

# Forward School

**Program Code: J620-002-4:2020**

**Program Name: FRONT-END SOFTWARE DEVELOPMENT**

**Title : Exe23 - Dimension Reduction Exercise**

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**Introduction :**

**Conclusion :**

## Dimension Reduction Exercise

### Principal Component Analysis

Let's begin by importing the various library

```
In [5]: import pandas as pd  
import numpy as np
```

```

In [130]: from sklearn.datasets import load_breast_cancer
breast = load_breast_cancer()
# Combine the data and target arrays
data = breast.data

# Create a dictionary with column names as keys and data arrays as values
data_dict = {name: data[:, i] for i, name in enumerate(breast.feature_names)}

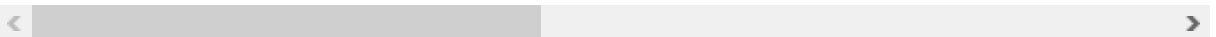
# Create the DataFrame
df = pd.DataFrame(data_dict)
df['target'] = breast.target
df

```

Out[130]:

	mean radius	mean texture	mean perimeter	mean area	mean smoothness	mean compactness	mean concavity	mean concave points	mean symmetry
0	17.99	10.38	122.80	1001.0	0.11840	0.27760	0.30010	0.14710	0.241
1	20.57	17.77	132.90	1326.0	0.08474	0.07864	0.08690	0.07017	0.181
2	19.69	21.25	130.00	1203.0	0.10960	0.15990	0.19740	0.12790	0.206
3	11.42	20.38	77.58	386.1	0.14250	0.28390	0.24140	0.10520	0.259
4	20.29	14.34	135.10	1297.0	0.10030	0.13280	0.19800	0.10430	0.180
...	...	...	...	...	...	...	...	...	...
564	21.56	22.39	142.00	1479.0	0.11100	0.11590	0.24390	0.13890	0.172
565	20.13	28.25	131.20	1261.0	0.09780	0.10340	0.14400	0.09791	0.175
566	16.60	28.08	108.30	858.1	0.08455	0.10230	0.09251	0.05302	0.159
567	20.60	29.33	140.10	1265.0	0.11780	0.27700	0.35140	0.15200	0.239
568	7.76	24.54	47.92	181.0	0.05263	0.04362	0.00000	0.00000	0.158

569 rows × 31 columns



In [131]: breast.keys()

Out[131]: dict\_keys(['data', 'target', 'frame', 'target\_names', 'DESCR', 'feature\_names', 'filename', 'data\_module'])

In [132]: df.info()

```
<class 'pandas.core.frame.DataFrame'>
RangeIndex: 569 entries, 0 to 568
Data columns (total 31 columns):
#   Column                                Non-Null Count  Dtype
---  -
0   mean radius                           569 non-null    float64
1   mean texture                           569 non-null    float64
2   mean perimeter                         569 non-null    float64
3   mean area                             569 non-null    float64
4   mean smoothness                        569 non-null    float64
5   mean compactness                       569 non-null    float64
6   mean concavity                          569 non-null    float64
7   mean concave points                    569 non-null    float64
8   mean symmetry                          569 non-null    float64
9   mean fractal dimension                  569 non-null    float64
10  radius error                           569 non-null    float64
11  texture error                           569 non-null    float64
12  perimeter error                         569 non-null    float64
13  area error                             569 non-null    float64
14  smoothness error                       569 non-null    float64
15  compactness error                      569 non-null    float64
16  concavity error                        569 non-null    float64
17  concave points error                    569 non-null    float64
18  symmetry error                          569 non-null    float64
19  fractal dimension error                 569 non-null    float64
20  worst radius                           569 non-null    float64
21  worst texture                           569 non-null    float64
22  worst perimeter                         569 non-null    float64
23  worst area                             569 non-null    float64
24  worst smoothness                       569 non-null    float64
25  worst compactness                      569 non-null    float64
26  worst concavity                         569 non-null    float64
27  worst concave points                    569 non-null    float64
28  worst symmetry                          569 non-null    float64
29  worst fractal dimension                 569 non-null    float64
30  target                                 569 non-null    int32
dtypes: float64(30), int32(1)
memory usage: 135.7 KB
```

In [133]:

In [135]: data.shape

Out[135]: (569, 30)

```
In [158]: from sklearn.preprocessing import StandardScaler
X = df.iloc[:, 0:30].values
scaler = StandardScaler()
scaled_data = scaler.fit_transform(X)

pca = PCA(copy=True, iterated_power='auto', n_components=2, random_state=None,
pca.fit(scaled_data)
x_pca = pca.fit_transform(scaled_data)
x_pca.shape
```

Out[158]: (569, 2)

