# Programming 3

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# **Import Packages**

```
In [39]: import numpy as np
    import matplotlib.pyplot as plt
    from mpl_toolkits.mplot3d import Axes3D
    from sklearn.cluster import KMeans
    from sklearn import datasets
    from sklearn.preprocessing import normalize
    from copy import deepcopy

    iris = datasets.load_iris()
    #X = iris.data
    X = norm(iris.data)
    y = iris.target
```

# KMeans.py

```
>>> Select "k" random points from "data" as the initial centroids.
    11 11 11
   n_samples, n_features = np.shape(data)
    centroids = np.zeros((k, n_features))
    for i in range(k):
        centroid = data[np.random.choice(range(n_samples))]
        centroids[i] = centroid
    print(">>> initial centroids")
    print(centroids)
    return centroids
def converged(centroids1, centroids2):
    >>> Function you need to write
    >>> check whether centroids1==centroids
    >>> add proper code to handle infinite loop if it never converges
    diff = np.sum(np.abs(np.sum(centroids1 - centroids2, axis=1)), axis=0)
    if np.equal(centroids1, centroids2).all():
        return True
    elif diff < THRESHOLD:
       return True
    else:
       return False
def euclidean_dist(x1, x2):
    return np.sqrt(np.sum(np.square(x1 - x2), axis=1))
def closest_centroid(val, centroids):
    return np.argmin(np.sqrt(np.sum(np.square(val - centroids), axis=1)))
def update_centroids(data, centroids, k=3):
    >>> Function you need to write
    >>> Assign each data point to its nearest centroid based on the Euclidean distance
    >>> Update the cluster centroid to the mean of all the points assigned to that clu
   n_samples = np.shape(data)[0]
   n_features = np.shape(data)[1]
    clusters = [[] for _ in range(k)]
    labels = np.zeros((n_samples))
    for idx, val in enumerate(data):
        val_label = closest_centroid(val, centroids)
```

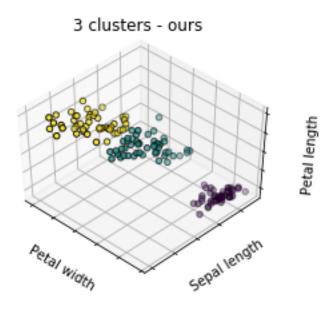
```
clusters[val_label].append(val)
        labels[idx] = val_label
    centroids = np.zeros((k, n_features))
    for idx, cluster_val in enumerate(clusters):
        centroid = np.mean(cluster_val, axis=0)
        centroids[idx] = centroid
    return centroids, labels
def kmeans(data, k=3):
    11 11 11
    >>> Function you should not touch
    # step 1:
    centroids = rand_center(data, k)
    converge = False
    iteration = 0
    while not converge:
        old_centroids = np.copy(centroids)
        # step 2 & 3
        centroids, label = update_centroids(data, old_centroids)
        converge = converged(old_centroids, centroids)
        iteration += 1
    print('number of iterations to converge: ', iteration)
    print(">>> final centroids")
    print(centroids)
    return centroids, label
def evaluation(predict, ground_truth):
    >>> use F1 and NMI in scikit-learn for evaluation
    f1 = f1_score(y_true=ground_truth, y_pred=predict, average='weighted')
    nmi = normalized_mutual_info_score(labels_true=ground_truth, labels_pred=predict)
    return f1, nmi
def gini(predict, ground_truth):
    >>> use the ground truth to do majority vote to assign a flower type for each clus
    >>> accordingly calculate the probability of missclassifiction and correct classif
    >>> finally, calculate gini using the calculated probabilities
    labels = np.unique(ground_truth)
    num_labels = len(labels)
    cluster_p, cluster_g = [0 for _ in range(num_labels)],
```

```
[0 for _ in range(num_labels)]
   gini_index = 0
    for i in labels:
        cluster_p[i], cluster_g[i] = np.array(predict[predict==i].shape[0]),
                                 np.array(ground_truth[ground_truth==i].shape[0])
        if cluster_p[i] < cluster_g[i]:</pre>
            correct_prob = cluster_p[i]/cluster_g[i]
            incorrect_prob = (cluster_g[i] - cluster_p[i])/cluster_g[i]
            gini_index += 1 - np.square(correct_prob) - np.square(incorrect_prob)
   gini_index /= num_labels
   print('Gini Index :', gini_index)
   return gini_index
def SSE(centroids, data):
   >>> Calculate the sum of squared errors for each cluster
    clusters = [[] for _ in centroids]
   num_centroids = len(centroids)
   for val in data:
        clusters[closest_centroid(val, centroids)].append(val)
    sse_each_cluster = [np.sum(np.sum(np.square(clusters[i] - centroids[i]), axis=1))
   print('SSE_each_cluster: ', sse_each_cluster)
   return sse_each_cluster
def plot_result(model):
   fig = plt.figure(1, figsize=(4, 3))
   ax = Axes3D(fig, rect=[0, 0, .95, 1], elev=48, azim=134)
   ax.scatter(X[:, 3], X[:, 0], X[:, 2],
               c=label.astype(np.float), edgecolor='k')
   ax.w_xaxis.set_ticklabels([])
   ax.w_yaxis.set_ticklabels([])
   ax.w_zaxis.set_ticklabels([])
   ax.set_xlabel('Petal width')
   ax.set_ylabel('Sepal length')
   ax.set_zlabel('Petal length')
   ax.set_title('3 clusters - {}'.format(model))
   ax.dist = 12
   plt.show()
```

