

Experiment No: 03

● <u>Aim</u>: To implement PCA / SVD / LDA.

●Theory:

Principal Component Analysis (PCA):

Principal Component Analysis is an unsupervised learning algorithm that is used for the dimensionality reduction in machine learning. It is a statistical process that converts the observations of correlated features into a set of linearly uncorrelated features with the help of orthogonal transformation. These new transformed features are called the Principal Components. It is one of the popular tools that is used for exploratory data analysis and predictive modeling. It is a technique to draw strong patterns from the given dataset by reducing the variances.

PCA works by considering the variance of each attribute because the high attribute shows the good split between the classes, and hence it reduces the dimensionality. Some real-world applications of PCA are image processing, movie recommendation system, optimizing the power allocation in various communication channels. It is a feature extraction technique, so it contains the important variables and drops the least important variable.

Steps for PCA algorithm:

Getting the dataset

Firstly, we need to take the input dataset and divide it into two subparts X and Y, where X is the training set, and Y is the validation set.

Representing data into a structure

Now we will represent our dataset into a structure. Such as we will represent the two-dimensional matrix of independent variable X. Here each row corresponds to the data items, and the column corresponds to the Features. The number of columns is the dimensions of the dataset.

Standardizing the data

In this step, we will standardize our dataset. Such as in a particular column, the features with high variance are more important compared to the features with lower variance.

If the importance of features is independent of the variance of the feature, then we will divide each data item in a column with the standard deviation of the column. Here we will name the matrix as Z.

Calculating the Covariance of Z

To calculate the covariance of Z, we will take the matrix Z, and will transpose it. After transpose, we will multiply it by Z. The output matrix will be the Covariance matrix of Z.



Calculating the Eigen Values and Eigen Vectors

Now we need to calculate the eigenvalues and eigenvectors for the resultant covariance matrix Z. Eigenvectors or the covariance matrix are the directions of the axes with high information. And the coefficients of these eigenvectors are defined as the eigenvalues.

Sorting the Eigen Vectors

In this step, we will take all the eigenvalues and will sort them in decreasing order, which means from largest to smallest. And simultaneously sort the eigenvectors accordingly in matrix P of eigenvalues. The resultant matrix will be named as P*.

Calculating the new features Or Principal Components

Here we will calculate the new features. To do this, we will multiply the P* matrix to the Z. In the resultant matrix Z*, each observation is the linear combination of original features. Each column of the Z* matrix is independent of each other.

Remove less or unimportant features from the new dataset.

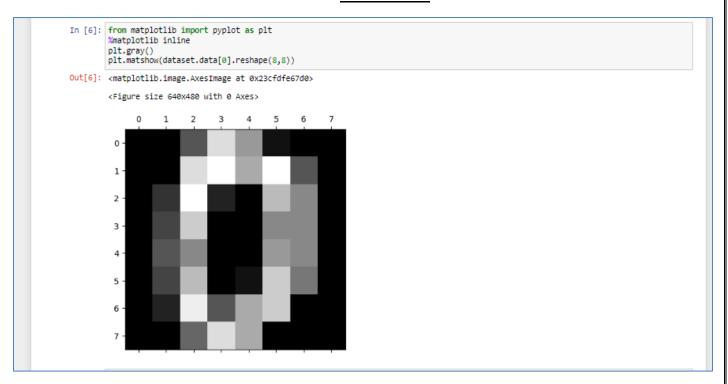
The new feature set has occurred, so we will decide here what to keep and what to remove. It means, we will only keep the relevant or important features in the new dataset, and unimportant features will be removed out.

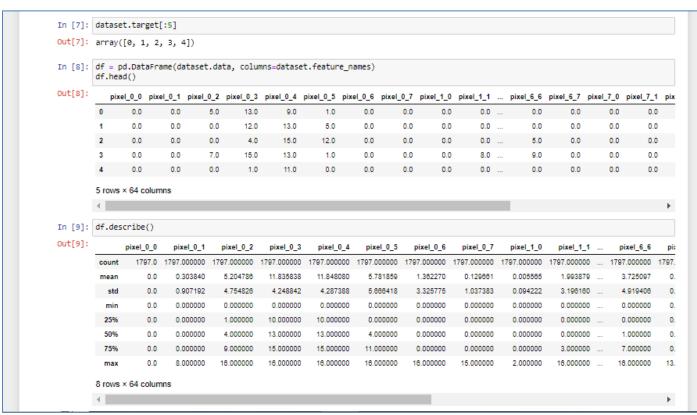
Applications of Principal Component Analysis

