1 Linear regression and correlation

The following dataset is a data frame with 7 economical variables, observed yearly from 1947 to 1962 (n=16).

- GNP.deflator: GNP implicit price deflator (1954=100)
- GNP: Gross National Product.
- Unemployed: number of unemployed.
- Armed Forces: number of people in the armed forces.
- Population: 'noninstitutionalized' population ≥ 14 years of age.
- Year: the year (time).
- Employed: number of people employed.

Source J. W. Longley (1967) An appraisal of least-squares programs from the point of view of the user. Journal of the American Statistical Association 62, 819–841.

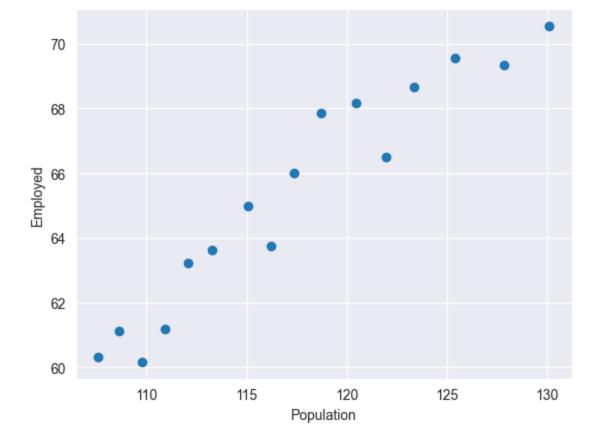
```
In [1]: import pandas as pd
    from pandas import read_csv
    from matplotlib import pyplot as plt

# load the dataset
url = 'https://raw.githubusercontent.com/jbrownlee/Datasets/master/longley.csv'
dataframe = read_csv(url, header=None)
data = dataframe.values

# choose the input and output variables
x, y = data[:, 4], data[:, -1]

# plot input vs output
plt.scatter(x, y)
plt.xlabel('Population')
plt.ylabel('Employed')
```

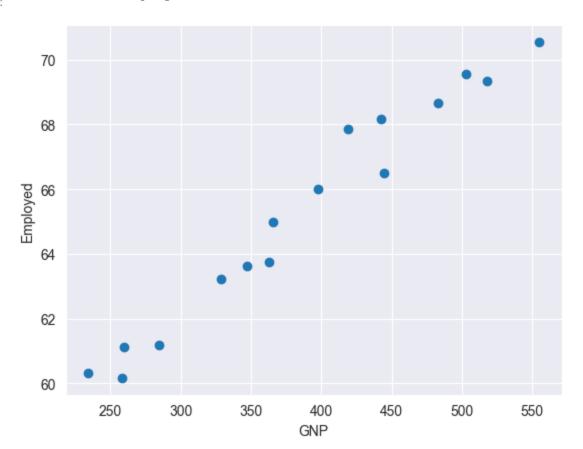
Out[1]: Text(0, 0.5, 'Employed')



```
In [2]: #GNP and employed
x, y = data[:, 1], data[:, -1]

# plot input vs output
plt.scatter(x, y)
plt.xlabel('GNP')
plt.ylabel('Employed')
```

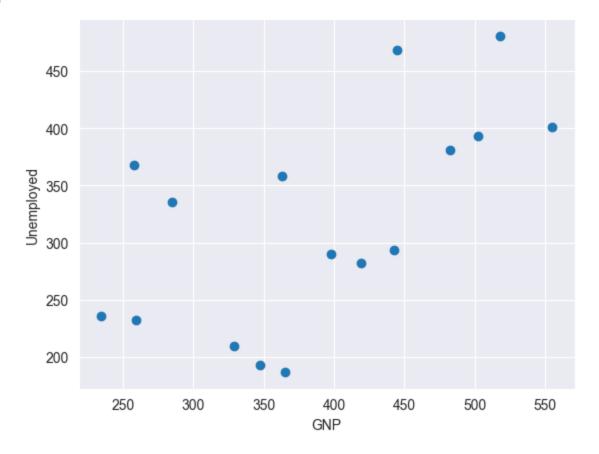
Out[2]: Text(0, 0.5, 'Employed')



```
In [3]: #GNP and UNemployed
x, y = data[:, 1], data[:, 2]

# plot input vs output
plt.scatter(x, y)
plt.xlabel('GNP')
plt.ylabel('Unemployed')
```

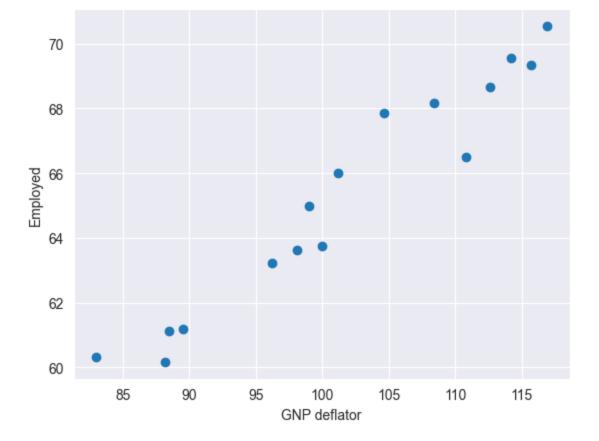
Out[3]: Text(0, 0.5, 'Unemployed')



```
In [4]: #GNP deflator and Employed
x, y = data[:, 0], data[:, -1]

# plot input vs output
plt.scatter(x, y)
plt.xlabel('GNP deflator')
plt.ylabel('Employed')
```

Out[4]: Text(0, 0.5, 'Employed')



- Looking at the figures, which variables exhibit a positive correlation index?
- Chose one set of varibles that are linearly correlated and construct a linear regression model to fit them.
- Choose a next value in the future and predict the corresponding value of your model.

(The topics related to this example can be found in: Example 5.4.3 and Chapter 4.6 from Meyer, "Matrix Analysis and Applied Linear Algebra").

Conduct of the exercise n.1

First point

Looking only at the graphs, we can say that those in which there might be a positive correlation index between the two relative variables represented are:

- **First graph** representing the correlation between population and number of employed. As the population increases, the number of employed increases.
- **Second graph** representing the correlation between GNP and number of employed. Knowing that GNP is how much the country has earned from the total production of goods and services (at current prices), we can say that as GNP increases, the number of employed also increases.
- **Fourth graph** representing the correlation between GNP deflator and number of employed. Knowing that the GNP deflator is a measure of price changes in an economy, we can say that as the GNP deflator increases, the number of employed also increases. A deflator greater than 100 indicates that prices have increased compared to the reference year. (In the trace the value 100 is reached in 1954)

All three graphs above, show a positive correlation index, compared to the third one (representing the correlation between GNP and number of unemployed) in which the data are scattered and show less correlation with each other.

