

# PROJECT TITLE : WATER QUALITY ANALYSIS

**Project Definition:** The project involves analyzing water quality data to assess the suitability of water for specific purposes, such as drinking. The objective is to identify potential issues or deviations from regulatory standards and determine water potability based on various parameters. This project includes defining analysis objectives, collecting water quality data, designing relevant visualizations, and building a predictive model.

## **Design Thinking:**

Before diving into the analysis, it's crucial to understand the context and the stakeholders' needs:

1. Identify the stakeholders: Who are the end-users of this analysis (e.g., regulatory bodies, public health agencies, local communities)?

2. Understand their needs and concerns: What are the specific water quality standards and regulations that need to be met? What are the potential health risks associated with poor water quality?

## **ANALYSIS PHASE:**

1. Determine specific goals: Are you primarily focused on assessing water potability, identifying deviations from standards, or both?

2. Define success criteria: What metrics or criteria will be used to measure the success of your analysis?

## **DATA COLLECTION:**

Identify the data sources and collection methods:

1. Identify the relevant parameters: List all the water quality parameters you have access to, such as pH, Hardness, Solids, Chlorine levels, etc.

2. Data collection plan: Determine how you will gather the data, whether it's through field measurements, historical records, or other sources.

3. Data quality assessment: Consider the reliability and completeness of the data. Are there any gaps or inconsistencies that need to be addressed?

## **VISUALIZATION STRATEGY:**

1. Choose visualization tools: Decide which tools or software will be best for creating visualizations (e.g., Python with libraries like Matplotlib and Seaborn).

2. Visualization types: Select appropriate visualization types for parameter distributions, correlations, and potability assessment. For example, histograms, scatter plots, and heatmaps can be useful.

3. Interactive dashboards: Consider creating interactive dashboards for stakeholders to explore the data themselves.

## **PREDICTIVE MODELING:**

1. Feature selection: Determine which water quality parameters are most relevant for predicting potability. This may involve feature engineering to create new variables.

2. Machine learning algorithms: Choose suitable algorithms for classification tasks (since you are predicting potability). Common choices include Decision Trees, Random Forests, Logistic Regression, or even more advanced methods like Neural Networks.

3. Model evaluation: Establish evaluation metrics (e.g., accuracy, precision, recall) to assess the performance of your predictive model.

By following the Design Thinking process, you can ensure that your water quality analysis project is not only technically sound but also addresses the needs and concerns of the stakeholders effectively.



# NAAN MUDHALVAN PROJECT PHASE 2: INNOVATION

PROJECT TITLE: **WATER QUALITY ANALYSIS**

## **DATA ANALYTICS OF WATER QUALITY ANALYSIS**

### VALUABLE INNOVATION STEPS:

#### **STEP1: DATA COLLECTIONS**

- Review the initial design concept to ensure it aligns with the identified problem.
- Gather feedback from stakeholders and subject matter experts for improvements.
- Incorporate necessary changes to enhance the design's effectiveness.

#### **STEP2: CLEANING DATA**

- Clean and reprocess the data to remove outliers, errors, and inconsistencies.
- Ensure data quality before analysis.
- Transformation data into proper format for further processes

#### **STEP3: EVALUATE AND ANALYSIS**

##### Machine Learning Models:

- Train machine learning models to predict water quality based on chemical components present in water.
- This can help pinpoint potential sources of water quality.

##### Cluster Analysis:

- Use clustering algorithms to group similar noise patterns together.
- This can help identify areas with distinct noise characteristics.

#### **STEP4: DATA VISUALIZATION**

- Create interactive maps and visualizations to communicate water quality patterns.
- Finding components present on water such as PH, sodium, carbon, hydrogen etc..

#### **STEP5: DISCRIBE RESULT (COMMUNICATOIN)**

- Result is predicted by using appropriate calculation method using statistic evaluation as per our data collections.
- Finally result of water quality is predicted.

## WATER QUALITY ANALYSIS DESIGN



# da-phase3

November 1, 2023

## 1 WATER QUALITY ANALYSIS

### 2 In this part we will begin building your project by loading

```
[3]: import pandas as pd
file_path = "/content/water_potability.csv"
data = pd.read_csv(file_path)
```

```
[4]: print(data.head())
```

	ph	Hardness	Solids	Chloramines	Sulfate	Conductivity \
0	NaN	204.890455	20791.318981	7.300212	368.516441	564.308654
1	3.716080	129.422921	18630.057858	6.635246	NaN	592.885359
2	8.099124	224.236259	19909.541732	9.275884	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	86.990970	2.963135	0
1	15.180013	56.329076	4.500656	0
2	16.868637	66.420093	3.055934	0
3	18.436524	100.341674	4.628771	0
4	11.558279	31.997993	4.075075	0

```
[6]: selected_columns = data[['ph', 'Hardness', 'Solids', 'Chloramines']]
print(selected_columns)
```

	ph	Hardness	Solids	Chloramines
0	NaN	204.890455	20791.318981	7.300212
1	3.716080	129.422921	18630.057858	6.635246
2	8.099124	224.236259	19909.541732	9.275884
3	8.316766	214.373394	22018.417441	8.059332
4	9.092223	181.101509	17978.986339	6.546600
...	...	...	...	...
3271	4.668102	193.681735	47580.991603	7.166639
3272	7.808856	193.553212	17329.802160	8.061362
3273	9.419510	175.762646	33155.578218	7.350233
3274	5.126763	230.603758	11983.869376	6.303357

```
3275  7.874671  195.102299  17404.177061      7.509306
```

```
[3276 rows x 4 columns]
```