- PCA is mainly used as the dimensionality reduction technique in various AI applications such as computer vision, image compression, etc.
- It can also be used for finding hidden patterns if data has high dimensions. Some fields where PCA is used are Finance, data mining, Psychology, etc

Implementation of Principal Component Analysis (PCA):







```
In [10]: X = df
    y = dataset.target

In [11]: y
Out[11]: array([0, 1, 2, ..., 8, 9, 8])
```



```
In [12]: from sklearn.preprocessing import StandardScaler
            scaler = StandardScaler()
            X_scaled = scaler.fit_transform(X)
            X_scaled
                       0. , -0.33501649, -0.04308102, ..., -1.14664746,
-0.5056698 , -0.19600752],
Out[12]: array([[ 0.
                       -0.595698 , -0.3591649, -1.09493684, ..., 0.54856067, -0.5056698 , -0.19600752], [ 0. , -0.33501649, -1.09493684, ..., 1.56568555, 1.6951369 , -0.19600752],
                     [ 0.
                     [ 0. , -0.33501649, -0.88456568, ..., -0.12952258, -0.5056698 , -0.19600752], [ 0. , -0.33501649, -0.67419451, ..., 0.8876023 ,
                       -0.5056698 , -0.19600752],
                       0. , -0.33501649,
-0.26113572, -0.19600752]])
                                                         1.00877481, ..., 0.8876023 ,
                     [ 0.
In [13]: from sklearn.model_selection import train_test_split
            X_train, X_test, y_train, y_test = train_test_split(X_scaled, y, test_size=0.2, random_state=30)
In [14]: from sklearn.linear model import LogisticRegression
            model = LogisticRegression()
            model.fit(X_train, y_train)
model.score(X_test, y_test)
Out[14]: 0.97222222222222
In [20]: X.shape
Out[20]: (1797, 64)
```

```
Use PCA to reduce dimensions
In [15]: from sklearn.decomposition import PCA
            pca = PCA(0.95)
            X_pca = pca.fit_transform(X)
            X_pca.shape
Out[15]: (1797, 29)
In [21]: pca.explained_variance_ratio_
Out[21]: array([0.14890594, 0.13618771, 0.11794594, 0.08409979, 0.05782415, 0.0491691 , 0.04315987, 0.03661373, 0.03353248, 0.03078806,
                     0.02372341, 0.02272697, 0.01821863, 0.01773855, 0.01467101,
                    0.01409716, 0.01318589, 0.01248138, 0.01017718, 0.00905617, 0.00889538, 0.00797123, 0.00767493, 0.00722904, 0.00695889, 0.00596081, 0.00575615, 0.00515158, 0.0048954 ])
In [22]: pca.n_components_
Out[22]: 29
            You can see that both combined retains 0.14+0.13=0.27 or 27% of important feature information
In [24]: X_train_pca, X_test_pca, y_train, y_test = train_test_split(X_pca, y, test_size=0.2, random_state=30)
In [25]: model = LogisticRegression(max_iter=1000)
           model.fit(X_train_pca, y_train)
           model.score(X_test_pca, y_test)
Out[25]: 0.96944444444444444
```



```
Let's now select only two components
In [26]: pca = PCA(n_components=2)
         X_pca = pca.fit_transform(X)
        X_pca.shape
Out[26]: (1797, 2)
In [27]: X_pca
[ 10.80130987, -6.96024689],
[ -4.87210787, 12.42395562],
                [ -0.34434859, 6.36555467]])
In [29]: pca.explained_variance_ratio_
Out[29]: array([0.14890594, 0.13618771])
         You can see that both combined retains 0.14+0.13=0.27 or 27% of important feature information
In [30]: X_train_pca, X_test_pca, y_train, y_test = train_test_split(X_pca, y, test_size=0.2, random_state=30)
         model = LogisticRegression(max_iter=1000)
         model.fit(X_train_pca, y_train)
         model.score(X_test_pca, y_test)
Out[30]: 0.60833333333333333
```

Singular Value Decomposition (SVD):

The Singular Value Decomposition (SVD) of a matrix is a factorization of that matrix into three matrices. It has some interesting algebraic properties and conveys important geometrical and theoretical insights about linear transformations. It also has some important applications in data science. In this article, I will try to explain the mathematical intuition behind SVD and its geometrical meaning.

Mathematics behind SVD:

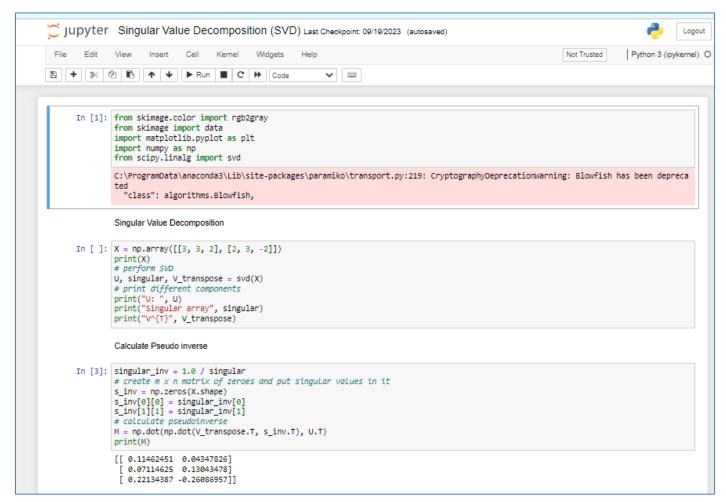
The SVD of mxn matrix A is given by the formula $A = U \setminus Sigma V^T$

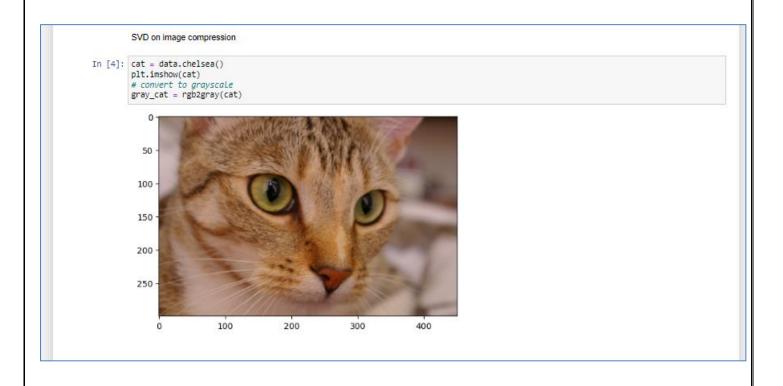
where:

- U: mxm matrix of the orthonormal eigenvectors of AA[^]{T}
- VT: transpose of a nxn matrix containing the orthonormal eigenvectors of A^TA.
- \Sigma: diagonal matrix with r elements equal to the root of the positive eigenvalues of AA^T or A^T A (both matrics have the same positive eigenvalues anyway).



Implementation of Singular Value Decomposition (SVD):



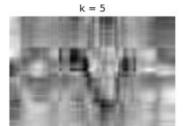




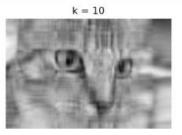
Calculate the SVD and plot the image

```
In [6]: U, S, V_T = svd(gray_cat, full_matrices=False)
S = np.diag(S)
fig, ax = plt.subplots(5, 2, figsize=(8, 20))

curr_fig = 0
for r in [5, 10, 70, 100, 200]:
    cat_approx = U[:, :r] @ S[0:r, :r] @ V_T[:r, :]
    ax[curr_fig][0].imshow(cat_approx, cmap='gray')
    ax[curr_fig][0].set_title("k = "+str(r))
    ax[curr_fig, 0].axis('off')
    ax[curr_fig][1].set_title("Original Image")
    ax[curr_fig][1].imshow(gray_cat, cmap='gray')
    ax[curr_fig, 1].axis('off')
    curr_fig = 1
plt.show()
```





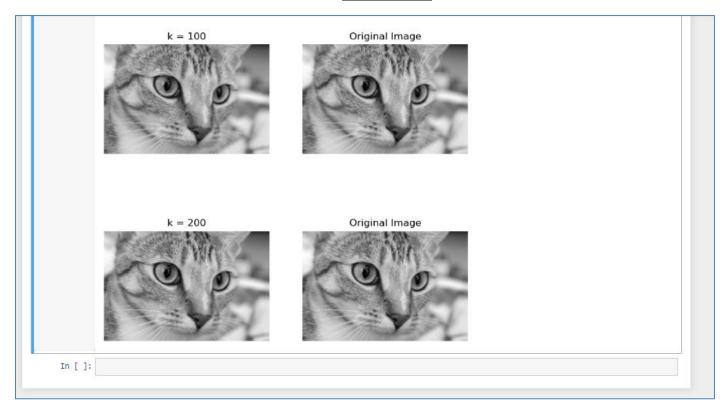












• Conclusion:

Both PCA and SVD are valuable tools for simplifying and analysing high-dimensional data, which can lead to improved data understanding and more efficient machine learning models. The choice between PCA and SVD depends on the specific problem and the mathematical and computational tools available.