# 2602063913\_Bernardus Ignasio\_Assignment 5 Machine Learning

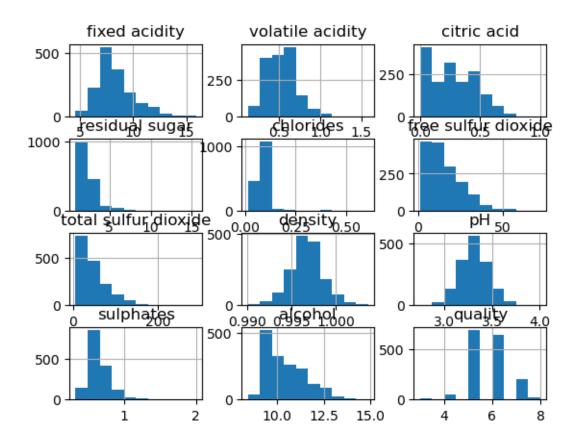
June 28, 2024

## 1 Assignment 5 Machine Learning

NIM: 2602063913

```
NAME : Bernardus Ignasio
      CLASS: LA01
      Assignment Number: 5
[378]: import pandas as pd
       import matplotlib.pyplot as plt
       from sklearn.metrics import silhouette score,
        →adjusted_rand_score,davies_bouldin_score
[379]: dataset=pd.read_csv('https://raw.githubusercontent.com/joh4nnes/MLData/main/
        ⇔winequality-red.csv')
       dataset.head()
[379]:
          fixed acidity volatile acidity citric acid residual sugar
                                                                         chlorides \
                    7.4
                                      0.70
                                                   0.00
                                                                     1.9
                                                                              0.076
                    7.8
                                                   0.00
                                                                    2.6
       1
                                      0.88
                                                                              0.098
       2
                    7.8
                                     0.76
                                                   0.04
                                                                    2.3
                                                                              0.092
       3
                   11.2
                                      0.28
                                                   0.56
                                                                     1.9
                                                                              0.075
       4
                    7.4
                                                   0.00
                                                                     1.9
                                      0.70
                                                                              0.076
          free sulfur dioxide total sulfur dioxide density
                                                                     sulphates
                                                                 рΗ
       0
                                                34.0
                                                       0.9978
                                                               3.51
                                                                           0.56
                         11.0
                         25.0
                                                67.0
       1
                                                       0.9968
                                                               3.20
                                                                           0.68
       2
                         15.0
                                                54.0
                                                       0.9970
                                                               3.26
                                                                           0.65
                                                                           0.58
       3
                         17.0
                                                60.0
                                                       0.9980
                                                               3.16
                         11.0
                                                34.0
                                                       0.9978 3.51
                                                                           0.56
          alcohol quality
       0
              9.4
              9.8
                         5
       1
       2
              9.8
                         5
       3
              9.8
                         6
              9.4
                         5
```

```
[403]: dataset.isna().sum()
[403]: fixed acidity
                               0
      volatile acidity
                               0
       citric acid
                               0
       residual sugar
                               0
       chlorides
                               0
       free sulfur dioxide
                               0
       total sulfur dioxide
       density
                               0
      рΗ
                               0
       sulphates
       alcohol
                               0
       quality
                               0
       dtype: int64
[404]: dataset.hist()
[404]: array([[<Axes: title={'center': 'fixed acidity'}>,
               <Axes: title={'center': 'volatile acidity'}>,
               <Axes: title={'center': 'citric acid'}>],
              [<Axes: title={'center': 'residual sugar'}>,
               <Axes: title={'center': 'chlorides'}>,
               <Axes: title={'center': 'free sulfur dioxide'}>],
              [<Axes: title={'center': 'total sulfur dioxide'}>,
               <Axes: title={'center': 'density'}>,
               <Axes: title={'center': 'pH'}>],
              [<Axes: title={'center': 'sulphates'}>,
               <Axes: title={'center': 'alcohol'}>,
               <Axes: title={'center': 'quality'}>]], dtype=object)
```

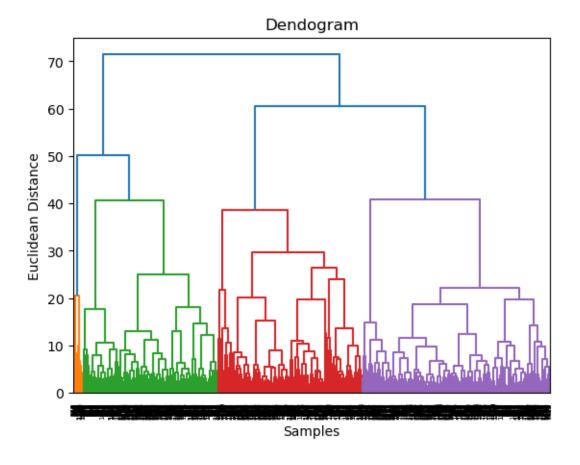


```
Extracting matrix of features
[381]: x=dataset.iloc[:,:11].values
      х
[381]: array([[ 7.4 , 0.7 , 0. , ..., 3.51 , 0.56 , 9.4 ],
            [7.8, 0.88, 0., ..., 3.2, 0.68, 9.8],
            [7.8, 0.76, 0.04, ..., 3.26, 0.65, 9.8],
            [6.3, 0.51, 0.13, ..., 3.42, 0.75, 11.
            [5.9, 0.645, 0.12, ..., 3.57, 0.71, 10.2],
            [6., 0.31, 0.47, ..., 3.39, 0.66, 11.
                                                          ]])
[382]: from sklearn.preprocessing import StandardScaler
      scale=StandardScaler()
      x=scale.fit_transform(x)
      х
[382]: array([[-0.52835961, 0.96187667, -1.39147228, ..., 1.28864292,
             -0.57920652, -0.96024611],
             [-0.29854743, 1.96744245, -1.39147228, ..., -0.7199333,
```

```
0.1289504 , -0.58477711],
[-0.29854743, 1.29706527, -1.18607043, ..., -0.33117661, -0.04808883, -0.58477711],
...,
[-1.1603431 , -0.09955388, -0.72391627, ..., 0.70550789, 0.54204194, 0.54162988],
[-1.39015528, 0.65462046, -0.77526673, ..., 1.6773996 , 0.30598963, -0.20930812],
[-1.33270223, -1.21684919, 1.02199944, ..., 0.51112954, 0.01092425, 0.54162988]])
```

### Find Optimal number of cluster with dendogram

```
[383]: import scipy.cluster.hierarchy as shc
  dendro=shc.dendrogram(shc.linkage(x,method='ward'))
  plt.title('Dendogram')
  plt.xlabel('Samples')
  plt.ylabel('Euclidean Distance')
  plt.show()
```

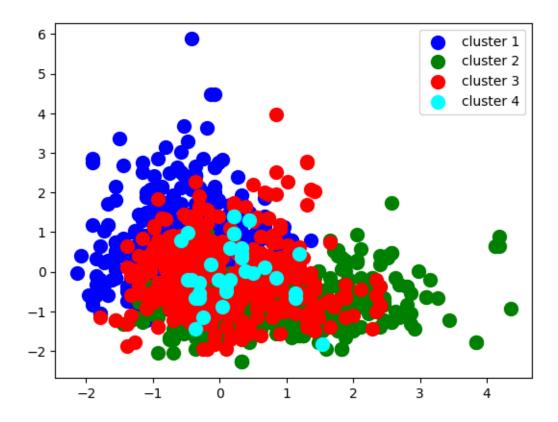


by slicing horizontally, optimal number of cluster is between 3 and 4

# Clustering Dataset with agglomerative clustering and optimal cluster from sklearn.cluster import AgglomerativeClustering hc=AgglomerativeClustering(n\_clusters=4,affinity='euclidean',linkage='ward') y\_predict=hc.fit\_predict(x) y\_predict C:\Users\Ignas\anaconda3\Lib\sitepackages\sklearn\cluster\\_agglomerative.py:1005: FutureWarning: Attribute 'affinity' was deprecated in version 1.2 and will be removed in 1.4. Use 'metric' instead warnings.warn( [384]: array([0, 0, 0, ..., 2, 2, 1], dtype=int64) [385]: plt.scatter(x[y\_predict==0,0],x[y\_predict==0,1], s=100,c='blue',label='cluster\_u=1')

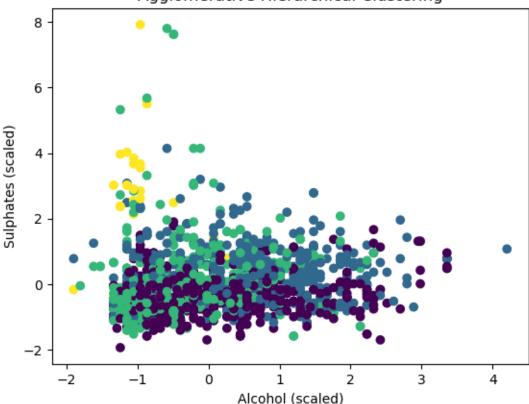
```
plt.scatter(x[y_predict==0,0],x[y_predict==0,1], s=100,c='blue',label='cluster_u
41')
plt.scatter(x[y_predict==1,0],x[y_predict==1,1], s=100,c='green',label='cluster_u
42')
plt.scatter(x[y_predict==2,0],x[y_predict==2,1], s=100,c='red',label='cluster_u
43')
plt.scatter(x[y_predict==3,0],x[y_predict==3,1], s=100,c='cyan',label='cluster_u
44')
#plt.scatter(x[y_predict==4,0],x[y_predict==4,1],u
4s=100,c='magenta',label='cluster_5')

plt.legend()
plt.show()
```



```
[386]: plt.scatter(x[:,10], x[:,9], c=y_predict,cmap='viridis',marker='o')
    plt.title('Agglomerative Hierarchical Clustering')
    plt.xlabel('Alcohol (scaled)')
    plt.ylabel('Sulphates (scaled)')
    plt.show()
```





```
[387]: print(f"Silhouette Score: {silhouette_score(x,hc.labels_,metric='euclidean')}") print(f"Davies Bouldin Index (DBI): {davies_bouldin_score(x,hc.labels_)}")
```

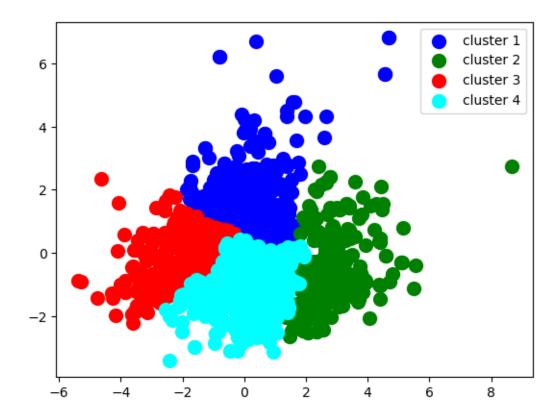
Silhouette Score: 0.17141826625204853 Davies Bouldin Index (DBI): 1.6648528437215329

from the visualization and evaluation metrics, we can see that the clustering is still poorly matched and the distance between every cluster cannot be seen, clusters are spreading and combining together which indicates poor clustering quality

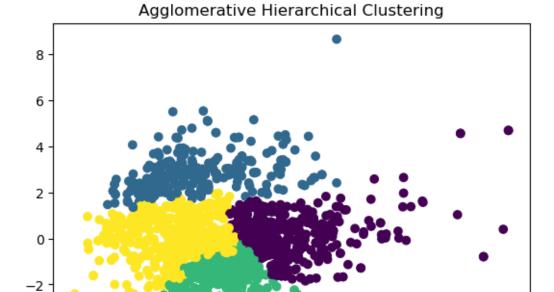
### Enhancing quality with PCA

```
[388]: from sklearn.decomposition import PCA
pca=PCA(n_components=2)
x=pca.fit_transform(x)
x
```

```
[-1.45612897, 0.31174559],
              [-2.27051793, 0.97979111],
              [-0.42697475, -0.53669021]])
[389]: hc=AgglomerativeClustering(n_clusters=4,affinity='euclidean',linkage='ward')
       y_predict=hc.fit_predict(x)
       y_predict
      C:\Users\Ignas\anaconda3\Lib\site-
      packages\sklearn\cluster\_agglomerative.py:1005: FutureWarning: Attribute
      `affinity` was deprecated in version 1.2 and will be removed in 1.4. Use
      `metric` instead
        warnings.warn(
[389]: array([2, 0, 0, ..., 2, 2, 3], dtype=int64)
[390]: plt.scatter(x[y_predict==0,0],x[y_predict==0,1], s=100,c='blue',label='cluster_
       <1¹)
       plt.scatter(x[y_predict==1,0],x[y_predict==1,1], s=100,c='green',label='cluster_
       plt.scatter(x[y_predict==2,0],x[y_predict==2,1], s=100,c='red',label='cluster_u
       plt.scatter(x[y_predict==3,0],x[y_predict==3,1], s=100,c='cyan',label='cluster_
        4')
       \#plt.scatter(x[y\_predict==4,0],x[y\_predict==4,1],
        ⇔s=100, c='magenta', label='cluster 5')
       plt.legend()
       plt.show()
```



```
[391]: plt.scatter(x[:,1], x[:,0], c=y_predict,cmap='viridis',marker='o') plt.title('Agglomerative Hierarchical Clustering') plt.show()
```



```
[392]: print(f"Silhouette Score: {silhouette_score(x,hc.labels_,metric='euclidean')}") print(f"Davies Bouldin Index (DBI): {davies_bouldin_score(x,hc.labels_)}")
```

0

Silhouette Score: 0.3413983711219637 Davies Bouldin Index (DBI): 0.9068520999649691

-2

after implementing pca and reducting columns until 2 columns left, we can see that the clusters separation can be seen clearly although there's no gap/distance between the clusters. also the evaluation metrics is better, indicating that the clustering is improving

2

4

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