Second point

To carry out the second point, let us actually see which variables, even among those not seen within the previous point, are linearly correlated with each other. To do this, we use the **coefficient of linear** correlation define like $\rho_{xy} = \cos\theta = \frac{\mathbf{z_x}^T\mathbf{z_y}}{\|\mathbf{z_x}\|\|\mathbf{z_y}\|}$, which exploits the cosine.

```
In [5]: #Renaming of features for easier reading of the code
dataframe.rename(columns={0: "GNPdeflator",1:"GNP",2:"Unemployed",3:"Armed Forces",4:"Po
dataframe
```

Out[5]:		GNPdeflator	GNP	Unemployed	Armed Forces	Population	Year	Employed
	0	83.0	234.289	235.6	159.0	107.608	1947	60.323
	1	88.5	259.426	232.5	145.6	108.632	1948	61.122
	2	88.2	258.054	368.2	161.6	109.773	1949	60.171
	3	89.5	284.599	335.1	165.0	110.929	1950	61.187
	4	96.2	328.975	209.9	309.9	112.075	1951	63.221
	5	98.1	346.999	193.2	359.4	113.270	1952	63.639
	6	99.0	365.385	187.0	354.7	115.094	1953	64.989
	7	100.0	363.112	357.8	335.0	116.219	1954	63.761
	8	101.2	397.469	290.4	304.8	117.388	1955	66.019
	9	104.6	419.180	282.2	285.7	118.734	1956	67.857
	10	108.4	442.769	293.6	279.8	120.445	1957	68.169
	11	110.8	444.546	468.1	263.7	121.950	1958	66.513
	12	112.6	482.704	381.3	255.2	123.366	1959	68.655
	13	114.2	502.601	393.1	251.4	125.368	1960	69.564
	14	115.7	518.173	480.6	257.2	127.852	1961	69.331
	15	116.9	554.894	400.7	282.7	130.081	1962	70.551

Also taking into account the analysis made in the previous point, the number of employed persons was identified as the target features. Consequently, with the aim of predicting the target y considering only highly correlated features, we proceed with the correlation calculation. As mentioned before, the cosine technique will be used, taking into account the mean, standard deviation and z-score of each feature.

```
In [6]: #Definition of independent features
    x1 = dataframe['GNPdeflator']
    x2 = dataframe['GNP']
    x3 = dataframe['Unemployed']
    x4 = dataframe['Armed Forces']
    x5 = dataframe['Population']
    x6 = dataframe['Year']

#Definition of target feature
    y = dataframe['Employed']
```

```
import numpy as np

#Average of all variables x
mean_x1 = np.mean(x1)
mean_x2 = np.mean(x2)
mean_x3 = np.mean(x3)
```

```
mean x4 = np.mean(x4)
mean x5 = np.mean(x5)
mean x6 = np.mean(x6)
#Standard deviation for each variable x
standardDev x1 = np.std(x1)
standardDev x2 = np.std(x2)
standardDev x3 = np.std(x3)
standardDev x4 = np.std(x4)
standardDev x5 = np.std(x5)
standardDev x6 = np.std(x6)
#Averange and Standard deviation for the y
mean y = np.mean(y)
standardDev y = np.std(y)
#z-score of all variables x
z \cdot scoreX1 = (x1 - mean x1) / standardDev x1
z_{scoreX2} = (x2 - mean x2) / standardDev x2
z \cdot scoreX3 = (x3 - mean x3) / standardDev x3
z \cdot scoreX4 = (x4 - mean x4) / standardDev x4
z \cdot scoreX5 = (x5 - mean x5) / standardDev x5
z \cdot scoreX6 = (x6 - mean x6) / standardDev x6
#z-score for y target
z scoreY = (y - mean y) / standardDev y
\#Compute the cosine for the correlation between y and x i
correlation y x1 = np.dot(z scoreX1, z scoreY) / (np.linalg.norm(z scoreX1) * np.linalg.
correlation_y_x2 = np.dot(z_scoreX2, z scoreY) / (np.linalg.norm(z scoreX2) * np.linalg.
correlation y x3 = np.dot(z scoreX3, z scoreY) / (np.linalg.norm(z scoreX3) * np.linalg.
correlation y x4 = np.dot(z scoreX4, z scoreY) / (np.linalg.norm(z scoreX4) * np.linalg.
correlation y x5 = np.dot(z scoreX5, z scoreY) / (np.linalg.norm(z scoreX5) * np.linalg.
correlation y x6 = np.dot(z scoreX6, z scoreY) / (np.linalg.norm(z scoreX6) * np.linalg.
print("Coefficient of linear correlation of NUMBER OF EMPLOYED with GNPdeflator ", corr
print("Coefficient of linear correlation of NUMBER OF EMPLOYED with GNP ", correlation
print("Coefficient of linear correlation of NUMBER OF EMPLOYED with UNEMPLOYED ", corre
print("Coefficient of linear correlation of NUMBER OF EMPLOYED with ARMED FORCES ", cor
print("Coefficient of linear correlation of NUMBER OF EMPLOYED with POPULATION ", corre
print("Coefficient of linear correlation of NUMBER OF EMPLOYED with YEARS ", correlatio
Coefficient of linear correlation of NUMBER OF EMPLOYED with GNPdeflator 0.97089852506
Coefficient of linear correlation of NUMBER OF EMPLOYED with GNP 0.9835516111796695
Coefficient of linear correlation of NUMBER OF EMPLOYED with UNEMPLOYED 0.502498083875
9943
Coefficient of linear correlation of NUMBER OF EMPLOYED with ARMED FORCES 0.4573073999
Coefficient of linear correlation of NUMBER OF EMPLOYED with POPULATION 0.960390571594
Coefficient of linear correlation of NUMBER OF EMPLOYED with YEARS 0.9713294591921189
```

We then obtained a measure of the degree of correlation between x and y. Accordingly, only the features with the highest correlation were chosen, namely: **GNPdeflator, GNP, Population, Year.**

Now, let us move on to the creation of the linear regression model. We know that the number of people employed depends on the GNP, the GNP deflactor, the population and the year. Therefore, to predict the number of employed knowing the value of these variables, we need a function (in this case linear) that best fits the data. Mainly we will have:

- A matrix A, which will represent the features x (+ bias)
- A vector b, which will represent the values of the y variables associated with each set of x variables.

