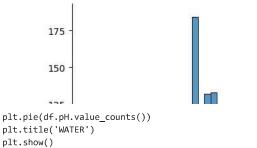
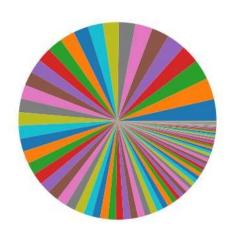
Kashish jain 21BCE3204

```
TASK 1
import pandas as pd
df=pd.read_csv('/content/winequality-red.csv')
df.info()
        <class 'pandas.core.frame.DataFrame'>
       RangeIndex: 1599 entries, 0 to 1598
       Data columns (total 12 columns):
                             Non-Null Count Dtype
         # Column
        0 fixed acidity 1599 non-null float64
1 volatile acidity 1599 non-null float64
2 citric acid 1599 non-null float64
3 residual sugar 1599 non-null float64
4 chlorides 1599 non-null float64
5 free sulfur dioxide 1599 non-null float64
6 total sulfur dioxide 1599 non-null float64
7 dossity 1599 non-null float64
                                                  -----
         7 density 1599 non-null float64
8 pH 1599 non-null float64
9 sulphates 1599 non-null float64
10 alcohol 1599 non-null float64
11 quality 1599 non-null int64
        dtypes: float64(11), int64(1)
       memory usage: 150.0 KB
df.shape
        (1599, 12)
TASK 2
Univariate
import pandas as pd
import numpy as np
import matplotlib.pyplot as plt
import seaborn as sns
sns.displot(df.pH)
```

<seaborn.axisgrid.FacetGrid at 0x7df31745cc40>



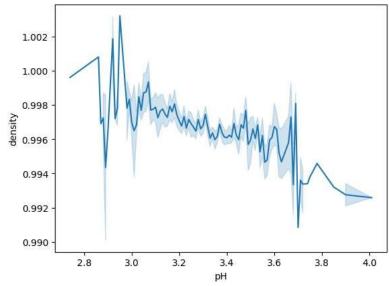
WATER



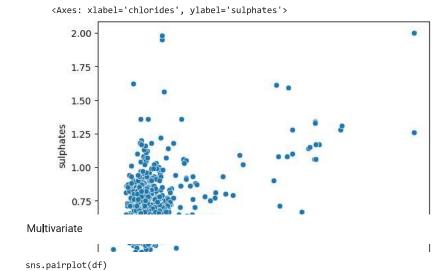
Bivariate

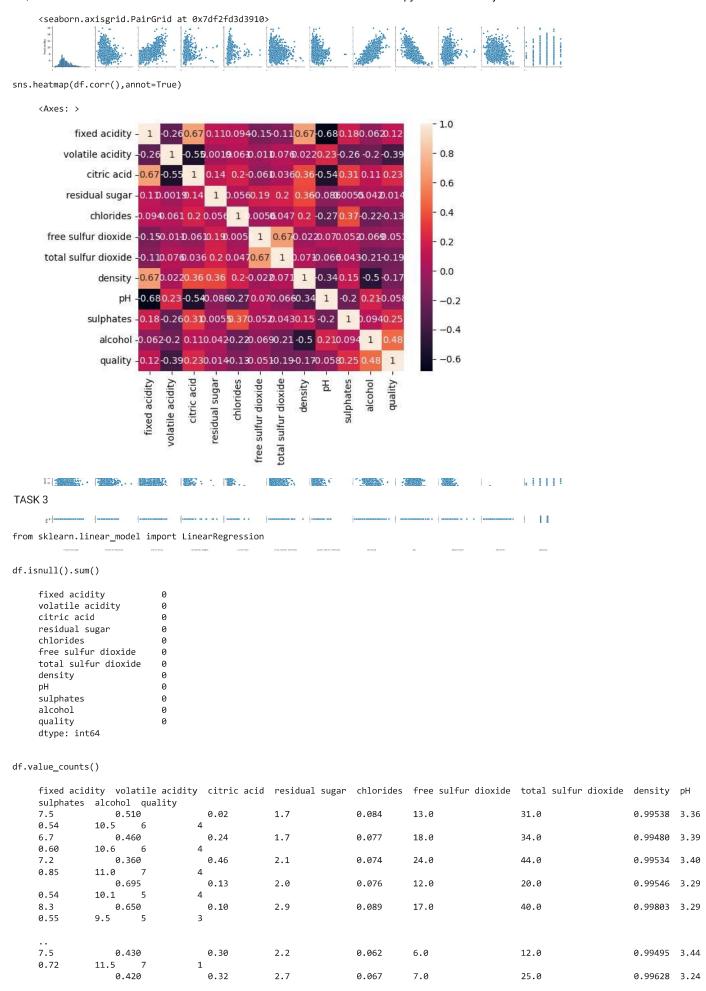
sns.lineplot(x = df.pH,y=df.density)

