CS 613 - Machine Learning

Assignment 3 - Dimensionality Reduction & Clustering

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CS 613 Machine Learning
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Introduction

In this assignment you'll work on visualizing data, reducing its dimensionality and clustering it.

You may not use any functions from machine learning library in your code, however you may use statistical functions. For example, if available you **MAY NOT** use functions like:

- pca
- · k-nearest neighbors functions

Unless explicitly told to do so.

But you MAY use basic statistical functions like:

- std
- mean
- COV
- eig
- svd

Submission

For your submission, upload to Blackboard a single zip file containing:

- 1. A LaTeX typeset PDF or Jupyter Notebook PDF containing:
 - (a) Part 1: Your answers to the theory questions.
 - 1.1a. The principle components are:
 - 1.1b. The data projected onto the to 1st principle component:
 - 1.2a.

- x1 information gain:
- x2 information gain:

1.2b.

1.2c. The directions corresponding to eignevectors are:

1.2d The data projected is:

1.2e



(b) Part 2: The visualization of the PCA result, KNN accuracies.

KNN score, k=1, no PCA:

KNN score, k=1, PCA100:

KNN score whitened:

Project PCA2 (image):



- (c) Part 3:
- (d) Part 4:
- (e) Source Code python notebook

```
In [2595]: import pandas as pd
           import numpy as np
           import random
           import matplotlib.pyplot as plt
           import numpy.matlib
           from sklearn.preprocessing import scale
           from sklearn.decomposition import PCA
           from sklearn.metrics import accuracy_score
           from scipy.spatial.distance import cdist
```

DataSets

Labeled Faces in the Wild Dataset. This dataset consists of celebrities download from the Internet from the early 2000s. We use the grayscale version from sklearn.datasets. We will download the images in a specific way as shown below. You will have 3,023 images, each 87x65 pixels large, belonging to 62 different people.



Part 1: Theory Questions

1.1 Consider the following data:

Out[2227]:

	x1	x2
0	-2.0	1.0
1	-5.0	-4.0
2	-3.0	1.0
3	0.0	3.0
4	-8.0	11.0
5	-2.0	5.0
6	1.0	0.0
7	5.0	-1.0
8	-1.0	-3.0
9	6.0	1.0

a. Find the principle components of the data (you must show the math, including how you compute the eivenvectors and eigenvalues). Make sure you standardize the data first and that your principle components are normalized to be unit length. As for the amount of detail needed in your work imagine that you were working on paper with a basic calculator. Show me whatever you would be writing on that paper. (5pts)

```
In [2228]: import math
            m1=np.sum(x1)/10
            var = np.sum((x1 - m1) ** 2) / 9
            std = var ** 0.5
            m2=np.sum(x2)/10
            var1= np.sum((x2 - m2) ** 2) / 9
            std1= var1 ** 0.5
            def Standardise(x):
                 return (x-m1)/std
            def Standardisel(x):
                 return (x-m2)/std1
            x1=Standardise(x1)
            x2=Standardise1(x2)
            x11=np.array([x1,x2])
            x11=x11.T
            print(x11)
            [[-0.26015724 -0.09359019]
             [-0.96967699 -1.2634676 ]
             [-0.49666382 -0.09359019]
             [ 0.21285592  0.37436077]
             [-1.67919674 2.24616461]
             [-0.26015724 \quad 0.84231173]
             [ 0.44936251 -0.32756567]
             [ 1.39538884 -0.56154115]
             [-0.02365066 -1.02949211]
             [ 1.63189542 -0.09359019]]
In [1605]: #Covariance
            A=(x11.T@x11)/(x11.shape[0]-1)
            df.cov()
            Α
Out[1605]: array([[ 1.
                                  , -0.40826218],
                    [-0.40826218, 1.
                                                ]])
In [1531]: 0.1666780076183524-1
Out[1531]: -0.8333219923816476
            (A - \lambda I)X=0
            [ [(1.-\lambda), -0.40826218]
              [(-0.40826218, (1.-\lambda)]] *X= 0
            Find the determinant of A-\lambda I
            Determinant = (1.-\lambda)*(1.-\lambda) - (-0.40826218*-0.40826218)
            1. -\lambda - \lambda + \lambda^2 -0.1666780076183524=0
            0.8333219923816476 -2\lambda + \lambda^2 = 0
            The equation to find the eigen value is
            \lambda^2 - 2\lambda + 0.8333219923816476 = 0
```

```
The equation to find X in a quadratic equation is x=-b \pm \text{sqrt}(b^2-4ac) / 2a In the above equation b=-2 a=1 c=0.8333219923816476 x1=(-(-2)+\text{math.sqrt}(((-2)**2)-(4*1*0.8333219923816476)))/2 \lambda 1=1.40826218 \ \lambda 2=0.5917378200000001 are the Eigen values
```

```
In [1535]: (2+math.sqrt(((-2)**2)-(4*1*0.8333219923816476)))/2
           (2-math.sqrt(((-2)**2)-(4*1*0.8333219923816476)))/2
           1.-1.40826218
           -0.4082621799999999*1/0.40826218
           -0.40826218*1/0.408262179999999
           el=np.array([[1],[0.99999999999999]])
           print(e1.T@e1)
           e11=np.array([[1],[1.0000000000000024]])
           print(e11.T@e11)
           vector=(1/math.sqrt(2))*e11
           vector
           [[2.]]
           [[2.]]
Out[1535]: array([[0.70710678],
                  [0.70710678]]
In [1507]: 0.4082621799999999*1/0.40826218
           -0.40826218*1/-0.4082621799999999
           el=np.array([[1],[0.99999999999999]])
           print(e1.T@e1)
           e11=np.array([[1],[1.0000000000000024]])
           print(e11.T@e11)
           1-0.591737820000001
```



```
Eigen Vector for \lambda 1=1.40826218
(A-\lambda I)X=0
Let X be [X1,X2]
([[1.,-0.40826218] [[1.40826218, 0]
                                           ([X1,
                                                     [0,
 [-0.40826218, 1.]] - [0, 1.40826218]]) * X2]) = 0]
[[1.-1.40826218, -0.40826218]
                                 [X1, [0,
 [-0.40826218, 1.-1.40826218] * X2] = 0]
[[-0.4082621799999999, -0.40826218] [X1, [0,
 [-0.40826218, -0.4082621799999999] * X2]= 0]
-0.408262179999999911 -0.40826218X2 = 0
-0.40826218X1 -0.4082621799999999X2 = 0
Lets consider X1 as 1 then
-0.4082621799999999*1=0.40826218X2
-0.40826218*1=0.408262179999999922
for X1=1, X1=1
X2=-0.9999999999999999999 X2=-1.0000000000000024
the matrix [X1,
            X2] is -0.999999999999999] , -1.0000000000000024]
So the eigenvector corresponding to eigenvalue \lambda 1 = 1/\sqrt{2} \times [1, 1]
-0.9999999999999991
```

```
= [0.70710678,
                                                     -0.70710678]
                                                or
                                                   = 1/\sqrt{2*[1,
-1.00000000000000241
                   Eigen Vector for \lambda 1 = 1.40826218 = [0.70710678]
                                                      -0.707106781
Eigen Vector for \lambda 2 = 0.5917378200000001
(A-\lambda I)X=0
Let X be [X1,X2]
([[1.,-0.40826218] [[0.5917378200000001, 0] ([X1, [0,
 [-0.40826218, 1.] - [0, 0.5917378200000001]) * X2]) = 0]
 [[0.4082621799999999, -0.40826218] [X1, [0,
 [-0.40826218, 0.4082621799999999]] * X2] = 0]
  0.408262179999999911 - 0.40826218X2 = 0
 -0.40826218X1 + 0.4082621799999999X2 = 0
Lets consider X1 as 1 then
 0.4082621799999999*1=0.40826218X2
 -0.40826218*1=-0.408262179999999922
for X1=1, X1=1
the matrix [X1, [1.,
            X2] is 0.999999999999999] , 1.000000000000024]
So the eigenvector corresponding to eigenvalue \lambda 1 = 1/\sqrt{2} \times [1, 1]
0.999999999999999
                                                   = [0.70710678]
                                                      0.70710678]
                                               or
                                                  = 1/\sqrt{2*[1,
1.00000000000000241
            Eigen Vector for \lambda 2=0.5917378200000001 = [0.70710678]
                                                      0.70710678]
```