• A vector x, which will represent the weights, the importance, given to each variable x in order to predict y.

In the next cell, an attempt will be made to obtain the vector x to predict y.

```
#Least square solition with pytorch
In [8]:
        import torch
        #Definition of the variables x and y, omitting the last line outside the fit in order to
        x values = dataframe[['GNPdeflator', 'GNP', 'Population', 'Year']].values
        b values = dataframe['Employed'].values
        x = x \text{ values } [:15,:]
        b = b values[:15]
        #Addition of column in matrix A to represent bias (column all to 1)
        (nx, ) = x.shape
        ones column = np.ones((nx, 1))
        A = np.hstack((ones column, x))
        AT = torch.tensor(A, dtype=torch.float64)
        bT = torch.tensor(b,dtype=torch.float64)
        X = torch.linalg.lstsq(AT, bT)
        print("Vector x:", np.asarray(X.solution))
        #print(torch.tensor(p)-X.solution)
        #print(X.rank)
        #print(X.residuals)
        #print(X.singular values)
        Vector x: [ 3.57927247e+02 -2.35671977e-01 9.16785742e-02 -3.41827722e-01
```

Then we can compose our linear function for the prediction of y:

$$f(x) = +3.5792(bias) - 2.36*GNP deflator + 9.17*GNP - 3.42*Population - 1.35*Year$$

Third point

-1.35059509e-011

Original y values: 70.551

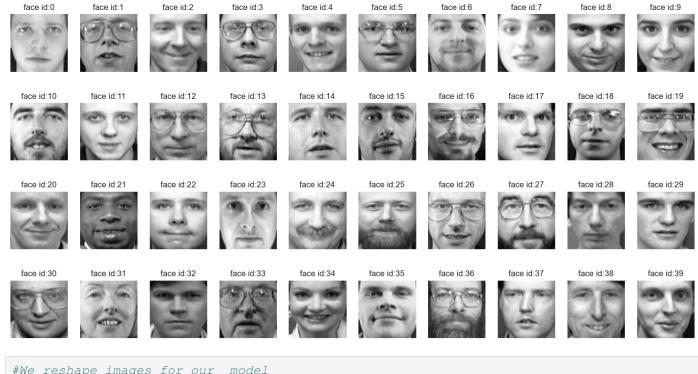
We now take the last row of the original dataset to predict its corresponding y using the linear function found earlier.

Predicted y values: 71.79703529095909

2 Eigenfaces

Download the files olivetti_faces.npy and olivetti_faces_target.npy .

```
import numpy as np
In [11]:
        data=np.load("olivetti faces.npy")
In [12]:
         target=np.load("olivetti faces target.npy")
In [13]: | print("There are {} images in the dataset".format(len(data)))
         print("There are {} unique targets in the dataset".format(len(np.unique(target))))
         print("Size of each image is {}x{}".format(data.shape[1],data.shape[2]))
         print("Pixel values were scaled to [0,1] interval. e.g:{}".format(data[0][0,:4]))
        There are 400 images in the dataset
        There are 40 unique targets in the dataset
        Size of each image is 64x64
        Pixel values were scaled to [0,1] interval. e.g:[0.30991736 0.3677686 0.41735536 0.4421
In [14]: def show_40_distinct people(images, unique ids):
             #Creating 4X10 subplots in 18x9 figure size
             fig, axarr=plt.subplots(nrows=4, ncols=10, figsize=(18, 9))
             #For easy iteration flattened 4X10 subplots matrix to 40 array
             axarr=axarr.flatten()
             #iterating over user ids
             for unique id in unique ids:
                 image index=unique id*10
                 axarr[unique id].imshow(images[image index], cmap='gray')
                axarr[unique id].set xticks([])
                 axarr[unique id].set yticks([])
                 axarr[unique id].set title("face id:{}".format(unique id))
             plt.suptitle("There are 40 distinct people in the dataset")
In [15]: show 40 distinct people(data, np.unique(target))
```



```
In [16]: #We reshape images for our model
    X=data.reshape((data.shape[0],data.shape[1]*data.shape[2]))
    print("X shape:",X.shape)
    X shape: (400, 4096)
```

```
In [17]: from sklearn.model_selection import train_test_split

X_train, X_test, y_train, y_test=train_test_split(X, target, test_size=0.3, stratify=tar print("X_train shape:",X_train.shape)
    print("y_train shape:{}".format(y_train.shape))

X_train shape: (280, 4096)
    y train shape: (280,)
```

- Check and compare all the studied techniques to find the minimal value of k, number of principal components that are useful to classify the faces.
- Classify all the faces in the testset and compute the confusion matrix. The element in position i, j of the confusion matrix contains the number of observations known to be in group i and predicted to be in group j. Use sklearn.metrics.confusion_matrix and compute also the classification report using sklearn.metrics.classification_report.

(The topics related to this exercise can be found in: eigenfaceforrecognition-TurkPentland-jcn1991.pdf Chapter 6 from Elden, "Numerical Linear Algebra and Applications in Data Mining")

Conduct of the exercise n.2

First point

In order to find the optimal k, we need the singular values of the train matrix. Accordingly, we centre the train matrix and decompose it into its corresponding SVD form. The SVD technique decomposes the original matrix into three matrices $U \sum V^T$ and the columns of the matrices U and V are called left and right singular vectors, respectively. While the diagonal elements of \sum are said singular values.

```
In [18]: | import warnings
        warnings.filterwarnings("ignore")
        from scipy.linalg import svd
        import numpy as np
        from numpy import linalg
        import matplotlib.pyplot as plt
        (nrc,nt) = X train.shape
        print(nrc,nt)
        def centralizing matrix(matrix, nrc, nt):
           matrix sum = np.sum(matrix,axis=1)/nt
            e = np.ones((1,nt))
            centralizing matrix =matrix-np.dot(matrix sum.reshape(nrc,1),e)
            return centralizing matrix
        X trainC = centralizing matrix(X train, nrc, nt)
        U, S, VT = svd(X trainC)
        singular values = S
        singular values
        280 4096
        array([115.65021921, 53.60780435, 37.87629115, 32.25059585,
Out[18]:
               29.44320185, 25.61298961, 23.36836107, 20.85217286,
               20.45953807, 19.14477133, 18.54698076, 17.59573116,
               16.8132345 , 15.99646094, 15.35254918, 14.89075181,
               14.24920778, 13.9137765, 13.32140692, 12.8026415,
               12.39373316, 11.76915229, 11.66594042, 11.54126058,
               11.24306775, 10.83865445, 10.66143974, 10.47269743,
               10.3278773 , 9.87578209, 9.67898009, 9.38421892,
                9.12079058, 8.8783905, 8.74233297, 8.48518612,
                8.40218073, 8.29560387,
                                         8.12899234, 8.06218092,
                7.93758782,
                            7.69444 ,
                                         7.59937322,
                                                       7.44696458,
                7.34172927, 7.25008474, 7.06275156, 7.02741961,
                6.96422957, 6.80067677, 6.66650756, 6.62012137,
                6.54565202, 6.49057137, 6.33531442, 6.3160628,
                6.29849895,
                           6.18796189, 6.17306985, 6.12036508,
                5.92912025, 5.82099695, 5.80345759, 5.74513523,
                5.73542557, 5.6541913,
                                         5.61675966, 5.52165666,
                                         5.42028324, 5.37966529,
                5.48922593,
                            5.43230059,
                           5.27091081, 5.21188211, 5.13311361,
                5.30740157,
                5.09346863, 5.06364799, 5.03487014, 4.96778415,
                4.90232077, 4.87078062, 4.84118112, 4.7695859,
                           4.72945653, 4.66720348, 4.62364966,
                4.74449241,
                4.58269281, 4.55214006, 4.52391342, 4.48863347,
                4.43537911, 4.41556807, 4.36536441, 4.32980339,
                                         4.23408067,
                                                      4.2257984 ,
                4.29025222,
                            4.25306565,
                           4.15406148, 4.10283054, 4.08193607,
                4.18748021,
                4.06859161, 4.03997539, 4.01249284, 3.9800224,
                3.96305615, 3.90575983, 3.8989024, 3.87266301,
                            3.78375972, 3.75519318,
                                                      3.7537432 ,
                3.82574286,
                3.71746202,
                             3.67509562, 3.67015683, 3.63098953,
                3.61571574, 3.60522938, 3.5661954, 3.55093532,
                3.52490792,
                            3.50502596, 3.4782899,
                                                      3.46356625,
                3.44994485,
                            3.40201945, 3.37281203, 3.36113016,
                3.34869014, 3.30467282, 3.2964765, 3.28046733,
                3.26091821, 3.23188743, 3.20805058, 3.19661851,
                3.17286409, 3.14733339, 3.12614944, 3.10340326,
                3.08739393,
                            3.0690069 , 3.05409411, 3.03151916,
                3.0013312 , 2.97721429,
                                        2.97527109, 2.95540494,
```

2.94119246, 2.89779459, 2.88206836,

2.75388891,

2.82979071, 2.79942112,

2.74375888,

2.86174459,

2.77237164,

2.94327065,

2.88116972, 2.78785888,

```
2.72222402, 2.70012489, 2.69487941, 2.68403112,
2.65514679, 2.65069271, 2.64428572, 2.6067529,
2.59783546, 2.58620802, 2.56034602, 2.54842995,
          2.50749469, 2.48889335,
                                    2.48052394,
2.51744979,
2.47305621, 2.44879463, 2.42454497, 2.41388977,
2.3931405 , 2.37768834, 2.36273907, 2.35291393,
2.34976477,
          2.33301571, 2.30551651, 2.29635522,
          2.27629985,
2.28419787,
                       2.26902961, 2.23703184,
2.23203051, 2.21265668, 2.20843622, 2.18825623,
2.17164128, 2.15784808, 2.14717903, 2.12445641,
2.12055466,
           2.09631843,
                       2.0802521 ,
                                    2.06456017,
          2.05312051, 2.04305902,
                                    2.03044364,
2.05384435,
2.01711034, 2.0052157, 1.99459029, 1.97152002,
1.96003176, 1.94298372, 1.93379292, 1.90527655,
          1.87618138, 1.86641563, 1.85992649,
1.89619983,
1.84704454, 1.83159518, 1.82019367, 1.80776214,
1.79281346, 1.77691586, 1.76482325, 1.75716607,
                                    1.70168225,
1.74955908,
           1.73101336, 1.71784734,
          1.68838132,
                       1.68743211, 1.6713523,
1.69910106,
1.66473587, 1.64370445, 1.6343436, 1.6154135,
1.605431 , 1.59892438, 1.58886199, 1.57595173,
          1.55260195, 1.52608531, 1.51924548,
1.55847573,
1.50639034,
           1.4715078 , 1.46013769, 1.45432827,
1.44277131, 1.42811794,
                       1.41587634, 1.40165562,
1.38571879, 1.36714703, 1.35627461, 1.35160018,
          1.32505853,
                       1.31804781, 1.29384458,
1.33397971,
1.26801461, 1.24668868, 1.22787691, 1.21047148,
1.1905029 , 1.18026676, 1.16448972, 1.14408813,
1.10116149, 1.05662922, 0.99259097, 0.92932449])
```

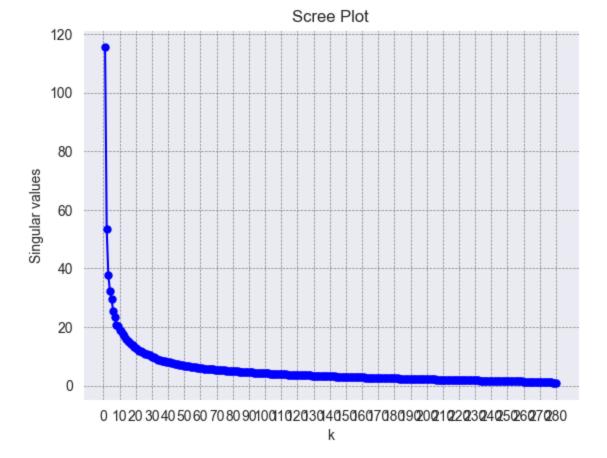
So let's start exploring the various techniques designed to define the best k.

Scree Plot

This first method is a visual and very often subjective method. It involves representing the singular values in a cartesian coordinate system and then selecting $\setminus (k \setminus)$ at the point where the graph shows a sharper curve, forming a kind of 'elbow'.

```
In [19]: plt.plot(np.arange(1, len(singular_values) + 1), singular_values, marker='o', linestyle=
    plt.xlabel('k')
    plt.ylabel('Singular values')
    plt.title('Scree Plot')
    plt.grid(True, linestyle='--', linewidth=0.5, color='gray')

plt.xticks(np.arange(0, len(singular_values) + 1, step=10))
    plt.grid(True)
    plt.show()
```



In our case, the point at which the graph shows a sharper curve is on a k of 30

In
$$[20]$$
: kScree = 30

Kaiser rule

Another method for finding how many principal components to keep is the Kaiser rule. In general, its basic version is that if a component has a singular value greater than 1, then it is kept because it is considered good as a summary of the data. However, this version has a flaw: it retains components with singular values even slightly greater than 1, which may not be representative. Consequently, the singular mean value is calculated first and then retained all those singular values which, when squared, are greater than the mean value.

$$k = \sigma_i^2 \geq rac{1}{n} \sum_{j=1}^n \sigma_j^2$$

Cumulative Percentage of Total Variance

Within this method, we consider a function S_k which is the percentage of variation accumulated by the k points:

$$s_k = 100 \cdot rac{\sum_{i=1}^k \sigma_i^2}{\sum_{i=1}^r \sigma_i^2}$$

It is then necessary to choose a threshold value S^* , generally between 70% and 80%, which allows us to return only those k's with $S_k > S^*$

```
In [22]:
        def compute sk(k):
             k sum = 0
             for i in range (1, k):
                k sum+= singular values[i]**2
             rank = np.linalg.matrix rank(X train)
             r sum = 0
             for i in range(1, rank):
                 r sum += singular values[i] **2
             s k = 100 * (k sum/r sum)
             return s k
         s cutoff = (compute sk(len(singular values)) * 80)/100
         kc = 0
         for i in range(1,100):
             sk = compute sk(i)
             if sk > s cutoff:
                 kc = i
         pass
         print("k = " + str(kc))
```

k = 99 **Entropy**

This method defines the contribution that each individual value makes to the component via the function:

$$f_k = rac{\sigma_k^2}{\sum_{i=1}^r \sigma_i^2}$$

Thus entropy, or how informative a component is, is given by:

$$E = rac{1}{\log(2)} \sum_{k=1}^r f_k \log(f_k)$$

```
In [23]: t = pd.DataFrame(columns=['Tollerance', 'k'])

tol=[1e-1,7e-2,5e-2,1e-2,5e-3,1e-3,0]
k_entropy = []
for ind in tol:
    f=s**2. / np.sum(S**2.)
    r = S.shape
    entropy = (-1/np.log(r))*np.sum(f*np.log(f))
    ks = int(r*entropy)
    perc=1-ind
    nc = int(r*entropy*perc)
    k_entropy.append(int(nc))

for i in range(len(tol)):
    t.loc[len(t)] = [tol[i],k_entropy[i]]
```

Out[23]:		Tollerance	k
	0	0.100	126.0
	1	0.070	130.0

```
2 0.050 133.0
3 0.010 139.0
4 0.005 140.0
5 0.001 140.0
6 0.000 140.0
```

The relative error inside the approssimation the matrix

In this method, however, reference is made to the Eckert-Young theorem.

```
In [24]: t = pd.DataFrame(columns=['Tollerance','k'])

tol=[1e-1,7e-2,5e-2,1e-2,5e-3,1e-3,0]
k_rb = []
for ind in tol:
    f= (S)/S[0]
    nc = np.sum(f >= ind)
    k_rb.append(int(nc))

for i in range(len(tol)):
    t.loc[len(t)] = [str(tol[i]),k_rb[i]]

t
```

Out[24]:		Tollerance	k
	0	0.1	23
	1	0.07	39
	2	0.05	63
	3	0.01	275
	4	0.005	280
	5	0.001	280
	6	0	280

Having arrived at this point, we do not just have a single value of k, but rather a set of k's which, when taken in this way, do not answer the question 'how many k's do I need to maintain in order to have a matrix that is reduced but still allows me to maintain a fair accuracy?' Consequently, I define a function that, based on the Euclidean norm, calculates the relative error between the original matrix and the reconstructed matrix having used the k principal values:

```
In [25]: def relative_error(original_Matrix, reconstruction_matrix):
    normNum = linalg.norm(original_Matrix - reconstruction_matrix)
    normDen = linalg.norm(original_Matrix)

    error = normNum/normDen

    return error
```

```
In [26]: from sklearn.decomposition import PCA

def get_reduction_PCA_matrix(original_matrix, k_components):
    pca = PCA(copy=True, n_components=k_components, whiten = True)
    reduction_matrix = pca.fit_transform(original_matrix)
```

```
reconstruction matrix = pca.inverse transform(reduction matrix)
return reduction matrix, reconstruction matrix
```

```
In [27]: t = pd.DataFrame(columns=['Method', 'Error', 'Tollerance', 'k'])
        method k finder = ['Scree-Plot','Kaiser-Rule','Cumulative-Percentage-of-Total-Variance',
         k finded = {'Scree-Plot': kScree,
                     'Kaiser-Rule': k kaiser rule,
                     'Cumulative-Percentage-of-Total-Variance': kc,
                     'Entropy': k entropy,
                     'Relative-error-based': k rb}
         for ,method in enumerate(method k finder):
             if method != 'Entropy' and method != 'Relative-error-based':
                 k = k finded[f'{method}']
                 reducted matrix, reconstruction matrix = get reduction PCA matrix(X trainC, k)
                 error = relative error(X trainC, reconstruction matrix)
                 t.loc[len(t)] = [f'{method}', f'{error:.9f}', '-', f'{k}']
             else:
                 n k = len(k finded[f'{method}'])
                 for n in range(0, n k):
                     k = k finded[f'{method}'][n]
                     reducted matrix, reconstruction matrix = get reduction PCA matrix(X trainC,k)
                     error = relative error(X trainC, reconstruction matrix)
                     t.loc[len(t)] = [f'\{method\}', f'\{error:.9f\}', tol[n], k]
         t
```

Ou	t	Γ	2	7	1	
		ь.			-1	

	Method	Error	Tollerance	k
0	Scree-Plot	0.360760781	-	30
1	Kaiser-Rule	0.365321329	-	29
2	Cumulative-Percentage-of-Total-Variance	0.201761148	-	99
3	Entropy	0.165131766	0.1	126
4	Entropy	0.160129287	0.07	130
5	Entropy	0.156567640	0.05	133
6	Entropy	0.149446571	0.01	139
7	Entropy	0.148455600	0.005	140
8	Entropy	0.148311544	0.001	140
9	Entropy	0.148187104	0	140
10	Relative-error-based	0.395734582	0.1	23
11	Relative-error-based	0.327572649	0.07	39
12	Relative-error-based	0.265534855	0.05	63
13	Relative-error-based	0.011904778	0.01	275
14	Relative-error-based	0.000000000	0.005	280
15	Relative-error-based	0.000000000	0.001	280
16	Relative-error-based	0.000000000	0	280

From this last table we can see that both entropy and relative error inside the approximation the matrix obtain the lowest error values, thus reducing the original matrix with one not too far from it. But, in order to maintain the right compromise between the number of k and the error obtained, we will choose k obtained through the entropy formula, i.e. k = 140.

We then set the PCA with a number of k equal to 140 and apply it to the train set of images.

```
In [28]: from sklearn.decomposition import PCA

k = k_entropy[6]

pca = PCA(copy=True, n_components=k,whiten = True).fit(X_trainC)

IM_pca = pca.transform(X_trainC)

IM_pca = IM_pca/np.sqrt(nrc)

print("IM_pca: " + str(IM_pca.shape))

IM_pca: (280, 140)
```

We assign the train set images processed by the PCA their own unique class:

```
In [29]: cli = {}
    for i in range(40):
        cli[str(i)] = []

unique_classes = np.unique(y_train)

for c in unique_classes:
    for index,val in enumerate(y_train):
        if val == c:
            PT1 = X_trainC[index]
            Pimg = pca.transform([PT1])
            cli[str(c)].append(Pimg)
        cli[str(c)] = np.asarray(cli[str(c)])

om = {}
    for i in range(0,40):
        om[i]=np.mean(cli[str(i)],axis=0)
```

Second point

To carry out the second point, we start by centralising the test set:

We project the images of the test set in the vector space established in the previous point:

```
In [31]: testProjector = []

for i in range(0,120):
    PT1_t = X_testC[i]
    Pimg_t = pca.transform([PT1_t])
    testProjector.append(Pimg_t)

len(testProjector)
```

And in order to associate the new class with the examples of the test set, we calculate the distance between each new element of the test set and the elements of the train set, so that the class of the train element that is closest to the set data is assigned:

```
In [32]: from numpy.linalg import norm

predicted_class = []

for projectionT in testProjector:
    distance_min = float('inf')
    predict_class = -1
    for class_id, mean_comp in om.items():
        distance = norm(mean_comp-projectionT,2)
        if distance < distance_min:
            distance_min = distance
            predict_class = class_id
        predicted_class.append(predict_class)</pre>
```

The predicted classes for the 120 images in the test set are as follows:

```
In [33]: print(predicted class)
         [39, 10, 33, 2, 37, 10, 28, 4, 27, 20, 29, 7, 28, 6, 13, 3, 35, 26, 9, 13, 2, 27, 38, 2
         3, 16, 35, 14, 2, 34, 37, 11, 32, 12, 22, 28, 3, 39, 7, 33, 9, 5, 14, 39, 4, 6, 33, 20,
         24, 34, 8, 36, 31, 17, 21, 18, 20, 35, 22, 39, 36, 16, 29, 25, 30, 32, 38, 17, 36, 8, 1
         1, 10, 25, 38, 9, 26, 15, 17, 2, 6, 11, 29, 18, 16, 24, 20, 13, 38, 32, 8, 21, 31, 5, 1,
         19, 23, 27, 37, 1, 29, 30, 19, 3, 22, 14, 19, 15, 39, 15, 23, 24, 20, 29, 18, 24, 30, 4,
         23, 1, 0, 12]
In [34]:
         #Visual example
         images = X test[:20].reshape(-1, 64, 64)
         labels = predicted class[:20]
         fig, axes = plt.subplots(2, 10, figsize=(20, 4))
         for i in range(2):
             for j in range(10):
                ax = axes[i, j]
                ax.imshow(images[i*10+j], cmap='grey')
                ax.set title(f"Predicted label: {labels[i*10+j]}")
                 ax.axis('off')
         plt.tight layout()
         plt.show()
```























The real classes of the test set are:

```
In [35]: print(y test)
        [ 4 10 33 7 37 10 28 4 27
                                      5
                                         7 28
                                               6 13
                                                     3 35 26
                                   2
                                                             9 13 25 27 38
         16 35 14 2 34 37 11 32 12 22 28
                                         3
                                           0
                                               7 33 9 5 14 39 34
                                                                  6 33 20 24
         34 8 36 31 17 21 18 20 35 22 39 36 16 29 25 30 32 21 17 36 8 11 10 25
         38 9 26 15 17 2 6 11 29 18 16 31 12 13 38 32 8 21 31 5 1 19 23 27
         37 1 29 30 19 3 22 14 19 15 39 15 23 24 20 26 18 24 30
                                                                4 23 1 0 121
```

```
In [36]: #Visual example
images = X_test[:20].reshape(-1, 64, 64)
labels = y_test[:20]

fig, axes = plt.subplots(2, 10, figsize=(20, 4))

for i in range(2):
    for j in range(10):
        ax = axes[i, j]
        ax.imshow(images[i*10+j], cmap='grey')
        ax.set_title(f"Label: {labels[i*10+j]}")
        ax.axis('off')

plt.tight_layout()
plt.show()
```



We now construct the confusion matrix and the report to also see the accuracy of the model.

```
from sklearn.metrics import confusion matrix, classification report
In [37]:
         import seaborn as sns
         def create confusionMatrix andReport(predicted class):
            conf matrix = confusion matrix(y test, predicted class)
            class report = classification report(y test, predicted class)
            print("CONFUSION MATRIX:")
            plt.figure(figsize=(10, 8))
            sns.heatmap(conf matrix, annot=True, fmt='d', cmap='Blues', cbar=False)
            plt.xlabel('Predicted Labels')
            plt.ylabel('True Labels')
            plt.title('Confusion Matrix')
            plt.show()
            print("\nCLASSIFICATION REPORT :")
            print(class report)
         create confusionMatrix andReport(predicted class)
```

CONFUSION MATRIX:

Confusion Matrix

0 0 Ō 0 12 13 14 15 16 18 19 20 21 22 23 24 25 26 27 28 29 30 31 32 33 34 35 36 37 38 0 0 0 0 0 3 0 0 3 0 0 0 0 0 0 0 0 0 0 0 0 0 0 Ō 0 0 0 0 0 0 0 3 0 Ō 0 0 0 3 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21 22 23 24 25 26 27 28 29 30 31 32 33 34 35 36 37 38 39

Predicted Labels

CI	7\	C	Q I	ΓF	T	$C \lambda$	т	Т	M	D	F	D	$\neg \neg \neg$	١.