## Q1

1. Run your own k-Means multiple times (at least 10 times) to find the best three clustering based on SSE (sum of squared errors) and compare it with the result obtained from k-Means in scikit-learn for comparison. Show both clustering results using matplotlib.

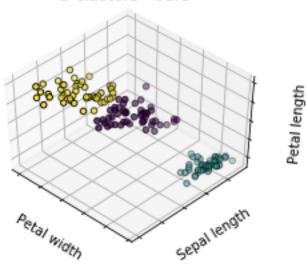
```
In [145]: for _ in range(10):
              # this is the function you are supposed to implement in kmeans.py
              centroids, label = kmeans(X,3)
              # this is the function you are supposed to implement in kmeans.py
              evaluation(label, y)
              gini(label, y)
              SSE(centroids, X)
              plot_result('ours')
>>> initial centroids
[[0.72151899 0.59090909 0.50724638 0.4
                                             ]
                                             ٦
 [0.69620253 0.54545455 0.53623188 0.4
                                             ]]
 [0.84810127 0.68181818 0.75362319 0.92
number of iterations to converge: 4
>>> final centroids
[[0.63367089 0.77681818 0.21217391 0.0976
                                             1
 [0.74853944 0.62237762 0.62263099 0.53
                                             ]
 [0.84045359 0.68560606 0.80676329 0.8225
                                             11
Gini Index: 0.025600000000000008
SSE_each_cluster: [0.5860971086851517, 1.0302179676457113, 1.2157725785900486]
```



SSE\_each\_cluster: [0.5860971086851517, 1.0302179676457113, 1.2157725785900486]

# 3 clusters - ours

#### >>> initial centroids [[0.84810127 0.56818182 0.84057971 0.72 ] [0.64556962 0.75 0.24637681 0.2 1 [0.78481013 0.77272727 0.7826087 0.92 ]] number of iterations to converge: 7 >>> final centroids [[0.74853944 0.62237762 0.62263099 0.53 ] [0.63367089 0.77681818 0.21217391 0.0976 ] [0.84045359 0.68560606 0.80676329 0.8225 11 Gini Index: 0.025600000000000008 SSE\_each\_cluster: [1.0302179676457113, 0.5860971086851517, 1.2157725785900486]



```
>>> initial centroids
```

- [[0.62025316 0.70454545 0.2173913 0.04 ]
- [0.72151899 0.65909091 0.60869565 0.52 ]
- [0.60759494 0.68181818 0.20289855 0.04 ]]

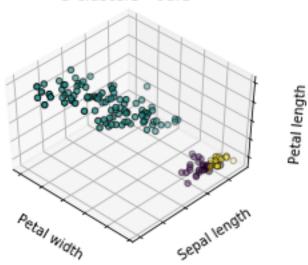
number of iterations to converge: 7

#### >>> final centroids

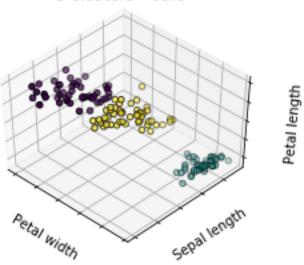
- [[0.66229656 0.8336039 0.21532091 0.11428571]
- [0.79265823 0.65272727 0.71101449 0.6704
- [0.5972382 0.70454545 0.20816864 0.07636364]]

Gini Index : 0.3285333333333333

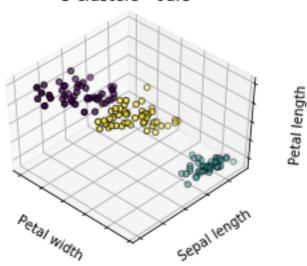
SSE\_each\_cluster: [0.21307645052258836, 5.538389212891668, 0.09732480213098582]



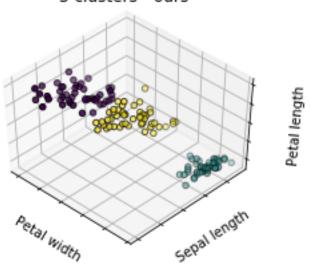
```
>>> initial centroids
[[0.86075949 0.72727273 0.85507246 0.92
                                             ]
[0.82278481 0.68181818 0.75362319 0.8
                                             ]
                                             ]]
 [0.87341772 0.70454545 0.73913043 0.92
number of iterations to converge: 10
>>> final centroids
[[0.84045359 0.68560606 0.80676329 0.8225
                                             ]
 [0.63367089 0.77681818 0.21217391 0.0976
                                             ]
 [0.74853944 0.62237762 0.62263099 0.53
                                             ]]
Gini Index: 0.025600000000000008
SSE_each_cluster: [1.2157725785900486, 0.5860971086851517, 1.0302179676457113]
```



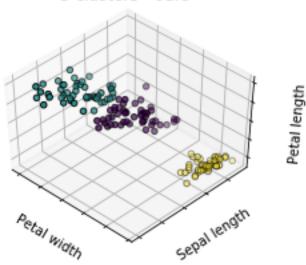
```
>>> initial centroids
[[0.91139241 0.68181818 0.84057971 0.64
                                             ]
[0.72151899 1.
                        0.2173913 0.16
                                             ]
 [0.70886076 0.65909091 0.52173913 0.52
                                             ]]
number of iterations to converge: 4
>>> final centroids
[[0.84045359 0.68560606 0.80676329 0.8225
                                             ]
[0.63367089 0.77681818 0.21217391 0.0976
                                             ]
 [0.74853944 0.62237762 0.62263099 0.53
                                             ]]
Gini Index: 0.025600000000000008
SSE_each_cluster: [1.2157725785900486, 0.5860971086851517, 1.0302179676457113]
```



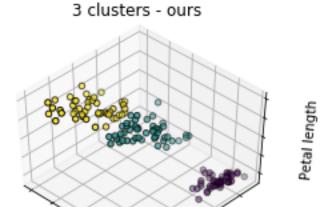
```
>>> initial centroids
[[0.84810127 0.68181818 0.75362319 0.92
                                             ]
[0.63291139 0.77272727 0.2173913 0.08
                                             ]
                                             ]]
 [0.83544304 0.68181818 0.63768116 0.56
number of iterations to converge: 4
>>> final centroids
[[0.84045359 0.68560606 0.80676329 0.8225
                                             ]
 [0.63367089 0.77681818 0.21217391 0.0976
                                             ]
 [0.74853944 0.62237762 0.62263099 0.53
                                             ]]
Gini Index: 0.025600000000000008
SSE_each_cluster: [1.2157725785900486, 0.5860971086851517, 1.0302179676457113]
```



```
>>> initial centroids
[[0.72151899 0.65909091 0.60869565 0.52
                                             ]
[0.84810127 0.70454545 0.8115942 0.96
                                             ]
                                             ]]
 [0.6835443 0.77272727 0.24637681 0.08
number of iterations to converge: 4
>>> final centroids
[[0.74853944 0.62237762 0.62263099 0.53
                                             ]
 [0.84045359 0.68560606 0.80676329 0.8225
                                             ]
 [0.63367089 0.77681818 0.21217391 0.0976
                                             ]]
Gini Index: 0.025600000000000008
SSE_each_cluster: [1.0302179676457113, 1.2157725785900486, 0.5860971086851517]
```

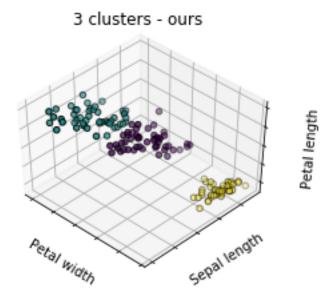


```
>>> initial centroids
[[0.69620253 0.52272727 0.57971014 0.52
                                             ]
[0.79746835 0.52272727 0.63768116 0.52
                                             ]
                                             ]]
 [0.83544304 0.68181818 0.63768116 0.56
number of iterations to converge: 5
>>> final centroids
[[0.63367089 0.77681818 0.21217391 0.0976
                                             ]
 [0.74853944 0.62237762 0.62263099 0.53
                                             ]
 [0.84045359 0.68560606 0.80676329 0.8225
                                             ]]
Gini Index: 0.025600000000000008
SSE_each_cluster: [0.5860971086851517, 1.0302179676457113, 1.2157725785900486]
```



```
>>> initial centroids
[[0.73417722 0.61363636 0.73913043 0.76
                                             ]
 [0.86075949 0.72727273 0.85507246 0.92
                                             ]
                                             ]]
 [0.60759494 0.77272727 0.27536232 0.08
number of iterations to converge: 7
>>> final centroids
[[0.74853944 0.62237762 0.62263099 0.53
                                             ]
 [0.84045359 0.68560606 0.80676329 0.8225
                                             ]
 [0.63367089 0.77681818 0.21217391 0.0976
                                             ]]
Gini Index: 0.025600000000000008
SSE_each_cluster: [1.0302179676457113, 1.2157725785900486, 0.5860971086851517]
```

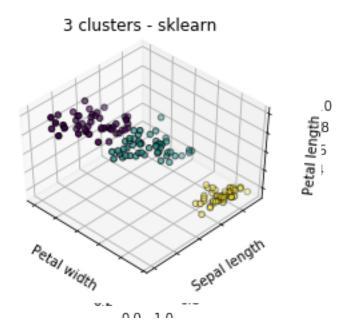
Petal Width



# sklearn

```
est = KMeans(n_clusters=3)
fig = plt.figure(1, figsize=(4, 3))
ax = Axes3D(fig, rect=[0, 0, .95, 1], elev=48, azim=134)
est.fit(X)
labels = est.labels_

# this is the function you are supposed to implement in kmeans.py
evaluation(labels, y)
gini(labels, y)
plot_result('sklearn')
Gini Index : 0.0256000000000000000
```



## Q2

2. Use the results obtained in (1) to explain why it's important to choose proper initial centroids.

From the result of question 1 we can observe in one case that when we choose the initial centroids as :

```
[[0.62025316 0.70454545 0.2173913 0.04 ]
[0.72151899 0.65909091 0.60869565 0.52 ]
[0.60759494 0.68181818 0.20289855 0.04 ]]
```

the evaluation parameters we get are:

- Gini Index: 0.3285333333333333
- SSE\_each\_cluster:
  - \* [0.21307645052258836, 5.538389212891668, 0.09732480213098582]

There is no feasible way to guarantee finding the optimal solution because the clustering analysis is a NP hard problem, heuristics approaches are needed to get an outcome that's more likely to be optimal. So there's a chance that with poorly chosen initial centroids, the results might be suboptimal.

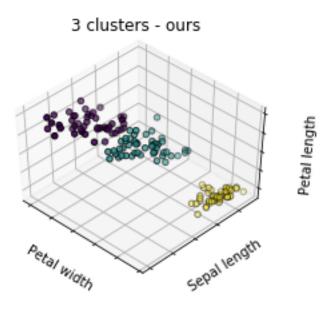
### Q3

3. Between the best clusterings you obtained via scikit-learn and your own k-Means implementation, which clustering is better? Please make comparison in terms of Impurity (Gini

Index) of the clusterings. To compute. Gini Index for a clustering, use the ground truth (i.e., flower types of data points) to do majority vote in order to assign a flower type for each cluster. Accordingly, calculate the probability of misclassification and correct classification. Finally, calculate Gini Index for a clustering using the calculated probabilities.

#### our Kmeans

```
In [147]: # this is the function you are supposed to implement in kmeans.py
          centroids, label = kmeans(X,3)
          # this is the function you are supposed to implement in kmeans.py
          evaluation(label, y)
          gini(label, y)
          SSE(centroids, X)
          plot_result('ours')
>>> initial centroids
[[0.70886076 0.68181818 0.65217391 0.6
                                             ]
                                             1
 [0.6835443  0.68181818  0.65217391  0.6
 [0.72151899 1.
                        0.2173913 0.16
                                             ]]
number of iterations to converge: 6
>>> final centroids
[[0.84045359 0.68560606 0.80676329 0.8225
                                             ]
 [0.74853944 0.62237762 0.62263099 0.53
                                             ]
 [0.63367089 0.77681818 0.21217391 0.0976
                                             ]]
Gini Index: 0.025600000000000008
SSE_each_cluster: [1.2157725785900486, 1.0302179676457113, 0.5860971086851517]
```

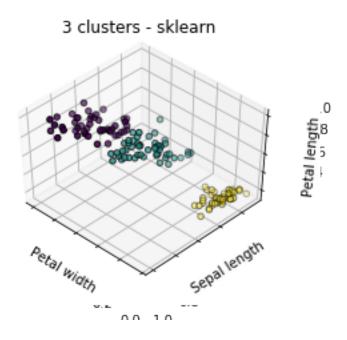


#### sklearn

```
In [152]: est = KMeans(n_clusters=3)
    fig = plt.figure(1, figsize=(4, 3))
    ax = Axes3D(fig, rect=[0, 0, .95, 1], elev=48, azim=134)
    est.fit(X)
    labels = est.labels_

# this is the function you are supposed to implement in kmeans.py
    evaluation(labels, y)
    gini(labels, y)
    plot_result('sklearn')
```

Gini Index: 0.025600000000000008



#### Discussion

Comparing the evaluation of our kmeans and the kmeans built in the sklearn package, the Gini Index of the best case in the 10 tests run with our kmeans is equivalent to the Gini Index calculated from sklearn.

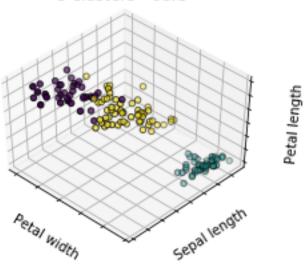
#### **O**4

4. In lab3.py, the data points are normalized (see line 12) by default as a data preprocessing step. What happens if you use the raw data (line 11) without any data preprocessing? Between normalized and unnormalized datasets, which one obtains better clustering? Please make comparison in terms of Impurity (Gini Index) of the clusterings and computational cost (number of iterations to converge).

#### un-normalize

```
In [158]: X = iris.data
In [161]: # this is the function you are supposed to implement in kmeans.py
          centroids, label = kmeans(X,3)
          # this is the function you are supposed to implement in kmeans.py
          evaluation(label, y)
          gini(label, y)
          SSE(centroids, X)
         plot_result('ours')
>>> initial centroids
[[7.1 3. 5.9 2.1]
[6. 2.9 4.5 1.5]
[6.5 3. 5.5 1.8]]
number of iterations to converge: 8
>>> final centroids
ΓΓ6.85
             3.07368421 5.74210526 2.07105263]
                     1.464
[5.006
             3.418
                                  0.244
 [5.9016129 2.7483871 4.39354839 1.43387097]]
Gini Index: 0.1216
SSE_each_cluster: [23.87947368421052, 15.24040000000001, 39.82096774193548]
```

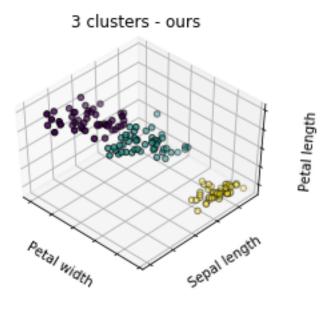
## 3 clusters - ours



#### normalize

In [162]: X = norm(iris.data)

```
In [169]: # this is the function you are supposed to implement in kmeans.py
          centroids, label = kmeans(X,3)
          # this is the function you are supposed to implement in kmeans.py
          evaluation(label, y)
          gini(label, y)
          SSE(centroids, X)
          plot_result('ours')
>>> initial centroids
[[0.73417722 0.61363636 0.73913043 0.76
                                             ]
 [0.83544304 0.68181818 0.63768116 0.56
                                             ]
                                             ]]
 [0.70886076 0.65909091 0.52173913 0.52
number of iterations to converge: 5
>>> final centroids
[[0.84045359 0.68560606 0.80676329 0.8225
                                             ]
 [0.74853944 0.62237762 0.62263099 0.53
                                             ]
 [0.63367089 0.77681818 0.21217391 0.0976
                                             ]]
Gini Index: 0.025600000000000008
SSE_each_cluster: [1.2157725785900486, 1.0302179676457113, 0.5860971086851517]
```



#### Discussion

From the evaluation results, we can see that the normalized dataset obtains better clustering, reducing the Gini Index by approximately 0.096 from 0.1216 to 0.0256. And with regard to the

computational cost, the number of iterations consumed to for the algorithm trained by unnormalized dataset to converge is 8 and the number of iterations for the algorithm trained by normalized dataset is 5 (however by running the test for a few times, the numbers of iterations took for both unnormalized and normalized dataset to converge are mostly similar and not significantly different)

#### Q5

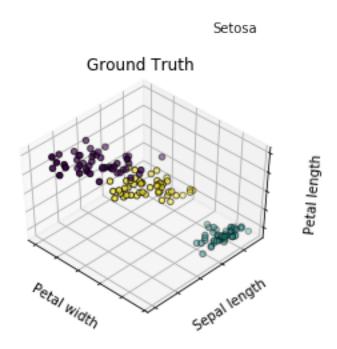
5. By comparing the clusterings in (3) and (4) against the ground truth, explain whether Impurity is a reasonable quality measure for clustering?

#### ground truth visualization

```
In [170]: fig = plt.figure(1, figsize=(4, 3))
          ax = Axes3D(fig, rect=[0, 0, .95, 1], elev=48, azim=134)
          for name, label in [('Setosa', 0),
                              ('Versicolour', 1),
                              ('Virginica', 2)]:
              ax.text3D(X[y == label, 3].mean(),
                        X[y == label, 0].mean(),
                        X[y == label, 2].mean() + 2, name,
                        horizontalalignment='center',
                        bbox=dict(alpha=.2, edgecolor='w', facecolor='w'))
          y = np.choose(y, [1, 2, 0]).astype(np.float)
          ax.scatter(X[:, 3], X[:, 0], X[:, 2], c=y, edgecolor='k')
          ax.w_xaxis.set_ticklabels([])
          ax.w_yaxis.set_ticklabels([])
          ax.w_zaxis.set_ticklabels([])
          ax.set_xlabel('Petal width')
          ax.set_ylabel('Sepal length')
          ax.set_zlabel('Petal length')
          ax.set_title('Ground Truth')
          ax.dist = 12
          plt.show()
```

# Virginica

#### Versicolour



#### Discussion

Impurity is a reasonable quality measure for clustering from the clusterings we got from (3) and (4). Because:

- Impurity (in this case, the Gini Index) could be used as evaluation criteria for the algorithm to determine its optimality:
  - \* In question 3, Gini Index is used for comparison of clusterings obtained between our KMeans and the one from sklearn package;
  - \* In question 4, Gini Index is used for comparison of clusterings obtained between KMeans models trained with normalized and unnormalized dataset