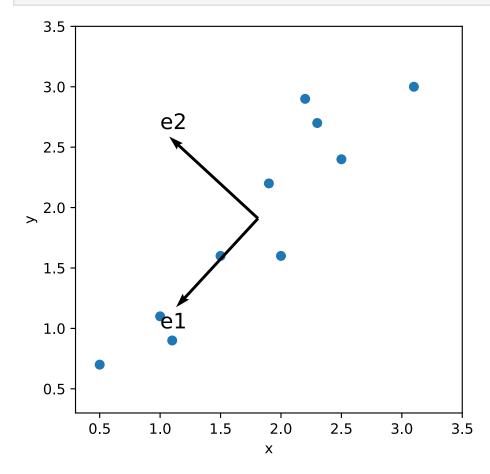
Q1: PCA

(a) Principle Component and Transformed space

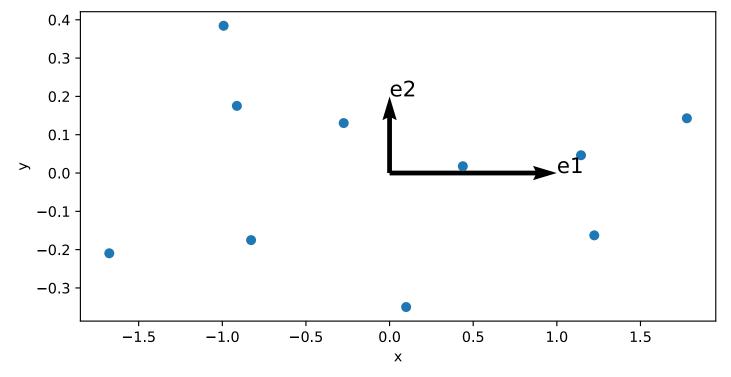
1.

```
In [115...
          import numpy as np
          import matplotlib.pyplot as plt
In [116...
         data = np.loadtxt('qladata.txt')
          mean_a = np.mean(data,axis = 0)
          D = data - mean_a
          n = len(data)
          cov_a = D.T@D
          v,d = np.linalg.eig(cov_a)
          idx = v.argsort()[::-1]
          v = v[idx]
          d = d[:,idx]
          print("eigenvalues: ",v)
          print("eigenvector: ",d)
         eigenvalues: [11.55624941 0.44175059]
         eigenvector: [[-0.6778734 -0.73517866]
          [-0.73517866 0.6778734 ]]
In [117... fig,ax=plt.subplots(figsize = (5,5))
          ax.set_aspect('equal', adjustable='box')
          ax.scatter(data[:,0],data[:,1])
          e1 = d[:,0]
          e2 = d[:,1]
          a = ax.quiver(mean_a[0], mean_a[1], e1[0], e1[1], angles='xy', scale_units='xy', scale=1)
          b = ax.quiver(mean_a[0], mean_a[1], e2[0], e2[1], angles='xy', scale_units='xy', scale=1)
          ax.set xlabel("x")
          ax.set_ylabel("y")
          ax.set_xlim(0.3,3.5)
          ax.set_ylim(0.3,3.5)
          ax.text(1,2.65,"e2",size = 15)
          ax.text(1,1,"e1",size = 15)
          plt.show()
```



2.

```
In [118... A = D@d A
```



3.

Out[120... 3.453381743924969

Out[118... array([[-0.82797019, -0.17511531],

I think e1 is the optimal one-dimentional representation of the data because it can capture more information of the data. And the range of data is 3.453.

(b)

1.

```
data_b = np.loadtxt('q1bdata.txt')
In [121...
          mean_b = np.mean(data_b,axis = 0)
          D_b = data_b - mean_b
          n = len(data b)
          cov_b = D_b@D_b.T
          v,d = np.linalg.eig(cov_b)
          idx = v.argsort()[::-1]
          v = v[idx]
          d = d[:,idx]
          print("covariance: ",cov_b)
          print("eigenvalues: ",v)
          print("eigenvector: ",d)
         covariance: [[ 69.875 -18.875 -26.375 -24.625]
          [-18.875 121.375 -53.125 -49.375]
          [-26.375 -53.125 98.375 -18.875]
          [-24.625 -49.375 -18.875 92.875]]
         eigenvalues: [175.55118219 114.31705238 92.63176543 0.
         eigenvector: [[-0.06628148 0.04124587 -0.86249959 -0.5
          [-0.79038331 - 0.06822502 0.34733208 - 0.5]
          [ 0.47285044 -0.69123739  0.22046165 -0.5
          [ 0.38381435  0.71821654  0.29470586 -0.5
                                                          ]]
```

There are three non-zero eigenvalues which means there are three effective eigenvectors.

2.

```
In [122... u = D b.T@d[:,0:3]
          # normalize u
          u = u/np.linalg.norm(u,axis=0)
          Projection = D_b@u
          print("projection: ",Projection)
         projection: [[ -0.8782013
                                     0.44099733 -8.3011616 ]
          [-10.47224127 -0.72945617 3.34291139]
          [ 6.26506632 -7.39065157 2.12184196]
          [ 5.08537624 7.67911041 2.83640825]]
        3.
          omega = mean b.reshape(1,-1) + Projection@u.T
In [123...
          mse = np.mean((data_b-omega)**2,axis = 1)
          print("MSE: ",mse)
         MSE: [6.31380655e-31 2.30252020e-30 3.01077587e-30 3.12689931e-30]
        4.
In [124... u2 = D_b.T@d[:,0:2]
          # normalize u
          u2 = u2/np.linalg.norm(u2,axis=0)
          Projection2 = D_b@u2
          print("projection2: ",Projection2)
          omega2 = mean_b.reshape(1,-1) + Projection2@u2.T
          mse2 = np.mean((data b-omega2)**2,axis = 1)
          print("MSE: ",mse2)
         projection2: [[ -0.8782013
                                        0.44099733]
          [-10.47224127 -0.72945617]
          [ 6.26506632 -7.39065157]
          [ 5.08537624 7.67911041]]
         MSE: [3.62680441 0.58816087 0.2369586 0.4234322 ]
        5.
         Y = np.array([1,3,0,3,-2,2,4,1,3,0,-2,0,1,1,-3,0,1,-2,-3])
In [125...
In [126...
          ED = np.linalg.norm(Y-omega,axis = 1)
In [127...
Out[127... array([12.92284798, 5.65685425, 16.79285562, 15.77973384])
```

The second sample is the most similar one to this new vector.

6.

```
In [128... ED2 = np.linalg.norm(Y-data_b,axis = 1)
    ED2
Out[128... array([12.92284798, 5.65685425, 16.79285562, 15.77973384])
```

This result is the same as part(5). Although there are 19 dimensions in original space, there are only three informative eigenvectors, we can reduce dimensions to 3 without loss.

Q2: Eigenface with PCA

(a)

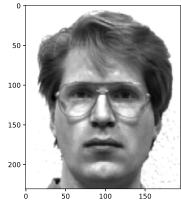
```
import scipy.io
from PIL import Image
mat = scipy.io.loadmat("data-1.mat")
print(mat.keys())

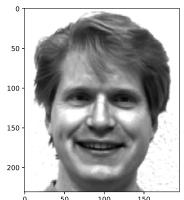
dict_keys(['__header__', '__version__', '__globals__', 'Y', 'X', 'testimages', 'trainimages'])

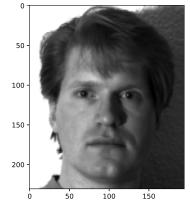
In [130... X = mat["X"]
```

```
fig,axs = plt.subplots(1,5,figsize = (25,5))
for i in range(5):
    face = np.transpose(X[:,i].reshape(195,231))
    axs[i].imshow(face,cmap = "gray")
```





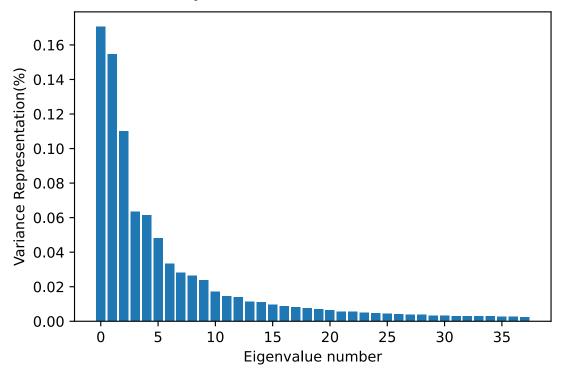






(b)

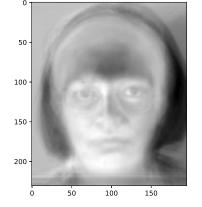
```
In [132... from sklearn.decomposition import PCA
    mean = np.mean(X,axis = 1)
    D = X-mean.reshape(-1,1)
    pca = PCA(n_components=0.9)
    pca.fit(X.T)
    print(pca.explained_variance_ratio_)
    m = np.arange(0,pca.n_components_,1)
    plt.bar(m,pca.explained_variance_ratio_)
    plt.xlabel("Eigenvalue number")
    plt.ylabel("Variance Representation(%)")
    plt.show()
```

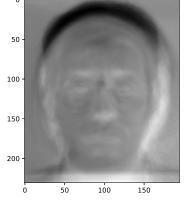


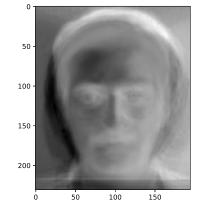
```
In [133... A = pca.transform(X.T)

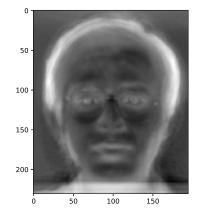
u = D@A
```

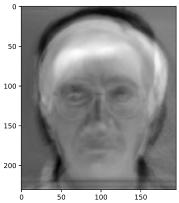
fig,axs = plt.subplots(1,5,figsize = (25,5))
for i in range(5):
 face = np.transpose(u[:,i].reshape(195,231))
 axs[i].imshow(face,cmap = "gray")











(c)

```
In [135...
          fig,axs = plt.subplots(1,6,figsize = (30,5))
          face = np.transpose(X[:,0].reshape(195,231))
          axs[0].imshow(face,cmap = "gray")
          axs[0].set title("original")
          eigen_num = [10, 20, 30, 40, 50]
          for i in eigen_num:
               pca = PCA(n_components = i)
              pca.fit(X.T)
              A = pca.transform(X.T)
               approx = pca.inverse_transform(A)
               face = np.transpose(approx.T[:,0].reshape(195,231))
               axs[eigen_num.index(i)+1].imshow(face,cmap = "gray")
               axs[eigen_num.index(i)+1].set_title(f"eigenfaces num = {i}")
                                    eigenfaces num = 10
                                                                             eigenfaces num = 30
                  original
                                                        eigenfaces num = 20
                                                                                                                       eigenfaces num = 50
                                                                                                  eigenfaces num = 40
         (d)
          from sklearn.neighbors import KNeighborsClassifier
In [136...
          train = mat['trainimages'].ravel()-1
          test = mat['testimages'].ravel()-1
          Y = mat['Y'].ravel()
          pca = PCA(n components = 30)
In [137...
          A1 = pca.fit transform(X[:, train].T)
          train_X = pca.inverse_transform(A1)
          A2 = pca.fit_transform(X[:, test].T)
          test_X = pca.inverse_transform(A2)
          score_30 = KNeighborsClassifier(n_neighbors=1).fit(train_X, Y[train]).score(test_X, Y[test])
In [138...
          score_original = KNeighborsClassifier(n_neighbors=1).fit(X[:,train].T, Y[train]).score(X[:,test].T, Y[test])
          print("accuracy of reduced dimension space: ",score_30)
          print("accuracy of original dimension space: ",score_original)
```

Yes, this is what I expected. We reduce dimensions from 45045 to 30, but the accuracy only decreases 3.33%, which means these 30 dimensions capture most information of the data.

accuracy of reduced dimension space: 0.8

accuracy of original dimension space: 0.8333333333333333