<Axes: xlabel='pH', ylabel='density'>



sns.scatterplot(x = df.chlorides,y=df.sulphates)





0.44	10.4 5	1					
		0.31	1.6	0.080	15.0	42.0	0.99780 3.31
0.64	9.0 5	1					
	0.410	0.15	3.7	0.104	29.0	94.0	0.99786 3.14
0.58	9.1 5	1					
15.9	0.360	0.65	7.5	0.096	22.0	71.0	0.99760 2.98
0.84	14.9 5	1					
Length: 1	1359, dtype: int64						

df.nunique()

fixed acidity	96
volatile acidity	143
citric acid	80
residual sugar	91
chlorides	153
free sulfur dioxide	60
total sulfur dioxide	144
density	436
pH	89
sulphates	96
alcohol	65
quality	6
dtype: int64	

y = df['pH']
y.head()

0 3.51 1 3.20 2 3.26 3 3.16 4 3.51

Name: pH, dtype: float64

X= df.drop(columns = ['quality'],axis =1)
X.head()

	fixed acidity	volatile acidity	citric acid	residual sugar	chlorides	free sulfur dioxide	total sulfur dioxide	density	рН	sulphates	alcohol
0	7.4	0.70	0.00	1.9	0.076	11.0	34.0	0.9978	3.51	0.56	9.4
1	7.8	0.88	0.00	2.6	0.098	25.0	67.0	0.9968	3.20	0.68	9.8
2	7.8	0.76	0.04	2.3	0.092	15.0	54.0	0.9970	3.26	0.65	9.8
3	11.2	0.28	0.56	1.9	0.075	17.0	60.0	0.9980	3.16	0.58	9.8
4	7.4	0.70	0.00	1.9	0.076	11.0	34.0	0.9978	3.51	0.56	9.4

from sklearn.preprocessing import StandardScaler
scale =StandardScaler()

X_scaled =pd.DataFrame(scale.fit_transform(X),columns = X.columns)
X_scaled.head()

	fixed acidity	volatile acidity	citric acid	residual sugar	chlorides	free sulfur dioxide	total sulfur dioxide	density	рН	sulphates
0	-0.528360	0.961877	-1.391472	-0.453218	-0.243707	-0.466193	-0.379133	0.558274	1.288643	-0.579207
1	-0.298547	1.967442	-1.391472	0.043416	0.223875	0.872638	0.624363	0.028261	-0.719933	0.128950
2	-0.298547	1.297065	-1.186070	-0.169427	0.096353	-0.083669	0.229047	0.134264	-0.331177	-0.048089
3	1.654856	-1.384443	1.484154	-0.453218	-0.264960	0.107592	0.411500	0.664277	-0.979104	-0.461180
4	-0.528360	0.961877	- 1.391472	-0.453218	-0.243707	-0.466193	-0.379133	0.558274	1.288643	-0.579207

from sklearn.model_selection import train_test_split
x_train,x_test,y_train,y_test = train_test_split(X_scaled,y,test_size =0.2,random_state =0)

x_train.shape

(1279, 11)

0.60

-0.75

1.07

-1.40

-0.16

```
x_test.shape
```

x_test.head()

