```
In [1]: from sklearn.datasets import load_iris
        import pandas as pd
In [4]: | iris = load_iris()
        iris
Out[4]: {'data': array([[5.1, 3.5, 1.4, 0.2],
                [4.9, 3., 1.4, 0.2],
                [4.7, 3.2, 1.3, 0.2],
                [4.6, 3.1, 1.5, 0.2],
                [5., 3.6, 1.4, 0.2],
                [5.4, 3.9, 1.7, 0.4],
                [4.6, 3.4, 1.4, 0.3],
                [5., 3.4, 1.5, 0.2],
                [4.4, 2.9, 1.4, 0.2],
                [4.9, 3.1, 1.5, 0.1],
                [5.4, 3.7, 1.5, 0.2],
                [4.8, 3.4, 1.6, 0.2],
                [4.8, 3., 1.4, 0.1],
                [4.3, 3., 1.1, 0.1],
                [5.8, 4., 1.2, 0.2],
                [5.7, 4.4, 1.5, 0.4],
                [5.4, 3.9, 1.3, 0.4],
                [5.1, 3.5, 1.4, 0.3],
                [5.7, 3.8, 1.7, 0.3],
```

Out[5]:

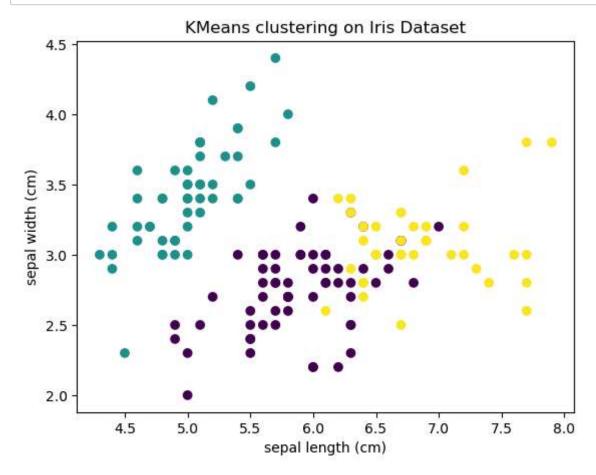
	sepal length (cm)	sepal width (cm)	petal length (cm)	petal width (cm)
0	5.1	3.5	1.4	0.2
1	4.9	3.0	1.4	0.2
2	4.7	3.2	1.3	0.2
3	4.6	3.1	1.5	0.2
4	5.0	3.6	1.4	0.2
145	6.7	3.0	5.2	2.3
146	6.3	2.5	5.0	1.9
147	6.5	3.0	5.2	2.0
148	6.2	3.4	5.4	2.3
149	5.9	3.0	5.1	1.8

150 rows × 4 columns

In [8]: from sklearn.cluster import KMeans
import matplotlib.pyplot as plt

```
In [9]:
         kmeans = KMeans(n clusters=3,random state=42)
         kmeans.fit(x)
         C:\Users\999ra\anaconda3\Lib\site-packages\sklearn\cluster\ kmeans.py:1412: FutureWarning: The default valu
         e of `n_init` will change from 10 to 'auto' in 1.4. Set the value of `n_init` explicitly to suppress the wa
         rning
           super(). check params vs input(X, default n init=10)
         C:\Users\999ra\anaconda3\Lib\site-packages\sklearn\cluster\ kmeans.py:1436: UserWarning: KMeans is known to
         have a memory leak on Windows with MKL, when there are less chunks than available threads. You can avoid it
         by setting the environment variable OMP NUM THREADS=1.
           warnings.warn(
 Out[9]:
                          KMeans
          KMeans(n_clusters=3, random_state=42)
In [11]: x['cluster'] = kmeans.labels_
         x['cluster']
Out[11]: 0
                1
         1
                1
         2
                1
         3
                1
         4
                1
         145
                2
         146
                0
         147
                2
         148
                2
         149
                0
         Name: cluster, Length: 150, dtype: int32
```

```
In [13]: plt.scatter(x.iloc[:,0], x.iloc[:,1], c=x['cluster'], cmap='viridis')
    plt.title('KMeans clustering on Iris Dataset')
    plt.xlabel(iris.feature_names[0])
    plt.ylabel(iris.feature_names[1])
    plt.show()
```



In []: #Description of KMeans Clustering

#KMeans is a centroid-based clustering algorithm that works by partitioning data into k distinct clusters.
#It tries to minimize the distance between the points and their assigned cluster centroids. The steps involve

- #1. Randomly selecting k centroids.
- #2. Assigning each data point to the nearest centroid.
- #3. Updating the centroids based on the points assigned to each cluster.
- #4. Repeating the process until the centroids do not change or a maximum number of iterations is reached.