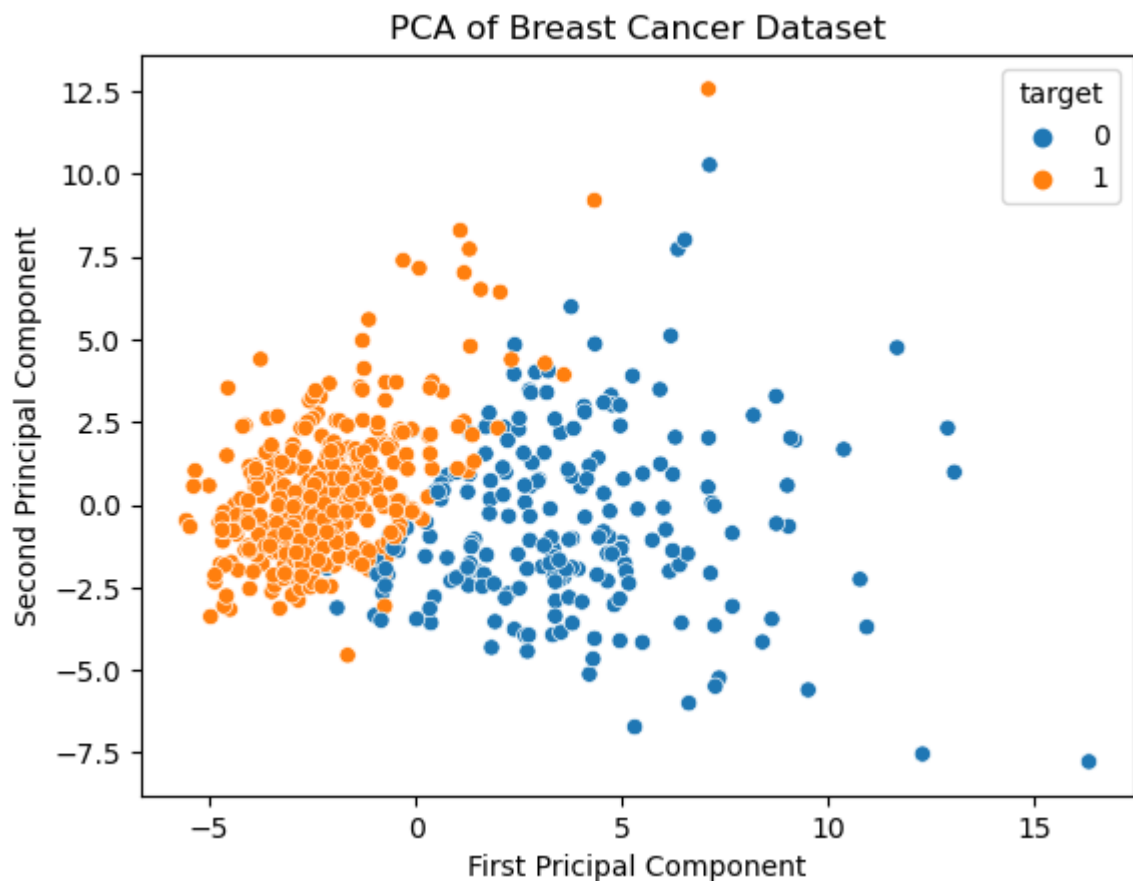
```
In [159]: import seaborn as sns
import matplotlib.pyplot as plt

pca_df = pd.DataFrame(x_pca,
                      columns=['First Principal Component',
                              'Second Principal Component'])

pca_df['target'] = breast.target

sns.scatterplot(x='First Principal Component',
                y='Second Principal Component',
                hue='target',
                data=pca_df)

plt.title('PCA of Breast Cancer Dataset')
plt.show()
```

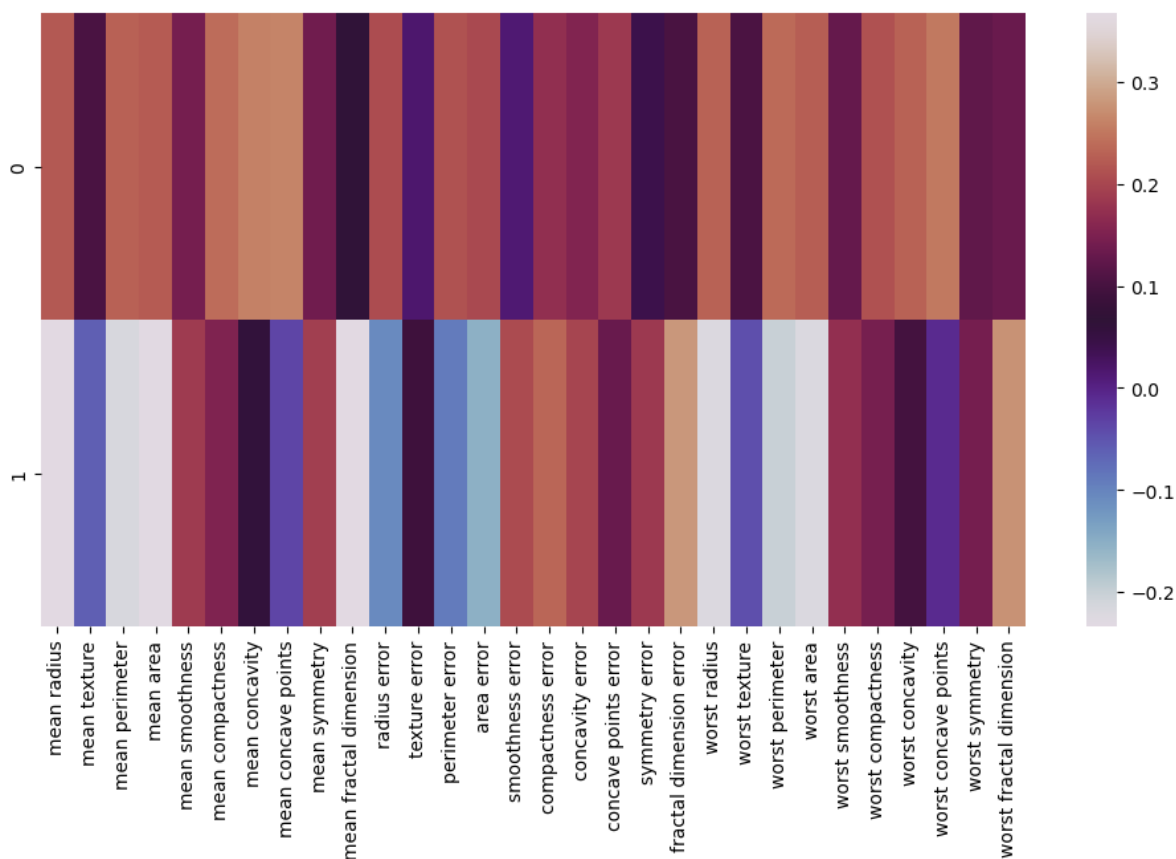


```
In [161]: pca.components_
```

```
Out[161]: array([[ 0.21890244,  0.10372458,  0.22753729,  0.22099499,  0.14258969,
                   0.23928535,  0.25840048,  0.26085376,  0.13816696,  0.06436335,
                   0.20597878,  0.01742803,  0.21132592,  0.20286964,  0.01453145,
                   0.17039345,  0.15358979,  0.1834174 ,  0.04249842,  0.10256832,
                   0.22799663,  0.10446933,  0.23663968,  0.22487053,  0.12795256,
                   0.21009588,  0.22876753,  0.25088597,  0.12290456,  0.13178394],
                  [-0.23385713, -0.05970609, -0.21518136, -0.23107671,  0.18611302,
                   0.15189161,  0.06016536, -0.0347675 ,  0.19034877,  0.36657547,
                  -0.10555215,  0.08997968, -0.08945723, -0.15229263,  0.20443045,
                   0.2327159 ,  0.19720728,  0.13032156,  0.183848 ,  0.28009203,
                  -0.21986638, -0.0454673 , -0.19987843, -0.21935186,  0.17230435,
                   0.14359317,  0.09796411, -0.00825724,  0.14188335,  0.27533947]])
```

```
In [165]: map_df = pd.DataFrame(pca.components_, columns=breast['feature_names'])
plt.figure(figsize=(12,6))
sns.heatmap(map_df,cmap='twilight')
```

```
Out[165]: <Axes: >
```





