```
import numpy as np
In [159...
          import pandas as pd
          import matplotlib.pyplot as plt
          import seaborn as sns
          import scipy
          import statsmodels as stm
          from sklearn.datasets import load_breast_cancer
          from sklearn.neighbors import KDTree, BallTree, KNeighborsClassifier, RadiusNeighbor
          %matplotlib inline
```

Before we work with KNN and other distance dependent algorithms lets understand the various distance metrics. My motive in understanding the distances because there is no one metric that can be used for every problem.

```
CASE-I: Entire variable as vectors
         X = np.array([[1,2,3]])
In [3]:
         Y = np.array([[3,4,5]])
         X, Y
Out[3]: (array([[1, 2, 3]]), array([[3, 4, 5]]))
         X.shape, Y.shape, X.ndim, Y.ndim
In [4]:
Out[4]: ((1, 3), (1, 3), 2, 2)
In [5]:
         eu_dist = DistanceMetric.get_metric('euclidean')
         man_hat_dist = DistanceMetric.get_metric('manhattan')
In [6]:
In [7]:
         eu dist.pairwise(X,Y)
Out[7]: array([[3.46410162]])
In [8]:
         man hat dist.pairwise(X,Y)
Out[8]: array([[6.]])
        CASE-II: Different shape vectors
         XX = X.copy().reshape(3,1)
In [9]:
         YY = Y.copy().reshape(3,1)
         XX,YY
```

```
In [10]:
Out[10]: (array([[1],
                  [2],
                  [3]]),
           array([[3],
                  [4],
                  [5]]))
          eu_dist.pairwise(XX,YY)
In [12]:
Out[12]: array([[2., 3., 4.],
                 [1., 2., 3.],
                 [0., 1., 2.]])
```

localhost:8888/lab 1/13

Out[44]:

```
In [13]:
          man hat dist.pairwise(XX,YY)
Out[13]: array([[2., 3., 4.],
                 [1., 2., 3.],
                 [0., 1., 2.]])
         CASE-III
           shoe_size = np.array([7,8,9,10,6,5,7])
In [23]:
           body_mass = np.array([23,24,29,35,21,20,25])
           shoe_size, body_mass, shoe_size.shape, shoe_size.ndim, body_mass.shape, body_mass.nd
          (array([7, 8, 9, 10, 6, 5, 7]),
array([23, 24, 29, 35, 21, 20, 25]),
Out[23]:
           (7,),
           1,
           (7,),
           1)
          sns.scatterplot(x=shoe_size,y=body_mass)
In [26]:
          <AxesSubplot:>
Out[26]:
          34
          32
          30
          28
          26
          24
          22
          20
               5
                                          8
                                                   9
                                                            10
          [shoe_size[5],body_mass[5]], [shoe_size[4],body_mass[4]]
In [35]:
          ([5, 20], [6, 21])
Out[35]:
           eu_dist.pairwise([[shoe_size[5],body_mass[5]]], [[shoe_size[4],body_mass[4]]])
In [37]:
         array([[1.41421356]])
Out[37]:
In [38]:
          man_hat_dist.pairwise([[shoe_size[5],body_mass[5]]], [[shoe_size[4],body_mass[4]]])
Out[38]: array([[2.]])
         CASE-IV
           chby_dist = DistanceMetric.get_metric('chebyshev')
In [40]:
In [42]:
           chby_dist.pairwise([[shoe_size[5],body_mass[5]]], [[shoe_size[4],body_mass[4]]])
          array([[1.]])
Out[42]:
In [44]:
           chby_dist.pairwise(shoe_size.reshape((1,7)),body_mass.reshape((1,7)))
          array([[25.]])
```

