# DCS 630 Predictive Analytics (DSC630-T302 2231-1)

# **Bellevue University**

**Term Project: Prediction of Water Quality** 

Milestone 4: Finalizing the Results

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For Milestone 4 Predictive Model code, see the Predictive Modeling and Evaluation section.

#### **Dataset Source**

The objective of this project is to construct a model to predict whether water is potable based on water quality measurements. The dataset used for this analysis can be found through the link below:

**Prediction of Water Quality Dataset** 

# **Import Necessary Libraries**

```
In [1]:
        Import the necessary libraries for the analysis.
        import numpy as np
        import pandas as pd
        import seaborn as sns
        import scipy.stats
        import sklearn
        import operator
        import matplotlib
        import matplotlib.pyplot as plt
        from scipy import stats
        from collections import OrderedDict
        from scipy.stats import shapiro
        from scipy.stats import pointbiserialr
        from sklearn.model selection import train test split
        from sklearn.decomposition import PCA
        from sklearn.linear model import LogisticRegression
        from sklearn.neighbors import KNeighborsClassifier
        from sklearn.tree import DecisionTreeClassifier
        from sklearn.tree import plot tree
        from sklearn.ensemble import RandomForestClassifier, GradientBoostingClassifier, BaggingCl
        from sklearn.svm import SVC
        from sklearn.preprocessing import StandardScaler
        from sklearn.preprocessing import MinMaxScaler
        from sklearn.metrics import roc auc score
        from sklearn.model selection import cross val score
        from sklearn.metrics import precision score, recall score, f1 score
        from sklearn.metrics import confusion matrix, accuracy score, classification report
```

```
Check the versions of the packages.

'''

print('numpy version:', np.__version__)

print('pandas version:', pd.__version__)

print('seaborn version:', sns.__version__)

print('matplotlib version:', matplotlib.__version__)

print('sklearn:', sklearn.__version__)
```

numpy version: 1.20.3 pandas version: 1.3.4 seaborn version: 0.11.2 matplotlib version: 3.4.3

sklearn: 0.24.2

#### **Overview of Dataset**

Out[4]:		ph	Hardness	Solids	Chloramines	Sulfate	Conductivity	Organic_carbon	Trihalomethanes	Tur
	0	NaN	204.890456	20791.31898	7.300212	368.516441	564.308654	10.379783	86.990970	2.9
	1	3.716080	129.422921	18630.05786	6.635246	NaN	592.885359	15.180013	56.329076	4.5
	2	8.099124	224.236259	19909.54173	9.275884	NaN	418.606213	16.868637	66.420093	3.0
	3	8.316766	214.373394	22018.41744	8.059332	356.886136	363.266516	18.436525	100.341674	4.6
	4	9.092223	181.101509	17978.98634	6.546600	310.135738	398.410813	11.558279	31.997993	4.0
	5	5.584087	188.313324	28748.68774	7.544869	326.678363	280.467916	8.399735	54.917862	2.5
	6	10.223862	248.071735	28749.71654	7.513408	393.663395	283.651634	13.789695	84.603556	2.6
	7	8.635849	203.361523	13672.09176	4.563009	303.309771	474.607645	12.363817	62.798309	4.4
	8	NaN	118.988579	14285.58385	7.804174	268.646941	389.375566	12.706049	53.928846	3.5
	9	11.180284	227.231469	25484.50849	9.077200	404.041635	563.885481	17.927806	71.976601	4.3

There are 3276 rows and 10 columns in this dataset.

```
In [6]:
    Display the total size of this dataset.
    '''
    print('This dataset contains {} records.'.format(df.size))
```

This dataset contains 32760 records. In [7]: 1.1.1 Find the type of data within each column initially. df.dtypes Out[7]: ph float64 float64 Hardness float64 float64 Solids Chloramines Sulfate float64
Conductivity float64
Organic\_carbon float64
Trihalomethanes float64 float64 Turbidity Potability int64 dtype: object In [8]: Understand if there are any missing values in the dataset. df.isna().sum().sort values(ascending = False) 781 Sulfate Out[8]: 491 Trihalomethanes 162 Hardness Solids Chloramines Conductivity Organic carbon Turbidity 0 Potability dtype: int64 In [9]:

This dataset contains 1434 total missing values.

The missing values are located within Sulfate, ph, and Trihalomethane features. There are 781, 491, and 162 missing values in each respective feature. This will need to be addressed during the Data Preparation step.

print('This dataset contains {} total missing values.'.format(missing values))

## **Highlights for Dataset Overview:**

missing values = df.isna().sum().sum()

- There are 3276 records and 10 features in the dataset.
- There are 1434 missing values within Sulfate, ph, and Trihalomethanes features.

Understand how many missing values are in the dataset initially.

## **Data Understanding - Preliminary Preparation**

```
Out[11]:
                            ph
                                    hardness
                                                     solids chloramines
                                                                                sulfate conductivity organic_carbon trihalometha
           count 2785.000000
                                 3276.000000
                                               3276.000000
                                                             3276.000000
                                                                           2495.000000
                                                                                         3276.000000
                                                                                                          3276.000000
                                                                                                                            3114.000
                                                                                                                              66.396
                      7.080795
                                  196.369496
                                              22014.092526
                                                                7.122277
                                                                            333.775777
                                                                                          426.205111
                                                                                                            14.284970
            mean
                      1.594320
                                               8768.570828
              std
                                   32.879761
                                                                1.583085
                                                                             41.416840
                                                                                           80.824064
                                                                                                             3.308162
                                                                                                                              16.175
                      0.000000
                                   47.432000
                                                 320.942611
                                                                            129.000000
                                                                                                             2.200000
                                                                                                                               0.738
             min
                                                                0.352000
                                                                                          181.483754
             25%
                      6.093092
                                  176.850538 15666.690300
                                                                6.127421
                                                                            307.699498
                                                                                          365.734414
                                                                                                            12.065801
                                                                                                                              55.844
             50%
                      7.036752
                                  196.967627
                                              20927.833605
                                                                7.130299
                                                                            333.073546
                                                                                          421.884968
                                                                                                            14.218338
                                                                                                                              66.622
             75%
                      8.062066
                                  216.667456 27332.762125
                                                                8.114887
                                                                            359.950170
                                                                                          481.792305
                                                                                                            16.557652
                                                                                                                              77.337
             max
                     14.000000
                                  323.124000 61227.196010
                                                               13.127000
                                                                            481.030642
                                                                                          753.342620
                                                                                                            28.300000
                                                                                                                             124.000
```

```
In [12]:
    Review the percentage of missing features for sulfate, ph, and trihalomethanes.
    '''
    perc_missing = round(df.isnull().sum() * 100 / len(df),2)
    df_missing = pd.DataFrame({'feature': df.columns, 'percent_missing': perc_missing})
    df_missing.sort_values(by= 'percent_missing', ascending = False)
```

	feature	percent_missing
sulfate	sulfate	23.84
ph	ph	14.99
trihalomethanes	trihalomethanes	4.95
hardness	hardness	0.00
solids	solids	0.00
chloramines	chloramines	0.00
conductivity	conductivity	0.00
organic_carbon	organic_carbon	0.00
turbidity	turbidity	0.00
potability	potability	0.00

333.389426

Out[12]:

Identify a plan on how to handle the missing values in sulfate, ph, and trihalomethane columns. Evaluate whether to replace the missing values (with mean or median) or remove the values from the data set.

```
331.838167
         Name: sulfate, dtype: float64
In [14]:
         Evaluate the mean for the missing values in the sulfate column for potable and non-potable
         df.groupby('potability')['sulfate'].mean()
         potability
Out[14]:
             334.56429
              332.56699
         Name: sulfate, dtype: float64
In [15]:
          1.1.1
         Evaluate the median for the missing values in the ph column for potable and non-potable we
         df.groupby('potability')['ph'].median()
         potability
Out[15]:
             7.035456
             7.036752
         Name: ph, dtype: float64
In [16]:
         Evaluate the mean for the missing values in the sulfate column for potable and non-potable
          df.groupby('potability')['ph'].mean()
         potability
Out[16]:
             7.085378
              7.073783
         Name: ph, dtype: float64
In [17]:
         Evaluate the median for the missing values in the trihalomethanes column for potable and
         df.groupby('potability')['trihalomethanes'].median()
         potability
Out[17]:
             66.542198
              66.678214
         Name: trihalomethanes, dtype: float64
In [18]:
         Evaluate the mean for the missing values in the trihalomethanes column for potable and nor
         df.groupby('potability')['trihalomethanes'].mean()
         potability
Out[18]:
             66.303555
              66.539684
         Name: trihalomethanes, dtype: float64
        Based on the descriptive statistics above, the mean or median may be a suitable choice to replace the features
        for each respective variables. As a result, I'll replace the sulfate, ph, and trihalomethane values with the median
        of each respective feature.
```

Replace the NaN values for sulfate, trihalomethanes, and ph with the median values of each

In [19]:

1.1.1

Handle the missing values as described above.