In [6]: `data.describe()`

Out[6]:

	Alcohol	Malic_Acid	Ash	Ash_Alcanity	Magnesium	Total_Phenols	Flavanoids
count	178.000000	178.000000	178.000000	178.000000	178.000000	178.000000	178.000000
mean	13.000618	2.336348	2.366517	19.494944	99.741573	2.295112	2.029270
std	0.811827	1.117146	0.274344	3.339564	14.282484	0.625851	0.998859
min	11.030000	0.740000	1.360000	10.600000	70.000000	0.980000	0.340000
25%	12.362500	1.602500	2.210000	17.200000	88.000000	1.742500	1.205000
50%	13.050000	1.865000	2.360000	19.500000	98.000000	2.355000	2.135000
75%	13.677500	3.082500	2.557500	21.500000	107.000000	2.800000	2.875000
max	14.830000	5.800000	3.230000	30.000000	162.000000	3.880000	5.080000

In [7]: *#2. Split the dataset into the Training and the Test set. Set the test set to 0.3*  
`from sklearn.model_selection import train_test_split`  
`X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.3, random_state=42)`

In [106]: *#3. Scale the train and test set using the StandardScaler*  
`from sklearn.preprocessing import StandardScaler`  
`model = StandardScaler()`  
`X_train_scaled = model.fit_transform(X_train)`  
`X_test_scaled = model.transform(X_test)`

In [107]: `from sklearn.decomposition import PCA`  
`pca = PCA(n_components=2)`  
`X_train_pca = pca.fit_transform(X_train_scaled)`  
`X_test_pca = pca.transform(X_test_scaled)`

In [108]: *#4. Apply the PCA function to both the test and train set, to extract the first two principal components*  
`pca.explained_variance_ratio_`

Out[108]: `array([0.36196226, 0.18763862])`

In [114]: `from sklearn.linear_model import LogisticRegression`  
`logr = LogisticRegression()`  
`logr.fit(X_train_pca, y_train)`

Out[114]: `LogisticRegression()`

**In a Jupyter environment, please rerun this cell to show the HTML representation or trust the notebook.**

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```
In [115]: #5. Create a Logistic Regression based on the training set
```

```
In [116]: #6. Apply the created LR model onto the test data
# Predicting the test set result using
# predict function under LogisticRegression
from sklearn.linear_model import LogisticRegression
y_pred = logr.predict(X_test_pca)
y_pred
```

```
Out[116]: array([1, 1, 3, 1, 2, 1, 2, 3, 2, 3, 2, 3, 1, 2, 1, 2, 2, 2, 1, 2, 1, 2,
                2, 3, 3, 3, 2, 2, 2, 1, 1, 2, 3, 1, 1, 1, 3, 3, 2, 3, 2, 2, 2, 2,
                3, 1, 2, 2, 3, 1, 2, 1, 1, 3], dtype=int64)
```

```
In [117]: #7. Create a confusion matrix to score the prediction performed
from sklearn.metrics import confusion_matrix
confusion_matrix = confusion_matrix(y_test, y_pred)
print(confusion_matrix)
```

```
[[17  2  0]
 [ 0 21  0]
 [ 0  0 14]]
```



```

In [119]: from matplotlib.colors import ListedColormap

X_set, y_set = X_train_pca, y_train
X1, X2 = np.meshgrid(np.arange(start = X_set[:, 0].min() - 1,
                               stop = X_set[:, 0].max() + 1, step = 0.01),
                     np.arange(start = X_set[:, 1].min() - 1,
                               stop = X_set[:, 1].max() + 1, step = 0.01))

plt.contourf(X1, X2, logr.predict(np.array([X1.ravel(),
                                             X2.ravel()])).T.reshape(X1.shape), alpha = 0.75,
              cmap = ListedColormap(('yellow', 'white', 'aquamarine'))))

plt.xlim(X1.min(), X1.max())
plt.ylim(X2.min(), X2.max())

for i, j in enumerate(np.unique(y_set)):
    plt.scatter(X_set[y_set == j, 0], X_set[y_set == j, 1],
                c = ListedColormap(('red', 'green', 'blue'))(i), label = j)