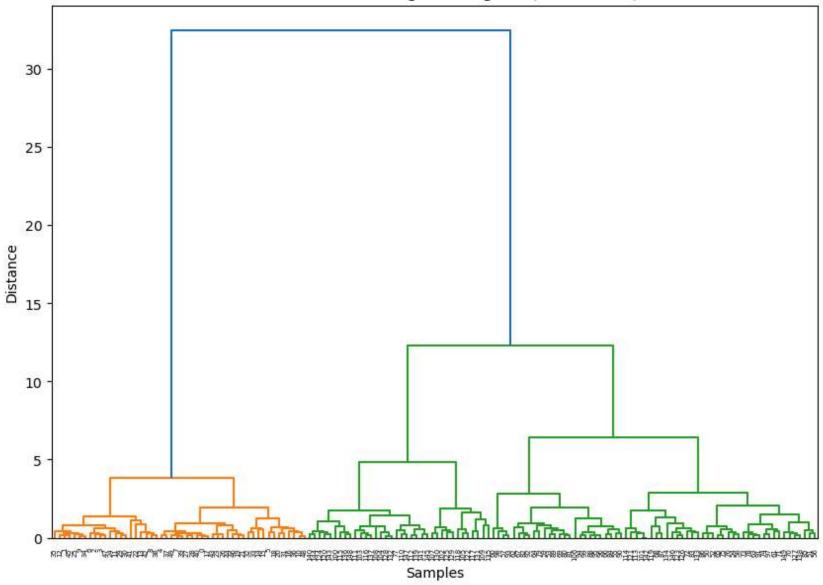
#Why KMeans is suitable for Iris dataset

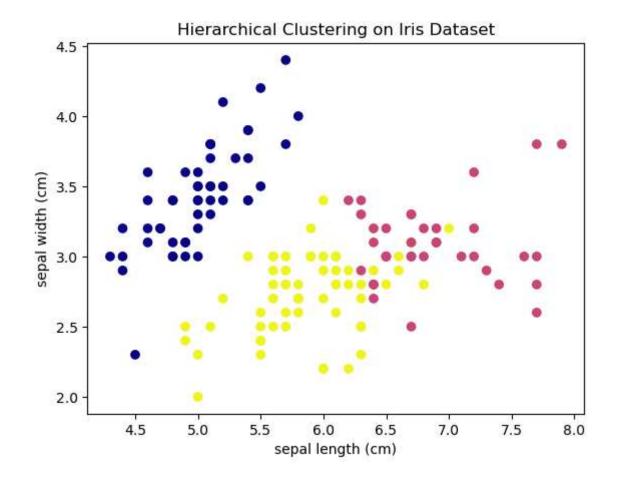
#KMeans is suitable for the Iris dataset because it's a simple, well-separated dataset.

#There are three known classes (species), and KMeans can cluster based on the inherent patterns in the featur #without knowing the actual species labels. The continuous numeric features (sepal and petal lengths/widths) #a good candidate for this dataset.

```
In [15]: from scipy.cluster.hierarchy import dendrogram, linkage
         import seaborn as sns
         # Hierarchical Clustering
         Z = linkage(x.iloc[:, :-1], method='ward')
         # Plot the dendrogram
         plt.figure(figsize=(10, 7))
         dendrogram(Z)
         plt.title('Hierarchical Clustering Dendrogram (Iris Dataset)')
         plt.xlabel('Samples')
         plt.ylabel('Distance')
         plt.show()
         # To create clusters from hierarchical clustering (choose 3 clusters)
         from scipy.cluster.hierarchy import fcluster
         # Cut the dendrogram to form 3 clusters
         x['h cluster'] = fcluster(Z, 3, criterion='maxclust')
         # Visualize the clusters (using only the first two features for easy visualization)
         plt.scatter(x.iloc[:, 0], x.iloc[:, 1], c=x['h_cluster'], cmap='plasma')
         plt.title('Hierarchical Clustering on Iris Dataset')
         plt.xlabel(iris.feature names[0])
         plt.ylabel(iris.feature_names[1])
         plt.show()
```

Hierarchical Clustering Dendrogram (Iris Dataset)





In []: #Description of Hierarchical Clustering #Hierarchical clustering builds a hierarchy of clusters by either a bottom-up approach (agglomerative) or #a top-down approach (divisive). In agglomerative clustering, each point starts as its own cluster, and #the algorithm recursively merges the closest clusters until only a single cluster or the desired number of c #The clustering process can be visualized using a dendrogram.

#Why Hierarchical Clustering is suitable for Iris dataset
#Hierarchical clustering can reveal the nested structure of the data, which is useful when the number of clus
#not known a priori. It allows us to observe how the clusters merge at different levels, which could provide
#the relationships between different species in the Iris dataset.