683765, 0.05822343], dtype=float32)
[50]: # initialize the idicies
      indices = np.argsort(xgimportantfeatures)
      # Defining the size of the plot
      fig, ax = plt.subplots(figsize = (10,8))
      # Plotting the graph
      ax.barh(range(len(xgimportantfeatures)), xgimportantfeatures[indices])
      # Y-ticks
      ax.set_yticks(range(len(xgimportantfeatures)))
      _ = ax.set_yticklabels(np.array(X_train.columns)[indices])
      # X-label
      plt.xlabel("Feature Importance", fontsize = 10)
      # Y - label
      plt.ylabel("Features", fontsize = 10)
      # Title of the plot
      plt.title("Feature Importance of XGB Classifier", fontsize = 20)
[50]: Text(0.5, 1.0, 'Feature Importance of XGB Classifier')
```



0.0.14 Model Comparisions

```
[51]: # Let's form arrays for the calculated accuracy score for the train and test \Box
      ⇔datasets, Precision,
      # Recall and F1 Score for the above three models.
      KNN_classifier = {'Model':'KNN Classifier',
                       'Accuracy (test)':knnclass_accuracy,
                    'Accuracy (train)':knnclass_accu_train,
                    'Precision score' :knnclass_precision,
                    'Recall score'
                                      :knnclass_recall,
                     'F1 Score':knnclass_f1score,}
      RandomForestclassifier = {'Model':'RandomForestClassifier',
                               'Accuracy (test)':rdreg_accuracy,
                                'Accuracy (train)':rdreg_accu_train,
                               'Precision score' :rdreg_precision,
                                'Recall score' :rdreg_recall,
                                'F1 Score':rdreg_f1score,}
      XGBClassifier = {'Model':'XGBClassifier',
```

```
'Accuracy (test)':xg_accuracy,
    'Accuracy (train)':xg_accu_train,
'Precision score' :xg_precision,
'Recall score' :xg_recall,
    'F1 Score':xg_f1score,}
```

[52]:		KNN	Classifier	${\tt RandomForest}$	Classifier	XGB Classifier
M	Model	KNN	${\tt Classifier}$	RandomForest	tClassifier	XGBClassifier
A	Accuracy (test)		0.53811		0.807927	0.801829
A	Accuracy (train)		0.714504		1.0	1.0
P	recision score		0.335		0.811	0.761
R	Recall score		0.246		0.631	0.68
F	71 Score		0.284		0.71	0.719