b. Project the data onto the principal component corresponding to the largest eigenvalue found in the previous part (3pts)

```
The largest eigen value is \lambda 1 = 1.40826218 The corresponding eigen vector is [0.70710678, -0.70710678]
```

1.2. Consider the following data

```
In [2328]: td2.groupby(['y', 'x1', 'x2']).first()
```

Out[2328]:

Count

у	x1	x2	
0	-8.0	11.0	1
	-5.0	-4.0	1
	-3.0	1.0	1
	-2.0	1.0	1
	0.0	3.0	1
1	-2.0	5.0	1
	-1.0	-3.0	1
	1.0	0.0	1
	5.0	-1.0	1
	6.0	1.0	1

```
In [2329]: m1=np.sum(x1)/10
           m2=np.sum(x2)/10
           def threshold(x):
               for i in range(0,len(x)):
                    if x[i]<m1:
                        x[i]=0
                    else:
                        x[i]=1
                    i+=1
               return x
           def threshold1(x):
               for i in range(0,len(x)):
                    if x[i]<m2:
                        x[i]=0
                    else:
                        x[i]=1
                    i+=1
               return x
           x11=threshold(x1)
           x21=threshold1(x2)
           m2
```

Out[2329]: 1.4

```
In [2330]: td3 = pd.DataFrame({
    'x11': x11,
    'x21': x21,
    'y1': y,
    'Count1': 1
})
td3
```

Out[2330]:

	x11	x21	у1	Count1
0	0	0	0	1
1	0	0	0	1
2	0	0	0	1
3	1	1	0	1
4	0	1	0	1
5	0	1	1	1
6	1	0	1	1
7	1	0	1	1
8	0	0	1	1
9	1	0	1	1

(a) Compute the information gain for each feature. You could standardize the data overall, although it won't make a difference. (5pts).

```
In [2331]: hy = -5/10*np.log2(5/10)-5/10*np.log2(5/10)
hx1 = (6/10)*(-(4/6)*np.log2(4/6)-(2/6)*np.log2(2/6))+(4/10)*(-(1/4)*np.log
hx2 = (7/10)*(-(3/7)*np.log2(3/7)-(4/7)*np.log2(4/7))+(3/10)*(-(2/3)*np.log
infox1=hy-hx1
infox2=hy-hx2
print(infox1,infox2)
```

0.12451124978365313 0.034851554559677034

```
In [2332]: (6/10)*(-(4/6)*np.log2(4/6)-(2/6)*np.log2(2/6))+(4/10)*(-(1/4)*np.log2(1/4))
```

Out[2332]: 0.8754887502163469

```
In [2333]: print("information gain of x1=",infox1,"information gain of x2=",infox2)
```

information gain of x1=0.12451124978365313 information gain of x2=0.034851554559677034

(b) Which feature is more discriminating based on results in part a (1pt)?