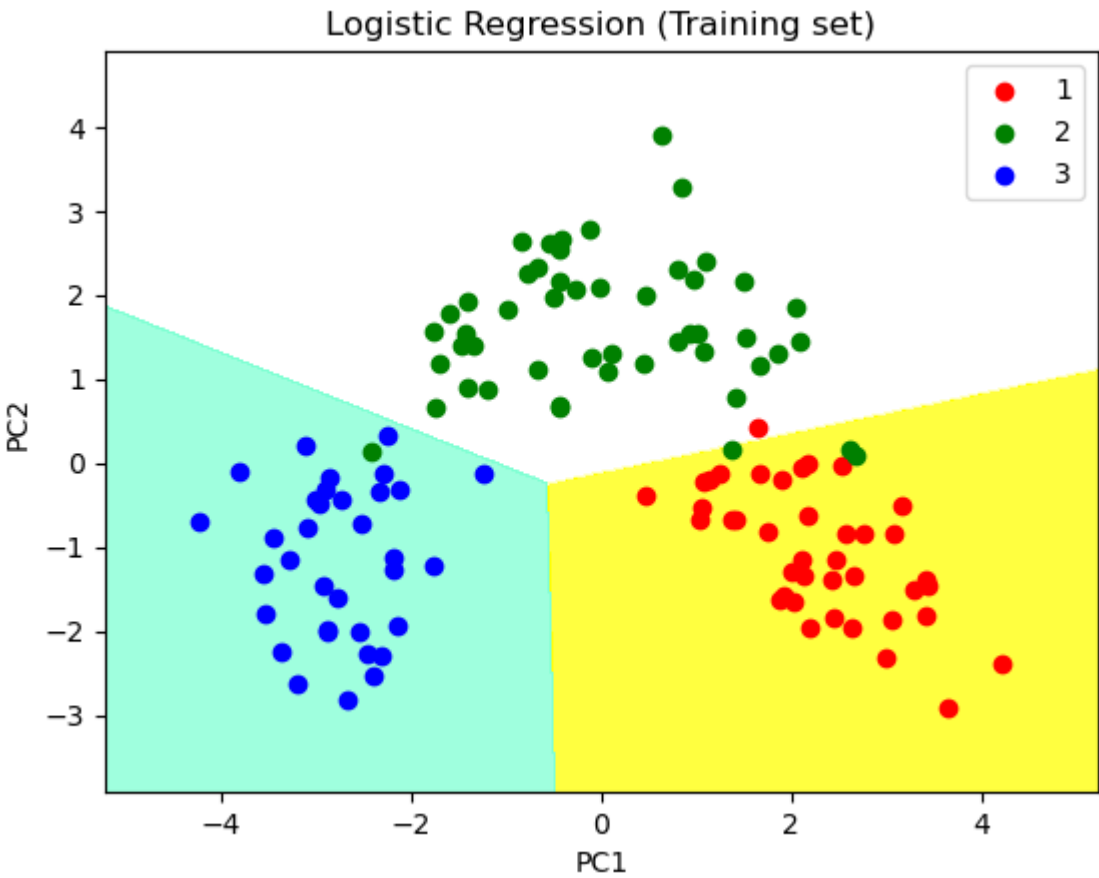
plt.title('Logistic Regression (Training set)')
plt.xlabel('PC1') # for Xlabel
plt.ylabel('PC2') # for Ylabel
plt.legend() # to show legend

# show scatter plot
plt.show()

```

C:\Users\Xiang Ze\AppData\Local\Temp\ipykernel\_11056\2160364856.py:17: UserWarning: \*c\* argument looks like a single numeric RGB or RGBA sequence, which should be avoided as value-mapping will have precedence in case its length matches with \*x\* & \*y\*. Please use the \*color\* keyword-argument or provide a 2D array with a single row if you intend to specify the same RGB or RGBA value for all points.

```
plt.scatter(X_set[y_set == j, 0], X_set[y_set == j, 1],
```



In [121]: *# Visualising the Test set results through scatter plot*

```

from matplotlib.colors import ListedColormap

X_set, y_set = X_test_pca, y_test

X1, X2 = np.meshgrid(np.arange(start = X_set[:, 0].min() - 1,
                               stop = X_set[:, 0].max() + 1, step = 0.01),
                     np.arange(start = X_set[:, 1].min() - 1,
                               stop = X_set[:, 1].max() + 1, step = 0.01))

plt.contourf(X1, X2, logr.predict(np.array([X1.ravel(),
                                             X2.ravel()]).T).reshape(X1.shape), alpha = 0.75,
             cmap = ListedColormap(('yellow', 'white', 'aquamarine'))))

plt.xlim(X1.min(), X1.max())
plt.ylim(X2.min(), X2.max())

for i, j in enumerate(np.unique(y_set)):
    plt.scatter(X_set[y_set == j, 0], X_set[y_set == j, 1],
                c = ListedColormap(('red', 'green', 'blue'))(i), label = j)

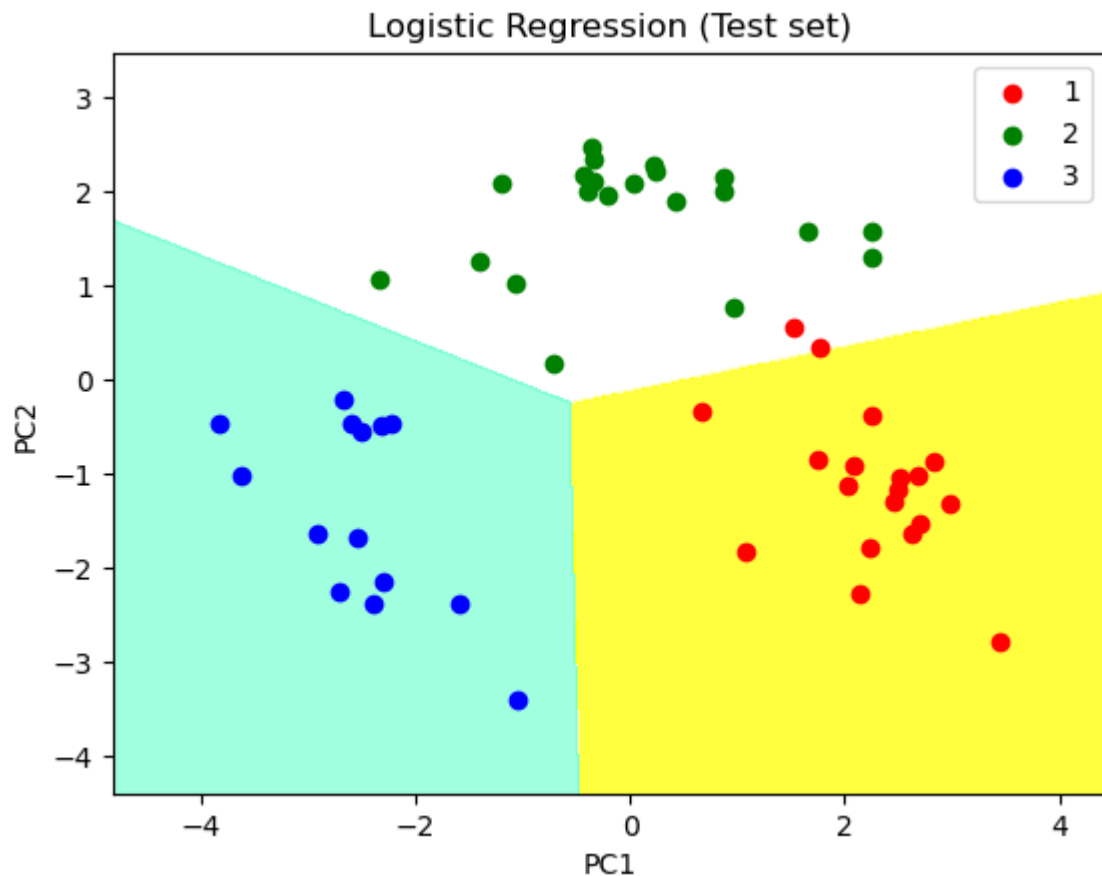
# title for scatter plot
plt.title('Logistic Regression (Test set)')
plt.xlabel('PC1') # for XLabel
plt.ylabel('PC2') # for YLabel
plt.legend()

