Assignment 3

```
In [1]: import pandas as pd
   import numpy as np
   import seaborn as sns
   import matplotlib.pyplot as plt
   from sklearn.neighbors import KNeighborsClassifier
```

```
In [2]: import pymc as pm
```

WARNING (pytensor.configdefaults): g++ not available, if using conda: `conda install m2w64-toolchain` WARNING (pytensor.configdefaults): g++ not detected! PyTensor will be unable to compile C-implementations and will d efault to Python. Performance may be severely degraded. To remove this warning, set PyTensor flags cxx to an empty st ring.

WARNING (pytensor.tensor.blas): Using NumPy C-API based implementation for BLAS functions.

```
In [3]: # import warnings filter
from warnings import simplefilter
# ignore all future warnings
simplefilter(action='ignore', category=FutureWarning)
```

Exercise 1 Bayesian exercises - Theory

Answer:

- 1. $logp_{\theta}(x)$ is the prior probability of the data. Since it does not dependent on z, it is essentially a normalization constant. In addition, E(a) = a, if a is a constant. So we could write as line 1.
- 2. the product rule is used here. p(z, x) = p(z|x)p(x)
- 3. Multiply both the numerator and denominator by $q_{\phi}(z|x)$
- 4. The Logarithmic Formula and the properties of Expectation are used here. $log_a(MN) = log_aM + log_aN$ and E(X+Y) = E(X) + E(Y)

Exercise 2 Bayesian exercises - Practical

Answer:

- 1. $x = pm.Data('x', x_data)$ should be corrected as $x = pm.Data('x', x_data, mutable=True)$ Because in this way, the x is keeping updated to predict y.
- 2. s = pm.Normal("sigma", sigma=0.001) should be corrected as s = pm.HalfNormal("sigma", sigma=0.001) Because the sigma should be a non-negative value.
- 3. likelihood = pm.Normal("y", mu=a*x + b, sigma=s) should be corrected as likelihood = pm.Normal("y", mu=a*x + b, sigma=s, observed=y_data) Because the information of the observed data is needed to define the likelihood function.
- 4. trace = pm.sample(1000, tune=1000, init=None, step=step, cores=2) should be corrected as trace = pm.sample(1000, tune=1000, init=None, step=step, chains=2) The "cores" means the number of chains to run in parallel processing tasks. We need to define "chains" in the MCMC method here.

```
p.s.
y = pm.Data('y_obs', y_data) is not necessary.

a = pm.Normal("slope", mu=100, sigma=100)
b = pm.Normal("intercept", mu=100, sigma=100)
s = pm.Normal("sigma", sigma=0.001) The value of mu and sigma is weird, but it still could work.
```

Exercise 3 Clustering I

```
In [4]: from sklearn.cluster import KMeans
```

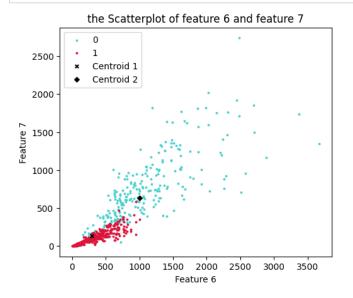
```
In [5]: import numpy as np
         # read in the data
         dataTrain = np.loadtxt('IDSWeedCropTrain.csv', delimiter=',')
         dataTest = np.loadtxt('IDSWeedCropTest.csv', delimiter=',')
         # split input variables and labels
         XTrain = dataTrain[:, :-1]
         YTrain = dataTrain[:, -1]
         XTest = dataTest[:, :-1]
         YTest = dataTest[:, -1]
In [6]: from sklearn.cluster import KMeans
         startingPoint = np.vstack((XTrain[0,], XTrain[1,]))
         kmeans = KMeans(2, n_init = 1, init= startingPoint, algorithm = "full").fit(XTrain) #set k=2
         print(kmeans.cluster_centers_)
         [[5.69426752e+00 4.93800425e+01 7.91594480e+02 3.84771338e+03
           3.38588535e+03 1.35988535e+03 2.93734607e+02 1.31609342e+02
           7.07282378e+01 3.96433121e+01 1.94437367e+01 4.23566879e+00
           4.41613588e-011
          [2.19092628e+00 1.37315690e+01 1.70943289e+02 1.39436484e+03
           3.18853497e+03 2.62461815e+03 1.00372023e+03 6.32814745e+02
           4.95829868e+02 2.95400756e+02 1.45809074e+02 2.91984877e+01
           2.83742911e+00]]
         C:\Users\dell\anaconda3\envs\IDS-A1\lib\site-packages\sklearn\cluster\_kmeans.py:1036: UserWarning: KMeans is known t
         o have a memory leak on Windows with MKL, when there are less chunks than available threads. You can avoid it by sett
         ing the environment variable OMP_NUM_THREADS=4.
           warnings.warn(
In [7]: df = pd.DataFrame()
         df["feature6"] = XTest[:, 6]
         df["feature7"] = XTest[:, 7]
df["feature1"] = XTest[:, 1]
         df["feature2"] = XTest[:, 2]
         df["label"] = YTest
         type0 = df[df['label'].isin([0])]
         type1 = df[df['label'].isin([1])]
         df_type0 = pd.DataFrame()
         df_type0["feature6"] = type0["feature6"]
        df_type0["feature7"] = type0["feature7"]
df_type0["feature1"] = type0["feature1"]
df_type0["feature2"] = type0["feature2"]
         df_type0["label"]= type0["label"]
         df_type1 = pd.DataFrame()
         df_type1["feature6"] = type1["feature6"]
        df_type1["feature7"] = type1["feature7"]
df_type1["feature1"] = type1["feature1"]
df_type1["feature2"] = type1["feature2"]
```