Feature x1 is more discriminating as it has higher information gain

(c) Using LDA, find the direction of projection (you must show the math, however for this one you don't have to show the computation for finding the eigenvalues and eigenvectors). Normalize this vector to be unit length (5pts).

```
In [2439]: import math
           m12=np.sum(x1)/10
           var12 = np.sum((x1 - m1) ** 2) / 9
           std12 = var ** 0.5
           m22=np.sum(x2)/10
           var13 = np.sum((x2 - m2) ** 2) / 9
           std13= var1 ** 0.5
           def Standardise(x):
               return (x-m12)/std12
           def Standardisel(x):
               return (x-m22)/std13
           x1=Standardise(x1)
           x2=Standardise1(x2)
           x11=np.array([x1,x2])
           x11=x11.T
           print(m12, m22, std13, std12)
```

-0.9 1.4 4.273952113286562 4.2282121254470875

```
In [2441]: m_class1=np.mean(x11[:5,:],axis=0)
           m class2=np.mean(x11[5:,:],axis=0)
           print(m class1,m class2)
           c1=x11[:5,:]-m_class1
           c2=x11[5:,:]-m_class2
           print(c1)
           print(c2)
           [-0.63856777 \quad 0.23397548] \quad [\quad 0.63856777 \quad -0.23397548]
           [[ 0.37841053 -0.32756567]
            [-0.33110922 -1.49744308]
            [ 0.14190395 -0.32756567]
                           0.140385291
            [ 0.8514237
            [-1.04062896 2.01218913]]
           [[-0.89872501 1.07628721]
            [-0.18920527 -0.09359019]
            [ 0.75682106 -0.32756567]
            [-0.66221843 - 0.79551663]
            [ 0.99332765  0.14038529]]
In [2442]: covc1=4*(c1.T@c1/4)
           covc1
Out[2442]: array([[ 2.08079553, -1.64903489],
                   [-1.64903489, 6.52554745]])
In [2443]: covc2=4*(c2.T@c2/4)
           covc2
Out[2443]: array([[ 2.84151647, -0.53123272],
                   [-0.53123272, 1.9270073]]
In [2444]: sw=covc1+covc2
           SW
Out[2444]: array([[ 4.922312 , -2.18026761],
                   [-2.18026761, 8.45255474]])
In [2445]: mean1=m class1.reshape(2,1)
           mean2=m class2.reshape(1,2)
           sb=mean1@mean2
Out[2445]: array([[-0.4077688 , 0.1494092 ],
                   [0.1494092, -0.05474453]]
In [2446]: | swinv=np.linalg.inv(sw)
           swinv
Out[2446]: array([[0.22936149, 0.05916193],
                   [0.05916193, 0.13356776]])
```

(d) Project the data onto the principal component found in the previous part (3pts).

The largest eigen Value is 0 hence the eigen vector chosen for the projection of data is V[:,1]

```
In [2448]: print(V[:,1])
    print(m_class1@V[:,1])
    print(m_class2@V[:,1])

        [-0.34403943 -0.9389552 ]
        8.326672684688674e-17
        -8.326672684688674e-17

In [2449]: print(x11@V[:,1])
        [ 0.17738135    1.51994659    0.25874894 -0.42473882 -1.53133805 -0.70138863
```

0.15297107 0.0471932

(e) Does the projection you performed in the previous part seem to provide good class separation? Why or why not (1pt)?

0.97478373 - 0.47355938

LDA provides a better class seperation than PCA as the difference between the class values is higher.

Part 2: Dimensionality Reduction via PCA

Import the data as shown above. This the labeled faces in the wild dataset. Verify that you have the correct number of people and classes

```
In [2455]: print("people.images.shape: {}".format(people.images.shape))
    print("Number of classes: {}".format(len(people.target_names)))
    people.images.shape: (2537, 87, 65)
    Number of classes: 45
```

This dataset is skewed toward George W. Bush and Colin Powell as you can verify here:

```
In [2456]: # count how often each target appears
    counts = np.bincount(people.target)
    # print counts next to target names
    for i, (count, name) in enumerate(zip(counts, people.target_names)):
        print("{0:25} {1:3}".format(name, count), end=' ')
        if (i + 1) % 3 == 0:
            print()
```