Drop the remaining missing values from the sulfate column.

df['trihalomethanes'].fillna(value=df['trihalomethanes'].median(), inplace = True)

df['sulfate'].fillna(value=df['sulfate'].median(), inplace = True)

df['ph'].fillna(value=df['ph'].median(), inplace = True)

Out[20]: feature percent\_missing 0.0 ph ph hardness hardness 0.0 solids solids 0.0 chloramines chloramines 0.0 sulfate sulfate 0.0 conductivity conductivity 0.0 organic\_carbon organic\_carbon 0.0 trihalomethanes trihalomethanes 0.0 turbidity turbidity 0.0

potability

potability

```
In [21]:
    Show the revised dataframe after preliminary prepearation steps.
    iii
    df.head()
```

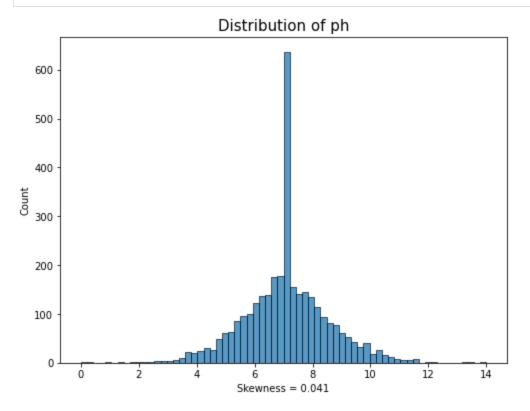
0.0

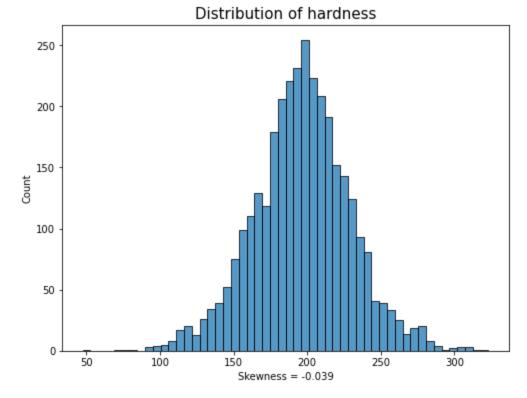
Out[21]:		ph	hardness	solids	chloramines	sulfate	conductivity	organic_carbon	trihalomethanes	turbic
	0	7.036752	204.890456	20791.31898	7.300212	368.516441	564.308654	10.379783	86.990970	2.963
	1	3.716080	129.422921	18630.05786	6.635246	333.073546	592.885359	15.180013	56.329076	4.500
	2	8.099124	224.236259	19909.54173	9.275884	333.073546	418.606213	16.868637	66.420093	3.055!
	3	8.316766	214.373394	22018.41744	8.059332	356.886136	363.266516	18.436525	100.341674	4.628
	4	9.092223	181.101509	17978.98634	6.546600	310.135738	398.410813	11.558279	31.997993	4.0750

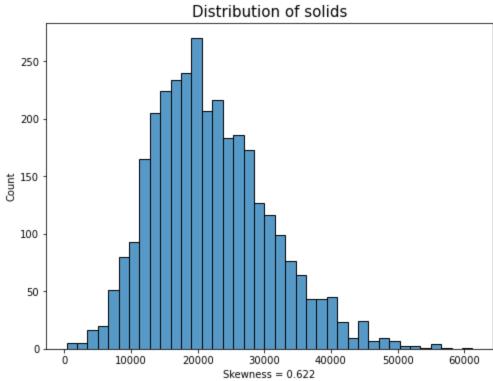
# **Highlights for Preliminary Data Preparation:**

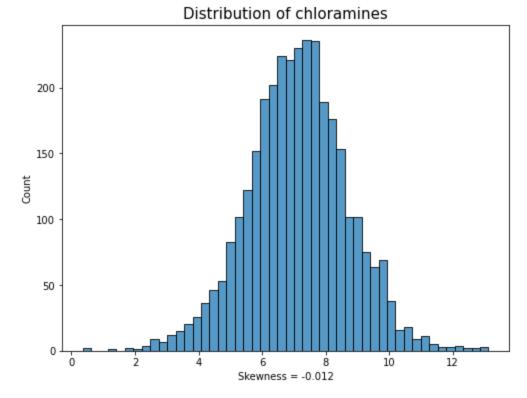
- Columns were renamed for convenience of analysis.
- Descriptive statistics of each numeric variable were displayed.
- Missing values for sulfate,ph, and trihalomethanes were replaced with the median values for each respective feature.

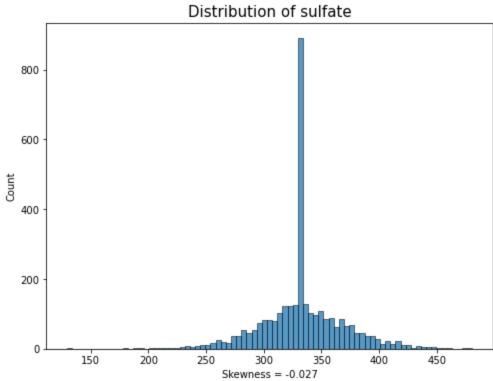
# **Data Understanding - Univariate Analysis**