### 3 In this part we begin with the preprocessing of datasets

```
[7]: import pandas as pd
      from sklearn.model_selection import train_test_split
      from sklearn.preprocessing import StandardScaler
      from sklearn.preprocessing import LabelEncoder
```

```
[8]: data = pd.read_csv('/content/water_potability.csv')
```

```
[9]: data['ph'].fillna(data['ph'].mean(), inplace=True)
```

```
[10]: label_encoder = LabelEncoder()
      data['Hardness'] = label_encoder.fit_transform(data['Hardness'])
```

```
[11]: X = data.drop('Potability', axis=1)
      y = data['Potability']
```

```
[12]: X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.2,
      random_state=42)
```

```
[13]: scaler = StandardScaler()
      X_train = scaler.fit_transform(X_train)
      X_test = scaler.transform(X_test)
```

### 4 In this part we begin with the exploratory data analysis

```
[22]: import pandas as pd
      import numpy as np
      import matplotlib.pyplot as plt
      import seaborn as sns
```

```
[17]: data = pd.read_csv('/content/water_potability.csv')
```

```
[18]: print("First 5 rows of the dataset:")
      print(data.head())
```

First 5 rows of the dataset:

	ph	Hardness	Solids	Chloramines	Sulfate	Conductivity \
0	NaN	204.890455	20791.318981	7.300212	368.516441	564.308654
1	3.716080	129.422921	18630.057858	6.635246	NaN	592.885359
2	8.099124	224.236259	19909.541732	9.275884	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
```

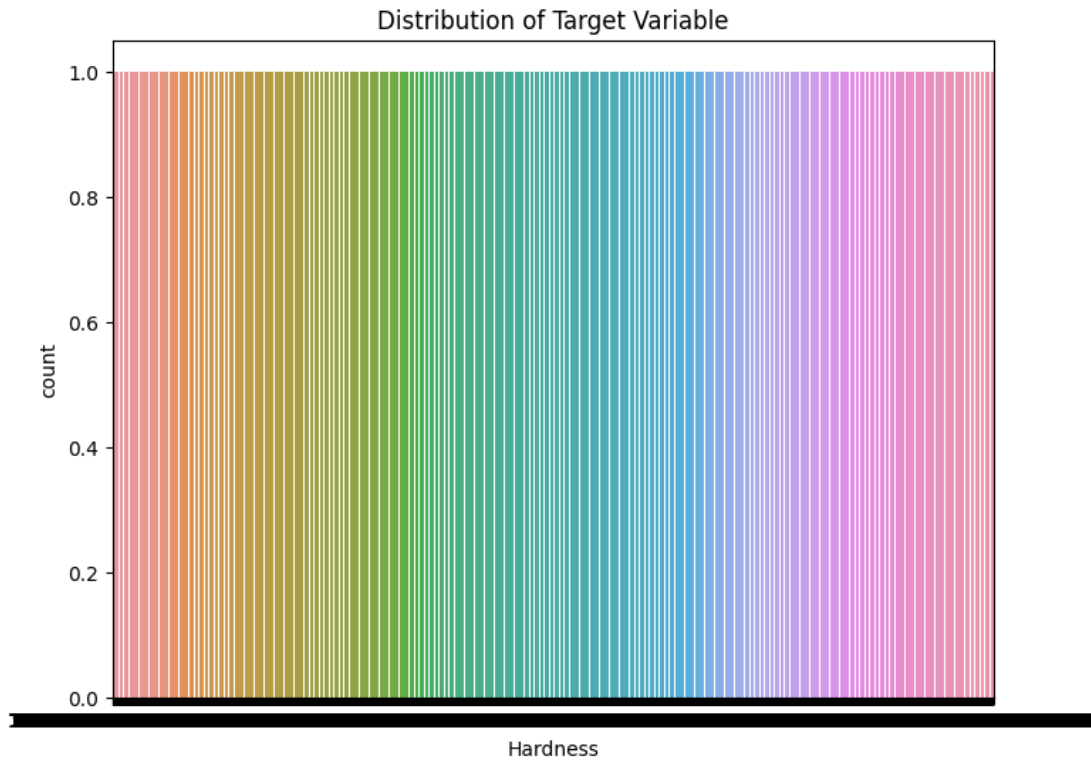
	Organic_carbon	Trihalomethanes	Turbidity	Potability
0	10.379783	86.990970	2.963135	0
1	15.180013	56.329076	4.500656	0
2	16.868637	66.420093	3.055934	0
3	18.436524	100.341674	4.628771	0
4	11.558279	31.997993	4.075075	0