# show scatter plot
plt.show()
# show scatter plot

```

C:\Users\Xiang Ze\AppData\Local\Temp\ipykernel\_11056\1590122172.py:20: UserWarning: \*c\* argument looks like a single numeric RGB or RGBA sequence, which should be avoided as value-mapping will have precedence in case its length matches with \*x\* & \*y\*. Please use the \*color\* keyword-argument or provide a 2D array with a single row if you intend to specify the same RGB or RGBA value for all points.

```
plt.scatter(X_set[y_set == j, 0], X_set[y_set == j, 1],
```



## Exercise 1. Take home test / challenge test

Repeat this exercise but this time do it without dimension reduction. What do you get?

```
In [10]: from sklearn.tree import DecisionTreeClassifier
from sklearn.metrics import accuracy_score
model = DecisionTreeClassifier()
model.fit(X_train, y_train)
y_pred = model.predict(X_test)
accuracy_rf = accuracy_score(y_test, y_pred)
print("Random Forest Accuracy:", accuracy_rf)
```

Random Forest Accuracy: 0.9444444444444444

## Random Forest Feature Selection

Let's do a guided walkthrough for a RF feature selection )

([https://chrisalbon.com/machine\\_learning/trees\\_and\\_forests/feature\\_selection\\_using\\_random\\_for](https://chrisalbon.com/machine_learning/trees_and_forests/feature_selection_using_random_for)  
([https://chrisalbon.com/machine\\_learning/trees\\_and\\_forests/feature\\_selection\\_using\\_random\\_for](https://chrisalbon.com/machine_learning/trees_and_forests/feature_selection_using_random_for)

```
In [19]: import numpy as np
from sklearn.ensemble import RandomForestClassifier
from sklearn import datasets
from sklearn.model_selection import train_test_split
from sklearn.feature_selection import SelectFromModel
from sklearn.metrics import accuracy_score
```

```
In [20]: # Load the iris dataset
iris = datasets.load_iris()

# Create a List of feature names
feat_labels = ['Sepal Length', 'Sepal Width', 'Petal Length', 'Petal Width']

# Create X from the features
X = iris.data

# Create y from output
y = iris.target
```

```
In [21]: # View the features
X[:5]
```

```
Out[21]: array([[5.1, 3.5, 1.4, 0.2],
               [4.9, 3. , 1.4, 0.2],
               [4.7, 3.2, 1.3, 0.2],
               [4.6, 3.1, 1.5, 0.2],
               [5. , 3.6, 1.4, 0.2]])
```

```
In [22]: # View the target data
iris.target
```

```
Out[22]: array([0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0,
                0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0, 0,
                0, 0, 0, 0, 0, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1,
                1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1,
                1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 2, 2, 2, 2, 2, 2, 2, 2,
                2, 2, 2, 2, 2, 2, 2, 2, 2, 2, 2, 2, 2, 2, 2, 2, 2, 2, 2, 2,
                2, 2, 2, 2, 2, 2, 2, 2, 2, 2, 2, 2, 2, 2, 2, 2, 2, 2, 2, 2])
```

```
In [ ]:
```

```
In [23]: # Split the data into 40% test and 60% training
from sklearn.model_selection import train_test_split
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.4, random_state=42)
```

```
In [24]: from sklearn.ensemble import RandomForestClassifier
# Create a random forest classifier
model = RandomForestClassifier(criterion="gini", max_depth = 2)

# Train the classifier
model.fit(X_train, y_train)

gini_importance = model.feature_importances_
print('Sepal Length', gini_importance[0])
print('Sepal Width', gini_importance[1])
print('Petal Length', gini_importance[2])
print('Petal Width', gini_importance[3])
```

```
Sepal Length 0.12883574204615036
Sepal Width 0.0027025199853343575
Petal Length 0.43459171536344887
Petal Width 0.4338700226050663
```

In [ ]:

```
In [25]: # Create a selector object that will use the random forest classifier to ident
# features that have an importance of more than 0.15
feature_selector = SelectFromModel(estimator=model, threshold=0.15)

# Train the selector
feature_selector.fit(X_train, y_train)
```

Out[25]: SelectFromModel(estimator=RandomForestClassifier(max\_depth=2), threshold=0.15)

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```
In [26]: # Print the names of the most important features
feature_selector.fit(X_train, y_train)
selected_feature_names = [iris.feature_names[i] for i in feature_selector.get_support()]
print(selected_feature_names)

['petal length (cm)', 'petal width (cm)']
```

```
In [27]: # Transform the data to create a new dataset containing only the most important features
# Note: We have to apply the transform to both the training X and test X data.
X_train_selected = feature_selector.transform(X_train)
X_test_selected = feature_selector.transform(X_test)
```

```
In [28]: # Create a new random forest classifier for the most important features
new_model = RandomForestClassifier(n_estimators=10000, random_state=0, n_jobs=

# Train the new classifier on the new dataset containing the most important fe
new_model.fit(X_train_selected, y_train)
```

Out[28]: RandomForestClassifier(n\_estimators=10000, n\_jobs=-1, random\_state=0)

**In a Jupyter environment, please rerun this cell to show the HTML representation or trust the notebook.**

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```
In [29]: # Apply the limited classifier to the test data
y_pred = model.predict(X_test)

# View The Accuracy Of Our Full Feature (4 Features) Model
accuracy_score(y_test, y_pred)
```

Out[29]: 0.8833333333333333