Alejandro Toledo		39	Alvaro Uribe	35	Amelie Ma
uresmo	21	2.6	nuicl Chance	77	ntol Dibo
Andre Agassi	2.4	36	Ariel Sharon	77	Atal Biha
ri Vajpayee	24	20	a.1'. p11	226	D. 11 D.
Bill Clinton	2.1	29	Colin Powell	236	David Bec
kham	31	101	G D. l	2.2	G
Donald Rumsfeld	5 20	121	George Robertson	22	George W
	530	100			
Gerhard Schroeder		109	Gloria Macapagal Arroyo	44	Gray Davi
S	26	0.0		2.0	01
Hamid Karzai		22	Hans Blix	39	Hugo Chav
ez	71	•			_
Igor Ivanov		20	Jack Straw	28	Jacques C
hirac	52				
Jean Chretien		55	Jennifer Aniston	21	Jennifer
Capriati	42				
Jennifer Lopez		21	John Ashcroft	53	Juan Carl
os Ferrero	28				
Junichiro Koizumi		60	Kofi Annan	32	Laura Bus
h	41				
Lindsay Davenport		22	Lleyton Hewitt	41	Megawati
Sukarnoputri	33				
Michael Bloomberg		20	Pete Sampras	22	Ricardo L
agos	27				
Rudolph Giuliani		26	Saddam Hussein	23	Serena Wi
lliams	52				
Tiger Woods		23	Tom Daschle	25	Tom Ridge
33					
Tony Blair		144	Vicente Fox	32	Vladimir
Putin	49				

To make the data less skewed, we will only take up to 50 images of each person (otherwise, the feature extraction would be overwhelmed by the likelihood of George W. Bush):

```
In [2457]: mask = np.zeros(people.target.shape, dtype=np.bool)
    for target in np.unique(people.target):
        mask[np.where(people.target == target)[0][:50]] = 1

X_people = people.data[mask]
    y_people = people.target[mask]

# scale the grayscale values to be between 0 and 1
# instead of 0 and 255 for better numeric stability
X_people = X_people / 255.
```

<ipython-input-2457-0f53c27b3db9>:1: DeprecationWarning: `np.bool` is a d
eprecated alias for the builtin `bool`. To silence this warning, use `boo
l` by itself. Doing this will not modify any behavior and is safe. If you
specifically wanted the numpy scalar type, use `np.bool_` here.
Deprecated in NumPy 1.20; for more details and guidance: https://numpy.or
g/devdocs/release/1.20.0-notes.html#deprecations (https://numpy.org/devdo
cs/release/1.20.0-notes.html#deprecations)
mask = np.zeros(people.target.shape, dtype=np.bool)

```
In [2458]: y_people.shape
Out[2458]: (1577,)
```

We are now going to compute how well a KNN classifier does using just the pixels alone.

```
In [2459]: from sklearn.neighbors import KNeighborsClassifier
    from sklearn.model_selection import train_test_split
    # split the data into training and test sets
    X_train, X_test, y_train, y_test = train_test_split(X_people, y_people, str
    # build a KNeighborsClassifier using one neighbor
    knn = KNeighborsClassifier(n_neighbors=1)
    knn.fit(X_train, y_train)
    print("Test set score of 1-nn: {:.10f}".format(knn.score(X_test, y_test)))
```

Test set score of 1-nn: 0.2784810127

You should have an accuracy around 23% - 27%.

Once you have your setup complete, write a script to do the following:

- 1. Write your own version of KNN (k=1) where you use the SSD (sum of squared differences) to compute similarity
- 2. Verify that your KNN has a similar accuracy as sklearn's version
- 3. Standardize your data (zero mean, divide by standard deviation)
- 4. Reduces the data to 100D using PCA
- 5. Compute the KNN again where K=1 with the 100D data. Report the accuracy
- 6. Compute the KNN again where K=1 with the 100D Whitened data. Report the accuracy
- 7. Reduces the data to 2D using PCA
- 8. Graphs the data for visualization

Recall that although you may not use any package ML functions like *pca*, you may use statistical functions like *eig* or *svd*.

Your graph should end up looking similar to Figure 1 *refer to the PDF for the figure* (although it may be rotated differently, de-pending how you ordered things)

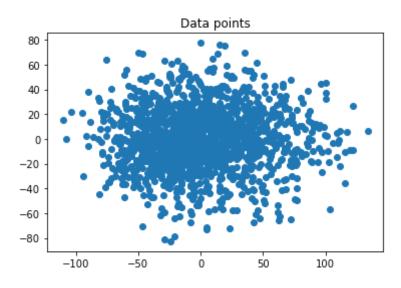
```
In [2460]: | df = pd.DataFrame(data=X people)
           def Standardise(string1):
               m1=np.mean(string1)
               std1=np.std(string1)
               string2=(string1-m1)/std1
               return string2
           df1=df.apply(Standardise)
           f=df1.to_numpy()
In [2461]: e1=df1.T.to_numpy()
           el.shape
Out[2461]: (5655, 1577)
In [2462]: | v=np.cov(e1)
In [2463]: v.shape
Out[2463]: (5655, 5655)
In [2464]: from numpy.linalg import eig
           val, vec=eig(v)
In [2301]: vec.shape
Out[2301]: (5655, 5655)
In [2465]: sorted index = np.argsort(val)[::-1]
           sorted eigenvalue = val[sorted index]
           #similarly sort the eigenvectors
           sorted eigenvectors = vec[:,sorted index]
In [2466]: n components = 100 #you can select any number of components.
           eigenvector subset = sorted eigenvectors[:,0:n components]
           eigenvector subset.shape
Out[2466]: (5655, 100)
In [2467]: X reduced = np.dot(eigenvector subset.transpose(),f.transpose()).transpose(
In [2468]: dff y=pd.DataFrame(data=y people)
In [2469]: X_train, X_test, y_train, y_test = train_test_split(X_reduced, y_people, st
```

```
In [2470]: from scipy.stats import mode
           def sumofsquare distance(row1, row2):
               dist = np.sum((row1-row2)**2)
               return dist
           '''def mode(label,val):
               for i in range(0,len(label)):
                   dict=dict(label[i],)
               if label in val:
                   dict'''
           def predict_classification(train, test,input1, num):
               output=[]
               value=[]
               for i in input1:
                   distances =list()
                   for j in range(len(train)):
                       distance = sumofsquare_distance(np.array(train[j,:]) , i)
                        distances.append(distance)
                   dist = np.argsort(distances)[:num]
                   labels = test[dist]
                   value.append(labels)
                   lab = mode(labels)
                   lab = lab.mode[0]
                   output.append(lab)
               return output
In [2471]: def accuracy metric(act, pred):
               correct = 0
               for i in range(0,act.shape[0]):
                   if act[i] == pred[i]:
                        correct += 1
                   i=i+1
               return correct / float(len(act)) * 100.0
In [2472]: Yn=predict_classification(X_train,y_train,X_test,1)
In [2473]: | accuracy metric(y test,Yn)
Out[2473]: 31.39240506329114
   In [ ]: #pcawhitening
```

```
In [2474]:
           sorted_index2 = np.argsort(val)[::-1]
           d=100
           vec_new=vec[:,sorted_index2]
           val_new=val[sorted_index2]
           vec new=vec[:,0:d]
           val new=val[0:d]
           temp=1/np.sqrt(val new)
           lambda1=np.diag(temp)
           eigenvector=vec new@lambda1
           pca_wht=f@eigenvector
           print(lambda1.shape, vec new.T.shape)
           print(pca_wht.shape)
           (100, 100) (100, 5655)
           (1577, 100)
In [2476]: X train, X test, y train, y test = train test split(pca wht, y people, stra
In [2477]: Yn1=predict_classification(X_train,y_train,X_test,1)
           accuracy_metric(y_test,Yn1)
Out[2477]: 37.9746835443038
In [2479]: n components1 = 2 #you can select any number of components.
           eigenvector subset1 = sorted eigenvectors[:,0:n components]
           X reduced1 = np.dot(eigenvector subset1.transpose(),f.transpose()).transpos
In [2480]: import matplotlib.pyplot as plt
           plt.figure()
           plt.scatter(X reduced1[:,0], X reduced1[:,1])
           plt.title("Data points")
           plt.show()
```

/Users/brindakulkarni/opt/anaconda3/lib/python3.8/site-packages/numpy/core/_asarray.py:171: ComplexWarning: Casting complex values to real discards the imaginary part

return array(a, dtype, copy=False, order=order, subok=True)



Part 3: Eigenfaces

Import the data as shown above. This the labeled faces in the wild dataset. Use the Xtrain data from above. Let's analyze the first and second principal components.

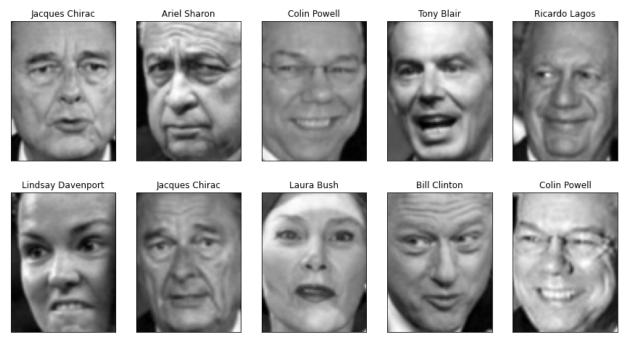
Write a script that:

- 1. Imports the data as mentioned above.
- 2. Standardizes the data.
- 3. Performs PCA on the data (again, although you may not use any package ML functions like pca, you may use statistical functions like eig). No need to whiten here.
- 4. Find the max and min image on PC1's axis. Find the max and min of PC2. Plot and report which faces these points correspond to, what variation do these components capture?
- 5. Visualizes the most important principle component as a 87x65 image (see Figure 2).
- 6. Reconstructs the Xtrain[0,:] image using the primary principle component. To best see the full reconstruction, "unstandardize" the reconstruction by multiplying it by the original standard deviation and adding back in the original mean.
- 7. Determines the number of principle components necessary to encode at least 95% of the information, k.
- 8. Reconstructs the Xtrain [0,:] image using the k most significant eigenvectors (found in the previous step, see Figure 4). For the fun of it maybe even look to see if you can perfectly reconstruct the face if you use all the eigen-vectors! Again, to best see the full reconstruction, "unstandardize" the reconstruction by multiplying it by the original standard deviation and adding back in the original mean.

Your principle eigenface should end up looking similar to Figure 2 of the PDF.

Your principal reconstruction should end up looking similar to Figure 3 of the PDF.

Your 95% reconstruction should end up looking similar to Figure 4 of the PDF.



```
In [2501]: print("people.images.shape: {}".format(people.images.shape))
    print("Number of classes: {}".format(len(people.target_names)))

    people.images.shape: (2537, 87, 65)
    Number of classes: 45
In [2502]: n_samples, h, w = people.images.shape
```

Alejandro Toledo		39	Alvaro Uribe	35	Amelie Ma
uresmo	21				
Andre Agassi		36	Ariel Sharon	77	Atal Biha
ri Vajpayee	24	2.0	0.11 p11	226	D. 11 D.
Bill Clinton kham	2.1	29	Colin Powell	236	David Bec
Donald Rumsfeld	31	121	George Robertson	22	George W
	530	121	George Robertson	22	George w
Gerhard Schroeder		109	Gloria Macapagal Arroyo	44	Gray Davi
S	26				0107 2011
Hamid Karzai		22	Hans Blix	39	Hugo Chav
ez	71				_
Igor Ivanov		20	Jack Straw	28	Jacques C
hirac	52				
Jean Chretien		55	Jennifer Aniston	21	Jennifer
Capriati	42				
Jennifer Lopez		21	John Ashcroft	53	Juan Carl
os Ferrero	28				
Junichiro Koizumi		60	Kofi Annan	32	Laura Bus
h	41			4.1	
Lindsay Davenport		22	Lleyton Hewitt	41	Megawati
Sukarnoputri	33	2.0	Dobo Compres	22	Ricardo L
Michael Bloomberg	27	20	Pete Sampras	22	RICALGO L
agos Rudolph Giuliani	21	26	Saddam Hussein	23	Serena Wi
lliams	52	20	baddam Husselli	23	berena wi
Tiger Woods	J -	23	Tom Daschle	25	Tom Ridge
33					
Tony Blair		144	Vicente Fox	32	Vladimir
Putin	49				

```
In [2504]: mask = np.zeros(people.target.shape, dtype=np.bool)
for target in np.unique(people.target):
    mask[np.where(people.target == target)[0][:50]] = 1

X_people = people.data[mask]
y_people = people.target[mask]

# scale the grayscale values to be between 0 and 1
# instead of 0 and 255 for better numeric stability
X_people = X_people / 255.
```

<ipython-input-2504-0f53c27b3db9>:1: DeprecationWarning: `np.bool` is a d
eprecated alias for the builtin `bool`. To silence this warning, use `boo
l` by itself. Doing this will not modify any behavior and is safe. If you
specifically wanted the numpy scalar type, use `np.bool_` here.
Deprecated in NumPy 1.20; for more details and guidance: https://numpy.or
g/devdocs/release/1.20.0-notes.html#deprecations (https://numpy.org/devdo
cs/release/1.20.0-notes.html#deprecations)

mask = np.zeros(people.target.shape, dtype=np.bool)



```
In [2508]: f=df1.to_numpy()
In [2509]: |e1=df1.T.to_numpy()
In [2510]: v=np.cov(e1)
In [2511]: v.shape
Out[2511]: (5655, 5655)
In [2512]: from numpy.linalg import eig
In [2513]: val, vec=eig(v)
In [2514]: print(vec.shape)
           print(val.shape)
           (5655, 5655)
           (5655,)
In [2515]: sorted_index = np.argsort(val)[::-1]
           sorted_eigenvalue = val[sorted_index]
           #similarly sort the eigenvectors
           sorted_eigenvectors = vec[:,sorted_index]
In [2516]: n components = 2 #you can select any number of components.
           eigenvector subset = sorted eigenvectors[:,0:n components]
           X reduced = np.dot(eigenvector subset.transpose(),f.transpose()).transpose(
           eigenvector subset.shape
Out[2516]: (5655, 2)
```

[195] (array([1220]),) (array([795]),) (array([1331]),)







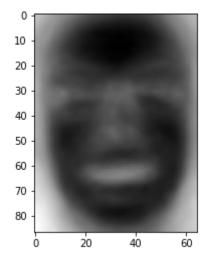


```
In [2499]: besteigenvec = eigenvector_subset[:,0]
    img1 = besteigenvec.reshape((87,65))
    img1 = img1.astype(np.float)
    print("primary principal component")
    plt.imshow(img1, cmap=cm.gray)
```

<ipython-input-2499-0be3715b7a8c>:3: DeprecationWarning: `np.float` is a
deprecated alias for the builtin `float`. To silence this warning, use `f
loat` by itself. Doing this will not modify any behavior and is safe. If
you specifically wanted the numpy scalar type, use `np.float64` here.
Deprecated in NumPy 1.20; for more details and guidance: https://numpy.or
g/devdocs/release/1.20.0-notes.html#deprecations (https://numpy.org/devdo
cs/release/1.20.0-notes.html#deprecations)

img1 = img1.astype(np.float)
<ipython-input-2499-0be3715b7a8c>:3: ComplexWarning: Casting complex valu
es to real discards the imaginary part
img1 = img1.astype(np.float)

Out[2499]: <matplotlib.image.AxesImage at 0x7f89e34bb400>



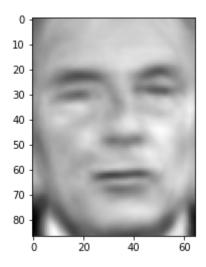
```
In [2542]: sorted_index_1 = np.argsort(val)[::-1]
    sorted_eigenvalue = val[sorted_index_1]
    #similarly sort the eigenvectors
    sorted_eigenvectors_1 = vec[:,sorted_index]
    n_components_1 = 100 #you can select any number of components.
    eigenvector_subset_1 = sorted_eigenvectors_1[:,0:n_components_1]
    X_reduced_1 = np.dot(eigenvector_subset_1.transpose(),f.transpose()).transpeigenvector_subset_1.shape
```

Out[2542]: (5655, 100)

```
In [2545]: mean23, std23 = np.mean(X_people, axis=0).reshape(-1,1), np.std(X_people,
    img_1 = X_reduced_1[0].reshape((1,-1))
    img_1 = img_1 @eigenvector_subset_1.T
    img_1= (img_1*std23.T)+mean23.T
    img_1 = img_1.reshape((87,65))
    plt.imshow(img_1.astype(float), cmap=cm.gray)
```

<ipython-input-2545-6c7d0637b341>:6: ComplexWarning: Casting complex valu
es to real discards the imaginary part
 plt.imshow(img_1.astype(float), cmap=cm.gray)

Out[2545]: <matplotlib.image.AxesImage at 0x7f8990fe3280>

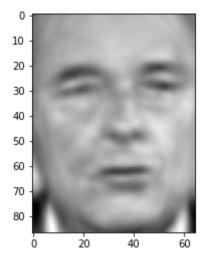


```
In [2549]:
           dim = 2
           var1 = np.var(f)
           sorted_index_2 = np.argsort(val)[::-1]
           vec_new = vec[:,sorted_index_2]
           val_new = val[sorted_index_2]
           while True:
               vec_temp1 = vec_new[:,0:dim]
               pca1= f@vec temp1
               reconstruction = pcal@vec_temp1.T
               pca_variance = np.var(reconstruction)
               diff = var1-pca variance
               if diff <= 0.05:
                   print(diff,dim)
                   break
               dim +=1
           img_f1 = pcal[0].reshape((1,-1))
           img_f1 = img_f1@vec_temp1.T
           img_f1= (img_f1*std23.T)+mean23.T
           img_f1 = img_f1.reshape((87,65))
           print(img f1.shape)
           plt.imshow(img_f1.astype(float), cmap=cm.gray)
           0.04999491436550518 186
```

(87, 65)

<ipython-input-2549-bb02a8b901eb>:21: ComplexWarning: Casting complex val ues to real discards the imaginary part plt.imshow(img f1.astype(float), cmap=cm.gray)

Out[2549]: <matplotlib.image.AxesImage at 0x7f89c3f7f3d0>

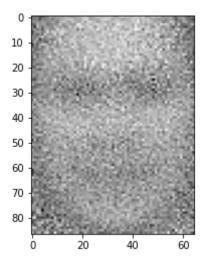


```
In [2610]: dim = X_people.shape[0]
    print(dim)
# meanX, stdX
    img_test = f[0].reshape((1,-1))
    img_test = img_test@vec_new.T
    img_test = (img_test*std23.T)+mean23.T
    img_test = img_test.reshape((87,65))
    print(img_test.shape)
    plt.imshow(img_test.astype(float), cmap=cm.gray)

1577
    (87, 65)

<ipython-input-2610-b9447190d7a1>:9: ComplexWarning: Casting complex values to real discards the imaginary part
    plt.imshow(img_test.astype(float), cmap=cm.gray)
```

Out[2610]: <matplotlib.image.AxesImage at 0x7f89e345b5b0>



Part 4: Clustering

Let's implement our own version of k-means to cluster data! Once you have your setup complete as shown above, write a script to do the following:

1. Write your own version of K-means clustering where you use the L2 distance to compute similarity

- 2. Standardize your data (zero mean, divide by standard deviation)
- 3. Reduces the data to 100D using PCA.
- 4. Run K-means clustering with K = 10.
- 5. Report the number of images within each cluster.
- 6. Reconstruct the cluster centers for each of the K clusters. You will have to rotate the cluster centers back to the original space to visualize. Report these images.
- 7. Find the image closest to the cluster center, and furthest from the cluster center and report these images. Again, you will have to rotate the images centers back to the original space to visualize.

