SESSION 5th

K-Nearest Neighbor(KNN) Algorithm for Machine Learning

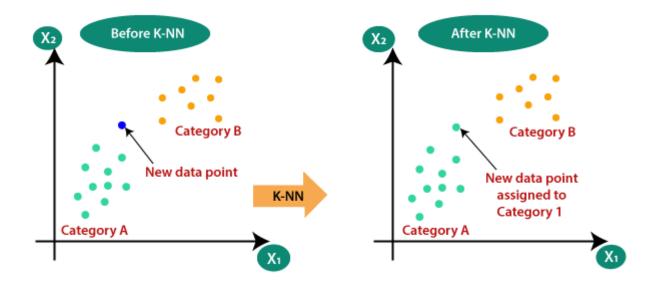
- K-Nearest Neighbour is one of the simplest Machine Learning algorithms based on Supervised Learning technique.
- K-NN algorithm assumes the similarity between the new case/data and available cases and put the new case into the category that is most similar to the available categories.
- K-NN algorithm stores all the available data and classifies a new data point based on the similarity. This means when new data appears then it can be easily classified into a well suite category by using K- NN algorithm.
- K-NN algorithm can be used for Regression as well as for Classification but mostly it is used for the Classification problems.
- K-NN is a non-parametric algorithm, which means it does not make any assumption on underlying data.
- It is also called a **lazy learner algorithm** because it does not learn from the training set immediately instead it stores the dataset and at the time of classification, it performs an action on the dataset.
- KNN algorithm at the training phase just stores the dataset and when it gets new data, then it classifies that data into a category that is much similar to the new data.
- Example: Suppose, we have an image of a creature that looks similar to cat and dog, but we want to know either it is a cat or dog. So for this identification, we can use the KNN algorithm, as it works on a similarity measure. Our KNN model will find the similar features of the new data set to the cats and dogs images and based on the most similar features it will put it in either cat or dog category.

KNN Classifier



Why do we need a K-NN Algorithm?

Suppose there are two categories, i.e., Category A and Category B, and we have a new data point x1, so this data point will lie in which of these categories. To solve this type of problem, we need a K-NN algorithm. With the help of K-NN, we can easily identify the category or class of a particular dataset. Consider the below diagram:



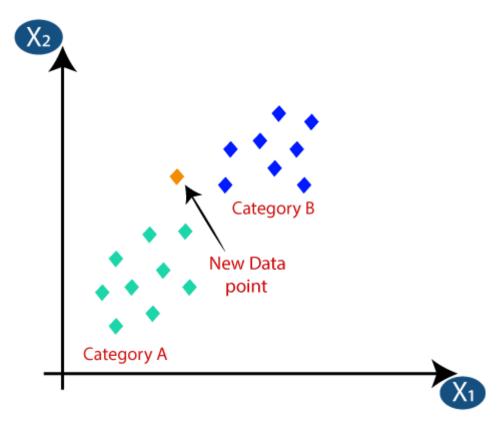
How does K-NN work?

The K-NN working can be explained on the basis of the below algorithm:

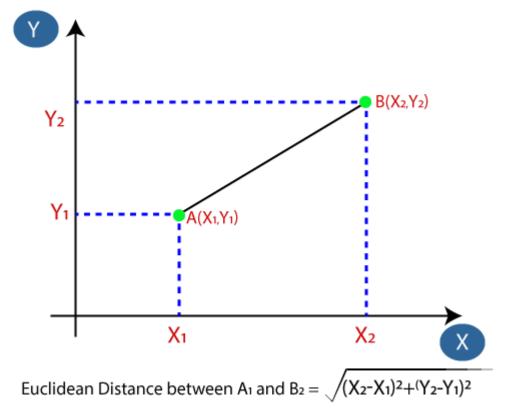
- **Step-1:** Select the number K of the neighbors
- Step-2: Calculate the Euclidean distance of K number of neighbors

- Step-3: Take the K nearest neighbors as per the calculated Euclidean distance.
- Step-4: Among these k neighbors, count the number of the data points in each category.
- Step-5: Assign the new data points to that category for which the number of the neighbor is maximum.
- Step-6: Our model is ready.