```
In [30]: y_pred_two = new_model.predict(X_test_selected)
accuracy_selected = accuracy_score(y_test, y_pred_two)
print("Accuracy with selected features:", accuracy_selected)
```

Accuracy with selected features: 0.9

```
In [31]: # Apply The Full Featured Classifier To The Test Data

# View The Accuracy Of Our Full Feature (4 Features) Model
```

```
In [32]: # Apply The Full Featured Classifier To The Test Data

# View The Accuracy Of Our Limited Feature (2 Features) Model
```

## Exercise 2. Now repeat RF feature selection but now use the wine dataset

```
In [36]: from sklearn.ensemble import RandomForestClassifier
data = pd.read_csv("./Wine.csv")
X = data.drop('Customer_Segment', axis = 1)
y = data['Customer_Segment']

X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.4, random_state=42)
model = RandomForestClassifier(criterion="gini", max_depth = 2)

# Train the classifier
model.fit(X_train, y_train)
feature_selector = SelectFromModel(estimator=model, threshold=0.15)

# Train the selector
feature_selector.fit(X_train, y_train)
X_train_selected = feature_selector.transform(X_train)
X_test_selected = feature_selector.transform(X_test)
new_model = RandomForestClassifier(n_estimators=10000, random_state=0, n_jobs=-1)

# Train the new classifier on the new dataset containing the most important features
new_model.fit(X_train_selected, y_train)

y_pred_two = new_model.predict(X_test_selected)
accuracy_selected = accuracy_score(y_test, y_pred_two)
print("Accuracy with selected features:", accuracy_selected)
```

Accuracy with selected features: 0.8888888888888888

## RFE Walkthrough

---



```

In [12]: # Automatically select the features
# automatically select the number of features for RFE
from numpy import mean
from numpy import std
from sklearn.datasets import make_classification
from sklearn.model_selection import cross_val_score
from sklearn.model_selection import RepeatedStratifiedKFold
from sklearn.feature_selection import RFECV
from sklearn.tree import DecisionTreeClassifier
from sklearn.pipeline import Pipeline
# define dataset
X, y = make_classification(n_samples=1000, n_features=10, n_informative=5, n_r

# create pipeline
model = DecisionTreeClassifier()
rfe = RFECV(estimator=model, step=1)
pipeline = Pipeline(steps=[('feature_selection', rfe), ('classifier', model)])

# evaluate model
cv = RepeatedStratifiedKFold(n_splits=10, n_repeats=3, random_state=1)
scores = cross_val_score(pipeline, X, y, scoring='accuracy', cv=cv)

# report performance
print("Accuracy:", mean(scores), (std(scores)))

```

Accuracy: 0.882 0.026255158223353628

```

In [30]: # report which features were selected by RFE
from sklearn.datasets import make_classification
from sklearn.feature_selection import RFE
# define dataset
X, y = make_classification(n_samples=1000, n_features=10, n_informative=5, n_r
# define RFE
rfe = RFE(estimator=DecisionTreeClassifier(), n_features_to_select=5)
# fit RFE
rfe.fit(X, y)
# summarize all features
for i in range(X.shape[1]):
    print('Column: %d, Selected %s, Rank: %.3f' % (i, rfe.support_[i], rfe.ra

```

```

Column: 0, Selected False, Rank: 5.000
Column: 1, Selected False, Rank: 4.000
Column: 2, Selected True, Rank: 1.000
Column: 3, Selected True, Rank: 1.000
Column: 4, Selected True, Rank: 1.000
Column: 5, Selected False, Rank: 6.000
Column: 6, Selected True, Rank: 1.000
Column: 7, Selected False, Rank: 2.000
Column: 8, Selected True, Rank: 1.000
Column: 9, Selected False, Rank: 3.000

```

**Exercise 3. Now repeat RF feature selection but now use the iris dataset**

```
In [38]: #Starter code for Exercise 3
# Load libraries

from sklearn import datasets
import matplotlib.pyplot as plt

# Load digits dataset
iris = datasets.load_iris()

# Create feature matrix
X = iris.data

# Create target vector
y = iris.target

# View the first observation's feature values
iris.data[0]
```

```
Out[38]: array([5.1, 3.5, 1.4, 0.2])
```

```
In [39]: X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.4, random_state=42)
model = RandomForestClassifier(criterion="gini", max_depth = 2)

# Train the classifier
model.fit(X_train, y_train)
feature_selector = SelectFromModel(estimator=model, threshold=0.15)

# Train the selector
feature_selector.fit(X_train, y_train)
X_train_selected = feature_selector.transform(X_train)
X_test_selected = feature_selector.transform(X_test)
new_model = RandomForestClassifier(n_estimators=10000, random_state=0, n_jobs=-1)

# Train the new classifier on the new dataset containing the most important features
new_model.fit(X_train_selected, y_train)

y_pred_two = new_model.predict(X_test_selected)
accuracy_selected = accuracy_score(y_test, y_pred_two)
print("Accuracy with selected features:", accuracy_selected)
```

Accuracy with selected features: 0.9

## t-SNE Walkthrough

(<https://cmdlinetips.com/2019/07/dimensionality-reduction-with-tsne/>)

(<https://cmdlinetips.com/2019/07/dimensionality-reduction-with-tsne/>)

```
In [40]: import matplotlib.pyplot as plt
import seaborn as sns
%matplotlib inline
import pandas as pd
```

```
In [41]: from sklearn.datasets import load_digits
digits = load_digits()
digits
```

```
Out[41]: {'data': array([[ 0.,  0.,  5., ...,  0.,  0.,  0.],
        [ 0.,  0.,  0., ..., 10.,  0.,  0.],
        [ 0.,  0.,  0., ..., 16.,  9.,  0.],
        ...,
        [ 0.,  0.,  1., ...,  6.,  0.,  0.],
        [ 0.,  0.,  2., ..., 12.,  0.,  0.],
        [ 0.,  0., 10., ..., 12.,  1.,  0.])),
  'target': array([0, 1, 2, ..., 8, 9, 8]),
  'frame': None,
  'feature_names': ['pixel_0_0',
    'pixel_0_1',
    'pixel_0_2',
    'pixel_0_3',
    'pixel_0_4',
    'pixel_0_5',
    'pixel_0_6',
    'pixel_0_7',
    'pixel_1_0',
    'pixel_1_1',
    ...]
```