localhost:8888/lab 2/13

```
np.square(np.corrcoef(shoe size,body mass))
In [47]:
                           , 0.8792535],
Out[47]: array([[1.
                [0.8792535, 1.
         CASE-V
          filepath = 'https://raw.githubusercontent.com/selva86/datasets/master/diamonds.csv'
In [327...
          df = pd.read_csv(filepath).iloc[0:500, [0,4,6]]
          df.shape, df.ndim, df.head()
Out[327... ((500, 3),
          2,
             carat depth price
              0.23
                     61.5
                              326
          a
              0.21
                     59.8
                              326
          1
          2
              0.23
                     56.9
                              327
          3
              0.29
                     62.4
                              334
              0.31
                     63.3
                              335)
          def mahalanobis(x=None, data=None, cov=None):
In [161...
              """Compute the Mahalanobis Distance between each row of x and the data
                   : vector or matrix of data with, say, p columns.
              data : ndarray of the distribution from which Mahalanobis distance of each obser
              cov : covariance matrix (p x p) of the distribution. If None, will be computed
              x_{minus_mu} = x - np.mean(data)
              if not cov:
                  cov = np.cov(data.values.T)
              inv_covmat = scipy.linalg.inv(cov)
              left_term = np.dot(x_minus_mu, inv_covmat)
              mahal = np.dot(left_term, x_minus_mu.T)
              return mahal.diagonal(), inv_covmat
          df_x = df[['carat', 'depth', 'price']].head(500)
          df_x['mahala'],VI = mahalanobis(x=df_x, data=df[['carat', 'depth', 'price']])
```

#### Out[161... carat depth price mahala 0.23 0 61.5 326 3.766898 1 0.21 59.8 326 5.040171 2 0.23 56.9 327 11.473766 3 0.29 62.4 334 3.876022 0.31 63.3 335 4.661105

df\_x.head()

## **Understanding the above function**

```
In [166... x_minus_mu = df - np.mean(df)
    cov = np.cov(df.values.T)
    inv_covmat = scipy.linalg.inv(cov)
    left_term = np.dot(x_minus_mu, inv_covmat)
    mahal = np.dot(left_term, x_minus_mu.T)
```

#### STEP-1

localhost:8888/lab 3/13

```
price 2233.49800
dtype: float64
```

```
In [181...
          x_minus_mu.head()
Out[181... carat
                      depth
                                 price
          0 -0.40612 -0.2128 -1907.498
          1 -0.42612 -1.9128 -1907.498
          2 -0.40612 -4.8128 -1906.498
          3 -0.34612  0.6872  -1899.498
          4 -0.32612 1.5872 -1898.498
In [184... | pd.DataFrame(df - np.mean(df)).head()
Out[184...
                      depth
           carat
                                 price
          0 -0.40612 -0.2128 -1907.498
          1 -0.42612 -1.9128 -1907.498
            -0.40612 -4.8128 -1906.498
          3 -0.34612  0.6872  -1899.498
          4 -0.32612 1.5872 -1898.498
         STEP-2
          pd.DataFrame(cov)
In [186...
Out[186...
                                            2
          0
              0.046833 0.022599
                                 197.305744
              0.022599 2.983724
                                    -5.602980
          2 197.305744 -5.602980 993181.561118
          df.cov().values
In [187...
Out[187... array([[ 4.68330116e-02, 2.25988617e-02, 1.97305744e+02],
                 [ 2.25988617e-02, 2.98372361e+00, -5.60298036e+00],
                 [ 1.97305744e+02, -5.60298036e+00, 9.93181561e+05]])
In [173...
          df.corr()
Out[173...
                            depth
                                       price
                    carat
           carat 1.000000
                          0.060455
                                    0.914849
          depth 0.060455
                          1.000000
                                   -0.003255
           price 0.914849 -0.003255
                                   1.000000
         np.corrcoef(df,rowvar=False)
In [175...
                              , 0.06045486, 0.91484888],
Out[175... array([[ 1.
                 [ 0.06045486, 1. , -0.00325481],
                 [ 0.91484888, -0.00325481, 1.
                                                          ]])
```