```
[5]: print(data.describe())
```

	ph	Hardness	Solids	Chloramines	Sulfate \
count	2785.000000	3276.000000	3276.000000	3276.000000	2495.000000
mean	7.080795	196.369496	22014.092526	7.122277	333.775777
std	1.594320	32.879761	8768.570828	1.583085	41.416840
min	0.000000	47.432000	320.942611	0.352000	129.000000
25%	6.093092	176.850538	15666.690297	6.127421	307.699498
50%	7.036752	196.967627	20927.833607	7.130299	333.073546
75%	8.062066	216.667456	27332.762127	8.114887	359.950170
max	14.000000	323.124000	61227.196008	13.127000	481.030642

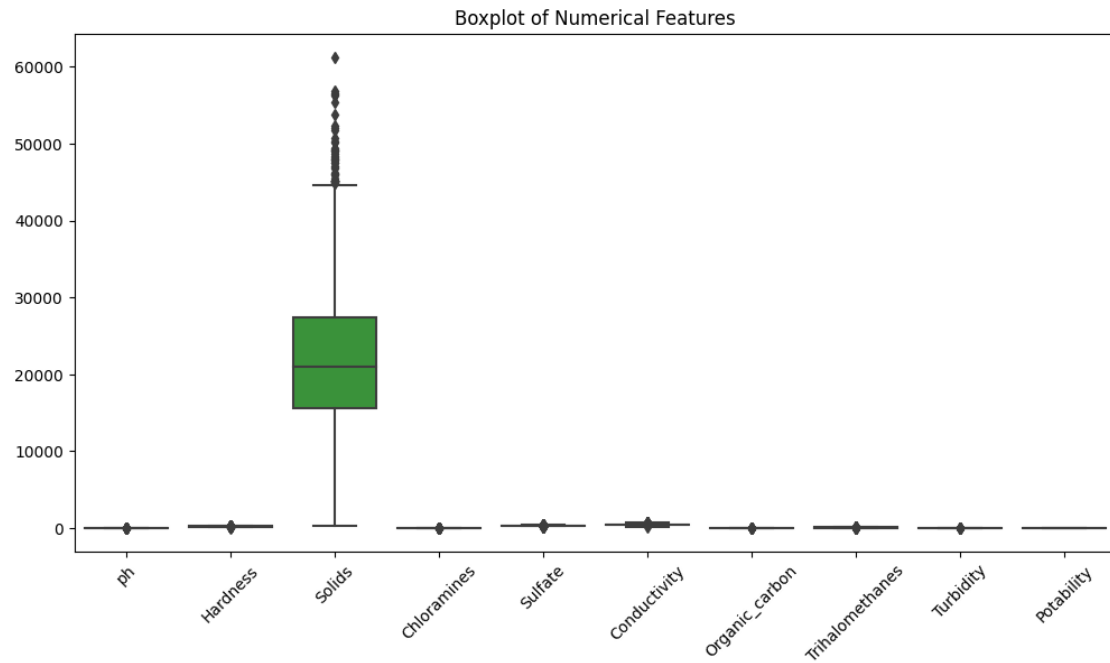
	Conductivity	Organic_carbon	Trihalomethanes	Turbidity	Potability
count	3276.000000	3276.000000	3114.000000	3276.000000	3276.000000
mean	426.205111	14.284970	66.396293	3.966786	0.390110
std	80.824064	3.308162	16.175008	0.780382	0.487849
min	181.483754	2.200000	0.738000	1.450000	0.000000
25%	365.734414	12.065801	55.844536	3.439711	0.000000
50%	421.884968	14.218338	66.622485	3.955028	0.000000
75%	481.792304	16.557652	77.337473	4.500320	1.000000
max	753.342620	28.300000	124.000000	6.739000	1.000000

