In [1]: import seaborn as sns
 data=sns.load_dataset('iris')
 data

Out[1]:		sepal_length	sepal_width	petal_length	petal_width	species
	0	5.1	3.5	1.4	0.2	setosa
	1	4.9	3.0	1.4	0.2	setosa
	2	4.7	3.2	1.3	0.2	setosa
	3	4.6	3.1	1.5	0.2	setosa
	4	5.0	3.6	1.4	0.2	setosa
	145	6.7	3.0	5.2	2.3	virginica
	146	6.3	2.5	5.0	1.9	virginica
	147	6.5	3.0	5.2	2.0	virginica
	148	6.2	3.4	5.4	2.3	virginica
	149	5.9	3.0	5.1	1.8	virginica

150 rows × 5 columns

In [2]: features=data.drop(columns='species',axis=0)
 features

Out[2]:		sepal_length	sepal_width	petal_length	petal_width
_	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 [3]: data.info()
       <class 'pandas.core.frame.DataFrame'>
       RangeIndex: 150 entries, 0 to 149
       Data columns (total 5 columns):
           Column
                        Non-Null Count Dtype
        0 sepal_length 150 non-null float64
        1 sepal width 150 non-null float64
            petal_length 150 non-null
                                       float64
        3 petal_width 150 non-null
                                       float64
        4 species
                       150 non-null
                                        object
       dtypes: float64(4), object(1)
       memory usage: 6.0+ KB
```

In [4]: features.describe()

Out[4]:

	sepal_length	sepal_width	petal_length	petal_width
count	150.000000	150.000000	150.000000	150.000000
mean	5.843333	3.057333	3.758000	1.199333
std	0.828066	0.435866	1.765298	0.762238
min	4.300000	2.000000	1.000000	0.100000
25%	5.100000	2.800000	1.600000	0.300000
50%	5.800000	3.000000	4.350000	1.300000
75%	6.400000	3.300000	5.100000	1.800000
max	7.900000	4.400000	6.900000	2.500000

KMeans Clustering

KMeans is an iterative clustering algorithm that divides the dataset into a predefined number of clusters (k). It randomly initializes k centroids, then assigns each point to the nearest centroid. Centroids are recalculated based on the mean of points assigned to them, and the process repeats until convergence.

KMeans is suitable for the Iris dataset because Iris has 3 species (which can be thought of as natural clusters). KMeans works well when the number of clusters is known, and Iris has clear boundaries between species based on feature values.

```
In [5]: from sklearn.preprocessing import StandardScaler
    standard=StandardScaler()
    scaled_features=standard.fit_transform(features)
    from sklearn.cluster import KMeans
```

```
In [6]: inertia=[]
k_values=range(1,11)
for k in k_values:
    kmeans=KMeans(n_clusters=k)
    kmeans.fit(scaled_features)
    inertia.append(kmeans.inertia_)
```

```
C:\Users\aathi\anaconda3\Lib\site-packages\sklearn\cluster\_kmeans.py:1412: FutureW
arning: The default value of `n_init` will change from 10 to 'auto' in 1.4. Set the
value of `n_init` explicitly to suppress the warning
  super(). check params vs input(X, default n init=10)
C:\Users\aathi\anaconda3\Lib\site-packages\sklearn\cluster\_kmeans.py:1436: UserWar
ning: KMeans is known to have a memory leak on Windows with MKL, when there are les
s chunks than available threads. You can avoid it by setting the environment variab
le OMP_NUM_THREADS=1.
 warnings.warn(
C:\Users\aathi\anaconda3\Lib\site-packages\sklearn\cluster\_kmeans.py:1412: FutureW
arning: The default value of `n_init` will change from 10 to 'auto' in 1.4. Set the
value of `n_init` explicitly to suppress the warning
  super()._check_params_vs_input(X, default_n_init=10)
C:\Users\aathi\anaconda3\Lib\site-packages\sklearn\cluster\_kmeans.py:1436: UserWar
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s chunks than available threads. You can avoid it by setting the environment variab
le OMP NUM THREADS=1.
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C:\Users\aathi\anaconda3\Lib\site-packages\sklearn\cluster\_kmeans.py:1412: FutureW
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  super()._check_params_vs_input(X, default_n_init=10)
C:\Users\aathi\anaconda3\Lib\site-packages\sklearn\cluster\_kmeans.py:1436: UserWar
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s chunks than available threads. You can avoid it by setting the environment variab

C:\Users\aathi\anaconda3\Lib\site-packages\sklearn\cluster_kmeans.py:1412: FutureW arning: The default value of `n_init` will change from 10 to 'auto' in 1.4. Set the

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C:\Users\aathi\anaconda3\Lib\site-packages\sklearn\cluster_kmeans.py:1412: FutureW arning: The default value of `n_init` will change from 10 to 'auto' in 1.4. Set the

C:\Users\aathi\anaconda3\Lib\site-packages\sklearn\cluster_kmeans.py:1436: UserWar ning: 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 variab

C:\Users\aathi\anaconda3\Lib\site-packages\sklearn\cluster_kmeans.py:1412: FutureW
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C:\Users\aathi\anaconda3\Lib\site-packages\sklearn\cluster_kmeans.py:1436: UserWar ning: 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 variab

C:\Users\aathi\anaconda3\Lib\site-packages\sklearn\cluster_kmeans.py:1412: FutureW arning: The default value of `n init` will change from 10 to 'auto' in 1.4. Set the

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value of `n_init` explicitly to suppress the warning super()._check_params_vs_input(X, default_n_init=10)

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warnings.warn(

```
value of `n_init` explicitly to suppress the warning
   super()._check_params_vs_input(X, default_n_init=10)
C:\Users\aathi\anaconda3\Lib\site-packages\sklearn\cluster\_kmeans.py:1436: UserWar
ning: KMeans is known to have a memory leak on Windows with MKL, when there are les
s chunks than available threads. You can avoid it by setting the environment variab
le OMP_NUM_THREADS=1.
   warnings.warn(
C:\Users\aathi\anaconda3\Lib\site-packages\sklearn\cluster\_kmeans.py:1412: FutureW
arning: The default value of `n_init` will change from 10 to 'auto' in 1.4. Set the
value of `n_init` explicitly to suppress the warning
   super()._check_params_vs_input(X, default_n_init=10)
```

C:\Users\aathi\anaconda3\Lib\site-packages\sklearn\cluster_kmeans.py:1436: UserWar ning: KMeans is known to have a memory leak on Windows with MKL, when there are les s chunks than available threads. You can avoid it by setting the environment variab le OMP_NUM_THREADS=1.

warnings.warn(

C:\Users\aathi\anaconda3\Lib\site-packages\sklearn\cluster_kmeans.py:1412: FutureW arning: The default value of `n_init` will change from 10 to 'auto' in 1.4. Set the value of `n_init` explicitly to suppress the warning

super()._check_params_vs_input(X, default_n_init=10)

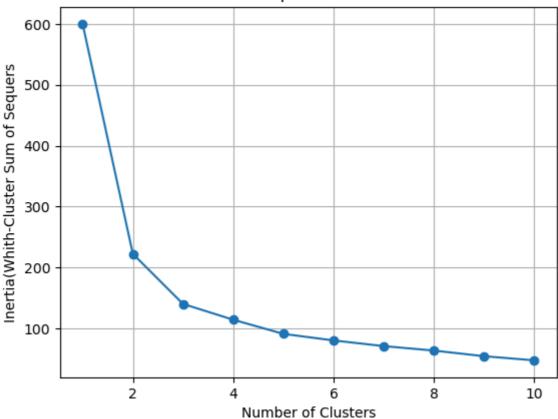
C:\Users\aathi\anaconda3\Lib\site-packages\sklearn\cluster_kmeans.py:1436: UserWar ning: KMeans is known to have a memory leak on Windows with MKL, when there are less chunks than available threads. You can avoid it by setting the environment variab le OMP_NUM_THREADS=1.

warnings.warn(

In [7]: inertia

```
In [8]: import matplotlib.pyplot as plt
plt.plot(k_values,inertia,marker='o')
plt.xlabel('Number of Clusters')
plt.ylabel('Inertia(Whith-Cluster Sum of Sequers')
plt.title('Elbow Method for Optimal Number of Clusters')
plt.grid(True)
plt.show()
```

Elbow Method for Optimal Number of Clusters



```
In [9]: from sklearn.cluster import KMeans
    Kmeans=KMeans(n_clusters=3)
    kmeans.fit_predict(scaled_features)
    data['Cluster']=kmeans.labels_
```

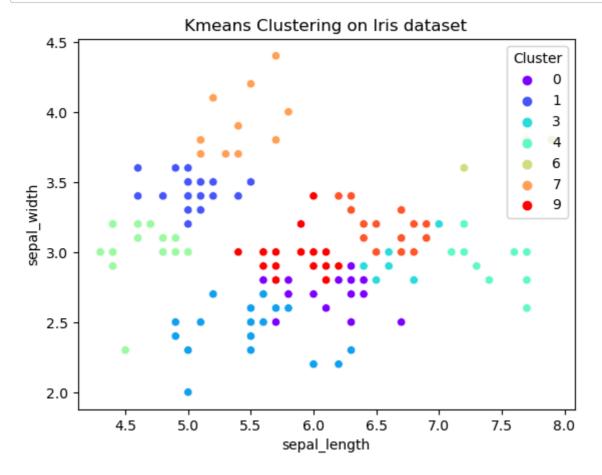
C:\Users\aathi\anaconda3\Lib\site-packages\sklearn\cluster_kmeans.py:1412: FutureW arning: The default value of `n_init` will change from 10 to 'auto' in 1.4. Set the value of `n_init` explicitly to suppress the warning

super()._check_params_vs_input(X, default_n_init=10)

C:\Users\aathi\anaconda3\Lib\site-packages\sklearn\cluster_kmeans.py:1436: UserWar ning: KMeans is known to have a memory leak on Windows with MKL, when there are les s chunks than available threads. You can avoid it by setting the environment variab le OMP_NUM_THREADS=1.

warnings.warn(

In [11]: import seaborn as sns
 sns.scatterplot(x=features['sepal_length'],y=features['sepal_width'],hue=data['Clust
 plt.title('Kmeans Clustering on Iris dataset')
 plt.show()



b) Hierarchial Clustering

Brief Description: Hierarchical clustering builds a hierarchy of clusters using either an agglomerative (bottom-up) or divisive (top-down) approach. In agglomerative clustering, it starts with each point as its own cluster and merges the closest pairs of clusters iteratively until a single cluster is formed. The results are often visualized using a dendrogram.

Why is Hierarchical clustering suitable for the Iris dataset? Hierarchical clustering is good for small datasets like Iris since it helps in visualizing how data points cluster at different levels of granularity. It does not require the number of clusters to be pre-specified, allowing for exploration.

```
In [18]: data=sns.load_dataset('iris')
```

C:\Users\aathi\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(

Out[13]:	sepal_length	sepal_width	petal_length	petal_width	species	cluster
	5.1	3.5	1.4	0.2	setosa	1
	4.9	3.0	1.4	0.2	setosa	1
	2 4.7	3.2	1.3	0.2	setosa	1
	3 4.6	3.1	1.5	0.2	setosa	1
	4 5.0	3.6	1.4	0.2	setosa	1
14	5 6.7	3.0	5.2	2.3	virginica	0
14	6.3	2.5	5.0	1.9	virginica	0
14	6.5	3.0	5.2	2.0	virginica	0
14	6.2	3.4	5.4	2.3	virginica	0

3.0

150 rows × 6 columns

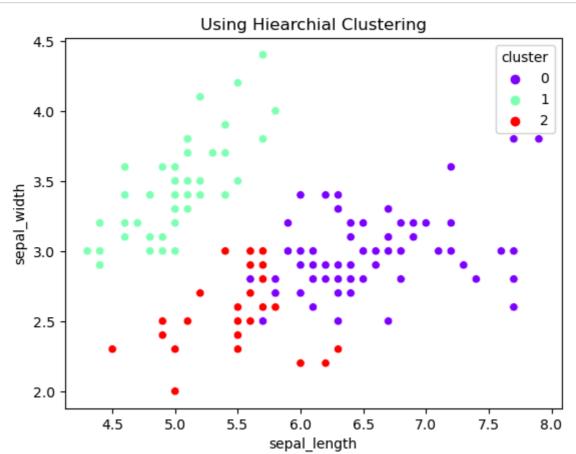
5.9

149

In [14]: sns.scatterplot(x='sepal_length',y='sepal_width',data=data,hue='cluster',palette='ra
plt.title('Using Hiearchial Clustering')
plt.show()

5.1

1.8 virginica



```
In [15]: from scipy.cluster.hierarchy import dendrogram,linkage
    z=linkage(scaled_features,method='ward')

In [16]: lab=data['species'].tolist()

In [17]: dendrogram(z,labels=lab,leaf_rotation=90)
    plt.title('Dendrogram of Agglomerative Hierarchical Clustering')
    plt.xlabel('Species')
    plt.ylabel('Distance')
    plt.show()
```