localhost:8888/lab 4/13

```
scipy.linalg.inv(df.cov().values)
In [189...
           array([[ 1.34268725e+02, -1.06705836e+00, -2.66798846e-02],
Out[189...
                    [-1.06705836e+00, 3.43635347e-01,
                                                             2.13920731e-04],
                   [-2.66798846e-02, 2.13920731e-04,
                                                             6.30830586e-06]])
In [177...
            inv_covmat
           array([[ 1.34268725e+02, -1.06705836e+00, -2.66798846e-02],
Out[177...
                    [-1.06705836e+00, 3.43635347e-01,
                                                              2.13920731e-04],
                   [-2.66798846e-02, 2.13920731e-04,
                                                             6.30830586e-06]])
          STEP-3
            pd.DataFrame(((df - np.mean(df)).values @ scipy.linalg.inv(df.cov().values)) @ ((df
In [203...
           (500, 500)
Out[203...
In [204..
            pd.DataFrame(((df - np.mean(df)).values @ scipy.linalg.inv(df.cov().values)) @ ((df
Out[204...
                        0
                                   1
                                              2
                                                         3
                                                                              5
                                                                                         6
                                                                                                    7
                                                                                                               8
             0
                 3.766898
                            3.916408
                                       3.985651
                                                  3.509290
                                                             3.396797
                                                                        3.658189
                                                                                  3.682101
                                                                                             3.631782
                                                                                                        3.615154
             1
                 3.916408
                            5.040171
                                       6.724389
                                                  3.101321
                                                             2.465016
                                                                        3.068993
                                                                                  3.374325
                                                                                             3.531883
                                                                                                        1.749026
             2
                 3.985651
                            6.724389
                                      11.473766
                                                  2.590670
                                                             1.152378
                                                                        1.861321
                                                                                  2.675488
                                                                                             3.354030
                                                                                                       -1.915498
             3
                 3.509290
                            3.101321
                                       2.590670
                                                  3.876022
                                                             4.122082
                                                                        3.776873
                                                                                  3.677306
                                                                                             3.664491
                                                                                                        4.162841
              4
                 3.396797
                            2.465016
                                       1.152378
                                                  4.122082
                                                             4.661105
                                                                        4.052604
                                                                                  3.808965
                                                                                             3.714524
                                                                                                        5.067003
           495
                -1.030892
                           -1.550898
                                       -2.805898
                                                 -1.077694
                                                            -0.864771
                                                                       -0.577531
                                                                                 -0.770679
                                                                                            -1.059940
                                                                                                        0.449597
                 -0.963937
                           -0.695969
                                                 -1.356482
                                                            -1.546962
                                                                       -1.124048
                                                                                 -1.076651
           496
                                       -0.526231
                                                                                            -1.143898
                                                                                                       -1.233603
           497
                 -0.877226
                            0.018579
                                       1.250222
                                                 -1.610250
                                                            -2.134672
                                                                       -1.516238
                                                                                 -1.285177
                                                                                            -1.208840
                                                                                                       -2.467239
           498
                 -0.901139
                           -0.286753
                                       0.436055
                                                 -1.510683
                                                            -1.891033
                                                                       -1.321054
                                                                                 -1.175901
                                                                                            -1.178855
                                                                                                       -1.866096
           499
                -1.741025 -3.158284
                                       -4.643153 -0.159136
                                                             0.751059
                                                                      -0.777156 -1.092296
                                                                                           -1.001465
                                                                                                        0.323868
          500 rows × 500 columns
          STEP-4
In [171...
            mahal.shape
Out[171...
           (500, 500)
In [170...
            pd.DataFrame(mahal)
                                   1
                                              2
                                                                    4
                                                                              5
                                                                                         6
                                                                                                    7
                                                                                                               3
Out[170...
                        0
                                                         3
             0
                 3.766898
                            3.916408
                                       3.985651
                                                  3.509290
                                                             3.396797
                                                                        3.658189
                                                                                  3.682101
                                                                                             3.631782
                                                                                                        3.615154
              1
                 3.916408
                            5.040171
                                       6.724389
                                                  3.101321
                                                             2.465016
                                                                        3.068993
                                                                                  3.374325
                                                                                             3.531883
                                                                                                        1.749026
             2
                 3.985651
                            6.724389
                                      11.473766
                                                  2.590670
                                                             1.152378
                                                                        1.861321
                                                                                  2.675488
                                                                                             3.354030
                                                                                                       -1.915498
             3
                 3.509290
                            3.101321
                                       2.590670
                                                  3.876022
                                                             4.122082
                                                                        3.776873
                                                                                  3.677306
                                                                                             3.664491
                                                                                                        4.162841
                 3.396797
                            2.465016
                                       1.152378
                                                  4.122082
                                                             4.661105
                                                                       4.052604
                                                                                  3.808965
                                                                                             3.714524
                                                                                                        5.067003
```