```
[19]: plt.figure(figsize=(8, 6))
sns.countplot(x='Hardness', data=data)
plt.title('Distribution of Target Variable')
plt.show()
```



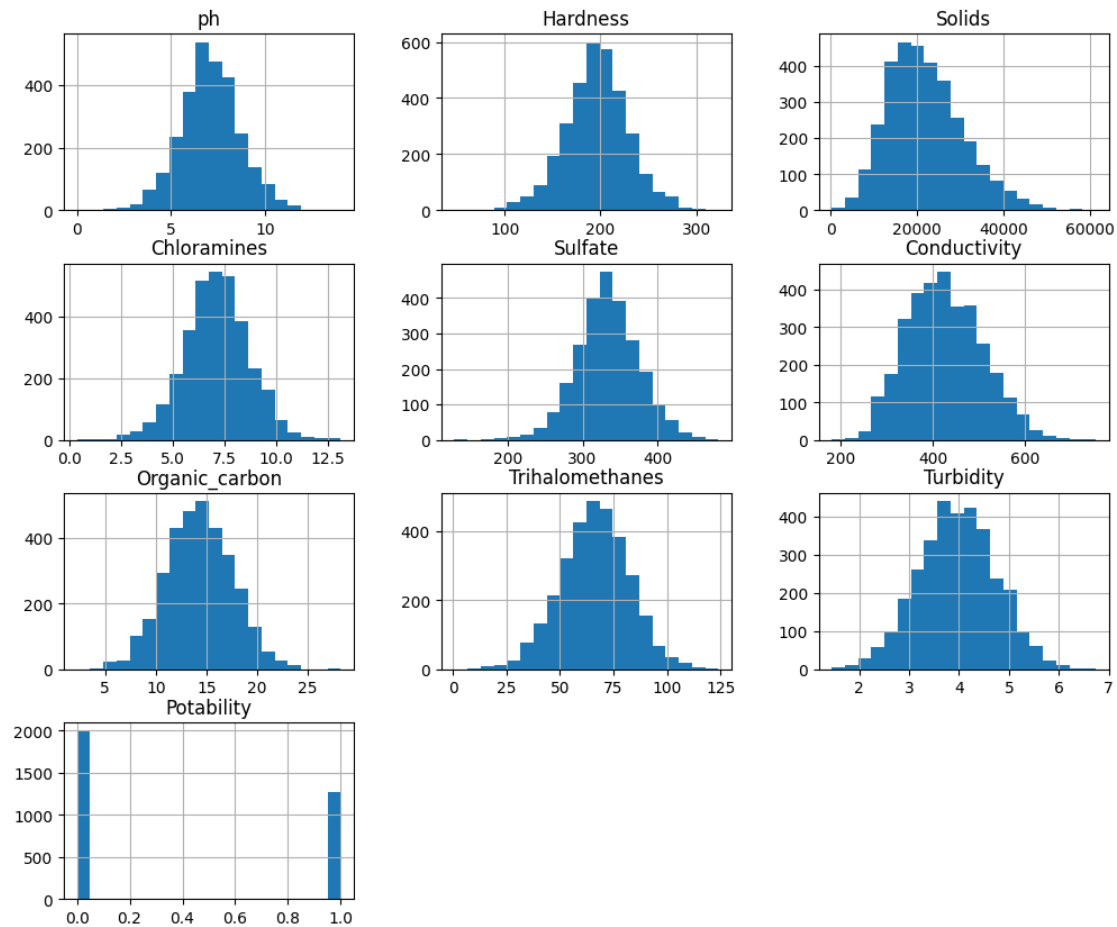
```
[20]: plt.figure(figsize=(12, 6))
sns.boxplot(data=data.select_dtypes(include=['float64', 'int64']))
plt.title('Boxplot of Numerical Features')
plt.xticks(rotation=45)
plt.show()
```





```
[21]: data.select_dtypes(include=['float64', 'int64']).hist(bins=20, figsize=(12, 10))  
plt.suptitle("Histograms of Numerical Features", y=1.02)  
plt.show()
```

## Histograms of Numerical Features