Implementation Details

- 1. Seed your random number generator with zero (do this right before running your k-means).
- 2. Randomly select k observations and use them for the initial reference vectors. I suggest you use randomize the indices and use the first k randomized indices to select the observations.
- Use the L2 distance (Euclidean) to measure the distance between observations and reference vectors.
- 4. Terminate the training process when the sum of magnitude of change of the cluster centers (from the previous iteration to the current one) is less than $\epsilon = 2^{-23}$. That is, when $\sum_{i=1}^k d(a_i(t-1), a_i(t)) < \epsilon$ where k is the number of clusters, $a_i(t)$ is the reference vector for cluster i at time t and d(x, y) is the L1 distance (Manhattan) between vectors x and y (as defined in the *Similarity and Distance Functions* link on BBlearn), or when you've hit 10,000 iterations.

```
In [2567]: df = pd.DataFrame(data=X_people)
    def Standardise(string1):
        ml=np.mean(string1)
        stdl=np.std(string1)
        string2=(string1-ml)/std1
        return string2
        df1=df.apply(Standardise)
In [2568]: X=df1.to_numpy()
    el=df1.T.to_numpy()
    el.shape
Out[2568]: (5655, 1577)
In [2569]: v=np.cov(e1)
    from numpy.linalg import eig
    val,vec=eig(v)
```

```
In [2618]: | sorted_index = np.argsort(val)[::-1]
           sorted eigenvalue = val[sorted index]
           #similarly sort the eigenvectors
           sorted_eigenvectors = vec[:,sorted_index]
           n_components = 100 #you can select any number of components.
           eigenvector_subset = sorted_eigenvectors[:,0:n_components]
           eigenvector_subset.shape
Out[2618]: (5655, 100)
In [2619]: X_reduced = np.dot(eigenvector_subset.transpose(),f.transpose()).transpose(
In [2620]: X_reduced
Out[2620]: array([[-35.54691031+0.j,
                                      12.54720679+0.j,
                                                         2.80998438+0.j, ...,
                    -3.42559574+0.j,
                                      0.16237098+0.j,
                                                         3.98125904+0.j],
                  [ 43.90277194+0.j, 16.55479901+0.j, -12.9888209 +0.j, ...,
                     5.07036923+0.j, -5.35488238+0.j,
                                                         3.80243584+0.j],
                    1.48052537+0.j, -17.624548 +0.j, 10.92830937+0.j, ...,
                    -3.94936438+0.j, -0.08731724+0.j,
                                                        1.51189737+0.j],
                  . . . ,
                  [-14.42154729+0.j,
                                      58.98607396+0.j, -14.48877391+0.j, ...,
                     4.60790643+0.j,
                                      -2.91598645+0.j,
                                                         0.34613491+0.j,
                                     18.84651195+0.j, -35.19563328+0.j, ...,
                  [-13.33391683+0.j,
                    -1.75510202+0.j,
                                      0.14249695+0.j, -0.81141118+0.j],
                  [-41.50046173+0.j, -9.87437792+0.j, 17.6243162 +0.j, ...,
                     3.00273369+0.j,
                                      6.28004404+0.j, 4.39189006+0.j]])
```

```
In [2621]: import numpy as np
           from scipy.spatial.distance import cdist
           import matplotlib.pyplot as plt
           #Defining our function
           def kmeans(x,k, no_of_iterations):
               idx = np.random.choice(len(x), k, replace=False)
               #Randomly choosing Centroids
               centroids = x[idx, :]
               print(centroids.shape)
               #finding the distance between centroids and all the data points
               distances = cdist(x, centroids , 'euclidean')
               #print(distances)
               #Centroid with the minimum Distance
               points = np.array([np.argmin(i) for i in distances]) #Step 3
               #Repeating the above steps for a defined number of iterations
               #Step 4
               for in range(no of iterations):
                   centroids = []
                   for idx in range(k):
                       #Updating Centroids by taking mean of Cluster it belongs to
                       temp cent = x[points==idx].mean(axis=0)
                       centroids.append(temp_cent)
                   centroids = np.vstack(centroids) #Updated Centroids
                   distances = cdist(x, centroids , 'euclidean')
                   points = np.array([np.argmin(i) for i in distances])
                   dist12 = np.array([[min(i), X reduced[ind]] for ind, i in enumerate(d
               return points,dist12
           #Applying our function
           label,centroids = kmeans(X reduced, 10, 1000)
           #Visualize the results
           print(label,centroids)
           (10, 100)
```

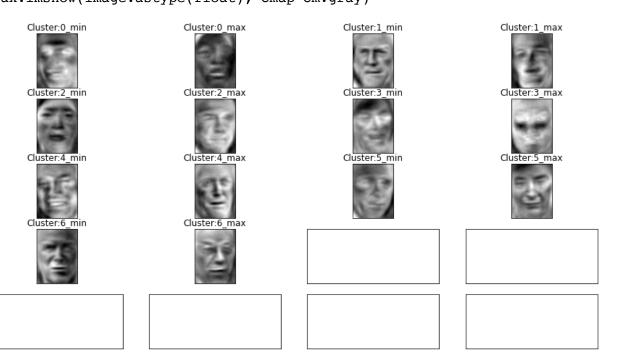
<ipython-input-2621-aa2f8aa008cb>:29: VisibleDeprecationWarning: Creating
an ndarray from ragged nested sequences (which is a list-or-tuple of list
s-or-tuples-or ndarrays with different lengths or shapes) is deprecated.
If you meant to do this, you must specify 'dtype=object' when creating th
e ndarray.

dist12 = np.array([[min(i),X_reduced[ind]] for ind,i in enumerate(dista
nces)])

```
[0 6 7 ... 2 2 1] [[46.30099222255128
             array([-3.55469103e+01+0.j, 1.25472068e+01+0.j, 2.80998438e+00+0.j,
In [2625]: clusters, i = np.unique(label, return_counts=True)
           print(np.array(( i)))
           [145 215 148 143 113 223 150 237 83 120]
```

```
In [2624]: images = []
           # print(dists clst.shape)
           fig, axes = plt.subplots(5,4, figsize=(15, 8),
                                     subplot_kw={'xticks': (), 'yticks': ()})
           # print(dists clst)
           faces = []
           c list = []
           for c in range(k):
               temporary = centroids[:,0][label==c]
               face1 = centroids[:,1][np.argmin(temporary)]
               face2 = centroids[:,1][np.argmax(temporary)]
               img_face = face1.reshape((1,-1))
               img face = img face@eigenvector subset.T
               img_face = img_min.reshape((87,65))
               faces.append(img min)
               c_list.append(f"{c}_min")
               img m = maxface.reshape((1,-1))
               img_m = img_m@eigenvector_subset.T
               img_m = img_m.reshape((87,65))
               faces.append(img m)
               c_list.append(f"{c}_max")
           for target, image, ax in zip(c_list, faces, axes.ravel()):
               ax.imshow(image.astype(float), cmap=cm.gray)
               ax.set title(f'Cluster:{target}')
```

<ipython-input-2624-318eda1764bb>:29: ComplexWarning: Casting complex val
ues to real discards the imaginary part
 ax.imshow(image.astype(float), cmap=cm.gray)



In []: