

## Assignment - 4

**TITLE:**

Statistical Processing .

**PROBLEM STATEMENT:**

Load Apply Basic PCA on the iris dataset.

- Describe the data set.
- Describe the structured of correlation among variables.
- Compute a PCA with maximum number of components.
- Compute commulative explained variance ratio.
- Determine number of component K by your computed values.
- Print K principle components and correlations of K principle component with original variables.
- Plot the samples projected into the K first PCs.
- Color sample by the species.

**OBJECTIVES:**

To be able to appropriately apply computational methodologies to real world statistical problems.

**OUTCOMES:**

Model and solve computing problem using correlation, and resampling using appropriate statistics algorithm.

**PREREQUISITE:**

→ Concept of data pre-processing.

## THEORY:

PCA finds the principle components of data. It is often useful to measure data in terms of its principle components rather than on normal  $x$ - $y$  axis. They are the directions where there is most variance the directions where data is most spread out.

Principle component Analysis (PCA) is unsupervised reduction algorithm. It identifies hyperplane that lies closest to the data, and then it projects the data onto it preserving the variance.

PCA finds a new set of dimensions such that all dimensions are orthogonal and ranked according to variance of data along them. It means more important principle axis occurs first.

How does PCA works?

- Calculate covariance matrix  $X$  of data points
- Calculate eigen vectors and corresponding eigen values.
- select eigen vector according to decreasing eigen values.
- Choose first  $k$  eigen vectors and that will be new  $k$  dimensions.
- Transform the original  $n$  dimensional data points into  $k$  dimensions.



## Implementing PCA on a 2-D dataset

→ Step-1: Normalise the data

First step is to normalise the data that we have so that PCA works properly.

→ Step-2: Calculate the covariance matrix

Since the dataset we took is 2-dimensional this will result in a  $2 \times 2$  covariance matrix

$$\text{Matrix (Covariance)} = \begin{bmatrix} \text{Var}[X_1] & \text{Cor}[X_1, X_2] \\ \text{Cor}[X_2, X_1] & \text{Var}[X_2] \end{bmatrix}$$

→ Step-3: Calculate the eigen values and eigen vectors.

$$\det(\lambda I - A) = 0$$

$$(\lambda I - A)v = 0$$

→ Step-4: Choosing components and forming a feature vector.

We do not lose out some information in the process, but if eigen values are small, we do not lose much.

### CONCLUSION:

Hence we conclude that, Principal Component Analysis (PCA) is useful to visualise the high-dimensional datasets, as it can compress it down to 2-dimensions.



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Apply Basic PCA on the iris dataset. The data set is available at: <https://raw.githubusercontent.com/neurospn/pystatsml/master/datasets/iris.csv>.

Describe the data set. Should the dataset been standardized? • Describe the structure of correlations among variables. • Compute a PCA with the maximum number of components. • Compute the cumulative explained variance ratio. Determine the number of components by your computed values. • Print the principal components directions and correlations of the principal components with the original variables. Interpret the contribution of the original variables into the PC. • Plot the samples projected into the first PCs. • Color samples by their species.

```
[ ] import numpy as np
import pandas as pd
import seaborn as sns
import matplotlib.pyplot as plt
```

```
[ ] iris_data = pd.read_csv("/content/drive/MyDrive/data/iris.csv")
iris_data.head()
```

	sepal_length	sepal_width	petal_length	petal_width	species
0	5.1	3.5	1.4	0.2	setosa
1	4.9	3.0	1.4	0.2	setosa
2	4.7	3.2	1.3	0.2	setosa
3	4.6	3.1	1.5	0.2	setosa

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[ ]	3	4.6	3.1	1.5	0.2	setosa
	4	5.0	3.6	1.4	0.2	setosa

Standardization 1)Each feature in the dataset has different mean and SD hence it is necessary to scale data for comparision with other features.

2)Standardization sets mean to zero and SD of 1 for all features.

3)As seen in the data below the SD and is different for each datasetalso the variation is different.

```
[ ] print("Sepal length range: [%s, %s]" % (min(iris_data["sepal_length"]),max(iris_data["sepal_length"])))
print("Sepal width range: [%s, %s]" % (min(iris_data["sepal_width"]),max(iris_data["sepal_width"])))
print("Petal length range: [%s, %s]" % (min(iris_data["petal_length"]),max(iris_data["petal_length"])))
print("Petal width range: [%s, %s]" % (min(iris_data["petal_width"]),max(iris_data["petal_width"])))

print("Sepal Length standard deviation :\t %f" %np.std(iris_data["sepal_length"]))
print("Sepal Width standard deviation :\t %f" %np.std(iris_data["sepal_width"]))
print("Petal Length standard deviation :\t %f" %np.std(iris_data["petal_length"]))
print("Petal Width standard deviation :\t %f" %np.std(iris_data["petal_width"]))
```

Sepal length range: [4.3, 7.9]

Sepal width range: [2.0, 4.4]

Petal length range: [1.0, 6.9]

Petal width range: [0.1, 2.5]

Sepal Length standard deviation : 0.825301

Sepal Width standard deviation : 0.435167

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Petal width standard deviation : 0.759693

Correlation Matrix :

```
[ ] corr = iris_data.corr(method = "pearson")
display(corr)
sns.heatmap(corr,cmap =sns.diverging_palette(10,0,as_cmap=True),annot= True,fmt = "f")
```

	sepal_length	sepal_width	petal_length	petal_width
sepal_length	1.000000	-0.117570	0.871754	0.817941
sepal_width	-0.117570	1.000000	-0.428440	-0.366126
petal_length	0.871754	-0.428440	1.000000	0.962865
petal_width	0.817941	-0.366126	0.962865	1.000000

<matplotlib.axes.\_subplots.AxesSubplot at 0x7f9ee30b5750>

	sepal_length	sepal_width	petal_length	petal_width
sepal_length	1.000000	-0.117570	0.871754	0.817941
sepal_width	-0.117570	1.000000	-0.428440	-0.366126
petal_length	0.871754	-0.428440	1.000000	0.962865
petal_width	0.817941	-0.366126	0.962865	1.000000

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1)Sepal\_length <-> Petal\_Length : 0.871754 (Pcsitive Correlation)

2)Sepal\_width <-> Petal\_Length : -0.428440 (Negative Correlation)

3)Petal\_length <-> Petal\_width : 0.962865 (Highly Positive Correlation)

4)Petal\_width <-> Sepal\_Length : 0.817941 (Positive Correlation)

PCA (Principal Ccmpnent Analysis):

```
[ ] from sklearn.decomposition import PCA
pca = PCA()
x_new1 =pca.fit_transform(iris_data.drop(["species"],axis =1))
x_new1[:5]
```

array([[ -2.68412563e+00, 3.19397247e-01, -2.79148276e-02,
 -2.26243707e-03],
 [ -2.71414169e+00, -1.77001225e-01, -2.10464272e-01,
 -9.90265503e-02],
 [ -2.88899057e+00, -1.44949426e-01, 1.79002563e-02,
 -1.99683897e-02],
 [ -2.74534286e+00, -3.18298979e-01, 3.15593736e-02,
 7.55758166e-02],
 [ -2.72871654e+00, 3.26754513e-01, 9.00792406e-02,
 6.12585926e-02]])

Explained Variance of PCA

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Explained Variance of R^2:

[ ] explained\_variance= pca.explained\_variance\_ratio\_  
  
explained\_variance  
  
array([0.92461872, 0.05306648, 0.01710261, 0.00521218])

[ ] plt.figure(figsize =(6,4))  
plt.bar(range(4),explained\_variance, alpha=0.5, align='center', label ="Individual Explained Variance", color ="darkblue")  
plt.ylabel("Explained Variance Ratio")  
plt.xlabel("Principal Component")  
plt.legend(loc ="best")  
plt.tight\_layout()



Principal Component	Explained Variance Ratio
1	0.92461872
2	0.05306648
3	0.01710261
4	0.00521218

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There are 3 Principal Components. Applying PCA

```
[ ] pca = PCA(n_components=3)
x_new =pca.fit_transform(iris_data.drop(['species'],axis =1))
x_new[:5]

array([[ -2.68412563,  0.31939725, -0.02791483],
       [ -2.71414169, -0.17700123, -0.21046427],
       [ -2.88899057, -0.14494943,  0.01790026],
       [ -2.74534286, -0.31829898,  0.03155937],
       [ -2.72871654,  0.32675451,  0.09007924]])
```

Correlation and Direction of PCA converting oategorical data to numerical

```
[ ] categ_num={"species":{"setosa":0,"versicolor":1,"virginica":2}}
iris_data1 = iris_data.replace(categ_num)
columns = list(iris_data.columns[:4])
```

Comparing Principal Components with original variables

```
[ ] fig, axes =plt.subplots(3,4,figsize =(15,10))
k= 0
for i in range(axes.shape[0]):
```

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```
[ ] k= 0
for i in range(axes.shape[0]):
    for j in range(axes.shape[1]):
        axes[i,j].scatter(x_new[:,i],iris_data[columns[j]],c =iris_data1["species"])
plt.show()
```

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2.5

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4.5

7

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