```
[ ]: data.fillna(data.mean(), inplace=True)
```

```
[25]: def handle_outliers_iqr(data, column):
        Q1 = data[column].quantile(0.25)
        Q3 = data[column].quantile(0.75)
        IQR = Q3 - Q1
        lower_bound = Q1 - 1.5 * IQR
        upper_bound = Q3 + 1.5 * IQR
        data = data[(data[column] >= lower_bound) & (data[column] <= upper_bound)]
        return data
```

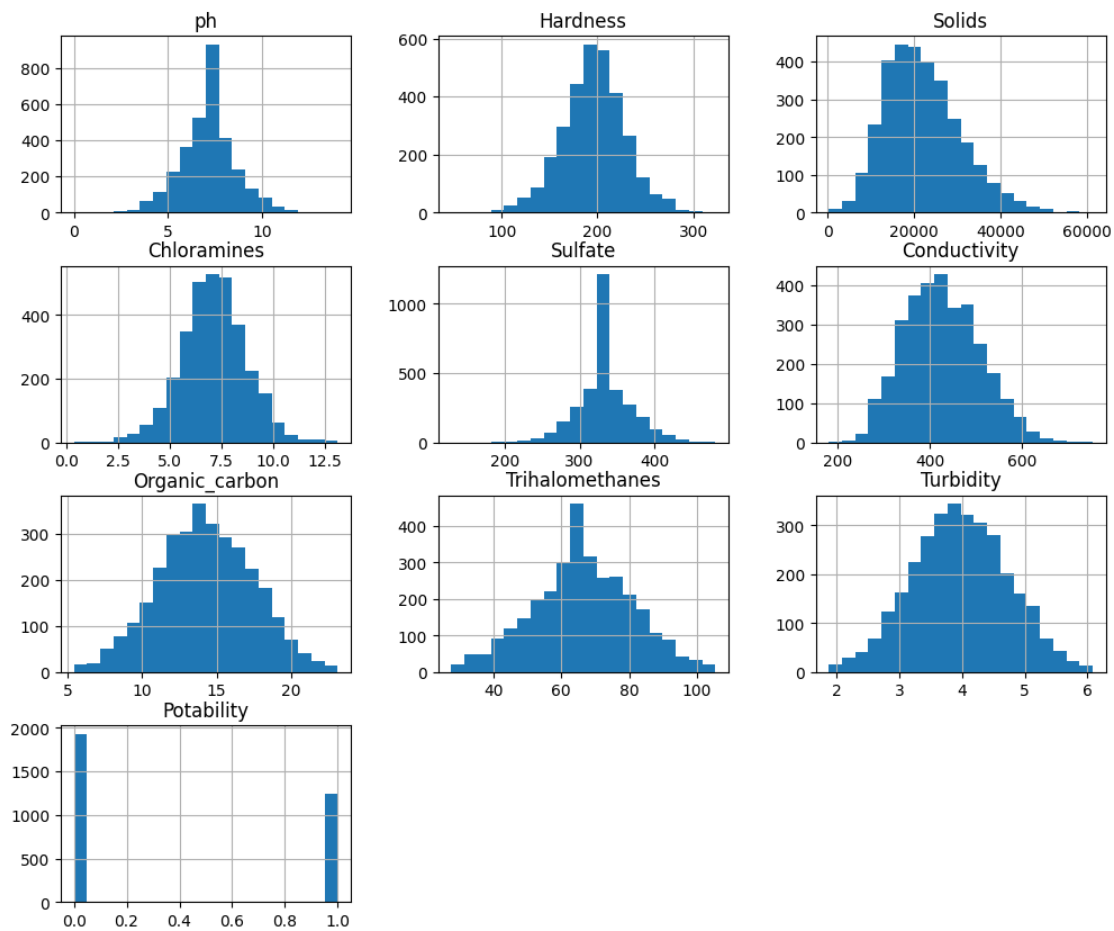
```
[26]: numeric_columns = ['Organic_carbon', 'Trihalomethanes', 'Turbidity']
```