True Labels

LASSIFICATIC	N REPORT :			
	precision	recall	f1-score	support
0	1.00	0.33	0.50	3
1	1.00	1.00	1.00	3
2	0.50	0.67	0.57	3
3	1.00	1.00	1.00	3
4	0.67	0.67	0.67	3
5	1.00	0.67	0.80	3
6	1.00	1.00	1.00	3
7	1.00	0.67	0.80	3
8	1.00	1.00	1.00	3
9	1.00	1.00	1.00	3
10	1.00	1.00	1.00	3
11	1.00	1.00	1.00	3
12	1.00	0.67	0.80	3
13	1.00	1.00	1.00	3
14	1.00	1.00	1.00	3
15	1.00	1.00	1.00	3
16	1.00	1.00	1.00	3
17	1.00	1.00	1.00	3
18	1.00	1.00	1.00	3
19	1.00	1.00	1.00	3
20	0.60	1.00	0.75	3
21	1.00	0.67	0.80	3
22	1.00	1.00	1.00	3
23	0.75	1.00	0.86	3
24	0.75	1.00	0.86	3
25	1.00	0.67	0.80	3

	26	1.00	0.67	0.80	3
	27	1.00	1.00	1.00	3
	28	1.00	1.00	1.00	3
	29	0.60	1.00	0.75	3
	30	1.00	1.00	1.00	3
	31	1.00	0.67	0.80	3
	32	1.00	1.00	1.00	3
	33	1.00	1.00	1.00	3
	34	1.00	0.67	0.80	3
	35	1.00	1.00	1.00	3
	36	1.00	1.00	1.00	3
	37	1.00	1.00	1.00	3
	38	0.75	1.00	0.86	3
	39	0.60	1.00	0.75	3
accura	acy			0.90	120
macro a	avg	0.93	0.90	0.90	120
weighted a	avg	0.93	0.90	0.90	120

This classification model achieves 90% accuracy.

3 Image compression

- How would you "compress" each image in the previous dataset by using any of the studied factorizations?
- Choose one image and plot the compressed one and determine the best k orthogonal bases for its range. The "best k" is computed such that the approximated image will have a relative distance in 2-norm from the original image less than 10^{-2} . Plot the approximated image.
- Use k principal directions for the range of all the images to repeat the classification task by using these directions rather than the whole images.

Conduct of the exercise n.3

First point

In order to compress each image in the dataset, a studied and useful technique is **URV-factorisation**, which reduces dimensionality while retaining the main information.

In URV, a matrix A (the image) is decomposed as:

$$A = URV^T$$

where:

- U and V are unitary matrices,
- *R* is a triangular matrix containing singular values.

To compress, we can truncate R, keeping only the most informative singular values. By doing so, we obtain an approximate version of A that retains the essential information but occupies less memory space.

Second point

```
plt.imshow(img, cmap='gray')
plt.grid(False)
plt.show()
```

```
0
10
20
30
40
50
60
0 20 40 60
```

```
In [39]: def relative_distance(original_Matrix, reconstruction_matrix):
    normNum = linalg.norm(original_Matrix - reconstruction_matrix)
    normDen = linalg.norm(original_Matrix)

    error = normNum/normDen

    return error
```

We apply URV factorisation to image:

```
In [41]: import numpy as np
import scipy.linalg as spl

#URV factorization, starting from QR fattorization
[Qimg,Rimg,Pimg]=spl.qr(img,pivoting=True)
[Ql,Rl]=spl.qr(Rimg.T)

ndi=(np.diag(Rl))
```

We calculate the best k taking into account a set of tolerances, but also the distance between the original and the reconstructed matrix, which, based on the Euclidean norm, must not exceed 1e-2:

```
In [42]:
        tol=[1e-1,5e-2,1e-2,5e-3,1e-3,1e-5,1e-8]
         imgs = []
         k = []
         distances = []
         k opt = []
         for ind in tol:
          nc=np.sum(abs(ndi)/max(abs(ndi)) >= ind)
          k.append(nc)
          QA=np.copy(Qimg[:,0:nc])
         # QA basis for the column space of A
          RA=np.copy(Rl[0:nc,0:nc].T)
          QLA=np.copy(Ql[:,0:nc].T)
          Apimg=np.dot(np.dot(QA,RA),QLA)
          PTimg = np.zeros(Pimg.shape, dtype=int)
          PTimg[Pimg] = range (Pimg.shape[0])
```

```
Apimg=Apimg[:,PTimg]
imgs.append(Apimg)

dist= relative_distance(img,Apimg)
if dist< 1e-2:
    k_opt.append(int(nc))
distances.append(dist)

fig, axes = plt.subplots(1, len(imgs), figsize=(20, 4))

for i in range(1):
    for j in range(len(imgs)):
        ax = axes[j]
        ax.imshow(imgs[i*10+j], cmap='grey')
        ax.set_title(f"k:{k[j]}\ndistance:{distances[j]}")
        ax.axis('off')

plt.tight_layout()
plt.show()</pre>
```



```
In [43]: print(f'The optimal k are: {k_opt}')
The optimal k are: [40, 63, 64]
```

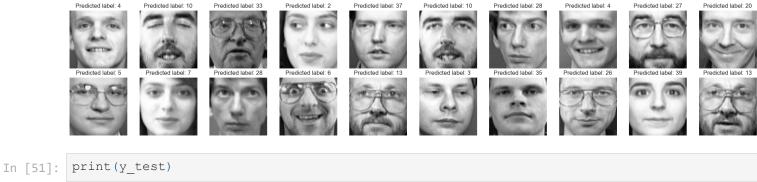
Since the objective is to compress the original image which is 64*64 in size, then we choose as best k, 40.

Third point

We now repeat the classification task considering 40 k:

```
from sklearn.decomposition import PCA
In [44]:
         k = 40
         pca = PCA(copy=True, n components=k, whiten = True).fit(X trainC)
         IM pca = pca.transform(X trainC)
         IM pca = IM pca/np.sqrt(nrc)
         print("IM_pca: " + str(IM_pca.shape))
        IM pca: (280, 40)
In [45]: cli = {}
         for i in range(40): # Da 0 a 40 inclusi
             cli[str(i)] = []
         unique classes = np.unique(y train)
         for c in unique classes:
             for index,val in enumerate(y train):
                 if val == c:
                     PT1 = X trainC[index]
                     Pimg = pca.transform([PT1])
                     cli[str(c)].append(Pimg)
             cli[str(c)] = np.asarray(cli[str(c)])
         om = {}
```

```
for i in range (0, 40):
             om[i]=np.mean(cli[str(i)],axis=0)
In [46]: (nrc,nt) = X test.shape
         X testC = centralizing matrix(X test,nrc,nt)
         X testC.shape
         (120, 4096)
Out[46]:
In [47]:
         testProjector = []
         for i in range (0, 120):
             PT1 t = X testC[i]
             Pimg t = pca.transform([PT1 t])
             testProjector.append(Pimg t)
         len(testProjector)
         120
Out[47]:
         from numpy.linalg import norm
In [48]:
         predicted class = []
         for projectionT in testProjector:
             distance min = float('inf')
             predict class = -1
             for class id, mean comp in om.items():
                 distance = norm(mean comp-projectionT,2)
                 if distance < distance min:</pre>
                     distance min = distance
                     predict class = class id
             predicted class.append(predict class)
In [49]: print(predicted class)
         [4, 10, 33, 2, 37, 10, 28, 4, 27, 20, 5, 7, 28, 6, 13, 3, 35, 26, 39, 13, 2, 27, 38, 20,
         16, 35, 14, 2, 34, 37, 11, 32, 12, 22, 28, 3, 20, 7, 33, 9, 5, 14, 39, 4, 6, 33, 20, 24,
         39, 8, 36, 31, 17, 21, 39, 20, 35, 22, 39, 36, 16, 29, 25, 30, 32, 38, 17, 36, 8, 11, 1
         0, 25, 38, 9, 26, 15, 17, 2, 6, 11, 5, 18, 16, 20, 14, 13, 38, 32, 8, 21, 31, 5, 1, 19,
         23, 27, 37, 1, 29, 5, 19, 3, 22, 14, 19, 15, 20, 15, 23, 24, 20, 26, 18, 24, 30, 4, 23,
         1, 0, 24]
In [50]: images = X test[:20].reshape(-1, 64, 64)
         labels = predicted class[:20]
         # Crea una figura e una griglia di subplot
         fig, axes = plt.subplots(2, 10, figsize=(20, 4))
         # Iterazione sui subplot e aggiunta delle immagini e delle label
         for i in range(2):
             for j in range(10):
                 ax = axes[i, j]
                 ax.imshow(images[i*10+j], cmap='grey') # Visualizza in scala di grigi
                 ax.set title(f"Predicted label: {labels[i*10+j]}")
                 ax.axis('off')
         plt.tight layout()
         plt.show()
```



```
[ 4 10 33 7 37 10 28 4 27 2 5 7 28 6 13 3 35 26 9 13 25 27 38 0
         16 35 14 2 34 37 11 32 12 22 28 3 0 7 33 9 5 14 39 34 6 33 20 24
         34 8 36 31 17 21 18 20 35 22 39 36 16 29 25 30 32 21 17 36 8 11 10 25
         38 9 26 15 17 2 6 11 29 18 16 31 12 13 38 32 8 21 31 5 1 19 23 27
         37 1 29 30 19 3 22 14 19 15 39 15 23 24 20 26 18 24 30 4 23 1 0 12]
In [51]: | images = X_test[:20].reshape(-1, 64, 64)
        labels = y test[:20]
         # Crea una figura e una griglia di subplot
        fig, axes = plt.subplots(2, 10, figsize=(20, 4))
         # Iterazione sui subplot e aggiunta delle immagini e delle label
        for i in range(2):
            for j in range(10):
                ax = axes[i, j]
                ax.imshow(images[i*10+j], cmap='grey') # Visualizza in scala di grigi
                ax.set title(f"Label: {labels[i*10+j]}")
                ax.axis('off')
        plt.tight layout()
        plt.show()
```



```
from sklearn.metrics import confusion matrix, classification report
In [52]:
         import seaborn as sns
         def create confusionMatrix andReport(predicted class):
             conf matrix = confusion_matrix(y_test,predicted_class)
             class report = classification report(y test,predicted class)
            print("CONFUSION MATRIX:")
            plt.figure(figsize=(10, 8))
            sns.heatmap(conf matrix, annot=True, fmt='d', cmap='Blues', cbar=False)
            plt.xlabel('Predicted Labels')
            plt.ylabel('True Labels')
            plt.title('Confusion Matrix')
            plt.show()
            print("\nCLASSIFICATION REPORT :")
            print(class report)
         create confusionMatrix andReport(predicted class)
```

Confusion Matrix

0 0 0 0 0 0 12 13 14 15 16 18 19 20 21 22 23 24 25 26 27 28 29 30 31 32 33 34 35 36 37 38 1 0 1 0 0 0 0 0 0 0 0 0 0 0 3 0 0 0 0 0 0 0 0 0 0 0 0 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21 22 23 24 25 26 27 28 29 30 31 32 33 34 35 36 37 38 39

Predicted Labels

CT.ASS	STFT	CATT	ON R	F.POR	г •

True Labels

LASSIFICATIO	N REPORT :			
	precision	recall	f1-score	support
0	1.00	0.33	0.50	3
1	1.00	1.00	1.00	3
2	0.50	0.67	0.57	3
3	1.00	1.00	1.00	3
4	0.75	1.00	0.86	3
5	0.75	1.00	0.86	3
6	1.00	1.00	1.00	3
7	1.00	0.67	0.80	3
8	1.00	1.00	1.00	3
9	1.00	0.67	0.80	3
10	1.00	1.00	1.00	3
11	1.00	1.00	1.00	3
12	1.00	0.33	0.50	3
13	1.00	1.00	1.00	3
14	0.75	1.00	0.86	3
15	1.00	1.00	1.00	3
16	1.00	1.00	1.00	3
17	1.00	1.00	1.00	3
18	1.00	0.67	0.80	3
19	1.00	1.00	1.00	3
20	0.38	1.00	0.55	3
21	1.00	0.67	0.80	3
22	1.00	1.00	1.00	3
23	1.00	1.00	1.00	3
24	0.75	1.00	0.86	3
25	1.00	0.67	0.80	3

2	6	1.00	1.00	1.00	3
2	7	1.00	1.00	1.00	3
2	8	1.00	1.00	1.00	3
2	9	1.00	1.00	1.00	3
3	0	1.00	0.67	0.80	3
3	1	1.00	0.67	0.80	3
3	2	1.00	1.00	1.00	3
3	3	1.00	1.00	1.00	3
3	4	1.00	0.33	0.50	3
3	5	1.00	1.00	1.00	3
3	6	1.00	1.00	1.00	3
3	7	1.00	1.00	1.00	3
3	8	0.75	1.00	0.86	3
3	9	0.40	0.67	0.50	3
accurac	У			0.88	120
macro av	g	0.93	0.88	0.88	120
weighted av	g	0.93	0.88	0.88	120

We achieved a slightly lower accuracy than the previous classifier.