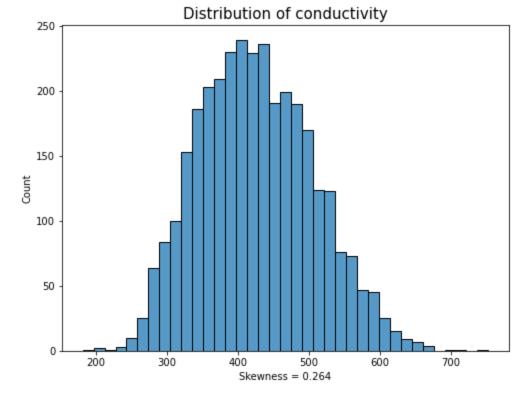


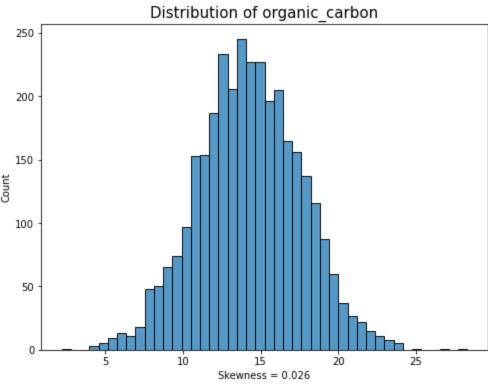




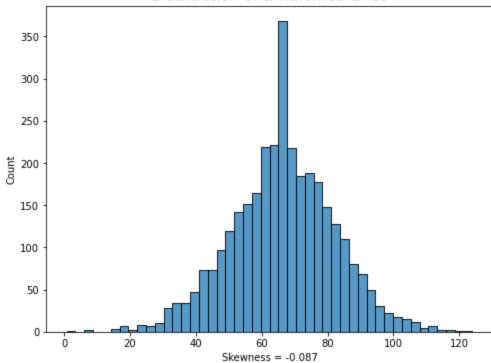


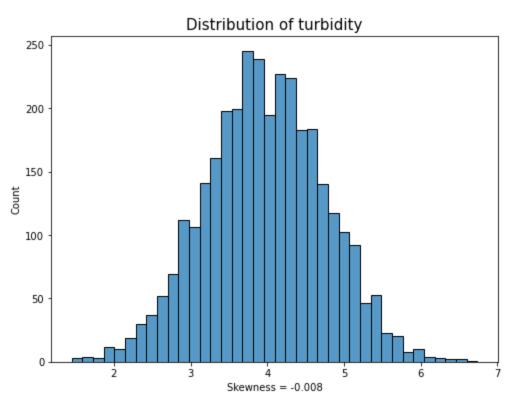






### Distribution of trihalomethanes

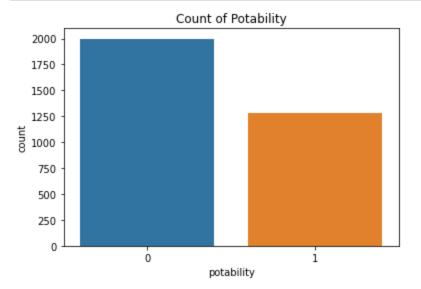


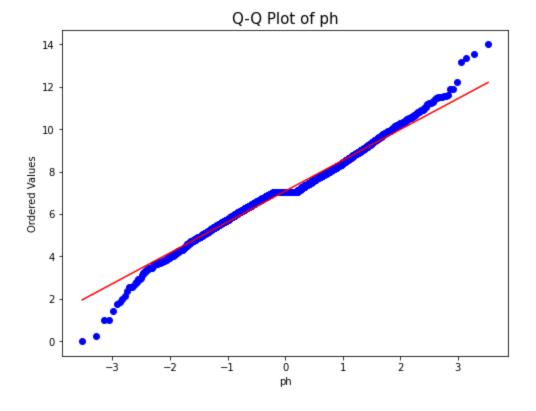


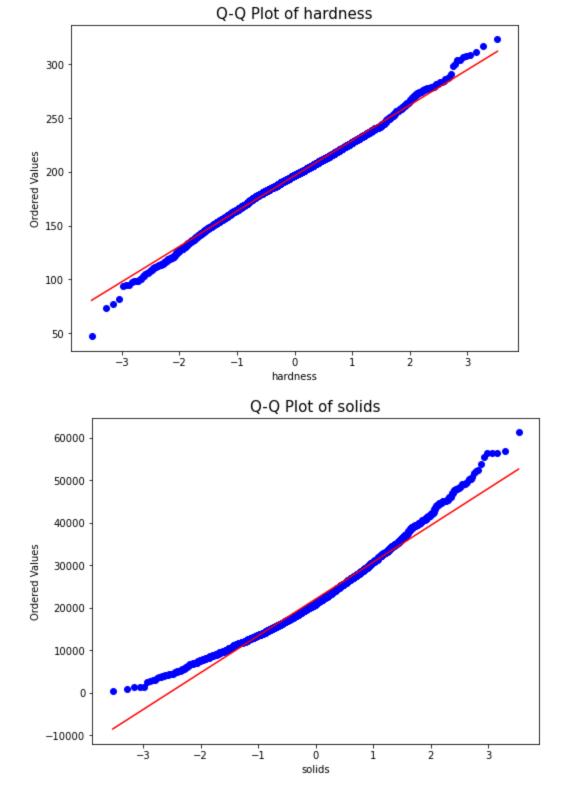
```
In [23]:
    Use describe() to understand the target variable descriptive statistics.
    df['potability'].value_counts()
```

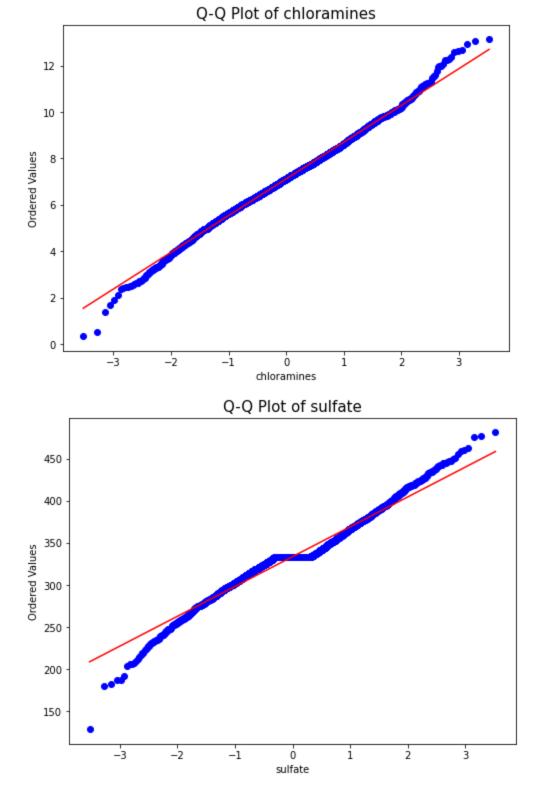
```
Out[23]: 0 1998
1 1278
Name: potability, dtype: int64
```

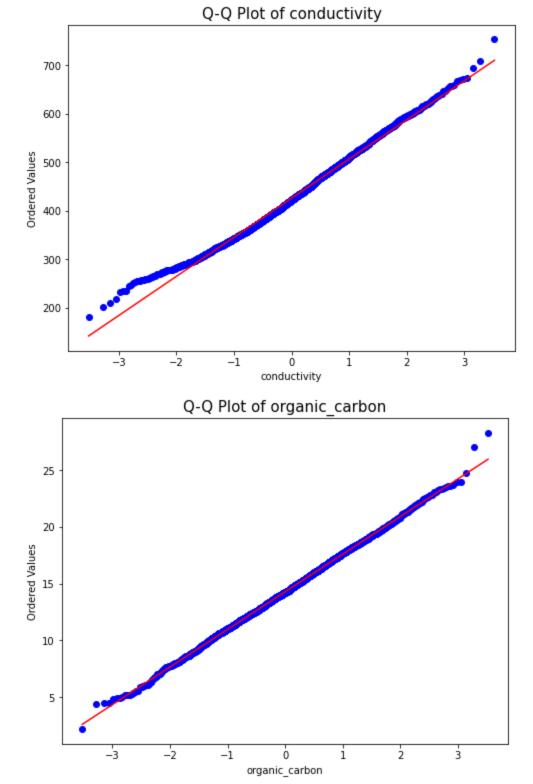
```
sns.countplot(x='potability', data = df).set(title = 'Count of Potability')
plt.show()
```

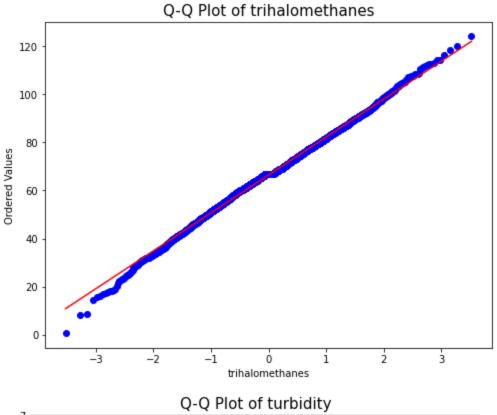


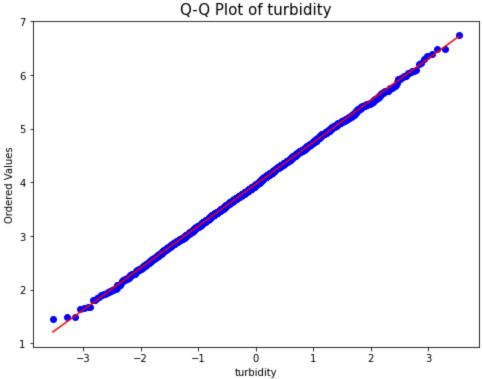






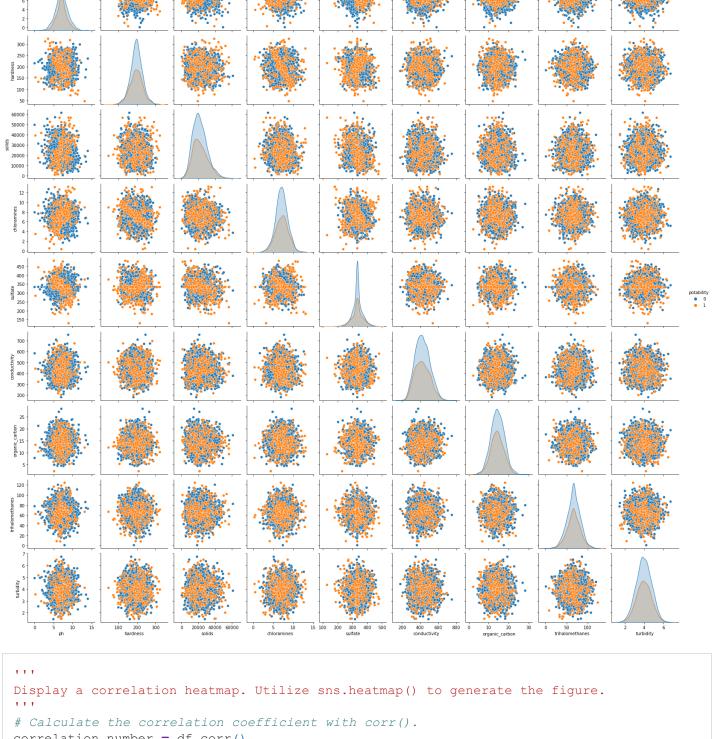




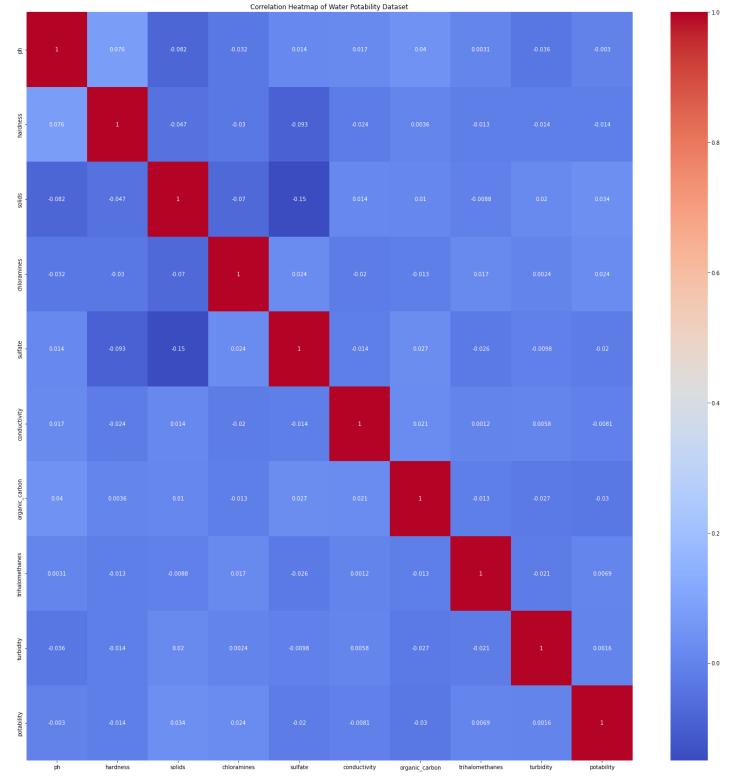


'conductivity test':shapiro test conductivity, 'oc test':shapiro test

## **Data Understanding - Multivariate Analysis**



Out[28]: Text(0.5, 1.0, 'Correlation Heatmap of Water Potability Dataset')



```
Check pointbiserial correlations for potability to see if there are any significant differ

"""

# Calculate the correlation coefficient for features with potability.

correlation_number_pb1 = pointbiserialr(df['potability'],df['ph'])

correlation_number_pb2 = pointbiserialr(df['potability'],df['hardness'])

correlation_number_pb3 = pointbiserialr(df['potability'],df['solids'])

correlation_number_pb4 = pointbiserialr(df['potability'],df['chloramines'])

correlation_number_pb5 = pointbiserialr(df['potability'],df['sulfate'])

correlation_number_pb6 = pointbiserialr(df['potability'],df['conductivity'])

correlation_number_pb7 = pointbiserialr(df['potability'],df['trihalomethanes'])

correlation_number_pb8 = pointbiserialr(df['potability'],df['turbidity'])

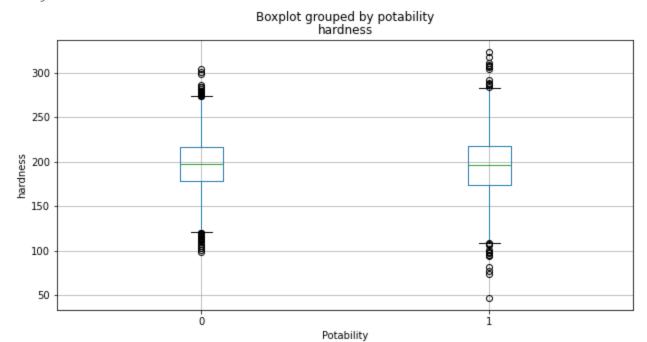
# Create a dictionary of the pointbiserial correlation values.
```

```
'solids corr':correlation number pb3, 'chloramines corr':correlation number pb4
                     'sulfate corr':correlation number pb5, 'conductivity':correlation number pb6,
                     'organic carbon corr':correlation number pb7, 'trihalomethanes corr':correlation
                     'turbidity corr':correlation number pb9}
         pb corr dict
        {'ph corr': PointbiserialrResult(correlation=-0.0030139288906728037, pvalue=0.863090956972
Out[29]:
          'hardness corr': PointbiserialrResult(correlation=-0.013836558096163432, pvalue=0.4285415
        8703761713),
         'solids corr': PointbiserialrResult(correlation=0.033743313390906485, pvalue=0.0534636382
        0366133),
         'chloramines corr': PointbiserialrResult(correlation=0.0237789720027849, pvalue=0.1736102
        245030519),
          'sulfate corr': PointbiserialrResult(correlation=-0.020476394607078924, pvalue=0.24132969
        39297576),
          'conductivity': PointbiserialrResult(correlation=-0.00812831972472755, pvalue=0.641884548
        4637925),
         'organic carbon corr': PointbiserialrResult(correlation=-0.030001366622918028, pvalue=0.0
        8599746151306269),
         'trihalomethanes corr': PointbiserialrResult(correlation=0.006886582826791652, pvalue=0.6
        935683946431569),
         'turbidity corr': PointbiserialrResult(correlation=0.0015806819408441912, pvalue=0.927939
        1626737951)}
In [30]:
         1.1.1
         Sort the dictionary by correlation value.
         sorted dict = sorted(pb corr dict.items(), key = operator.itemgetter(1))
         sorted dict
Out[30]: [('organic_carbon_corr',
          PointbiserialrResult(correlation=-0.030001366622918028, pvalue=0.08599746151306269)),
          ('sulfate corr',
          PointbiserialrResult(correlation=-0.020476394607078924, pvalue=0.2413296939297576)),
          ('hardness corr',
          PointbiserialrResult(correlation=-0.013836558096163432, pvalue=0.42854158703761713)),
          ('conductivity',
          PointbiserialrResult(correlation=-0.00812831972472755, pvalue=0.6418845484637925)),
          ('ph corr',
          PointbiserialrResult(correlation=-0.0030139288906728037, pvalue=0.8630909569723499)),
          ('turbidity corr',
          PointbiserialrResult(correlation=0.0015806819408441912, pvalue=0.9279391626737951)),
          ('trihalomethanes corr',
          PointbiserialrResult(correlation=0.006886582826791652, pvalue=0.6935683946431569)),
          ('chloramines corr',
          PointbiserialrResult(correlation=0.0237789720027849, pvalue=0.1736102245030519)),
          ('solids corr',
          PointbiserialrResult(correlation=0.033743313390906485, pvalue=0.05346363820366133))]
In [31]:
         Construct box plots to understand each feature distribution based on potability.
         1.1.1
         for col in df.columns[0:9]:
             plt.figure(figsize=(8,4))
             df.boxplot(column=col, by='potability', figsize=(10,5))
             plt.xlabel("Potability".format(col), fontsize=10)
             plt.ylabel("{}".format(col), fontsize=10)
             plt.show()
```

pb corr dict = {'ph corr':correlation number pb1, 'hardness corr':correlation number pb2,

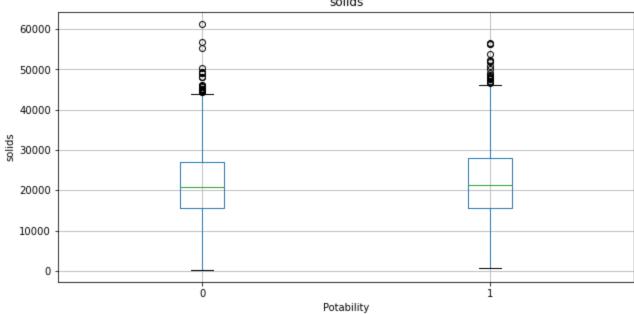
# Boxplot grouped by potability ph 듄

<Figure size 576x288 with 0 Axes>



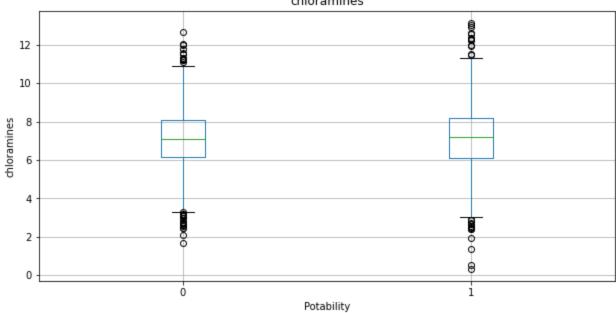
Potability

# Boxplot grouped by potability solids

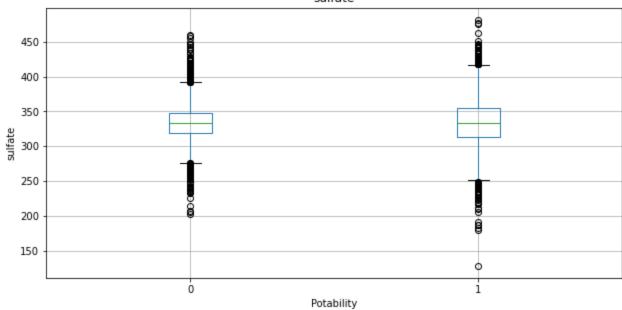


<Figure size 576x288 with 0 Axes>

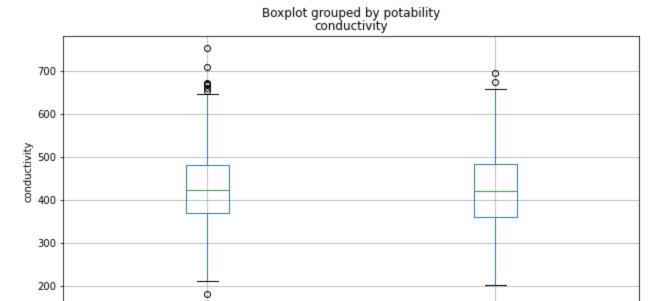
# Boxplot grouped by potability chloramines



# Boxplot grouped by potability sulfate



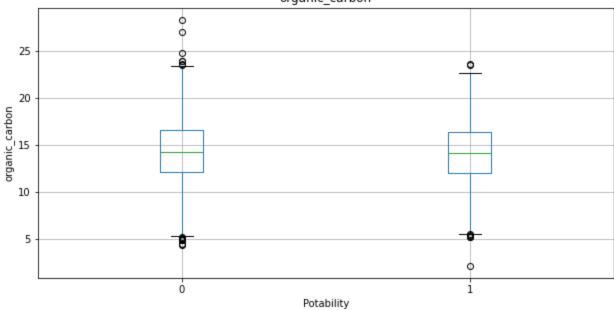
<Figure size 576x288 with 0 Axes>



Potability

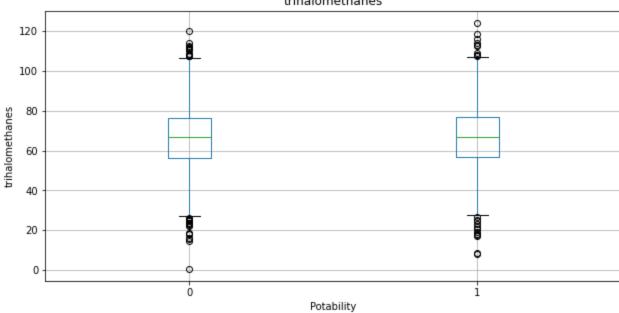
i

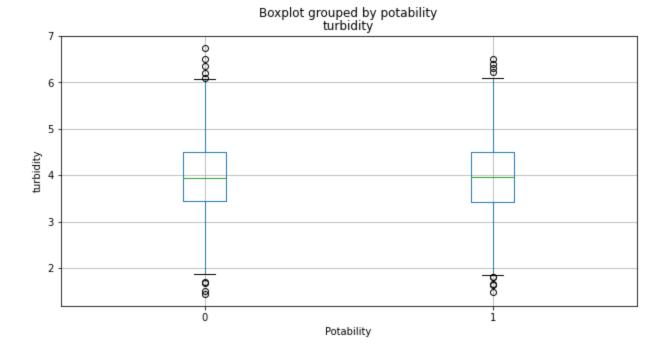
## Boxplot grouped by potability organic\_carbon



<Figure size 576x288 with 0 Axes>

# Boxplot grouped by potability trihalomethanes





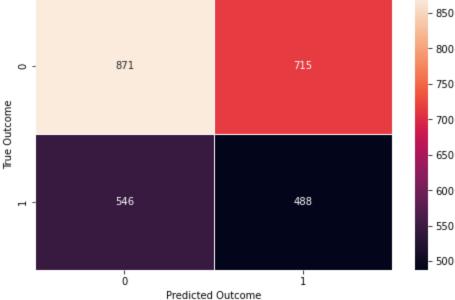
# **Data Preparation**

```
In [32]:
         Split the data into training and test set, where 'potability' is the target.
         X = df.drop('potability', axis = 1)
         y = df['potability']
In [33]:
         Split the data with train_test_split from sklearn.
         Use test size = 0.2 to split the data into 80% training and 20% testing data.
         X train, X test, y train, y test = train test split(X, y, test size = 0.2, random state =
In [34]:
         1.1.1
         Print out the shape of the resulting datasets for training and testing.
         print("The original data set shape was {} rows and {} columns.".format(df.shape[0],df.shape
         print("The X train shape is {} rows and {} columns.".format(X train.shape[0], X train.shape
         print("The y train shape is {} rows.".format(y train.shape[0]))
         print("The X test shape is {} rows and {} columns.".format(X test.shape[0], X test.shape[1]
         print("The y test shape is {} rows.".format(y test.shape[0]))
        The original data set shape was 3276 rows and 10 columns.
        The X train shape is 2620 rows and 9 columns.
        The y train shape is 2620 rows.
        The X test shape is 656 rows and 9 columns.
        The y test shape is 656 rows.
```

## **Predictive Modeling and Evaluation**

#### **Logistic Regression Model**

```
In [36]:
         Fit the Logistic Regression Classifier on the training dataset.
         lr classifier = lr.fit(X train, y train)
         lr classifier
         LogisticRegression(class weight='balanced')
Out[36]:
In [37]:
         Obtain the y prediction probabilities for each record in the training dataset.
         y predictions lr train = lr.predict(X train)
In [38]:
         Generate a Confusion Matrix for the Logistic Regression Classifier based on the training
         cm lr train = confusion matrix(y train, y predictions lr train)
In [39]:
         Plot the confusion matrix so it is clearly labelled and illustrated.
         f, ax = plt.subplots(figsize = (8,5))
         sns.heatmap(cm lr train, annot = True, linewidths = 0.5, fmt = ".0f", ax = ax)
         plt.xlabel('Predicted Outcome')
         plt.ylabel('True Outcome')
         plt.title('Confusion Matrix of Logistic Regression Classifier on Training Data')
         plt.show()
          Confusion Matrix of Logistic Regression Classifier on Training Data
```



In [41]: Generate a Confusion Matrix for the Logistic Regression Classifier based on the test datas



cm\_lr\_test = confusion\_matrix(y\_test, y\_predictions\_lr\_test)

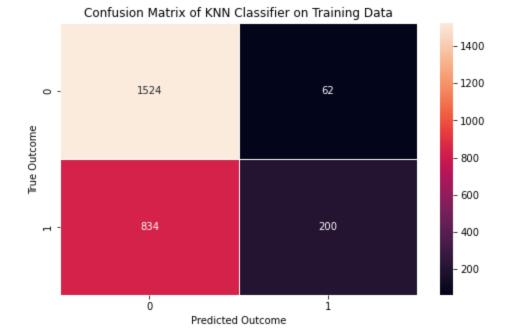
	precision	recall	f1-score	support
0	0.63	0.54	0.58 0.42	412 244
accuracy macro avg weighted avg	0.51 0.54	0.51 0.52	0.52 0.50 0.52	656 656 656

```
print("Recall of Logistic Regression Model on testing data is:{}".format(recall_lr))
print("F1 Score of Logistic Regression Model on testing data is:{}".format(f1_lr))
```

Accuracy of Logistic Regression Model on training data is:0.5187022900763358
Accuracy of Logistic Regression Model on testing data is:0.5152439024390244
Precision of Logistic Regression Model on testing data is:0.37748344370860926
Recall of Logistic Regression Model on testing data is:0.4672131147540984
F1 Score of Logistic Regression Model on testing data is:0.4175824175824176

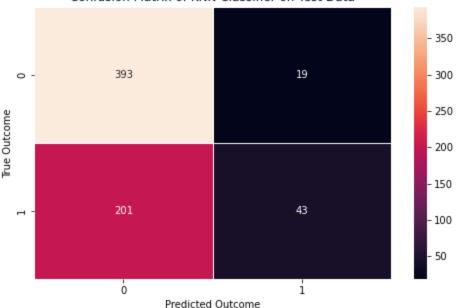
#### K-NN (K-Neareast Neighbor) Model

```
In [45]:
         Standardize the X train and X test datasets for the remainder of the models being evaluate
         sc = StandardScaler()
         X train = sc.fit transform(X train)
         X test = sc.transform(X test)
In [46]:
         Create the KNN Classifier.
         K = 38 based on Error Rate evaluation shown in this section.
         knn = KNeighborsClassifier(n neighbors = 38)
In [47]:
         Fit the KNN Classifier on the training dataset.
         knn classifier = knn.fit(X train, y train)
         knn classifier
        KNeighborsClassifier(n neighbors=38)
Out[47]:
In [48]:
         Obtain the y prediction probabilities for each record in the training dataset.
         y predictions knn train = knn.predict(X train)
In [49]:
         Generate a Confusion Matrix for the Logistic Regression Classifier based on the training
         cm knn train = confusion matrix(y train, y predictions knn train)
In [50]:
         1.1.1
         Plot the confusion matrix so it is clearly labelled and illustrated.
         f, ax = plt.subplots(figsize = (8,5))
         sns.heatmap(cm knn train, annot = True, linewidths = 0.5, fmt = ".0f", ax = ax)
         plt.xlabel('Predicted Outcome')
         plt.ylabel('True Outcome')
         plt.title('Confusion Matrix of KNN Classifier on Training Data')
         plt.show()
```



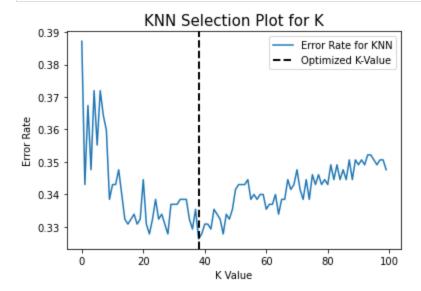
```
In [51]:
         Obtain the y predictions for the KNN Classifier.
         y predictions knn test = knn.predict(X test)
In [52]:
          1.1.1
         Generate a confusion matrix based on the test data set.
         cm_knn = confusion_matrix(y_test, y_predictions_knn_test)
         cm knn
         array([[393, 19],
Out[52]:
                [201, 43]], dtype=int64)
In [53]:
          1.1.1
         Plot the confusion matrix so it is clearly labelled and illustrated.
         f, ax = plt.subplots(figsize = (8,5))
         sns.heatmap(cm knn, annot = True, linewidths = 0.5, fmt = ".0f", ax = ax)
         plt.xlabel('Predicted Outcome')
         plt.ylabel('True Outcome')
         plt.title('Confusion Matrix of KNN Classifier on Test Data')
         plt.show()
```

#### Confusion Matrix of KNN Classifier on Test Data



	precision	recall	f1-score	support
0 1	0.66 0.69	0.95 0.18	0.78 0.28	412 244
accuracy	0.60	0 57	0.66	656
macro avg weighted avg	0.68 0.67	0.57 0.66	0.53	656 656

```
In [55]:
         Choose an optimal K value for KNN by plotting the error rate for the model against the dat
         # Create an empty list for error rate.
         error rate = []
         # Create a for loop to get error rates appended to the error rate list.
         for i in np.arange(1, 101):
             new model = KNeighborsClassifier(n neighbors = i)
             new model.fit(X train, y train)
             new predictions = new model.predict(X test)
             error rate.append(np.mean(new predictions != y test))
         # Create the plot to assist with selecting a K value for the KNN classifier.
         plt.plot(error rate, label = 'Error Rate for KNN')
         plt.title("KNN Selection Plot for K", fontsize=15)
         plt.xlabel("K Value", fontsize=10)
         plt.ylabel("Error Rate", fontsize=10)
         plt.xticks(fontsize=10)
         plt.yticks(fontsize=10)
         plt.axvline(x=pd.Series(error rate).idxmin(), linewidth = 2, color = 'k', linestyle = '--'
         plt.legend(loc = 'upper right')
         plt.show()
```



```
In [56]:
          1.1.1
         Show the optimum number of K based on the plot above.
         print('Optimum K-Value for KNN:{}'.format(pd.Series(error rate).idxmin()))
```

Optimum K-Value for KNN:38

```
In [57]:
         Calculate the accuracy for the model based on training and test data. Also, report the
         Precision, Recall, and F1 score for the model predications against the test data.
         accuracy knn train = accuracy score(y train, y predictions knn train)
         accuracy knn test = accuracy score(y test, y predictions knn test)
         precision_knn = precision_score(y_test, y_predictions_knn_test)
         recall knn = recall score(y test, y predictions knn test)
         f1 knn = f1 score(y test, y predictions knn test)
         print("Accuracy of KNN Model on training data is:{}".format(accuracy knn train))
         print("Accuracy of KNN Model on testing data is:{}".format(accuracy knn test))
         print("Precision of KNN Model on testing data is:{}".format(precision knn))
         print("Recall of KNN Model on testing data is:{}".format(recall knn))
         print("F1 Score of KNN Model on testing data is:{}".format(f1 knn))
```

Accuracy of KNN Model on training data is: 0.6580152671755726 Accuracy of KNN Model on testing data is:0.6646341463414634 Precision of KNN Model on testing data is:0.6935483870967742 Recall of KNN Model on testing data is:0.1762295081967213 F1 Score of KNN Model on testing data is:0.28104575163398693

#### **Decision Tree Model**

In [58]:

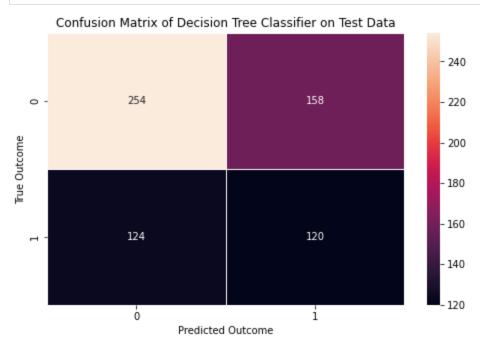
1.1.1

```
Create the Decision Tree Classifier.
         dt = DecisionTreeClassifier()
In [59]:
         Fit the Decision Tree Classifier on the training dataset.
         decision tree classifier = dt.fit(X train, y train)
         decision tree classifier
         DecisionTreeClassifier()
```

Out[59]:

```
In [60]:
          Obtain the y prediction probabilities for each record in the training dataset.
          y_predictions_dt_train = dt.predict(X train)
In [61]:
          Generate a Confusion Matrix for the Decision Tree Classifier based on the training dataset
          cm dt train = confusion matrix(y train, y predictions dt train)
In [62]:
          Plot the confusion matrix so it is clearly labelled and illustrated.
          f, ax = plt.subplots(figsize = (8,5))
          sns.heatmap(cm dt train, annot = True, linewidths = 0.5, fmt = ".0f", ax = ax)
          plt.xlabel('Predicted Outcome')
          plt.ylabel('True Outcome')
          plt.title('Confusion Matrix of Decision Tree Classifier on Training Data')
          plt.show()
             Confusion Matrix of Decision Tree Classifier on Training Data
                                                                    - 1400
                                                                    - 1200
                        1586
                                                  0
           0 -
                                                                    - 1000
         Frue Outcome
                                                                    - 800
                                                                    600
                                                                     400
                                                                     200
                                                  i
                         ò
                               Predicted Outcome
In [63]:
          Obtain the y predictions for the decision tree classifier.
          y predictions dt test = dt.predict(X test)
In [64]:
          1.1.1
          Generate a confusion matrix based on the test data set.
          cm dt = confusion matrix(y test, y predictions dt test)
          cm dt
         array([[254, 158],
Out[64]:
                [124, 120]], dtype=int64)
In [65]:
          Plot the confusion matrix so it is clearly labelled and illustrated.
```

```
f, ax = plt.subplots(figsize = (8,5))
sns.heatmap(cm_dt, annot = True, linewidths = 0.5, fmt = ".0f", ax = ax)
plt.xlabel('Predicted Outcome')
plt.ylabel('True Outcome')
plt.title('Confusion Matrix of Decision Tree Classifier on Test Data')
plt.show()
```



	precision	recall	f1-score	support
0	0.67 0.43	0.62	0.64	412 244
accuracy macro avg weighted avg	0.55 0.58	0.55 0.57	0.57 0.55 0.57	656 656 656

```
Calculate the accuracy for the model based on training and test data. Also, report the
Precision, Recall, and F1 score for the model predications against the test data.
'''
accuracy_dt_train = accuracy_score(y_train, y_predictions_dt_train)
accuracy_dt_test = accuracy_score(y_test, y_predictions_dt_test)
precision_dt = precision_score(y_test, y_predictions_dt_test)
recall_dt = recall_score(y_test, y_predictions_dt_test)
f1_dt = f1_score(y_test, y_predictions_dt_test)
print("Accuracy of Decision Tree Model on training data is:{}".format(accuracy_dt_train))
print("Precision of Decision Tree Model on testing data is:{}".format(precision_dt))
print("Recall of Decision Tree Model on testing data is:{}".format(precision_dt))
print("F1 Score of Decision Tree Model on testing data is:{}".format(f1_dt))
```

Accuracy of Decision Tree Model on training data is:1.0

Accuracy of Decision Tree Model on testing data is:0.5701219512195121

Precision of Decision Tree Model on testing data is:0.4316546762589928

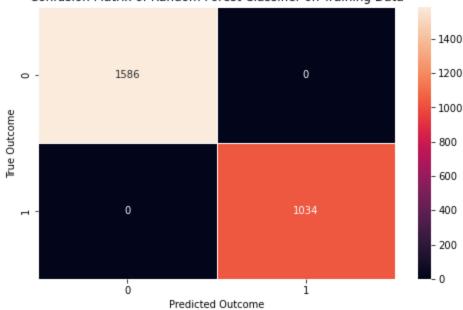
Recall of Decision Tree Model on testing data is:0.4918032786885246

F1 Score of Decision Tree Model on testing data is:0.45977011494252873

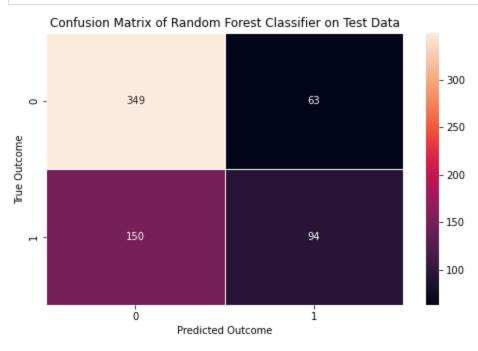
#### **Random Forest Model**

```
In [68]:
         Create the Random Forest Classifier.
         Select number of estimators at 31 based on error reduction plot for this model.
         rf = RandomForestClassifier(n estimators = 31)
In [69]:
         Fit the Random Forest Classifier on the training dataset.
         random forest classifier = rf.fit(X train, y train)
         random forest classifier
         RandomForestClassifier(n estimators=31)
Out[69]:
In [70]:
         Obtain the y prediction probabilities for each record in the training dataset.
         y predictions rf train = rf.predict(X train)
In [71]:
         Generate a Confusion Matrix for the Random Forest Classifier based on the training dataset
         cm rf train = confusion matrix(y train, y predictions rf train)
In [72]:
         Plot the confusion matrix so it is clearly labelled and illustrated.
         f, ax = plt.subplots(figsize = (8,5))
         sns.heatmap(cm rf train, annot = True, linewidths = 0.5, fmt = ".0f", ax = ax)
         plt.xlabel('Predicted Outcome')
         plt.ylabel('True Outcome')
         plt.title('Confusion Matrix of Random Forest Classifier on Training Data')
         plt.show()
```

#### Confusion Matrix of Random Forest Classifier on Training Data



```
Obtain the y predictions for the Random Forest classifier.
         y predictions rf test = rf.predict(X test)
In [74]:
         Generate a confusion matrix based on the test data set.
         cm rf = confusion matrix(y test, y predictions rf test)
         cm rf
         array([[349,
                      63],
Out[74]:
                [150, 94]], dtype=int64)
In [75]:
         Plot the confusion matrix so it is clearly labelled and illustrated.
         f, ax = plt.subplots(figsize = (8,5))
         sns.heatmap(cm rf, annot = True, linewidths = 0.5, fmt = ".0f", ax = ax)
         plt.xlabel('Predicted Outcome')
         plt.ylabel('True Outcome')
         plt.title('Confusion Matrix of Random Forest Classifier on Test Data')
         plt.show()
```



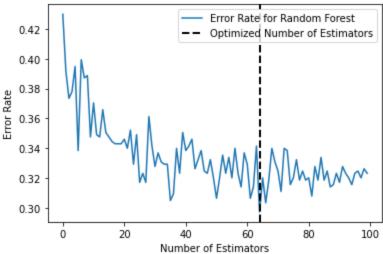
	precision	recall	f1-score	support
0	0.70	0.85	0.77	412
1	0.60	0.39	0.47	244
accuracy			0.68	656
macro avg	0.65	0.62	0.62	656
weighted avg	0.66	0.68	0.66	656

```
In [77]: ...

Choose an optimal number of estimators for Random Forest by plotting the error rate for the
```

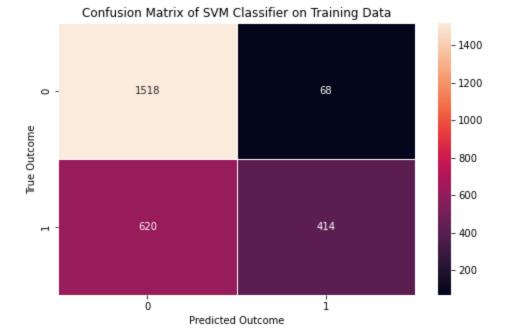
```
# Create an empty list for error rate.
error rate rf = []
# Create a for loop to get error rates appended to the error rate list.
for i in np.arange(1, 101):
    new model = RandomForestClassifier(n estimators = i)
    new model.fit(X train, y train)
    new predictions = new model.predict(X test)
    error rate rf.append(np.mean(new predictions != y test))
# Create the plot to assist with selecting a K value for the KNN classifier.
plt.plot(error rate rf, label = 'Error Rate for Random Forest')
plt.title("Random Forest Selection Plot for Number of Estimators", fontsize=15)
plt.xlabel("Number of Estimators", fontsize=10)
plt.ylabel("Error Rate", fontsize=10)
plt.xticks(fontsize=10)
plt.yticks(fontsize=10)
plt.axvline(x=pd.Series(error rate rf).idxmin(), linewidth = 2, color = 'k', linestyle =
            label = 'Optimized Number of Estimators')
plt.legend(loc = 'upper right')
plt.show()
```

#### Random Forest Selection Plot for Number of Estimators



Optimum number of estimators for Random Forest:64

```
In [80]:
         Create the SVM Classifier.
         svm = SVC()
In [81]:
         Fit the SVM Classifier on the training dataset.
         svm classifier = svm.fit(X train, y train)
         svm classifier
        SVC()
Out[81]:
In [82]:
         Obtain the y predictions for the SVM Classifier.
         y predictions svm = svm.predict(X test)
In [83]:
         Obtain the y prediction probabilities for each record in the training dataset.
         y predictions svm train = svm.predict(X train)
In [84]:
          1.1.1
         Generate a Confusion Matrix for the Support Vector Machine Model based on the training dat
         cm svm train = confusion matrix(y train, y predictions svm train)
In [85]:
         1.1.1
         Plot the confusion matrix so it is clearly labelled and illustrated.
         f, ax = plt.subplots(figsize = (8,5))
         sns.heatmap(cm svm train, annot = True, linewidths = 0.5, fmt = ".0f", ax = ax)
         plt.xlabel('Predicted Outcome')
         plt.ylabel('True Outcome')
         plt.title('Confusion Matrix of SVM Classifier on Training Data')
         plt.show()
```



```
In [86]:
         Obtain the y predictions for the SVM classifier.
         y predictions svm test = svm.predict(X test)
In [87]:
          1.1.1
         Generate a confusion matrix based on the test data set.
         cm_svm = confusion_matrix(y_test, y_predictions_svm_test)
         cm svm
         array([[375, 37],
Out[87]:
                [166,
                      78]], dtype=int64)
In [88]:
          1.1.1
         Plot the confusion matrix so it is clearly labelled and illustrated.
         f, ax = plt.subplots(figsize = (8,5))
         sns.heatmap(cm svm, annot = True, linewidths = 0.5, fmt = ".0f", ax = ax)
         plt.xlabel('Predicted Outcome')
         plt.ylabel('True Outcome')
         plt.title('Confusion Matrix of SVM Classifier on Test Data')
         plt.show()
```

## Confusion Matrix of SVM Classifier on Test Data 350 - 300 37 375 0 -- 250 Frue Outcome 200 - 150 166 78 - 100 Ò i Predicted Outcome

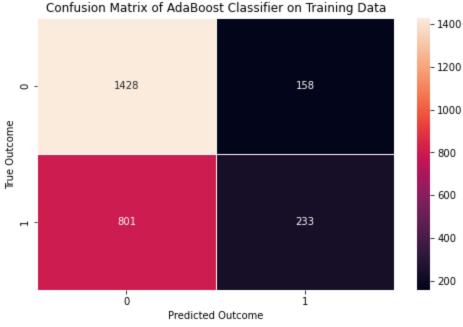
	precision	recall	f1-score	support
0	0.69	0.91	0.79	412
1	0.68	0.32	0.43	244
accuracy			0.69	656
macro avg	0.69	0.61	0.61	656
weighted avg	0.69	0.69	0.66	656

Accuracy of SVM Model on training data is:0.7374045801526717 Accuracy of SVM Model on testing data is:0.6905487804878049 Precision of SVM Model on testing data is:0.6782608695652174 Recall of SVM Model on testing data is:0.319672131147541 F1 Score of SVM Model on testing data is:0.43454038997214484

#### **Adaboost Model**

```
In [92]:
          1.1.1
         Fit the Adaboost Classifier on the training dataset.
         adaboost classifier = adaboost.fit(X train, y train)
         adaboost classifier
         AdaBoostClassifier(n estimators=20)
Out[92]:
In [93]:
         Obtain the y prediction probabilities for each record in the training dataset.
         y predictions adaboost train = adaboost.predict(X train)
In [94]:
          1.1.1
         Generate a Confusion Matrix for the AdaBoost Model based on the training dataset.
         cm adaboost train = confusion matrix(y train, y predictions adaboost train)
In [95]:
          1.1.1
         Plot the confusion matrix so it is clearly labelled and illustrated.
         f, ax = plt.subplots(figsize = (8,5))
         sns.heatmap(cm adaboost train, annot = True, linewidths = 0.5, fmt = ".0f", ax = ax)
         plt.xlabel('Predicted Outcome')
         plt.ylabel('True Outcome')
         plt.title('Confusion Matrix of AdaBoost Classifier on Training Data')
         plt.show()
```

adaboost = AdaBoostClassifier(n estimators=20)



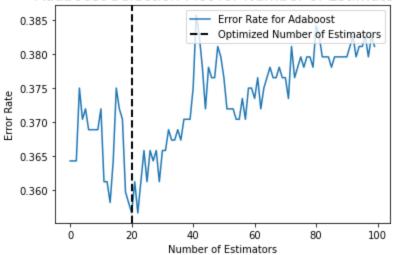
```
In [97]:
         Generate a confusion matrix based on the test data set.
         cm adaboost = confusion matrix(y test, y predictions adaboost test)
         cm adaboost
        array([[375,
                      37],
Out[97]:
                [198, 46]], dtype=int64)
In [98]:
          1.1.1
         Plot the confusion matrix so it is clearly labelled and illustrated.
         f, ax = plt.subplots(figsize = (8,5))
         sns.heatmap(cm adaboost, annot = True, linewidths = 0.5, fmt = ".0f", ax = ax)
         plt.xlabel('Predicted Outcome')
         plt.ylabel('True Outcome')
         plt.title('Confusion Matrix of Adaboost Classifier on Test Data')
         plt.show()
```

# Confusion Matrix of Adaboost Classifier on Test Data - 350 - 300 - 250 - 200 - 150 - 100 - 700

	precision	recall	f1-score	support
0	0.65 0.55	0.91	0.76 0.28	412 244
accuracy macro avg weighted avg	0.60 0.62	0.55	0.64 0.52 0.58	656 656 656

```
for i in np.arange(1, 101):
    new model = AdaBoostClassifier(n estimators = i)
    new model.fit(X train, y train)
    new predictions = new model.predict(X test)
    error rate adaboost.append(np.mean(new predictions != y test))
# Create the plot to assist with selecting a K value for the KNN classifier.
plt.plot(error rate adaboost, label = 'Error Rate for Adaboost')
plt.title("Adaboost Selection Plot for Number of Estimators", fontsize=15)
plt.xlabel("Number of Estimators", fontsize=10)
plt.ylabel("Error Rate", fontsize=10)
plt.xticks(fontsize=10)
plt.yticks(fontsize=10)
plt.axvline(x=pd.Series(error rate adaboost).idxmin(), linewidth = 2, color = 'k', linesty
            label = 'Optimized Number of Estimators')
plt.legend(loc = 'upper right')
plt.show()
```

#### Adaboost Selection Plot for Number of Estimators



Optimum Number of Estimators for AdaBoost Classifier:20

```
In [102...
Calculate the accuracy for the model based on training and test data. Also, report the
Precision, Recall, and F1 score for the model predications against the test data.
'''
accuracy_adaboost_train = accuracy_score(y_train, y_predictions_adaboost_train)
accuracy_adaboost_test = accuracy_score(y_test, y_predictions_adaboost_test)
precision_adaboost = precision_score(y_test, y_predictions_adaboost_test)
recall_adaboost = recall_score(y_test, y_predictions_adaboost_test)
f1_adaboost = f1_score(y_test, y_predictions_adaboost_test)
print("Accuracy of AdaBoost Model on training data is:{}".format(accuracy_adaboost_test))
print("Precision of SAdaBoost Model on testing data is:{}".format(precision_adaboost))
print("Recall of AdaBoost Model on testing data is:{}".format(precision_adaboost))
print("F1 Score of AdaBoost Model on testing data is:{}".format(f1_adaboost))
```

Accuracy of AdaBoost Model on testing data is:0.6417682926829268
Precision of SAdaBoostM Model on testing data is:0.5542168674698795
Recall of AdaBoost Model on testing data is:0.1885245901639344
F1 Score of AdaBoost Model on testing data is:0.28134556574923547

# Summary of Findings from Analysis and Model Deployment Recommendation

```
In [104...
         Create a dictionary summarizing the evaluation criteria for the six models.
         summary data = OrderedDict([('Model', ['Logistic Regression', 'KNN', 'Decision Tree', 'Rand
                                      ('Model Accuracy Test', [accuracy lr test, accuracy knn test, ac
                                                             accuracy rf test, accuracy svm test, accu
                                       ('Model Accuracy Training', [accuracy lr train, accuracy knn tr
                                                                 accuracy rf train, accuracy svm trai
                                      ('Model Precision Score', [precision lr, precision knn, precision]
                                                               precision rf, precision svm, precision &
                                      ('Model Recall Score', [recall lr, recall knn, recall dt,
                                                            recall_rf,recall_svm,recall adaboost]),
                                      ('Model F1 Score', [f1 lr,f1 knn,f1 dt,
                                                          f1 rf,f1 svm,f1 adaboost]),])
In [105...
         Display the summary evaluation metrics for the four models in a pandas dataframe.
         Sort the Models based on accuracy for each model from the test dataset.
```

summary df = pd.DataFrame(summary data, index = summary data['Model'])

display(summary df.sort values(by = 'Model Accuracy Test', ascending = False))

	Model	Model_Accuracy_Test	Model_Accuracy_Training	Model_Precision_Score	Model_
SVM	SVM	0.690549	0.737405	0.678261	
Random_Forest	Random_Forest	0.675305	1.000000	0.598726	
KNN	KNN	0.664634	0.658015	0.693548	
Adaboost	Adaboost	0.641768	0.633969	0.554217	
Decision_Tree	Decision_Tree	0.570122	1.000000	0.431655	
Logistic_Regression	Logistic_Regression	0.515244	0.518702	0.377483	

Thus far, the SVM Model provides the highest accuracy on the test dataset. To improve the accuracy, try hyperparameter tuning for these models to see if a significant difference can be acheived. At this stage in the analysis, even the best model (SVM) is not ready to be deployed and only has an accuracy of 69% on unseen data. I'd like to also look into cross-validation for these models to see which model(s) are the most repeatable.