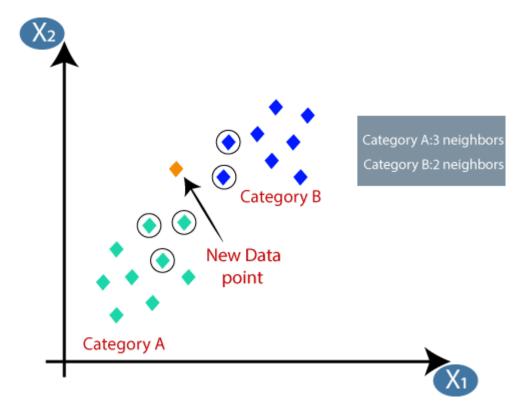
Suppose we have a new data point and we need to put it in the required category. Consider the below image:



- \circ Firstly, we will choose the number of neighbors, so we will choose the k=5.
- Next, we will calculate the **Euclidean distance** between the data points. The Euclidean distance is the distance between two points, which we have already studied in geometry. It can be calculated as:



 By calculating the Euclidean distance we got the nearest neighbors, as three nearest neighbors in category A and two nearest neighbors in category B.
 Consider the below image:



 As we can see the 3 nearest neighbors are from category A, hence this new data point must belong to category A.

How to select the value of K in the K-NN Algorithm?

Below are some points to remember while selecting the value of K in the K-NN algorithm:

- There is no particular way to determine the best value for "K", so we need to try some values to find the best out of them. The most preferred value for K is 5.
- A very low value for K such as K=1 or K=2, can be noisy and lead to the effects of outliers in the model.
- Large values for K are good, but it may find some difficulties.

Advantages of KNN Algorithm:

- o It is simple to implement.
- o It is robust to the noisy training data
- o It can be more effective if the training data is large.

Disadvantages of KNN Algorithm:

- o Always needs to determine the value of K which may be complex some time.
- The computation cost is high because of calculating the distance between the data points for all the training samples.

Python implementation of the KNN algorithm

To do the Python implementation of the K-NN algorithm, we will use the same problem and dataset which we have used in Logistic Regression. But here we will improve the performance of the model. Below is the problem description:

Problem for K-NN Algorithm: There is a Car manufacturer company that has manufactured a new SUV car. The company wants to give the ads to the users who are interested in buying that SUV. So for this problem, we have a dataset that contains multiple user's information through the social network. The dataset contains lots of information but the **Estimated Salary** and **Age** we will consider for the independent variable and the **Purchased variable** is for the dependent variable. Below is the dataset:

User ID	Gender	Age	EstimatedSalary	Purchased
15624510	Male	19	19000	0
15810944	Male	35	20000	0
15668575	Female	26	43000	0
15603246	Female	27	57000	0
15804002	Male	19	76000	0
15728773	Male	27	58000	0
15598044	Female	27	84000	0
15694829	Female	32	150000	1
15600575	Male	25	33000	0
15727311	Female	35	65000	0
15570769	Female	26	80000	0
15606274	Female	26	52000	0
15746139	Male	20	86000	0
15704987	Male	32	18000	0
15628972	Male	18	82000	0
15697686	Male	29	80000	0
15733883	Male	47	25000	1
15617482	Male	45	26000	1
15704583	Male	46	28000	1
15621083	Female	48	29000	1
15649487	Male	45	22000	1
15736760	Female	47	49000	1

Steps to implement the K-NN algorithm:

- Data Pre-processing step
- o Fitting the K-NN algorithm to the Training set
- o Predicting the test result
- Test accuracy of the result(Creation of Confusion matrix)
- Visualizing the test set result.

Data Pre-Processing Step:

The Data Pre-processing step will remain exactly the same as Logistic Regression. Below is the code for it:

- 1. # importing libraries
- 2. **import** numpy as nm
- 3. **import** matplotlib.pyplot as mtp
- 4. import pandas as pd
- 5. #importing datasets
- 6. data_set= pd.read_csv('user_data.csv')

- 7. #Extracting Independent and dependent Variable
- 8. $x = data_set.iloc[:, [2,3]].values$
- 9. y= data_set.iloc[:, 4].values

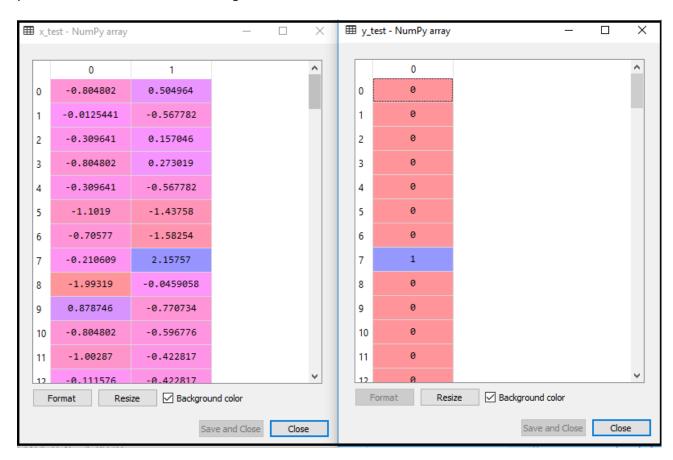
10.

- 11.# Splitting the dataset into training and test set.
- 12.from sklearn.model_selection import train_test_split
- 13.x_train, x_test, y_train, y_test= train_test_split(x, y, test_size= 0.25, random_state =0)

14.

- 15. #feature Scaling
- 16.from sklearn.preprocessing **import** StandardScaler
- 17.st_x= StandardScaler()
- 18.x_train = st_x.fit_transform(x_train)
- 19.x_test= st_x.transform(x_test)

By executing the above code, our dataset is imported to our program and well preprocessed. After feature scaling our test dataset will look like:



From the above output image, we can see that our data is successfully scaled.

Fitting K-NN classifier to the Training data:

Now we will fit the K-NN classifier to the training data. To do this we will import the **KNeighborsClassifier** class of **Sklearn Neighbors** library. After importing the class, we will create the **Classifier** object of the class. The Parameter of this class will be

- n_neighbors: To define the required neighbors of the algorithm.
 Usually, it takes 5.
- metric='minkowski': This is the default parameter and it decides the distance between the points.
- o **p=2:** It is equivalent to the standard Euclidean metric.

And then we will fit the classifier to the training data. Below is the code for it:

- 1. #Fitting K-NN classifier to the training set
- 2. from sklearn.neighbors import KNeighborsClassifier
- 3. classifier= KNeighborsClassifier(n_neighbors=5, metric='minkowski', p=2)
- classifier.fit(x_train, y_train)

Output: By executing the above code, we will get the output as:

- Predicting the Test Result: To predict the test set result, we will create
 a y_pred vector as we did in Logistic Regression. Below is the code for it:
- 1. #Predicting the test set result
- 2. y_pred= classifier.predict(x_test)

Output:

The output for the above code will be:



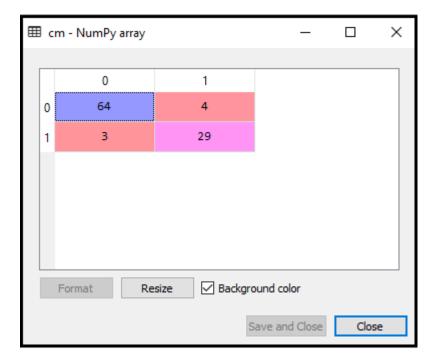
Creating the Confusion Matrix:

Now we will create the Confusion Matrix for our K-NN model to see the accuracy of the classifier. Below is the code for it:

- 1. #Creating the Confusion matrix
- 2. from sklearn.metrics **import** confusion_matrix
- 3. cm= confusion_matrix(y_test, y_pred)

In above code, we have imported the confusion_matrix function and called it using the variable cm.

Output: By executing the above code, we will get the matrix as below:



In the above image, we can see there are 64+29= 93 correct predictions and 3+4= 7 incorrect predictions, whereas, in Logistic Regression, there were 11 incorrect predictions. So we can say that the performance of the model is improved by using the K-NN algorithm.

o Visualizing the Training set result:

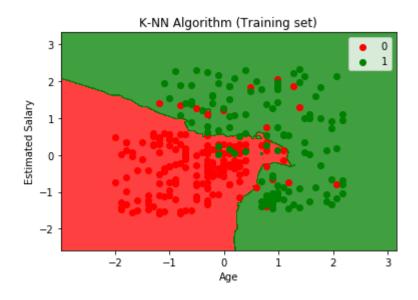
Now, we will visualize the training set result for K-NN model. The code will remain same as we did in Logistic Regression, except the name of the graph. Below is the code for it:

- 1. #Visulaizing the trianing set result
- 2. from matplotlib.colors **import** ListedColormap
- 3. x_{set} , $y_{set} = x_{train}$, y_{train}
- 4. x1, $x2 = nm.meshgrid(nm.arange(start = x_set[:, 0].min() 1, stop = x_set[:, 0]. max() + 1, step = 0.01),$
- 5. $nm.arange(start = x_set[:, 1].min() 1, stop = x_set[:, 1].max() + 1, step = 0.01)$
- mtp.contourf(x1, x2, classifier.predict(nm.array([x1.ravel(), x2.ravel()]).T).reshape(x1.shape),
- 7. alpha = 0.75, cmap = ListedColormap(('red','green')))
- 8. mtp.xlim(x1.min(), x1.max())
- 9. mtp.ylim(x2.min(), x2.max())
- 10.for i, j in enumerate(nm.unique(y_set)):
- 11.mtp.scatter($x_{set}[y_{set} == j, 0], x_{set}[y_{set} == j, 1],$

```
12.c = ListedColormap(('red', 'green'))(i), label = j)
13.mtp.title('K-NN Algorithm (Training set)')
14.mtp.xlabel('Age')
15.mtp.ylabel('Estimated Salary')
16.mtp.legend()
```

17.mtp.show() Output:

By executing the above code, we will get the below graph:



The output graph is different from the graph which we have occurred in Logistic Regression. It can be understood in the below points:

- As we can see the graph is showing the red point and green points.
 The green points are for Purchased(1) and Red Points for not
 Purchased(0) variable.
- The graph is showing an irregular boundary instead of showing any straight line or any curve because it is a K-NN algorithm, i.e., finding the nearest neighbor.
- The graph has classified users in the correct categories as most of the users who didn't buy the SUV are in the red region and users who bought the SUV are in the green region.
- The graph is showing good result but still, there are some green points in the red region and red points in the green region. But this is no big issue as by doing this model is prevented from overfitting issues.
- Hence our model is well trained.

Visualizing the Test set result:

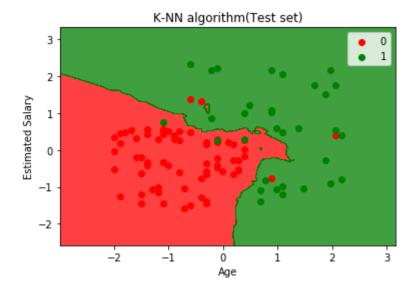
After the training of the model, we will now test the result by putting a new dataset, i.e., Test dataset. Code remains the same except some minor changes: such as **x_train and y_train** will be replaced by **x_test and y_test**.

Below is the code for it:

```
1. #Visualizing the test set result
2. from matplotlib.colors import ListedColormap
3. x_{set}, y_{set} = x_{test}, y_{test}
4. x1, x2 = nm.meshgrid(nm.arange(start = x_set[:, 0].min() - 1, stop = x_set[:, 0].
   max() + 1, step = 0.01),
5. nm.arange(start = x_set[:, 1].min() - 1, stop = x_set[:, 1].max() + 1, step = 0.01)
   )
6. mtp.contourf(x1, x2, classifier.predict(nm.array([x1.ravel(), x2.ravel()]).T).reshape(
   x1.shape),
7. alpha = 0.75, cmap = ListedColormap(('red','green')))
8. mtp.xlim(x1.min(), x1.max())
9. mtp.ylim(x2.min(), x2.max())
10.for i, j in enumerate(nm.unique(y_set)):
11.
      mtp.scatter(x_set[y_set == j, 0], x_set[y_set == j, 1],
12.
        c = ListedColormap(('red', 'green'))(i), label = j)
13.mtp.title('K-NN algorithm(Test set)')
14.mtp.xlabel('Age')
15.mtp.ylabel('Estimated Salary')
16.mtp.legend()
```

Output:

17.mtp.show()



The above graph is showing the output for the test data set. As we can see in the graph, the predicted output is well good as most of the red points are in the red region and most of the green points are in the green region.

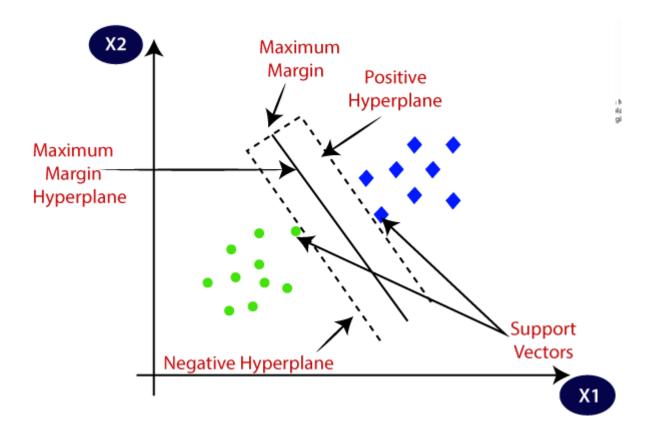
However, there are few green points in the red region and a few red points in the green region. So these are the incorrect observations that we have observed in the confusion matrix(7 Incorrect output).

Support Vector Machine Algorithm

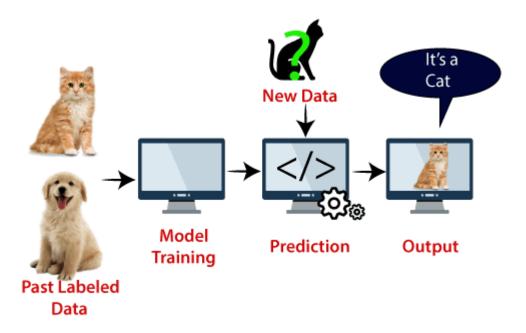
Support Vector Machine or SVM is one of the most popular Supervised Learning algorithms, which is used for Classification as well as Regression problems. However, primarily, it is used for Classification problems in Machine Learning.

The goal of the SVM algorithm is to create the best line or decision boundary that can segregate n-dimensional space into classes so that we can easily put the new data point in the correct category in the future. This best decision boundary is called a hyperplane.

SVM chooses the extreme points/vectors that help in creating the hyperplane. These extreme cases are called as support vectors, and hence algorithm is termed as Support Vector Machine. Consider the below diagram in which there are two different categories that are classified using a decision boundary or hyperplane:



Example: SVM can be understood with the example that we have used in the KNN classifier. Suppose we see a strange cat that also has some features of dogs, so if we want a model that can accurately identify whether it is a cat or dog, so such a model can be created by using the SVM algorithm. We will first train our model with lots of images of cats and dogs so that it can learn about different features of cats and dogs, and then we test it with this strange creature. So as support vector creates a decision boundary between these two data (cat and dog) and choose extreme cases (support vectors), it will see the extreme case of cat and dog. On the basis of the support vectors, it will classify it as a cat. Consider the below diagram:



SVM algorithm can be used for **Face detection**, **image classification**, **text categorization**, etc.

Types of SVM

SVM can be of two types:

- Linear SVM: Linear SVM is used for linearly separable data, which means if a
 dataset can be classified into two classes by using a single straight line, then
 such data is termed as linearly separable data, and classifier is used called as
 Linear SVM classifier.
- Non-linear SVM: Non-Linear SVM is used for non-linearly separated data, which means if a dataset cannot be classified by using a straight line, then such data is termed as non-linear data and classifier used is called as Non-linear SVM classifier.

Hyperplane and Support Vectors in the SVM algorithm:

Hyperplane: There can be multiple lines/decision boundaries to segregate the classes in n-dimensional space, but we need to find out the best decision boundary that helps to classify the data points. This best boundary is known as the hyperplane of SVM.

The dimensions of the hyperplane depend on the features present in the dataset, which means if there are 2 features (as shown in image), then hyperplane will be a straight line. And if there are 3 features, then hyperplane will be a 2-dimension plane.

We always create a hyperplane that has a maximum margin, which means the maximum distance between the data points.

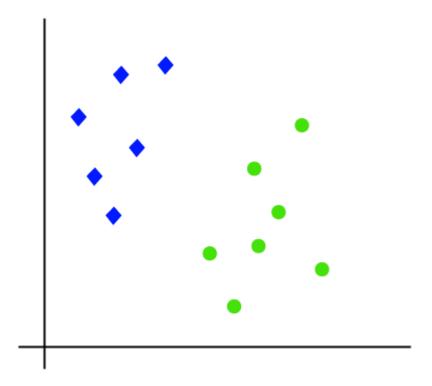
Support Vectors:

The data points or vectors that are the closest to the hyperplane and which affect the position of the hyperplane are termed as Support Vector. Since these vectors support the hyperplane, hence called a Support vector.

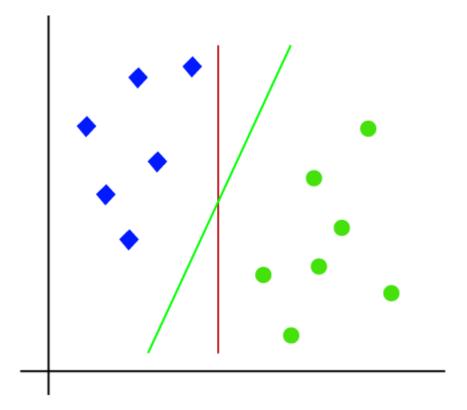
How does SVM works?

Linear SVM:

The working of the SVM algorithm can be understood by using an example. Suppose we have a dataset that has two tags (green and blue), and the dataset has two features x1 and x2. We want a classifier that can classify the pair(x1, x2) of coordinates in either green or blue. Consider the below image:

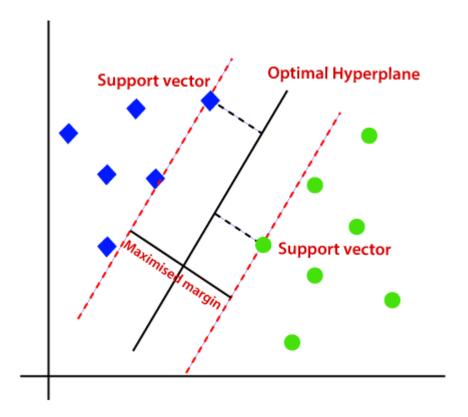


So as it is 2-d space so by just using a straight line, we can easily separate these two classes. But there can be multiple lines that can separate these classes. Consider the below image:



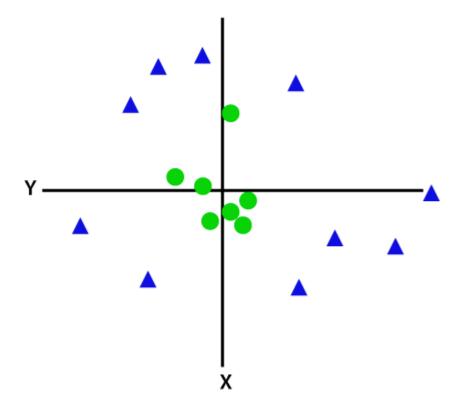
Hence, the SVM algorithm helps to find the best line or decision boundary; this best boundary or region is called as a **hyperplane**. SVM algorithm finds the closest point of the lines from both the classes. These points are called support vectors. The

distance between the vectors and the hyperplane is called as **margin**. And the goal of SVM is to maximize this margin. The **hyperplane** with maximum margin is called the **optimal hyperplane**.



Non-Linear SVM:

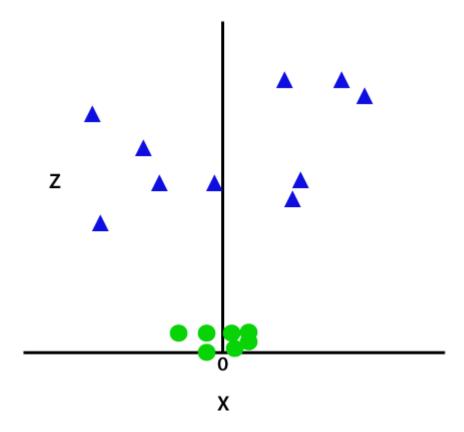
If data is linearly arranged, then we can separate it by using a straight line, but for non-linear data, we cannot draw a single straight line. Consider the below image:



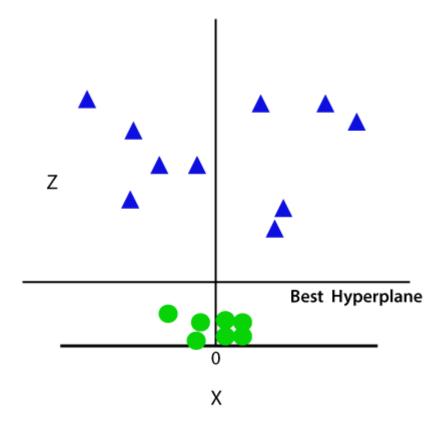
So to separate these data points, we need to add one more dimension. For linear data, we have used two dimensions x and y, so for non-linear data, we will add a third dimension z. It can be calculated as:

$$z=x^{2} + y^{2}$$

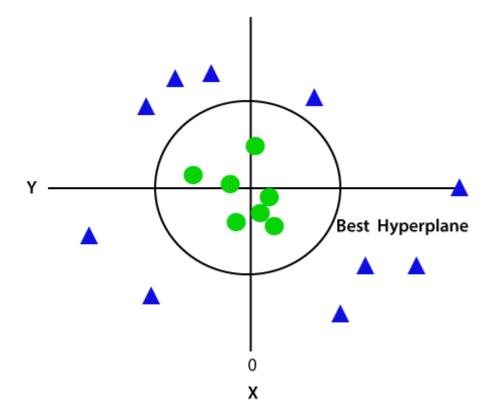
By adding the third dimension, the sample space will become as below image:



So now, SVM will divide the datasets into classes in the following way. Consider the below image:



Since we are in 3-d Space, hence it is looking like a plane parallel to the x-axis. If we convert it in 2d space with z=1, then it will become as:



Hence we get a circumference of radius 1 in case of non-linear data.

Python Implementation of Support Vector Machine

Now we will implement the SVM algorithm using Python. Here we will use the same dataset **user_data**, which we have used in Logistic regression and KNN classification.

Data Pre-processing step

Till the Data pre-processing step, the code will remain the same. Below is the code:

- 1. #Data Pre-processing Step
- 2. # importing libraries
- 3. **import** numpy as nm
- 4. **import** matplotlib.pyplot as mtp
- 5. import pandas as pd

6.

- 7. #importing datasets
- 8. data_set= pd.read_csv('user_data.csv')

9.

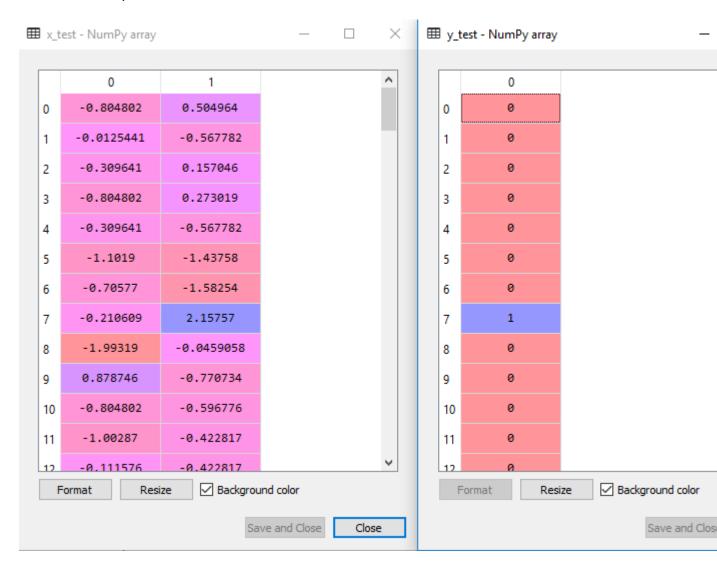
- 10. #Extracting Independent and dependent Variable
- 11.x= data_set.iloc[:, [2,3]].values

- 12.y= data_set.iloc[:, 4].values
- 13.
- 14.# Splitting the dataset into training and test set.
- 15.from sklearn.model_selection import train_test_split
- 16.x_train, x_test, y_train, y_test= train_test_split(x, y, test_size= 0.25, random_state =0)
- 17. #feature Scaling
- 18.from sklearn.preprocessing **import** StandardScaler
- 19.st_x= StandardScaler()
- 20.x_train= st_x.fit_transform(x_train)
- 21.x_test= st_x.transform(x_test)

After executing the above code, we will pre-process the data. The code will give the dataset as:

Index	User ID	Gender	Age	EstimatedSalary	Purchased	
0	15624510	Male	19	19000	0	
1	15810944	Male	35	20000	0	
2	15668575	Female	26	43000	0	
3	15603246	Female	27	57000	0	
4	15804002	Male	19	76000	0	
5	15728773	Male	27	58000	0	
6	15598044	Female	27	84000	0	
7	15694829	Female	32	150000	1	
8	15600575	Male	25	33000	0	
9	15727311	Female	35	65000	0	
10	15570769	Female	26	80000	0	
11	15606274	Female	26	52000	0	
12	15746139	Male	20	86000	0	
13	15704987	Male	32	18000	0	
14	15628972	Male	18	82000	0	

The scaled output for the test set will be:



Fitting the SVM classifier to the training set:

Now the training set will be fitted to the SVM classifier. To create the SVM classifier, we will import **SVC** class from **Sklearn.svm** library. Below is the code for it:

- 1. from sklearn.svm **import** SVC # "Support vector classifier"
- classifier = SVC(kernel='linear', random_state=0)
- 3. classifier.fit(x_train, y_train)

In the above code, we have used **kernel='linear'**, as here we are creating SVM for linearly separable data. However, we can change it for non-linear data. And then we fitted the classifier to the training dataset(x_train, y_train)

Output:

```
Out[8]:
SVC(C=1.0, cache_size=200, class_weight=None, coef0=0.0,
decision_function_shape='ovr', degree=3, gamma='auto_deprecated',
```

The model performance can be altered by changing the value of **C(Regularization factor)**, **gamma**, **and kernel**.

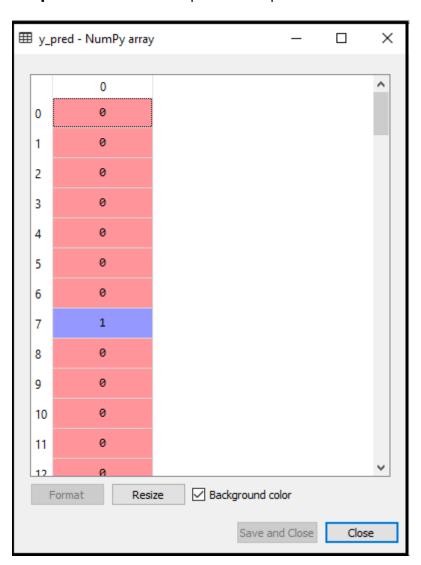
Predicting the test set result:

Now, we will predict the output for test set. For this, we will create a new vector y_pred. Below is the code for it:

- 1. #Predicting the test set result
- 2. y_pred= classifier.predict(x_test)

After getting the y_pred vector, we can compare the result of **y_pred** and **y_test** to check the difference between the actual value and predicted value.

Output: Below is the output for the prediction of the test set:

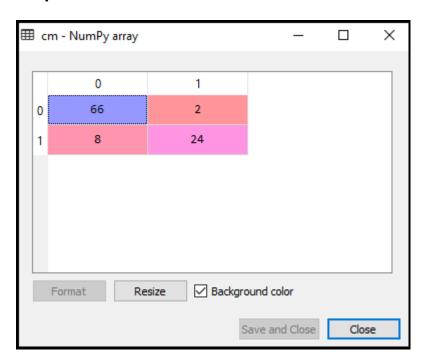


Creating the confusion matrix:

Now we will see the performance of the SVM classifier that how many incorrect predictions are there as compared to the Logistic regression classifier. To create the confusion matrix, we need to import the **confusion_matrix** function of the sklearn library. After importing the function, we will call it using a new variable **cm**. The function takes two parameters, mainly **y_true**(the actual values) and **y_pred** (the targeted value return by the classifier). Below is the code for it:

- 1. #Creating the Confusion matrix
- 2. from sklearn.metrics **import** confusion_matrix
- cm= confusion_matrix(y_test, y_pred)

Output:



As we can see in the above output image, there are 66+24=90 correct predictions and 8+2=10 correct predictions. Therefore we can say that our SVM model improved as compared to the Logistic regression model.

Visualizing the training set result:

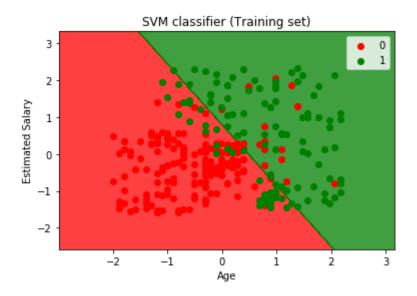
Now we will visualize the training set result, below is the code for it:

- 1. from matplotlib.colors **import** ListedColormap
- 2. x_set, y_set = x_train, y_train

```
3. x1, x2 = nm.meshgrid(nm.arange(start = x_set[:, 0].min() - 1, stop = x_set[:, 0].
   max() + 1, step = 0.01),
4. nm.arange(start = x_set[:, 1].min() - 1, stop = x_set[:, 1].max() + 1, step = 0.01)
   )
5. mtp.contourf(x1, x2, classifier.predict(nm.array([x1.ravel(), x2.ravel()]).T).reshape(
   x1.shape),
6. alpha = 0.75, cmap = ListedColormap(('red', 'green')))
7. mtp.xlim(x1.min(), x1.max())
8. mtp.ylim(x2.min(), x2.max())
9. for i, j in enumerate(nm.unique(y_set)):
      mtp.scatter(x_set[y_set == j, 0], x_set[y_set == j, 1],
10.
11.
        c = ListedColormap(('red', 'green'))(i), label = j)
12.mtp.title('SVM classifier (Training set)')
13.mtp.xlabel('Age')
14.mtp.ylabel('Estimated Salary')
15.mtp.legend()
16.mtp.show()
```

Output:

By executing the above code, we will get the output as:



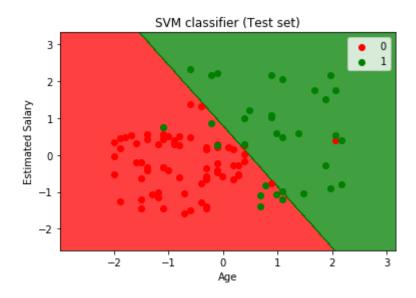
As we can see, the above output is appearing similar to the Logistic regression output. In the output, we got the straight line as hyperplane because we have **used a linear kernel in the classifier**. And we have also discussed above that for the 2d space, the hyperplane in SVM is a straight line.

o Visualizing the test set result:

- 1. #Visulaizing the test set result
- 2. from matplotlib.colors **import** ListedColormap
- 3. x_{set} , $y_{set} = x_{test}$, y_{test}
- 4. x1, $x2 = nm.meshgrid(nm.arange(start = x_set[:, 0].min() 1, stop = x_set[:, 0]. max() + 1, step = 0.01),$
- 5. $nm.arange(start = x_set[:, 1].min() 1, stop = x_set[:, 1].max() + 1, step = 0.01)$
- mtp.contourf(x1, x2, classifier.predict(nm.array([x1.ravel(), x2.ravel()]).T).reshape(x1.shape),
- 7. alpha = 0.75, cmap = ListedColormap(('red', 'green')))
- 8. mtp.xlim(x1.min(), x1.max())
- 9. mtp.ylim(x2.min(), x2.max())
- 10.**for** i, j in enumerate(nm.unique(y_set)):
- 11. mtp.scatter($x_set[y_set == j, 0]$, $x_set[y_set == j, 1]$,
- 12. c = ListedColormap(('red', 'green'))(i), label = j)
- 13.mtp.title('SVM classifier (Test set)')
- 14.mtp.xlabel('Age')
- 15.mtp.ylabel('Estimated Salary')
- 16.mtp.legend()
- 17.mtp.show()

Output:

By executing the above code, we will get the output as:



As we can see in the above output image, the SVM classifier has divided the users into two regions (Purchased or Not purchased). Users who purchased the SUV are in

the red region with the red scatter points. And users who did not purchase the SUV are in the green region with green scatter points. The hyperplane has divided the two classes into Purchased and not purchased variable.