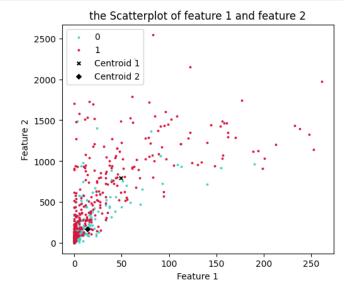
df_type1["label"]= type1["label"]

```
In [8]:
    fig = plt.figure(figsize=(12,10))
    ax1 = fig.add_subplot(2,2,1,title='the Scatterplot of feature 6 and feature 7',xlabel='Feature 6',ylabel='Feature 7')
    ax2 = fig.add_subplot(2,2,2,title='the Scatterplot of feature 1 and feature 2',xlabel='Feature 1',ylabel='Feature 2')
    ax1.scatter(data=df_type0,x='feature6',y='feature7',color='mediumturquoise',s=3,label='0')
    ax1.scatter(data=df_type0,x='feature6',y='feature7',color='crimson',s=3,label='1')
    ax1.scatter(x=293.73, y=131.61,s=15,marker='x',c='black',label='Centroid 1')
    ax1.scatter(x=1003.72, y=632.81,s=15,marker='D',c='black',label='Centroid 2')
    ax2.scatter(data=df_type0,x='feature1',y='feature2',color='mediumturquoise',s=3,label='0')
    ax2.scatter(data=df_type1,x='feature1',y='feature2',color='crimson',s=3,label='1')
    ax2.scatter(x=49.38, y=791.59,s=15,marker='x',c='black',label='Centroid 1')
    ax2.scatter(x=13.73, y=170.94,s=15,marker='D',c='black',label='Centroid 2')

ax1.legend(loc='upper left')
    fig.subplots_adjust(wspace=1)
    fig.tight_layout(pad=5)

plt.show()
```





1) The description of software used:

At first, two random points are chosen as the centroids, and the points are assigned to two centroids (two clusters formed). Then, the distances between points and centroids are calculated by iterating centroids and data assignment, until the sum of distances is to a minimum. In the end, the optimized centroids are found and the iteration stops.

2) Values of two cluster centers:

The first cluster center: [5.69, 49.38, 791.59, 3847.71, 3385.89, 1359.89, 293.73, 131.61, 70.73, 39.64, 19.44, 4.24, 0.44] The second cluster center: [2.19, 13.73, 170.94, 1394.36, 3188.53, 2624.62, 1003.72, 632.81, 495.83, 295.4, 145.8, 29.2, 2.84]

3) Two plots: The plots show different clusters of Test data (the blue and red "clusters" are defined by the labels, not by the algorithm). In the left plot, the two clusters have distinct characteristics. But in the right plot, two clusters are overlapped more. p.s. the centroids are plotted just for verifying. In the two plots, we could see the location of the centroids make sense (to some extent). So, the k-means algorithm works well here.