```
In [42]: digits.data.shape
```

```
Out[42]: (1797, 64)
```

```
In [43]: digits.target
```

```
Out[43]: array([0, 1, 2, ..., 8, 9, 8])
```

```
In [45]: data_X = digits.data[:600]
y = digits.target[:600]
```

```
In [17]: from sklearn.manifold import TSNE
tsne = TSNE(n_components=2, random_state=0)
tsne_obj = tsne.fit_transform(data_X)
tsne_obj
```

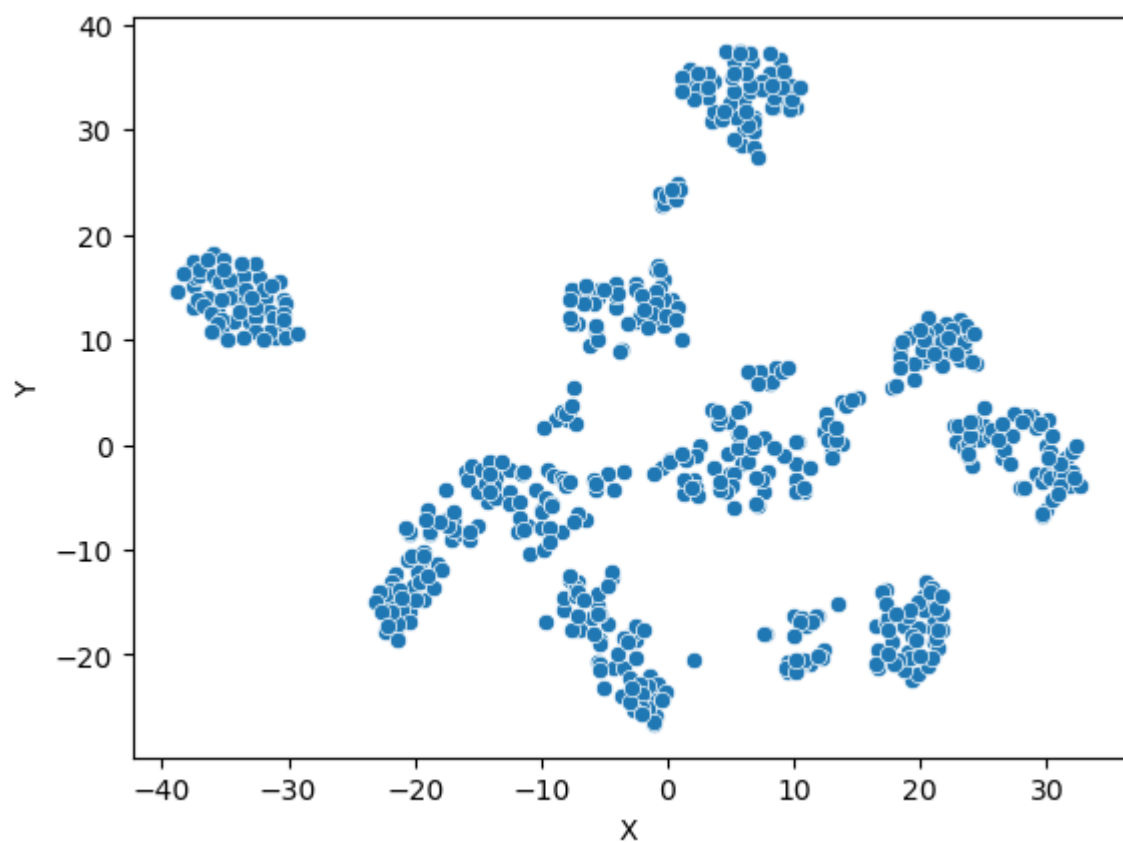
```
Out[17]: array([[ -33.064045 ,  11.775319 ],
        [ 17.76624 ,  5.4796257],
        [  8.055402 ,  5.8756485],
        ...,
        [ 18.094622 , -16.057703 ],
        [  5.6653833,  37.35514 ],
        [ -7.595555 ,  3.655271 ]], dtype=float32)
```

```
In [25]: tsne_df = pd.DataFrame({'X':tsne_obj[:,0], 'Y':tsne_obj[:,1], 'digit':y})  
tsne_df.head()
```

```
Out[25]:
```

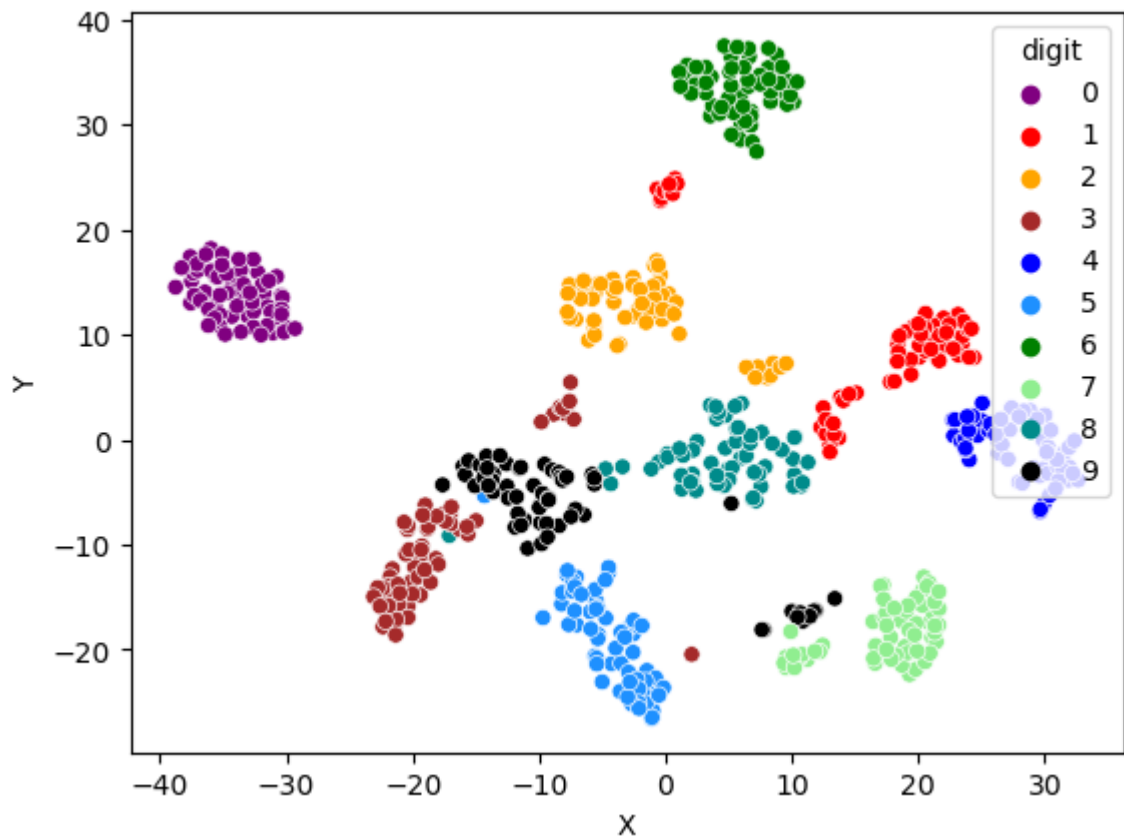
	X	Y	digit
0	-33.064045	11.775319	0
1	17.766239	5.479626	1
2	8.055402	5.875648	2
3	-18.861454	-8.514892	3
4	30.009453	2.323586	4