(320, 11)

```
free
                                                                              total
               fixed volatile
                                   citric residual
                                                     chlorides
                                                                   sulfur
                                                                             sulfur
                                                                                      density
                                                                                                      pH sulph
             acidity
                       acidity
                                     acid
                                              sugar
                                                                  dioxide
                                                                            dioxide
      1109
            1.425044 -0.323013 0.816598 -0.311323
                                                       1.775397
                                                                 1.063900
                                                                           0.593954
                                                                                     0.770280 -0.914312
                                                       0.160114 -1.039977 -0.987312
      1032
           -0.126188
                       1.632254 -1.391472
                                           1.107633
                                                                                     0.950485
                                                                                                0.316751
      1002
            0.448342 -1.328579 0.303093 -0.346797
                                                      -0.520005 -0.274931 -0.591995 -0.840962 -0.331177
      487
            1.080326
                       -0.732542 -1.039977 -0.987312
                                                                                     0.770280 -0.914312
      979
            2.229387 -0.434742 1.124700 -0.807957
                                                      -0.264960 -1.231239 -1.230584 0.081262 -1.173483
y_test
     1109
             3.17
     1032
             3.36
     1002
             3.26
     487
             3.17
     979
             3.13
     794
             3.17
     813
             3.44
     1322
             3.18
     704
             3.29
     1023
             3.27
     Name: pH, Length: 320, dtype: float64
from sklearn.linear model import LinearRegression
lr = LinearRegression()
lr.fit(x_train,y_train)
      ▼ LinearRegression
      LinearRegression()
y_predict = lr.predict(x_test)
y_predict
     array([3.17, 3.36, 3.26, 3.17, 3.13, 3.13, 3.12, 3.47, 3.3, 3.27, 3.2,
            3.14, 3.29, 3.74, 3.5, 3.34, 3.33, 3.15, 3.55, 3.38, 3.35, 3.27,
            3.36, 3.3 , 3.57, 3.27, 3.23, 3.35, 3.2 , 3.9 , 3.14, 3.31, 3.53,
            3.5 , 3.52, 3.26, 3.46, 3.69, 3.62, 3.35, 3.33, 3.37, 3.49, 3.02,
            3.3 , 3.19, 3.16, 3.3 , 3.58, 3.36, 3.41, 3.38, 3.38, 3.3 , 3.32,
            3.35,\ 3.3 , 3.09,\ 3.23,\ 3.27,\ 3.68,\ 3.32,\ 3.32,\ 3.56,\ 3.51,\ 3.42,
            3.29, 2.92, 3.19, 3.07, 3.3, 3.04, 3.38, 3.17, 3.34, 3.58, 3.46,
            3.45, 3.35, 3.3 , 3.1 , 3.39, 3.51, 3.14, 3.36, 3.4 , 3.44, 3.25,
            3.42, 3.44, 3.26, 3.22, 3.62, 3.35, 3.4, 3.3, 3.15, 3.19, 3.26,
            3.04, 3.23, 3.38, 3.29, 3.34, 3.46, 3.39, 3.46, 3.41, 3.42, 3.23,
            3.56, 3.1, 3.23, 3.29, 3.33, 3.37, 3.25, 3.27, 3.45, 3.36, 3.23,
            2.92, 3.03, 3.36, 3.27, 3.3 , 3.23, 3.13, 3.36, 3.06, 3.4 , 3.35,
            3.32, 3.22, 3.28, 3.29, 3.06, 3. , 2.88, 3.36, 3.39, 3.19, 3.16,
            3.67, 3.17, 3.3 , 3.23, 3.33, 3.18, 3.34, 3.42, 3.39, 3.41, 3.29,
            3.61, 3.46, 3.25, 3.28, 3.44, 3.38, 3.48, 3.42, 3.32, 3.5, 3.14,
            3.25, 3.54, 3.41, 3.18, 3.27, 3.1, 3.35, 3.48, 3.15, 3.52, 3.21,
            3.18, 3.34, 3.71, 3.16, 3.32, 3.38, 3.41, 3.2, 3.39, 3.2, 3.26,
            3.29, 3.25, 3.42, 3.33, 2.94, 3.18, 3.29, 3.46, 3.28, 3.2, 3.39,
            3.28, 3.16, 3.19, 3.41, 3.26, 3.31, 3.26, 3.59, 3.24, 3.38, 3.36,
            3.42, 3.29, 3.25, 3.11, 3.18, 3.49, 3.47, 3.04, 3.39, 3.38, 3.28,
            3.51, 3.16, 3.2, 3.03, 3.36, 3.12, 3.22, 3.56, 3.16, 3.39, 3.53,
            3.57, 3.35, 3.23, 3.13, 3.42, 3.07, 3.33, 3.36, 3.22, 3.48, 3.4,
            3.36, 3.39, 3.24, 3.26, 3.29, 3.29, 3. , 3.39, 3.19, 3.44, 3.23, 3.26, 3.4 , 3.36, 3.21, 3.11, 3.48, 3.07, 3.44, 3.68, 3.14, 3.08,
            3.34, 3.36, 3.6 , 3.28, 3.28, 3.33, 3.14, 3.17, 3.24, 3.9 , 3.22,
            3.39, 3.14, 3.21, 3.39, 3.05, 3.14, 3.27, 3.17, 3.26, 3.26, 3.55,
```

3.36, 3.27, 3.34, 3.15, 3.49, 3.5, 3.41, 3.19, 3.26, 3.38, 3.4, 3.29, 3.35, 3.15, 3.44, 3.56, 3.12, 3.31, 3.35, 3.61, 3.45, 3.48, 3.37, 3.39, 3.36, 3.42, 3.66, 3.19, 3.2, 3.17, 3.44, 3.18, 3.29,

3.27])

```
y_predict1 =lr.predict(x_train)
y_predict1
    array([3.39, 3.13, 3.26, ..., 3.29, 3.3 , 3.25])
```

profit =pd.DataFrame({'Actual_pH':y_test,'Predicted_pH':y_predict})
profit

	Actual_pH	Predicted_pH	Ħ
1109	3.17	3.17	ılı
1032	3.36	3.36	
1002	3,26	3.26	
487	3.17	3.17	
979	3,13	3.13	
794	3.17	3.17	
813	3.44	3.44	
1322	3.18	3.18	
704	3.29	3.29	
1023	3.27	3.27	

320 rows × 2 columns

Task 4

TASK 5

df.head()

	fixed acidity	volatile acidity	citric acid	residual sugar	chlorides	free sulfur dioxide	total sulfur dioxide	density	рН	sulph
0	7.4	0.70	0.00	1.9	0.076	11.0	34.0	0.9978	3.51	
1	7.8	0.88	0.00	2.6	0.098	25.0	67.0	0.9968	3.20	
2	7.8	0.76	0.04	2.3	0.092	15.0	54.0	0.9970	3.26	
3	11,2	0.28	0.56	1.9	0.075	17.0	60.0	0,9980	3.16	
4										-

 $lr.predict([[7.4,\ 0.70,\ 0.00,\ 1.9,\ \ 0.076,\ \ 11.0,\ 34.0,\ 0.9978,\ 3.51,\ 0.56,\ 9.4]])/100000$

/usr/local/lib/python3.10/dist-packages/sklearn/base.py:439: UserWarning: X does not have valid feature names, but LinearRegression was

+ Text

array([3.85284021e-05])

```
lr.predict([[7.8, 0.88, 0.00, 2.6, 0.098, 25.0, 67.0, 0.9968, 3.20, 0.68, 9.8]])/100000
```

/usr/local/lib/python3.10/dist-packages/sklearn/base.py:439: UserWarning: X does not have valid feature names, but LinearRegression was warnings.warn(array([3.80499538e-05])

lr.predict([[7.4, 0.70, 0.00, 1.9, 0.076, 11.0, 34.0, 0.9978, 3.51, 0.56, 9.4]])/100000

/usr/local/lib/python3.10/dist-packages/sklearn/base.py:439: UserWarning: X does not have valid feature names, but LinearRegression was warnings.warn(

array([3.85284021e-05])