```
[29]: for column in numeric_columns:
      data = handle_outliers_iqr(data, column)
```

## 5 Visualize Parameter Distributions

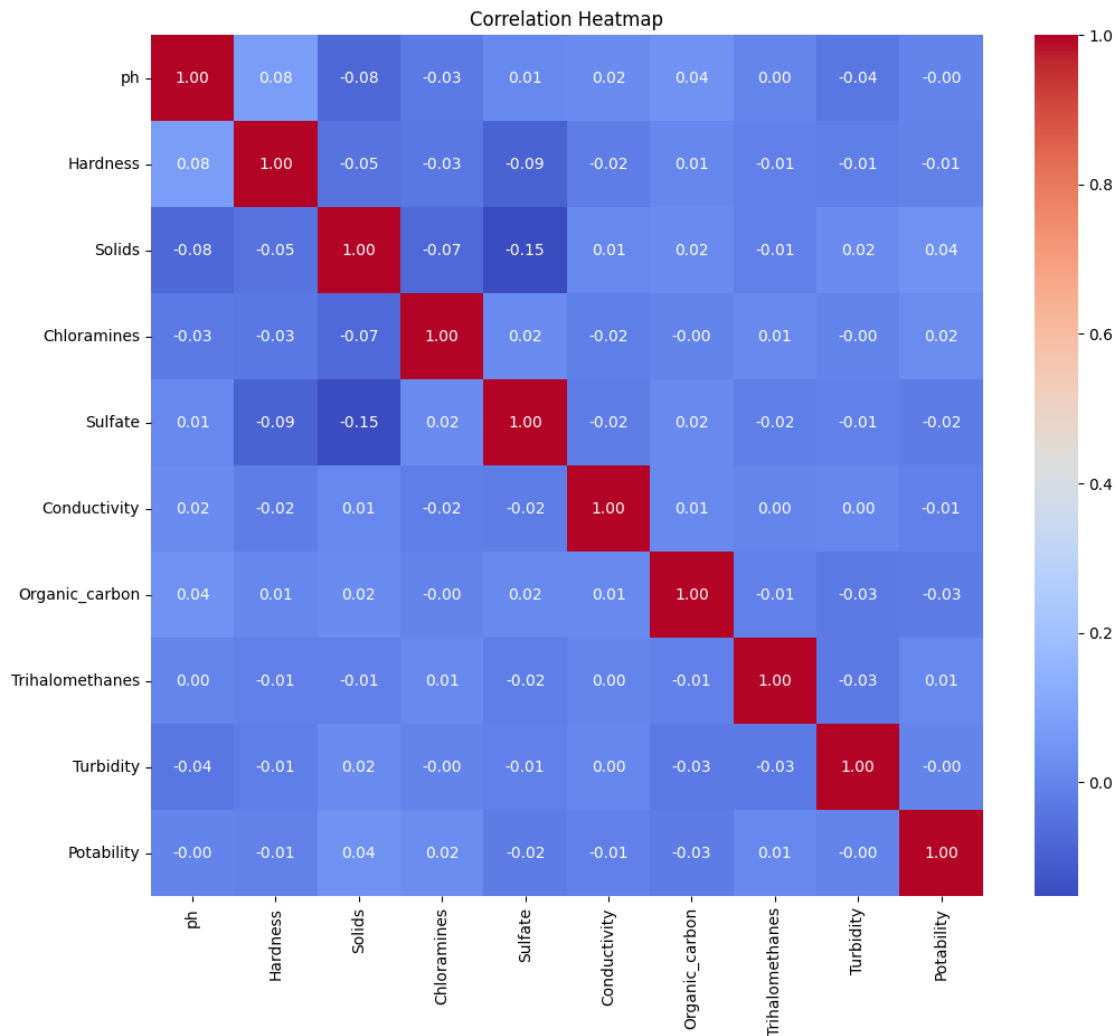
```
[30]: data.select_dtypes(include=['float64', 'int64']).hist(bins=20, figsize=(12, 10))
plt.suptitle("Histograms of Numerical Parameters", y=1.02)
plt.show()
```

Histograms of Numerical Parameters



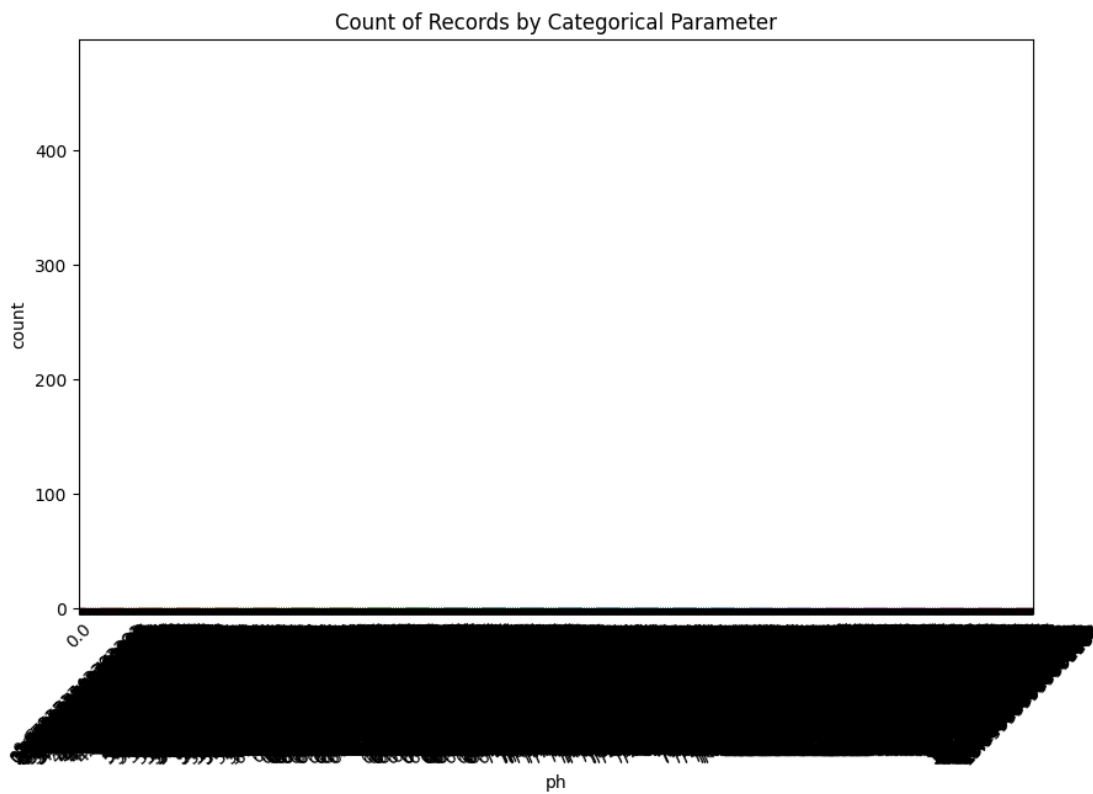
## 6 Visualize Correlations

```
[31]: correlation_matrix = data.corr()
# Plot correlation heatmap
plt.figure(figsize=(12, 10))
sns.heatmap(correlation_matrix, annot=True, cmap='coolwarm', fmt=".2f")
plt.title('Correlation Heatmap')
plt.show()
```



## 7 Identify Potential Deviations from Standards

```
[34]: plt.figure(figsize=(10, 6))
sns.countplot(x='ph', data=data)
plt.title('Count of Records by Categorical Parameter')
plt.xticks(rotation=45)
plt.show()
```



## **PROJECT TITLE: WATER QUALITY ANALYSIS**

### **PHASE 4: CREATING VISUALIZATIONS AND BUILDING A PREDICTIVE MODEL**

In this model we are going to create visualizations of the previous loaded dataset and to building a predictive model.

#### **STEPS FOLLOWED:**

##### **STEP 1 : LOAD THE DATASET**

Loading the given dataset from the source shared, by using the library functions.

##### **STEP 2: HANDLING THE MISSING VALUES**

After loading the dataset we have to identify the missing values by using the functions like `isnull()`

Drop the missing values based upon the nature of the dataset.

##### **STEP 3: CREATING VISUALIZATIONS**

After preprocessing the data, we have to create the visualizations of the given dataset.

To create the visualizations, use the library functions such as the matplotlib, seaborn. These libraries can be used to create histograms, scatter plots

##### **STEP 4: BUILDING A PREDICTIVE MODEL**

The machine learning models such as the logistic regression and the random forest to determine the water portability based upon the water quality parameters.

## **VISUALIZATION OF DATA**

Importing libraries:

Importing necessary libraries for loading the dataset.

```
In [1]: import pandas as pd
import numpy as np
import matplotlib.pyplot as plt
import seaborn as sns
```

Load the dataset:

```
In [2]: data=pd.read_csv("/kaggle/input/water-potability/water_potability.csv")
```

#### **DATA PREPROCESSING:**

Perform data cleaning and preprocessing. This may include handling missing values, converting data types, and ensuring data quality.

Here already the dataset is cleaned and loaded, so no preprocessing is needed.

```
In [3]: data.head()
```

```
Out[3]:
```

	ph	Hardness	Solids	Chloramines	Sulfate	Conductivity	Organic_carbon	Trihalomethanes	Turbidity	Potability
0	NaN	204.890455	20791.318981	7.300212	368.516441	564.308654	10.379783	86.990970	2.963135	0
1	3.716080	129.422921	18630.057858	6.635246	NaN	592.885359	15.180013	56.329076	4.500656	0
2	8.099124	224.236259	19909.541732	9.275884	NaN	418.606213	16.868637	66.420093	3.055934	0
3	8.316766	214.373394	22018.417441	8.059332	356.886136	363.266516	18.436524	100.341674	4.628771	0
4	9.092223	181.101509	17978.986339	6.546600	310.135738	398.410813	11.558279	31.997993	4.075075	0

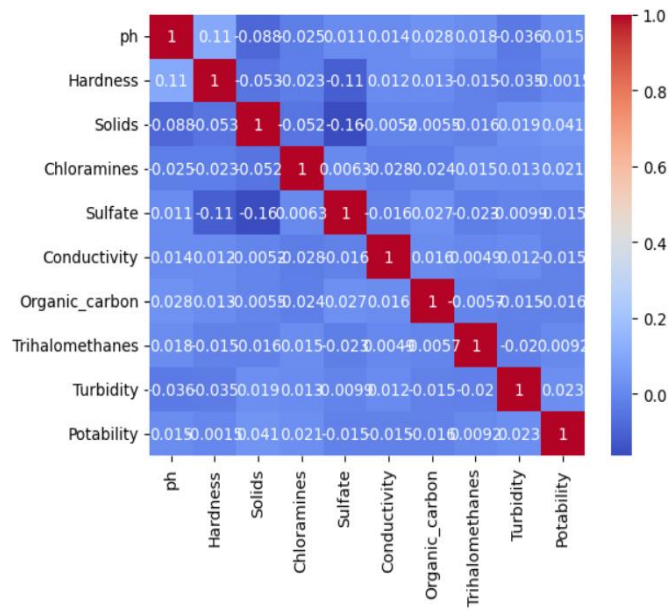
```
In [4]: data.info()
```

```
<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
```

```
In [7]: data.shape
```

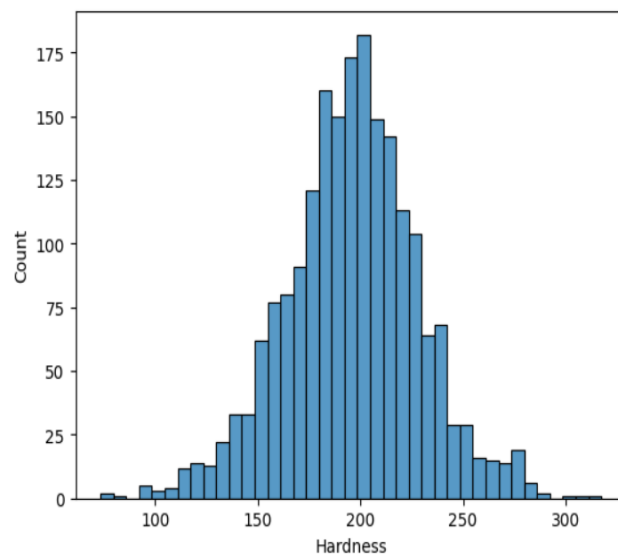
```
Out[7]: (3276, 10)
```

```
In [15]: import seaborn as sns
import matplotlib.pyplot as plt
sns.heatmap(cor,annot=True,cmap='coolwarm')
plt.show()
```



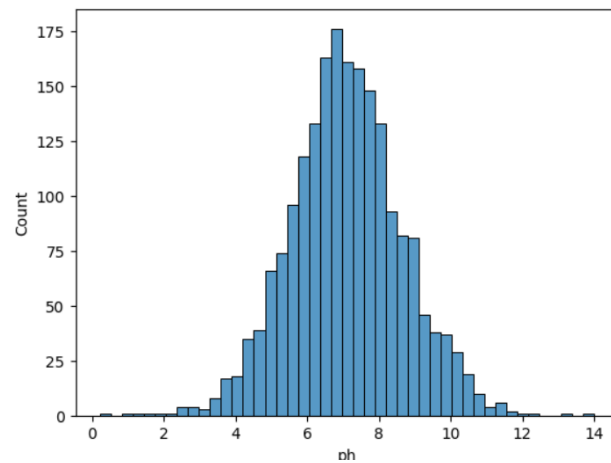
```
In [19]: sns.histplot(data['Hardness'])
```

```
Out[19]: <Axes: xlabel='Hardness', ylabel='Count'>
```



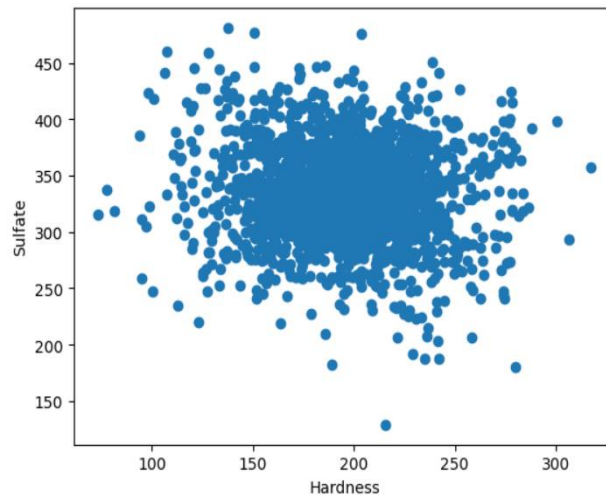
```
In [20]: sns.histplot(data['ph'])
```

```
Out[20]: <Axes: xlabel='ph', ylabel='Count'>
```





```
In [21]: gp = plt.scatter(data['Hardness'],data['Sulfate'])
plt.xlabel('Hardness')
plt.ylabel('Sulfate')
plt.show(gp)
```



## DATA NORMALIZATION AND STANDARDIZATION

```
In [17]: from sklearn.preprocessing import MinMaxScaler,StandardScaler
```

```
normalizer=MinMaxScaler()
standardizer=StandardScaler()
X= normalizer.fit_transform(X)
X=standardizer.fit_transform(X)
```

```
In [18]: from sklearn.model_selection import train_test_split
X_train,X_test,Y_train,Y_test=train_test_split(X,Y,test_size=0.2,random_state=62)
```

## MODEL BUILDING

```
In [19]: from sklearn.linear_model import LogisticRegression
from sklearn.naive_bayes import GaussianNB
from sklearn.svm import SVC
from sklearn.neighbors import KNeighborsClassifier
from sklearn.tree import DecisionTreeClassifier
from sklearn.tree import ExtraTreeClassifier
from sklearn.ensemble import RandomForestClassifier
from sklearn.ensemble import BaggingClassifier
from sklearn.ensemble import GradientBoostingClassifier
from sklearn.ensemble import AdaBoostClassifier
from sklearn.metrics import accuracy_score

models = {
    'Logistic Regression': LogisticRegression(),
    'Naive Bayes': GaussianNB(),
    'Support Vector Machine': SVC(),
    'K-Nearest Neighbors': KNeighborsClassifier(),
    'Decision Tree': DecisionTreeClassifier(),
    'Random Forest': RandomForestClassifier(),
    'Bagging': BaggingClassifier(),
    'AdaBoost': AdaBoostClassifier(),
    'Gradient Boosting': GradientBoostingClassifier(),
    'Extra Trees': ExtraTreeClassifier(),
}
```

```
for name, md in models.items():  
    md.fit(X_train,Y_train)  
    ypred = md.predict(X_test)  
  
    print(f"{name} with accuracy : {accuracy_score(Y_test,ypred)}")
```

Logistic Regression with accuracy : 0.6178660049627791  
Naive Bayes with accuracy : 0.6327543424317618  
Support Vector Machine with accuracy : 0.7245657568238213  
K-Nearest Neighbors with accuracy : 0.6550868486352357  
Decision Tree with accuracy : 0.6104218362282878  
Random Forest with accuracy : 0.7096774193548387  
Bagging with accuracy : 0.6650124069478908  
AdaBoost with accuracy : 0.6004962779156328  
Gradient Boosting with accuracy : 0.6898263027295285  
Extra Trees with accuracy : 0.5806451612903226