Principal Component Analysis

The goal of this question is to build a conceptual understanding of dimensionality reduction using PCA and implement it on a toy dataset. You'll only have to use numpy and matplotlib for this question.

```
In [62]:
```

```
import numpy as np
import matplotlib.pyplot as plt
# (a) Load data (features) and normalize the data
def load_data():
    data=np.load('features.npy')
    data_normalized=(data-np.mean(data,axis=0))/np.std(data,axis=0)
    return data_normalized
P=load_data()
```

In [63]:

```
# (b) Perform eigen decomposition and return eigen pairs in desecending order of eigen values
def eigendecomp(X):
          n=len(X)
         C = np.cov(X.T)
         print(C.shape)
          # Eigen decomposition
         eigenValues, eigenVectors = np.linalg.eig(C)
          idx = eigenValues.argsort()[::-1]
          sorted_eig_vals = eigenValues[idx]
         sorted eig vecs = eigenVectors[:,idx]
         return (sorted eig vals, sorted eig vecs)
sorted_eig_vals, sorted_eig_vecs=eigendecomp(P)
print("Sorted Eigen Values are :"+str(sorted eig vals))
print("Sorted Eigen Vectors are :"+str(sorted_eig_vecs))
(8, 8)
Sorted Eigen Values are: [ 4.74298961e+00 2.29585309e+00 7.76910512e-01 2.04172901e-01
    3.37651661e-02 8.98851439e-16 -1.31385926e-16 -6.12506839e-16]
Sorted Eigen Vectors are :[[-0.39124937  0.13884872 -0.46160937  0.58034539  0.24934936  0.09502813
       0.41458399 0.15875925]
   [ 0.11687696 - 0.4391715 - 0.78711289 - 0.2905579 - 0.12725786 0.15733318 ]
     -0.20945618 0.09463144]
   [-0.40655289 \quad 0.29080021 \quad -0.13961871 \quad -0.12636707 \quad -0.54994554 \quad 0.17331863 
      0.09114106 -0.61127232]
   [-0.39944906 \quad 0.26454833 \quad -0.16206048 \quad -0.54404218 \quad 0.49904279 \quad -0.43946377 ] 
     -0.05284949 -0.0484821 ]
  [-0.3778555 \quad -0.35426671 \quad 0.07790627 \quad 0.42060984 \quad 0.12822569 \quad -0.14932029
     -0.65144711 -0.24946272]
    [-0.09816172 \ -0.64299795 \ \ 0.11941452 \ -0.04972667 \ -0.0795516 \ \ -0.43247222 ] 
      0.57574617 -0.26011976]
    [-0.45509399 \ -0.03231459 \ \ 0.12200908 \ -0.08034689 \ -0.51935676 \ -0.19102316 ] 
    -0.10045113 0.673713911
  [-0.38587285 \ -0.30545597 \quad 0.29393481 \ -0.28457653 \quad 0.27864817 \quad 0.7051450681 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ -0.28457653 \ 
       0.08480007 0.07779233]]
```

In [64]:

```
return d
print("Variance gives the measure of contribution of each feature. When K is 3 it captures around
97% data. I would select k=3 as computational cost is reduced by lesser features")
```

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In [65]:

```
# (d) Visualize after projecting to 2-D space
def viz():
    eig vect=sorted eig vecs[:,0:2]
    datanew=P@eig vect
    datanew=datanew[:,:3]
    ynew=datanew.T
    l=np.load('labels.npy', allow_pickle=True)
    ynewplot1=ynew.T[:,0]
    ynewplot2=ynew.T[:,1]
    plt.title("Data in 2-D Space")
    unique = list(set(l))
    colors = [plt.cm.jet(float(i)/max(unique)) for i in unique]
    for i, u in enumerate(unique):
        xi = [ynewplot1[j] for j in range(len(ynewplot1)) if 1[j] == u]
yi = [ynewplot2[j] for j in range(len(ynewplot2)) if 1[j] == u]
         plt.scatter(xi, yi, c=colors[i], label=str(u))
    plt.legend()
    plt.show()
```

In [66]:

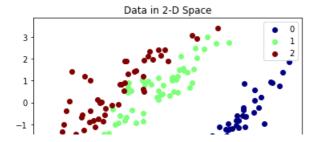
```
def main():
    eval()
    viz()

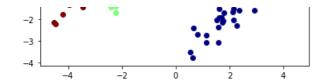
if __name__ == "__main__":
    main()
```

'c' argument looks like a single numeric RGB or RGBA sequence, which should be avoided as value-ma pping will have precedence in case its length matches with 'x' & 'y'. Please use a 2-D array with a single row if you really want to specify the same RGB or RGBA value for all points. 'c' argument looks like a single numeric RGB or RGBA sequence, which should be avoided as value-ma pping will have precedence in case its length matches with 'x' & 'y'. Please use a 2-D array with a single row if you really want to specify the same RGB or RGBA value for all points. 'c' argument looks like a single numeric RGB or RGBA sequence, which should be avoided as value-ma pping will have precedence in case its length matches with 'x' & 'y'. Please use a 2-D array with a single row if you really want to specify the same RGB or RGBA value for all points.

```
Variance
```

```
{'k0 = ': 0.5889212098295767, 'k1 = ': 0.8739896347022307, 'k2 = ': 0.9704560233211401, 'k3 = ': 0
.9958074918820436, 'k4 = ': 0.99999999999999, 'k5 = ': 1.0, 'k6 = ': 1.0, 'k7 = ': 1.0}
Eigenvalues
k= 1 = 4.74298960936572
k= 2 = 2.2958530862227167
k= 3 = 0.7769105123670564
k= 4 = 0.20417290116163855
k= 5 = 0.03376516605065379
k= 6 = 8.988514394516542e-16
k= 7 = -1.3138592633207446e-16
k= 8 = -6.125068394845793e-16
```





(e1): If the number of features is 1000 and the number of data points is 10, what will be the dimension of your covariance matrix? Can you suggest what can be changed to improve the performance?

Answer e1- If there are 1000 features and 10 data points the dimension of covariance matrix will be (1000,1000). The Performance can be improved by reducing the number of features by PCA, using only those features which can explain the data well. The performance can also be improved by collecting more data points.

(e2): Assume you have a dataset with the original dimensionality as 2 and you have to reduce it to 1. Provide a sample scatter plot of the original data (less than 10 datapoints) where PCA might produce misleading results. You can plot it by hand and then take a picture. In the next cell, switch to Markdown mode and use the command: ![title]()