localhost:8888/lab 5/13

3/15/2021 Revision\_14\_03\_21

	0	1	2	3	4	5	6	7	8		
•••								•••			
495	-1.030892	-1.550898	-2.805898	-1.077694	-0.864771	-0.577531	-0.770679	-1.059940	0.449597		
496	-0.963937	-0.695969	-0.526231	-1.356482	-1.546962	-1.124048	-1.076651	-1.143898	-1.233603		
497	-0.877226	0.018579	1.250222	-1.610250	-2.134672	-1.516238	-1.285177	-1.208840	-2.467239		
498	-0.901139	-0.286753	0.436055	-1.510683	-1.891033	-1.321054	-1.175901	-1.178855	-1.866096		
499	-1.741025	-3.158284	-4.643153	-0.159136	0.751059	-0.777156	-1.092296	-1.001465	0.323868		
500 r	500 rows × 500 columns										

# So, good here as everything matched!!

#### **Critical Value**

```
In [154... from scipy.stats import chi2
chi2.ppf((1-0.01), df=2)
```

Out[154... 9.21034037197618

#### **P-Value**

```
In [156... 1 - chi2.cdf(12.7,df=2)
```

Out[156... 0.0017467471362611064

## **Applying Mahalanobis Distance Metric on Breast Cancer**

Out[223...

٠		mean radius	mean texture	mean perimeter	mean area	mean smoothness	mean compactness	mean concavity	concave points	mean symmetry	diı
	0	17.99	10.38	122.80	1001.0	0.11840	0.27760	0.3001	0.14710	0.2419	
	1	20.57	17.77	132.90	1326.0	0.08474	0.07864	0.0869	0.07017	0.1812	
	2	19.69	21.25	130.00	1203.0	0.10960	0.15990	0.1974	0.12790	0.2069	
	3	11.42	20.38	77.58	386.1	0.14250	0.28390	0.2414	0.10520	0.2597	
	4	20.29	14.34	135.10	1297.0	0.10030	0.13280	0.1980	0.10430	0.1809	

5 rows × 30 columns

```
In [224... y = pd.DataFrame(cancer.target,columns=['Label'])
y.head()
```

Out[224... Label 0 0

localhost:8888/lab 6/13

```
      Label

      1
      0

      2
      0

      3
      0

      4
      0
```

```
In [226... cancer_df = pd.concat([cancer_df,y],axis=1).copy(deep=True)
    cancer_df
```

Out[226...

	mean radius	mean texture	mean perimeter	mean area	mean smoothness	mean compactness	mean concavity	mean concave points	mean symmetry
0	17.99	10.38	122.80	1001.0	0.11840	0.27760	0.30010	0.14710	0.2419
1	20.57	17.77	132.90	1326.0	0.08474	0.07864	0.08690	0.07017	0.1812
2	19.69	21.25	130.00	1203.0	0.10960	0.15990	0.19740	0.12790	0.2069
3	11.42	20.38	77.58	386.1	0.14250	0.28390	0.24140	0.10520	0.2597
4	20.29	14.34	135.10	1297.0	0.10030	0.13280	0.19800	0.10430	0.1809
•••									
564	21.56	22.39	142.00	1479.0	0.11100	0.11590	0.24390	0.13890	0.1726
565	20.13	28.25	131.20	1261.0	0.09780	0.10340	0.14400	0.09791	0.1752
566	16.60	28.08	108.30	858.1	0.08455	0.10230	0.09251	0.05302	0.1590
567	20.60	29.33	140.10	1265.0	0.11780	0.27700	0.35140	0.15200	0.2397
568	7.76	24.54	47.92	181.0	0.05263	0.04362	0.00000	0.00000	0.1587

569 rows × 31 columns

# Seggregating the positive and negative observations

#### Only positivies

localhost:8888/lab 7/13

Out[282...

	mean radius	mean texture	mean perimeter	mean area	mean smoothness	mean compactness	mean concavity	mean concave points	mean symmetry
525	8.571	13.10	54.53	221.3	0.10360	0.07632	0.025650	0.015100	0.1678
58	13.050	19.31	82.61	527.2	0.08060	0.03789	0.000692	0.004167	0.1819
342	11.060	14.96	71.49	373.9	0.10330	0.09097	0.053970	0.033410	0.1776
413	14.990	22.11	97.53	693.7	0.08515	0.10250	0.068590	0.038760	0.1944
559	11.510	23.93	74.52	403.5	0.09261	0.10210	0.111200	0.041050	0.1388

5 rows × 31 columns

Only negatives

Only negatives

Out[283... (127, 31)

In [284... train\_df\_neg.head()

Out[284...

		mean radius	mean texture	mean perimeter	mean area	mean smoothness	mean compactness	mean concavity	mean concave points	mean symmetry	(
	39	13.48	20.82	88.40	559.2	0.10160	0.12550	0.1063	0.05439	0.1720	
2	229	12.83	22.33	85.26	503.2	0.10880	0.17990	0.1695	0.06861	0.2123	
	40	13.44	21.58	86.18	563.0	0.08162	0.06031	0.0311	0.02031	0.1784	
1	172	15.46	11.89	102.50	736.9	0.12570	0.15550	0.2032	0.10970	0.1966	
2	213	17.42	25.56	114.50	948.0	0.10060	0.11460	0.1682	0.06597	0.1308	

5 rows × 31 columns

Distance calculation for positive observations

In [286... train\_df\_pos['mahal\_dist'] = train\_df\_pos\_mh\_dist.diagonal()

In [287... train\_df\_pos.head()

Out[287...

_		mean radius	mean texture	mean perimeter	mean area	mean smoothness	mean compactness	mean concavity	mean concave points	mean symmetry	
	525	8.571	13.10	54.53	221.3	0.10360	0.07632	0.025650	0.015100	0.1678	
	58	13.050	19.31	82.61	527.2	0.08060	0.03789	0.000692	0.004167	0.1819	
	342	11.060	14.96	71.49	373.9	0.10330	0.09097	0.053970	0.033410	0.1776	

localhost:8888/lab 8/13

	mean radius	mean texture	mean perimeter	mean area	mean smoothness	mean compactness	mean concavity	mean concave points	mean symmetry
413	14.990	22.11	97.53	693.7	0.08515	0.10250	0.068590	0.038760	0.1944
559	11.510	23.93	74.52	403.5	0.09261	0.10210	0.111200	0.041050	0.1388

5 rows × 32 columns

#### Distance calculation for negative observations

Out[290...

••		mean radius	mean texture	mean perimeter	mean area	mean smoothness	mean compactness	mean concavity	mean concave points	mean symmetry	(
	39	13.48	20.82	88.40	559.2	0.10160	0.12550	0.1063	0.05439	0.1720	
	229	12.83	22.33	85.26	503.2	0.10880	0.17990	0.1695	0.06861	0.2123	
	40	13.44	21.58	86.18	563.0	0.08162	0.06031	0.0311	0.02031	0.1784	
	172	15.46	11.89	102.50	736.9	0.12570	0.15550	0.2032	0.10970	0.1966	
	213	17.42	25.56	114.50	948.0	0.10060	0.11460	0.1682	0.06597	0.1308	

5 rows × 32 columns

**Hypothesis Testing** 

#### **CASE-I**

Here, first I'll perform the test on positive observations considering the null hypothesis as TRUE or POSITIVE and the alternate hypothesis as FALSE or negative.

localhost:8888/lab 9/13

mean

Out[300...

```
mean
               mean
                           mean
                                  mean
                                                mean
                                                              mean
                                                                         mean
                                                                                               mean
                                                                                 concave
      radius
                      perimeter
                                          smoothness
                                                      compactness
              texture
                                    area
                                                                     concavity
                                                                                          symmetry
                                                                                  points
525
       8.571
                13.10
                           54.53
                                   221.3
                                              0.10360
                                                            0.07632
                                                                      0.025650
                                                                                0.015100
                                                                                              0.1678
 58 13.050
                19.31
                           82.61
                                   527.2
                                              0.08060
                                                            0.03789
                                                                      0.000692 0.004167
                                                                                              0.1819
342 11.060
                14.96
                           71.49
                                   373.9
                                              0.10330
                                                            0.09097
                                                                      0.053970
                                                                                0.033410
                                                                                              0.1776
413 14.990
                22.11
                           97.53
                                   693.7
                                              0.08515
                                                            0.10250
                                                                      0.068590
                                                                                0.038760
                                                                                              0.1944
559 11.510
                23.93
                           74.52
                                  403.5
                                              0.09261
                                                            0.10210
                                                                      0.111200 0.041050
                                                                                              0.1388
5 rows × 33 columns
 train df pos.shape
 train_df_pos[train_df_pos['p_val'] < 0.01].shape</pre>
```

```
In [303...
Out[303... (214, 33)
          (26, 33)
Out[304...
          rej_null_hyp = train_df_pos[train_df_pos['p_val'] < 0.01]['mahal_dist'].values</pre>
In [308...
          rej_null_hyp
         array([ 81.02384852,
                                 72.07114302,
                                                52.08522329,
                                                              62.90183841,
Out[308...
                  69.0939475 ,
                                66.10649304,
                                               66.76169799,
                                                              76.36211933,
                 118.79639287, 145.94310336,
                                               59.52453926,
                                                              73.90722389,
                  68.33222003,
                                96.34523381,
                                               54.67357222, 154.13771741,
                  76.20390925, 100.12061258,
                                               74.13167215,
                                                              58.07212966,
                                 54.58637054, 129.41505176,
                 194.52697533,
                                                              60.14696292,
                  89.70806506,
                                 54.60903334])
          rej_null_hyp.min(), rej_null_hyp.max()
In [311..
Out[311... (52.085223289182096, 194.5269753331549)
In [325...
          train_df_pos[train_df_pos['p_val'] >= 0.01]['mahal_dist'].min(), train_df_pos[train_
Out[325... (6.167006044471601, 50.03579960395341)
```

So, the null hypothesis got rejected based on the above 26 observations where the mahalanobis distance is greater than the critical value thus the distance ended up in the rejection region.

#### CASE-II

Now, I'll perform the test on negative observations considering the null hypothesis as FALSE or NEGATIVE and the alternate hypothesis as TRUE or POSITIVE.

```
train_df_neg['p_val'] = 1 - chi2.cdf(train_df_neg['mahal_dist'], df=dof)
In [318...
           train df neg.head()
Out[318...
                                                                                    mean
                                                                  mean
                                                                           mean
                                  mean
                                        mean
                                                     mean
                                                                                               mean
                mean
                        mean
               radius texture
                              perimeter
                                         area
                                              smoothness compactness concavity
                                                                                           symmetry
                                                                                   points
```

localhost:8888/lab 10/13

	mean radius	mean texture	mean perimeter	mean area	mean smoothness	mean compactness	mean concavity	mean concave points	mean symmetry	•
39	13.48	20.82	88.40	559.2	0.10160	0.12550	0.1063	0.05439	0.1720	
229	12.83	22.33	85.26	503.2	0.10880	0.17990	0.1695	0.06861	0.2123	
40	13.44	21.58	86.18	563.0	0.08162	0.06031	0.0311	0.02031	0.1784	
172	15.46	11.89	102.50	736.9	0.12570	0.15550	0.2032	0.10970	0.1966	
213	17.42	25.56	114.50	948.0	0.10060	0.11460	0.1682	0.06597	0.1308	