```
In [20]: import seaborn as sns  
sns.scatterplot(x="X", y="Y", data=tsne_df);
```



```
In [28]: sns.scatterplot(x="X", y="Y", hue="digit",  
                        palette=['purple', 'red', 'orange', 'brown', 'blue', 'dodgerblue', '  
                        legend='full', data=tsne_df)
```

Out[28]: <Axes: xlabel='X', ylabel='Y'>



**Exercise 4. Now repeat t-SNE feature selection but now use the iris dataset**

```
In [47]: from sklearn.datasets import load_iris

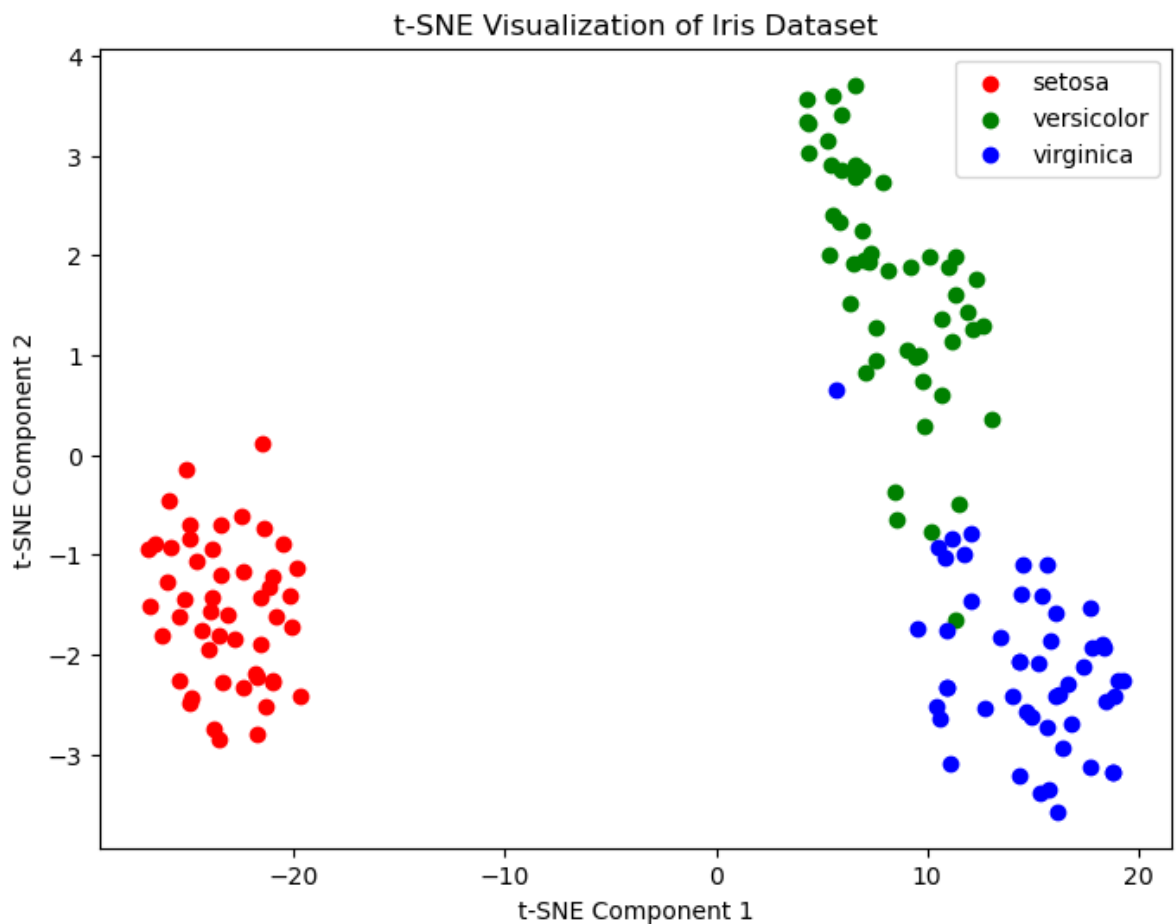
# Load the Iris dataset
iris = load_iris()
X = iris.data
y = iris.target

from sklearn.manifold import TSNE

tsne = TSNE(n_components=2, random_state=42)
X_tsne = tsne.fit_transform(X)

import matplotlib.pyplot as plt
plt.figure(figsize=(8, 6))
colors = ['red', 'green', 'blue']
for i in range(len(colors)):
    plt.scatter(X_tsne[y == i, 0], X_tsne[y == i, 1], c=colors[i], label=iris.

plt.xlabel('t-SNE Component 1')
plt.ylabel('t-SNE Component 2')
plt.title('t-SNE Visualization of Iris Dataset')
plt.legend()
plt.show()
```



In [ ]: