**Lab: Data Mining** 

Aim: To have a practical knowledge on the important concept of Data Mining

**Requirements: Google Colab** 

## 1. Clustering:

a. Install and Import libraries:

```
pip install palmerpenguins # (if it isn't working, use the next 2 line)
#import sys
#!{sys.executable} -m pip install palmerpenguins

from palmerpenguins import load_penguins # For penguins dataset
import pandas as pd # For dataframes
import matplotlib.pyplot as plt # For plotting functions
import seaborn as sns # For additional plotting functions
from sklearn.cluster import KMeans # For k-Means
from sklearn.model_selection import GridSearchCV # For grid search
from sklearn.metrics import silhouette_score # For metrics and scores
from sklearn.preprocessing import StandardScaler # For standardizing data
```

- b. Load and prepare data: Following steps are used to prepare the data:
  - i. Load the 'penguins' dataset in variable 'df'
  - ii. Remove the 'island', 'year', and 'sex' variables
  - iii. Rename the class variable 'species' as 'y'
  - iv. Drop all rows with 'NaN'
  - v. Display the first 5 rows of 'df'

c. Explore the data: Visualize various aspects of penguins dataset.

```
i. Bar Plot of Class Variable: sns.countplot(x='y', data=df)
```

ii. Scatter Plots and Density Plots for Feature Pairs:

```
# Creates a grid using Seaborn's PairGrid()
g = sns.PairGrid(
    df,
    vars=['bill_length_mm', 'bill_depth_mm', 'flipper_length_mm', 'body_mass_g']
,
    hue='y',
    diag_sharey=False,
    palette=["red", "green", "blue"])

# Adds histograms on the diagonal
g.map_diag(plt.hist)
# Adds density plots above the diagonal
g.map_upper(sns.kdeplot)
# Adds scatterplots below the diagonal
g.map_lower(sns.scatterplot)
# Adds a legend
```

```
g.add legend()
```

#### d. Prepare data:

#### e. k-MEANS: Train the Model

We'll set up a KMeans object with the following parameters:

- n clusters: Total number of clusters to make.
- random state: Set to one to reproduce these results.
- init: How to initialize the k-means centers; we'll use k-means++.
- n\_init: Number of times k-means would be run; the model returned would have the minimum value of inertia.

A few other attributes of the KMeans object:

- cluster centers\_: Stores the discovered cluster centers.
- labels : Label of each instance.
- inertia: Sum of square of distances of each instance from its corresponding center.
- n iter: Number of iterations run to find the centers.

```
# Sets up the kMeans object
km = KMeans(
    n_clusters=3,
    random_state=1,
    init='k-means++',
    n_init=10)

# Fits the model to the data
km.fit(df)

# Displays the parameters of the fitted model
km.get_params()
```

#### f. k-Means: Visualize the Clusters

```
# Creates a scatter plot
sns.scatterplot(
    x='bill_length_mm',
    y='bill_depth_mm',
    data=df,
    hue=y,
    style=km.labels_,
    palette=["orange", "green", "blue"])
```

```
# Adds cluster centers to the same plot
plt.scatter(
    km.cluster_centers_[:,0],
    km.cluster_centers_[:,1],
    marker='x',
    s=200,
    c='red')
```

## g. k-means: optimize via silhouette scores

The main challenge in k-means is to find the optimal number of clusters. We can set up a <code>GridSearchCV</code> object to search for the optimal parameters. For k-Mmeans, we require a custom scorer that computes the silhouette value for different number of clusters specified by n\_clusters. The custom scorer is called s2() in the code below, where it uses silhouette\_score() from the <code>sklearn.metrics</code> library to compute a score for an instance X.

A silhouette score is a value in [-1,+1]. It is a means for comparing how similar an instance is to its corresponding cluster compared to its similarity with other clusters. Formally, it takes into account cohesion and separation to compute a silhouette value. A +1 or close to this score value indicates better clusters.

```
# Sets up the custom scorer
def s2(estimator, X):
    return silhouette score(X, estimator.predict(X))
# List of values for the parameter `n_clusters`
param = range(2,10)
# KMeans object
km = KMeans(random state=0, init='k-means++')
# Sets up GridSearchCV object and stores in grid variable
grid = GridSearchCV(
    km,
    {'n clusters': param},
    scoring=s2,
    cv=2)
# Fits the grid object to data
grid.fit(df)
# Accesses the optimum model
best_km = grid.best_estimator_
# Displays the optimum model
best km.get params()
```

### h. Plot of Scores for Different Number of Clusters

```
# Plot mean_test_scores vs. n_clusters
plt.plot(
    param,
    grid.cv_results_['mean_test_score'])
# Draw a vertical line, where the best model is
plt.axvline(
    x=best_km.n_clusters,
    color='red',
    ls='--')
# Adds labels to the plot
```

```
plt.xlabel('Total Centers')
              plt.ylabel('Silhouette Score')
       i. Visualize the Best Model
           # Creates a scatter plot
           sns.scatterplot(
               x='bill length mm',
               y='bill depth mm',
               data=df,
               hue=y,
               style=best km.labels ,
               palette=['orange', 'green', 'blue'])
           # Adds cluster centers to the same plot
          plt.scatter(
               best km.cluster centers [:, 0],
               best km.cluster centers [:, 1],
               marker='x',
               s=200,
               c='red')
2. Classification
       a. Import libraries
           import pandas as pd
                                                                      # For dataframes
           import matplotlib.pyplot as plt
                                                                      # For plotting data
           import seaborn as sns
                                                                     # For plotting data
           from sklearn.model selection import train test split # For train/test splits
           from sklearn.model selection import GridSearchCV
                                                                    # For parameter optimization
           from sklearn.neighbors import KNeighborsClassifier
                                                                    # For kNN classification
           from sklearn.metrics import plot confusion matrix
                                                                    # Evaluation measure
       b. Load and prepare data
               i. Import Data
          df = pd.read csv('https://archive.ics.uci.edu/ml/machine-learning-
databases/spambase/spambase.data', header=None)
          df.head()
               ii. Rename Variables
                  • Assign a name to all attributes as X0, X1, ..., X56.
                  • Assign y to the class variable (the last column of df).
                  • Display the first 5 rows.
```

```
# Sequentially renames all attribute columns and renames the last column to 'y'
df.columns = ['X' + str(i) for i in range(0, len(df.columns) - 1)] + ['y']
# Shows the first few lines of the data
df.head()
iii. Split Data
```

- To prepare the dataset for classification, we have to split it into train and test sets.
- train test split() splits the data into train and test.
- In the arguments list, the data matrix consists of all attribute columns. Extract columns X0, X1, ..., X56 with df.filter(regex='\d'). The filter keeps only the names that have a numeric character in them.
- Specify the target variable as df.y.
- Set up trn and tst dataframes.

```
# Specifies X by filtering all columns with a number in name
          X_trn, X_tst, y_trn, y_tst = train_test_split(
               df.filter(regex='\d'),
               df.y,
              test size=0.30,
               random state=1)
           # Creates the training dataset, trn
          trn = X trn
          trn['y'] = y trn
           # Creates the testing dataset, tst
          tst = X tst
          tst['y'] = y tst
      iv. Explore training data

    Bar Plot of Class Variable

          sns.countplot(x='y', data=trn)
             Explore Attribute Variables: Select four arbitrary features and get paired plots
              # Creates a grid using Seaborn's PairGrid()
              g = sns.PairGrid(
                  trn,
                  vars=['X5', 'X20', 'X25', 'X53'],
                  hue='y',
                  diag sharey=False,
                  palette=['red', 'green'])
              # Adds histograms on the diagonal
              g.map diag(plt.hist)
              # Adds density plots above the diagonal
              g.map upper(sns.kdeplot)
              # Adds scatterplots below the diagonal
              g.map lower(sns.scatterplot)
              # Adds a legend
              g.add legend(title='Spam')
       v. Prepare data
              \# Separates the attributes X0-X56 into X trn
              X trn = trn.filter(regex='\d')
              # Separates the class variable into y_trn
              y trn = trn.y
              \# Separates the attributes X0-X56 into X tst
              X tst = tst.filter(regex='\d')
              # Separates the class variable into y tst
              y tst = tst.y
              # Class labels
              spam = ['Not Spam', 'Spam']
              trn.head()
c. kNN: train model
       # Sets up a kNN model and fits it to data
       knn = KNeighborsClassifier(n neighbors=5) \
```

.fit(X\_trn, y\_trn)

d. Mean Accuracy on Training Data

```
print('Accuracy on training data: ' + str("{:.2%}".format(knn.score(X trn, y trn))))
```

e. Optimize the kNN Model: The challenge in training a kNN model is to determine **the optimal number of neighbors.** To find the optimal parameters, GridSearchCV object can be used.

```
# Sets up the kNN classifier object
   knn = KNeighborsClassifier()
   # Search parameters
   param = range(3, 15, 2)
   # Sets up GridSearchCV object and stores it in grid variable
   grid = GridSearchCV(
       knn,{'n neighbors': param})
   # Fits the grid object and gets the best model
   best knn = grid \
       .fit(X_trn,y_trn) \
       .best estimator
   # Displays the optimum model
   best knn.get params()
f. Plot the Accuracy by Neighbors Parameter
       # Plots mean test scores vs. total neighbors
      plt.plot(
           param,
           grid.cv_results_['mean_test_score'])
       # Adds labels to the plot
      plt.xticks(param)
      plt.ylabel('Mean CV Score')
```

g. Test model: we'll evaluate the accuracy of the trained kNN model on the test set. A good evaluation measure is the confusion matrix that gives the fraction of true positives, true negatives, false positives, and false negatives.

```
plot_confusion_matrix(
    best_knn, X_tst, y_tst,
    display_labels=spam,
    normalize='true')
```

plt.xlabel('n\_neighbors')

x=best\_knn.n\_neighbors,

plt.axvline(

color='red',
ls='--')

h. Calculate Mean Accuracy on Testing Data

```
print( 'Accuracy on testing data: '
+ str("{:.2%}".format(best_knn.score(X_tst, y_tst))))
```

# Draws a vertical line where the best model is

# Your Turn!

- Do all the step that we discussed above for Clustering and Classification
  - **1.** Use Iris flower data and cluster them using K-Means
    - Download the Iris flower data: <a href="https://archive.ics.uci.edu/ml/datasets/iris">https://archive.ics.uci.edu/ml/datasets/iris</a>
  - 2. Use the breast cancer data and use the KNN determine whether they are benign or malignant.
    - Download the breast cancer data:
       <a href="https://archive.ics.uci.edu/ml/datasets/breast+cancer+wisconsin+(original)">https://archive.ics.uci.edu/ml/datasets/breast+cancer+wisconsin+(original)</a>