5 rows × 33 columns

```
train_df_neg.shape
In [319...
Out[319... (127, 33)
          train_df_neg[train_df_neg['p_val'] < 0.01].shape</pre>
In [320...
Out[320... (14, 33)
           rej_null_hyp_neg = train_df_neg[train_df_neg['p_val'] < 0.01]['mahal_dist'].values</pre>
In [321...
           rej_null_hyp_neg
Out[321... array([104.13997248,
                                 65.26593056, 57.53158832,
                                                              74.66179609,
                                79.93825875, 52.30118131, 67.52958184,
                  55.42755962,
                  58.82178158,
                                 52.2838042 , 61.00788787, 102.21571179,
                  60.01309311, 70.31700251])
          rej_null_hyp_neg.min(), rej_null_hyp_neg.max()
In [322...
Out[322... (52.28380420063802, 104.13997248029511)
          train_df_neg[train_df_neg['p_val'] >= 0.01]['mahal_dist'].min(), train_df_neg[train_
In [324...
Out[324... (8.361142675252562, 49.68938504959755)
```

So, the null hypothesis got rejected based on the above 14 observations where the mahalanobis distance is greater than the critical value thus the distance ended up in the rejection region.

**Euclidean Distance :: It performs the pairwise difference and not takes the distribution into account.** 

```
In [79]: from scipy.spatial.distance import pdist, squareform
In [163... distances = pdist(df.iloc[:,:], metric='euclidean')
    dist_matrix = pd.DataFrame(squareform(distances))
In [164... pd.DataFrame(dist_matrix.iloc[0:5,0])
Out[164... 0
```

localhost:8888/lab 11/13

```
0 0.000000
```

0

- **1** 1.700118
- **2** 4.707441
- **3** 8.050689
- **4** 9.178584

```
In [165... df.head(5)
```

# Out[165...

	carat	depth	price
0	0.23	61.5	326
1	0.21	59.8	326
2	0.23	56.9	327
3	0.29	62.4	334
4	0.31	63.3	335

## **VIF Calculation**

In [345... | from statsmodels.stats.outliers\_influence import variance\_inflation\_factor as vif

In [346... df.head()

#### Out[346...

```
carat depth price
0
    0.23
            61.5
                   326
1
    0.21
            59.8
                   326
2
    0.23
            56.9
                   327
3
    0.29
            62.4
                   334
            63.3
    0.31
                   335
```

```
In [350... vif_df = pd.DataFrame()
    vif_df['Values'] = [vif(df.values,i) for i in range(len(df.columns))]
    vif_df['Features'] = df.columns
```

In [351... vif\_df

#### Out[351...

	Values	Features
0	60.169063	carat
1	11.015691	depth
2	37.304660	price

#### **KNN**

```
In [405... knc = KNeighborsClassifier(n_neighbors=5, radius=0.7, weights='distance', algorithm='kd
```

localhost:8888/lab 12/13

```
In [406... | cancer_df.head()
```

Out[406...

	mean radius	mean texture	mean perimeter	mean area	mean smoothness	mean compactness	mean concavity	mean concave points	mean symmetry	diı
0	17.99	10.38	122.80	1001.0	0.11840	0.27760	0.3001	0.14710	0.2419	
1	20.57	17.77	132.90	1326.0	0.08474	0.07864	0.0869	0.07017	0.1812	
2	19.69	21.25	130.00	1203.0	0.10960	0.15990	0.1974	0.12790	0.2069	
3	11.42	20.38	77.58	386.1	0.14250	0.28390	0.2414	0.10520	0.2597	
4	20.29	14.34	135.10	1297.0	0.10030	0.13280	0.1980	0.10430	0.1809	

5 rows × 31 columns

```
In [407... knc.fit(cancer_df.iloc[:,0:-1],cancer_df.iloc[:,-1])
Out[407... KNeighborsClassifier(algorithm='kd_tree', weights='distance')
In [408... knc.predict(cancer_df.iloc[:,0:-1])
```

```
0, 1,
Out[408...
                 0, 0,
                      0, 0, 0, 0, 0, 0, 0, 0,
                                               0, 0, 0, 0, 0, 1, 0,
                                                                     0,
                   0,
                      1,
                          0,
                             1,
                                1,
                                   1,
                                      1,
                                         1,
                                            0,
                                               0,
                                                  1,
                                                     0, 0,
                                                           1,
                                                              1,
                                                                  1,
                                                                     1,
                                1,
                                   0,
                                      0,
                                         1,
                                            0,
                                               1,
                                                  0,
                                                     0,
                                                        1,
                                                           1,
                                                               1,
                                                                        1,
                                   0,
                                      0,
                                         1,
                                            1,
                                               1,
                                                  0,
                                                     0,
                                                         1,
                                                            1,
                                                                  1,
                                      0,
                                         0,
                                            0,
                                               1,
                                                  0,
                                                     0,
                                                         1,
                                                            1,
                                                                        1,
                                   0,
                                      1,
                                         1,
                                            0,
                                               1,
                                                  1,
                                                      1,
                                                        1,
                                                            0,
                                                               1,
                                                                  1,
                                                                     1,
                                                                        1,
                                                            0,
                                   1,
                                      0,
                                         0,
                                            1,
                                               0,
                                                  1,
                                                      1,
                                                         0,
                                                               1,
                                                                  1,
                                                                     0,
                                   0,
                                      1,
                                         0,
                                            1,
                                               0,
                                                  1,
                                                     1,
                                                         1,
                                                            0,
                                                               1,
                                                                  1,
                                                               0,
                                                                  0,
                                                                     0,
                                                                        1,
                                   1,
                                      0,
                                         1,
                                            0,
                                               1,
                                                  1,
                                                      0,
                                                        1,
                                                            0,
                                                                  0,
                                                                     0,
                                                                        1,
                                         1,
                                            0,
                                               0,
                                                  1,
                                                     1,
                                                         0,
                                                            1,
                                                               1,
                                                                  0,
                                   1,
                                      1,
                                         0,
                                            1,
                                               0,
                                                  0,
                                                      0,
                                                         0,
                                                            0,
                                   1,
                                      1,
                                         0,
                                            1,
                                               0,
                                                  1,
                                                      1,
                                                         0,
                                                            1,
                                                                     1,
                                   1,
                                      1,
                                         1,
                                            1,
                                               1,
                                                  0,
                                                      1,
                                                         1,
                                                            0,
                                                               1,
                                1,
                                   1,
                                      1,
                                         1,
                                            0,
                                               1,
                                                  1,
                                                     1,
                                                         0,
                                                            1,
                                   1,
                                      0,
                                         1,
                                            0,
                                               1,
                                                  1,
                                                     1,
                                                         0,
                                                            1,
                                   1,
                                      1,
                                         1,
                                            1,
                                               1,
                                                  1,
                                                      1,
                                                         0,
                                                            0,
                                                               1,
                                         1,
                                               1,
                                                  0,
                                   1,
                                      1,
                                            1,
                                                     1,
                                                         1,
                                                            1,
                                                               0,
                                      1,
                                         1,
                                               1,
                                                     0,
                             0,
                                   1,
                                            1,
                                                  1,
                                                        1,
                                                            1,
                                                               1,
                                                                     1,
                                      1,
                                         1,
                                            1,
                                               1,
                                   1,
                                                  1,
                                                      0,
                                                        1,
                                                            0,
                                                               0,
                                            0,
                                                  0,
                                      1,
                                               1,
                                1,
                                   0,
                                         1,
                                                     1,
                                                        1,
                                                            1,
                                                                     1,
                                   0,
                                      1,
                                         1,
                                            1,
                                                  1,
                          1,
                                1,
                                               1,
                                                     1,
                                                        1,
                                                           1,
                                                               1,
                                            1,
                                      1, 0,
                                               1,
                                                  1,
                          0,
                             1,
                                0,
                                   1,
                                                     1,
                                                        1, 0, 0,
                                                                  1,
                                                                        1,
                                1, 0, 1, 0, 1, 0, 0, 1, 1, 1, 0, 1,
                      1,
                         0,
                            1,
                                                                     1,
                                                                        1, 1,
                         1,
                 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 0, 0, 0, 0, 0, 0, 1])
```

localhost:8888/lab 13/13