Assessment

Machine Learning and Data Mining



MSc Data Science

Exploring Insights: Classification, Clustering, Text Classification, and Sentiment Analysis

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Part One: Classification Algorithms

Title: Identify if a mushroom is edible to eat or poisonous

1. Introduction

It is essential to distinguish between toxic and edible mushrooms in order to protect those who collect mushrooms for culinary use. Studying this dataset is driven by real-world issues around the gathering and eating of wild mushrooms. A great case study for assessing how well machine learning algorithms perform in categorization tasks is the Mushroom dataset. Based on observable traits, algorithms trained on this dataset can be used to automatically determine if mushrooms are edible.

In the field of machine learning and data mining, the Mushroom dataset is a well-known dataset that is frequently used for classification tasks. This dataset, which was initially taken from the UCI Machine Learning Repository, offers a thorough assortment of features that describe different aspects of mushrooms. Developing classification models that can correctly distinguish between edible and deadly mushrooms based on their observable properties is the main goal of using this dataset.

2. Research Objective

"Can machine learning algorithms effectively predict mushroom edibility

based on observable traits, and if yes, which classification model

performs best in discriminating between edible and dangerous mushrooms?"

We use at least two classification algorithms and perform a machine learning classification in both Python as well as Microsoft Azure to achieve the above objective. Finally, the results from algorithms are compared using different factors such as accuracy score, analyzing the confusion matrix etc.

3. Overview of dataset

The source of this dataset is https://archive.ics.uci.edu/dataset/73/mushroom. This data set contains descriptions of hypothetical samples belonging to 23 Agaricus and Lepiota Family gilled mushrooms (pp. 500-525). Each species is labeled as either definitely edible, definitely poisonous, or maybe edible but not recommended. This latter category was integrated with the toxic category. The Guide explains unequivocally that there is no easy rule for judging mushroom edibility; no rule like

"leaflets three, let it be" for Poisonous Oak and Ivy. The dataset have 8124 instances and 23 features. Below table describes the independent attributes.

| Attributes | Description |
|------------------------------|---|
| cap-shape | bell=b,conical=c,convex=x,flat=f, knobbed=k,sunken=s |
| cap-surface | fibrous=f,grooves=g,scaly=y,smooth=s |
| cap-color | brown=n,buff=b,cinnamon=c,gray=g,green=r, pink=p,purple=u,red=e,white=w,yellow=y |
| bruises? | bruises=t,no=f |
| odor | almond=a,anise=l,creosote=c,fishy=y,foul=f, musty=m,none=n,pungent=p,spicy=s |
| Gill-attachment | attached=a,descending=d,free=f,notched=n |
| gill-spacing | close=c,crowded=w,distant=d |
| gill-size | broad=b,narrow=n |
| gill-color | black=k,brown=n,buff=b,chocolate=h,gray=g, green=r,orange=o,pink=p,purple=u,red=e,white=w,yellow=y |
| stalk-shape | enlarging=e,tapering=t |
| stalk-root | bulbous=b,club=c,cup=u,equal=e, rhizomorphs=z,rooted=r,missing=? |
| stalk-surface- above-ring | fibrous=f,scaly=y,silky=k,smooth=s |
| stalk-surface- below-ring | fibrous=f,scaly=y,silky=k,smooth=s |
| stalk-color- above-ring | brown=n,buff=b,cinnamon=c,gray=g,orange=o, pink=p,red=e,white=w,yellow=y |
| stalk-color- below-ring | brown=n,buff=b,cinnamon=c,gray=g,orange=o, pink=p,red=e,white=w,yellow=y |
| veil-type | partial=p,universal=u |
| veil-color | brown=n,orange=o,white=w,yellow=y |

| ring-number | none=n,one=o,two=t |
|-----------------------|---|
| ring-type | cobwebby=c,evanescent=e,flaring=f,large=l, none=n,pendant=p,sheathing=s,zone=z |
| spore-print- color | black=k,brown=n,buff=b,chocolate=h,green=r, orange=o,purple=u,white=w,yellow=y |
| population | abundant=a,clustered=c,numerous=n, scattered=s,several=v,solitary=y |
| habitat | grasses=g,leaves=l,meadows=m,paths=p, urban=u,waste=w,woods=d |

The **target feature** is "class" and it has two elements in it edible=e and poisonous=p.

4. Exploratory Data Analysis

<u>Importing necessary libraries:</u> Below figure shows the necessary python libraries that we require to perform our classification task. We will make use some of the popular python libraries like pandas, numpy, sci-kit learn, matplotlib and seaborn. Apart from these we will also require some libraries for our classification which we will see in detail in the later sections.

```
In [9]: #importing libraries
import numpy as np
import pandas as pd
import sklearn as sk
import matplotlib.pyplot as plt
import seaborn as sns
```

<u>Loading the dataset:</u> using pandas' read_csv function, we load the dataset of csv format for analysis.

Load the dataset

```
In [10]: mushroom_dataset = pd.read_csv("mushrooms.csv")
```

<u>Dimensions of dataset:</u> shape is used to check the dimensions of dataset. This returns a tuple with the number of rows and columns that the dataset has. Our current dataset has 8124 rows and 23 columns.

```
In [11]: #check the number of rows and columns in the dataset
mushroom_dataset.shape
Out[11]: (8124, 23)
```

<u>Displaying the first few rows of dataset:</u> This is done by using the head() function which print the first 5 rows in the dataset. In the current dataset the target feature

column is the first column.



The target feature 'class' is the first column in this dataset

<u>Summary of the dataframe:</u> The info() method is used to print the concise summary of the dataframe. From the below figure we can interpret the following

- All the features in the dataset consists of categorical values
- Since all the feature columns have categorical values in it, we will not run a describe method to know the five-point summary of our dataset.

```
In [5]: mushroom_dataset.info()
                   <class
                                    'pandas.core.frame.DataFrame'>
                  RangeIndex: 8124 entries, 0 to 8
Data columns (total 23 columns):
# Column
                                                                                          Non-Null Count
                              class
cap-shape
cap-surface
                                                                                          8124 non-null
8124 non-null
8124 non-null
                                                                                                                               object
object
object
                               cap-color
                                                                                          8124
                                                                                                     non-null
                                                                                                                               object
                             bruises
odor
gill-attachment
gill-spacing
gill-size
gill-color
stalk-shape
stalk-root
stalk-surface-above-ring
stalk-surface-below-ring
stalk-color-above-ring
veil-type
                              bruises
                                                                                          8124 non-null
                                                                                                                               object
                                                                                          8124
8124
                                                                                          8124
8124
8124
8124
8124
8124
                                                                                                                               object
                                                                                          8124
                                                                                                      non-null
                                                                                                                               object
                                                                                          8124
                                                                                                     non-null
                                                                                          8124
                                                                                          8124
                              stalk-color-below-
veil-type
veil-color
ring-number
ring-type
spore-print-color
population
habitat
                                                                                          8124
                                                                                          8124 non-null
8124 non-null
8124 non-null
8124 non-null
8124 non-null
8124 non-null
                                                                                          8124 non-null
                  dtypes: object(23)
memory usage: 1.4+ MB
```

<u>Check for missing values:</u> isnull() is used to check if there is any null values in each of our column. Sum() counts the number of occurrences of null value in each of these column. From the below figure it is clear that there are no null values in our dataset.

<u>Checking for cardinality:</u> The number of distinct values in a column or feature of a dataset is referred to as its cardinality. For the current dataset I will drop all those features with cardinality less than 3 since they won't be of much importance for the analysis. Below figure shows the result of the cardinality.

```
In [6]: mushroom_dataset.select_dtypes('object').nunique()

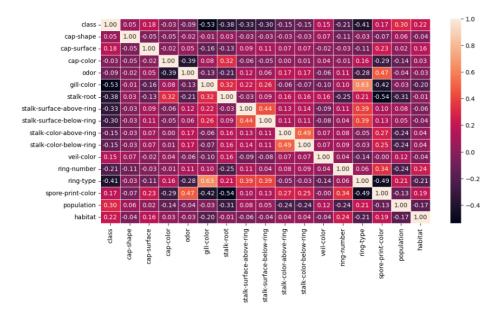
Out[6]: class 2
cap-shape 6
cap-surface 4
cap-color 10
bruises 2
odor 9
gill-spacing 2
gill-spacing 2
gill-spacing 2
gill-size 2
gill-color 12
stalk-shape 2
stalk-root 5
stalk-snore 5
stalk-surface-below-ring 4
stalk-surface-below-ring 9
stalk-color-below-ring 9
veil-type 1
veil-color 4
ring-number 3
ring-type 5
spore-print-color 9
population 6
habitat 7
dtype: int64

From the above result, we can interpret the following

- Feature columns bruises, gill-attachment, gill-spacing, gill-size, stalk-shape, veil-type have less cardinality.
- Action to be taken: drop these columns and train the model using remaining features
```

Based on the result from above figure we can observe that few features like bruises, gill-attachment, gill-spacing, gill-size, stalk-shape, veil-type have less cardinality. We will drop these features. To achieve this, we will make use of drop(). This method will drop the columns specified inside it. This is shown in below figure. The resultant dataset is stored as a new dataframe, mushroom new.

Correlation matrix: Figure below shows the correlation matrix visualized using the heatmap. Heatmap is a plotting technique in the seaborn library that is used to interpret the correlation. There is only a moderate correlation between our target variable and other feature variable. However, the negative correlation is bit high when compared to the positive ones. Gill color, stalk root, stalk surface above ring, stalk surface below ring and ring type has a higher negative correlation compared to other negatively correlated ones.



Visualizations:

Figure below shows a series of bar graphs that depict the distribution of the target variable ('class' - whether a mushroom is edible or poisonous) for each category feature in the dataset. For this we make use of the seaborn's barplot function with y axis displaying the count of occurances of the target variables for each element in every feature column.

A for loop is used to iterate through each categorical attribute in the dataframe. For each attribute, it calls the plotClass function that generate a barplot on the corresponding subplot. plt.tight_layout() is used to improve the appearance of the layout by preventing subplot overlaps. Finally, plt.show() is used to display the plots created.

```
In [35]: #Visualing the occurances of our target variables for each attribute

def plotClass(x,ax):
    group = mushroom_new.groupby([f'{x}','class'])['class'].count().reset_index(name='Count')
    sns.barplot(data=group,x=x,y='Count',hue='class',palette='viridis',ax=ax)
    ax.set_xticklabels(ax.get_xticklabels(), rotation=45, ha='right')
    ax.set_title(f'Count of {x}')
    cols = mushroom_new.columns.tolist()
    fig, axes = plt.subplots(8, 2, figsize=(12, 6 * 5))

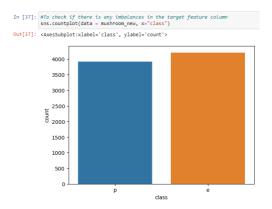
for index,column in enumerate(cols[1:]):
    row = index // 2
    col = index % 2
    ax = axes[row,col]
    plotClass(column, ax)

plt.tight_layout()
plt.show()
```

Figure below shows how the distribution of edible and deadly mushrooms vary for each categorical attribute in the dataset, revealing probable correlations between attributes and the target variable. From the plots it can be inferred that most of the mushrooms are edible for consuming.



<u>Check for class imbalance:</u> We use a countplot to visualize and check if the target feature column is balanced or not. From the below figure it can be interpreted that our target class is pretty much balanced since the imbalance is pretty minimal and we can proceed without any oversampling or undersampling methods.



5. Model Building Splitting dataset:

Before splitting the dataset, we need to segregate the independent and target columns. In the below code, we are selecting the features with more negative collinearity as our independent variables. The target column is placed in feature matrix y.

```
In [38]: # Arrange data into independent variables and target variables
X = mushroom_encoded.loc[: , ['cap-shape','stalk-surface-above-ring', 'habitat', 'spore-print-color']]
y=mushroom_encoded['class'] #is the target
```

Below code use the train_test_split function of scikit-learn for splitting the datasets to train and test sets. Here 30% of data will be used for testing and rest is training datasets. Value placed in random state parameter will help us ensure that we will get the same split when we use the same value in future. Training set(X_train and y_train) is used to train the machine learning model and Test set(X_test and y_test) is used to evaluate the performance.

```
X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.30, random_state=10)
```

<u>Label Encoding:</u> Label encoding is analogous to providing a dictionary to a computer that converts human language into a language that the computer understands. It's a basic yet efficient strategy for making machine learning algorithms more versatile and effective. In our dataset all the columns have categorical variables, so before training the model it is important to convert into numerical so that the algorithm can handle and process the data uniformly throughout the modeling process. mushroom_new.apply(lambda col: encoder.fit_transform(col)): Using the apply function to apply label encoding to each column (col) in the DataFrame mushroom_new. The lambda function uses the LabelEncoder's fit_transform method to each column independently.

Classification model1: Logistic Regression:

Below code shows the implementation of the logistic regression model using scikit-learn's LogisticRegression library. The solver option in scikit-learn's LogisticRegression class specifies the optimization algorithm to be utilized during logistic regression model training. We are using the liblinear value for solver since this value is basically used for small or mredium datasets.

In machine learning, "Fit (Train) the Model" entails training the model with a dataset to understand the patterns and correlations in the data. Here fit method is called on the logistic regression model with the training data.

Finally, the trained model is used to make predictions on the test data. This is done by the predict().

```
In [41]: from sklearn import metrics
    from sklearn.linear_model import LogisticRegression

# fitting the model on training
    log_r = LogisticRegression(solver="liblinear")
    log_r.fit(X_train, y_train)
    #predict on test
    y_predict = log_r.predict(X_test)
```

Below figure shows the code to detect the accuracy of how correctly our model has predicted the actual labels in test data. Then a confusion matrix is drawn to show the number of correct and wrong predictions against the actual data. Finally, it calculates the classification report for each class, which contains precision, recall, F1-score, and support. The overall accuracy of the model is 72% which means it correctly predicted the class for 72% of the instances in the test set. From confusion matrix we can interpret that True Positive is 775 i.e., the model predicted edible mushrooms as edible correctly. True Negative is 988 which means the model could predict 988 poisonous mushrooms correctly. Precision for 1 is 72% which means 72% of edible predictions were actually edible. Recall for edible is 67% which means that 67% of edible mushrooms were correctly predicted.

```
In [44]: from sklearn import metrics
          Accuracy_LR=metrics.accuracy_score(y_test,y_predict)
print('accuracy:%.2f\n\n'%(Accuracy_LR))
          LR_CM=metrics.confusion_matrix(y_test,y_predict)
print('Confusion Matrix:')
          print(LR_CM,'\n\n')
           result=metrics.classification_report(y_test,y_predict)
          print('Classification Report:\n')
          print(result)
           accuracy:0.72
          Confusion Matrix:
           [[988 288]
[387 775]]
          Classification Report:
                          precision recall f1-score
                                                              support
                                            0.77
                                      0.77
0.67
                                                                   1276
                                0.72
                                                     0.70
                                0.73
                                                                   1162
               accuracy
                                                       0.72
                                                                   2438
              macro avg
          weighted avg
                                0.72
                                            0.72
                                                        0.72
                                                                   2438
```

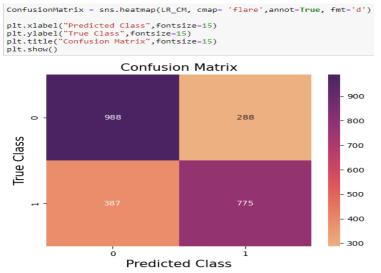
Confusion Matrix: Figure below shows the code for plotting the confusion matrix using seaborn's heatmap. From the confusion matrix the following can be interpreted:

True Positives (TP): 775

True Negatives (TN): 988

False Positives (FP): 288

False Negatives (FN): 387



Classification model2: KNeighbours Classifier:

The KNeighborsClassifier is a scikit-learn classification algorithm that implements the knearest neighbor's approach. Below code shows the implementation of the KNN model using scikit-learn's KNeighborsClassifier library. In the KNeighborsClassifier method we assign the number of neighbours to 2 which means that the algorithm will consider 2

neighbours while making predictions. The distance metric used here is Euclidean hence the p value is set to 2. Then we fit the model on the training dataset using fit(). Finally, the predict() predicts the labels for a test dataset (X_test) after training and keeps the predictions in y_pred.

Classification Model 2: K Nearest Neighbour

```
: from sklearn.neighbors import KNeighborsClassifier
classifier = KNeighborsClassifier(n_neighbors=2, metric='euclidean', p =2)
classifier.fit(X_train, y_train)
: KNeighborsClassifier(metric='euclidean', n_neighbors=2)
: #Evaluating the model
y_pred = classifier.predict(X_test)
C:\Users\vysha\anaconda3\lib\site-packages\sklearn\neighbors\classification
tions (e.g. `skew`, `kurtosis`), the default behavior of `mode` typically pr
is behavior will change: the default value of `keepdims` will become False,
e eliminated, and the value None will no longer be accepted. Set `keepdims`
mode, _ = stats.mode(_y[neigh_ind, k], axis=1)
```

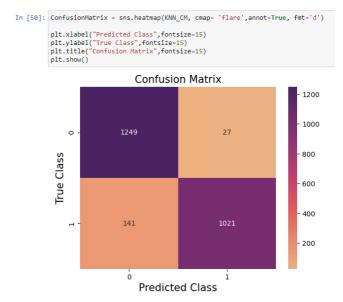
Below figure shows the code to detect the accuracy of how correctly our model has predicted the actual labels in test data. Then a confusion matrix is displayed to show the number of correct and wrong predictions against the actual data. Finally, it calculates the classification report for each class, which contains precision, recall, F1-score, and support. The overall accuracy of the model is 93% which means it correctly predicted the class for 93% of the instances in the test set. From confusion matrix we can interpret that True Positive is 1021 i.e., the model predicted edible mushrooms as edible correctly. True Negative is 1249 which means the model could predict 1249 poisonous mushrooms correctly.

Precision for 1 is 97% which means 97% of edible predictions were actually edible. Recall for edible is 92% which means that 92% of edible mushrooms were correctly predicted.

Confusion Matrix: Figure below shows the code for plotting the confusion matrix using seaborn's heatmap. From the confusion matrix the following can be interpreted:

• True Positives (TP): 1021

- True Negatives (TN): 1249
- False Positives (FP): 27
- False Negatives (FN): 141



Classification model2: Decision Tree Classifier:

The DecisionTreeClassifier is a scikit-learn classification algorithm that implements the decision tree algorithm for classification. Below code shows the implementation of the Decision tree model using scikit-learn's DecisionTreeClassifier library.

The criterion used here is entropy. Entropy is a measure of a node's disorder or impurity. As a result, a node with a more variable composition, such as 2Pass and 2Fail, is thought to have higher Entropy than a node with only pass or only fail. Then we fit the model on the training dataset using fit(). Finally, the predict() predicts the labels for a test dataset (X_test) after training and keeps the predictions in y_pred.

```
In [51]: from sklearn.tree import DecisionTreeClassifier
    classifier = DecisionTreeClassifier(criterion= 'entropy', random_state = 0)
    classifier.fit(X_train, y_train)

Out[51]: DecisionTreeClassifier(criterion='entropy', random_state=0)

In [52]: y_pred = classifier.predict(X_test)
```

Below figure shows the code to detect the accuracy of how correctly our model has predicted the actual labels in test data. Then a confusion matrix is displayed to show the number of correct and wrong predictions against the actual data. Finally, it calculates the classification report for each class, which contains precision, recall, F1-score, and support. The overall accuracy of the model is 94% which means it correctly predicted the class for 94% of the instances in the test set. From confusion matrix we can interpret that True Positive is 1056 i.e., the model predicted edible mushrooms as edible correctly. True

Negative is 1243 which means the model could predict 1243 poisonous mushrooms correctly.

Precision for 1 is 97% which means 97% of edible predictions were actually edible. Recall for edible is 91% which means that 91% of edible mushrooms were correctly predicted.

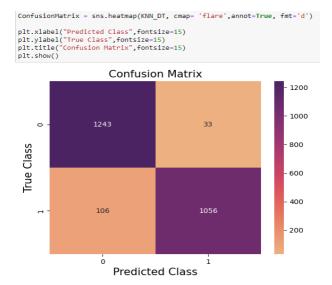
Confusion Matrix: Figure below shows the code for plotting the confusion matrix using seaborn's heatmap. From the confusion matrix the following can be interpreted:

True Positives (TP): 1056

True Negatives (TN): 1243

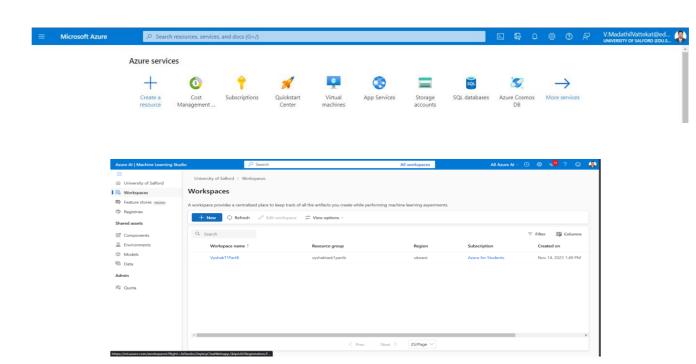
• False Positives (FP): 33

• False Negatives (FN): 106

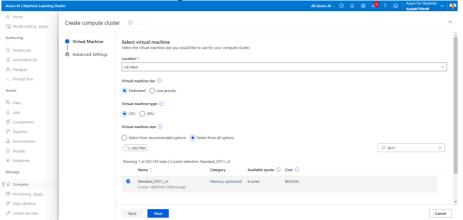


PART B: Machine Learning using Azure

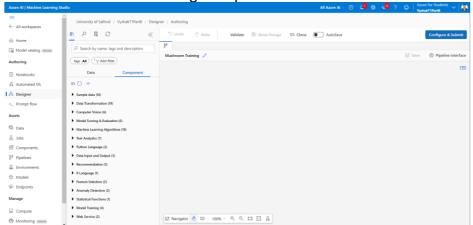
After creating an Azure Machine Learning Resource, fill in the necessary settings and click on create to get the workspace up and running.



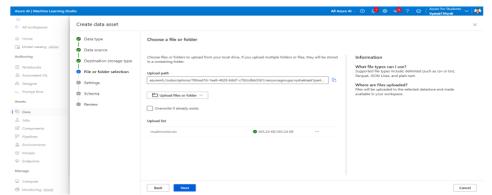
A new compute cluster is created using Standard_DS11_v2 virtual machine size



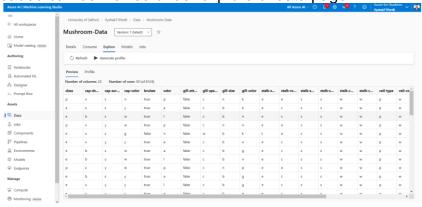
A new pipeline is created from the designer option in left menu.



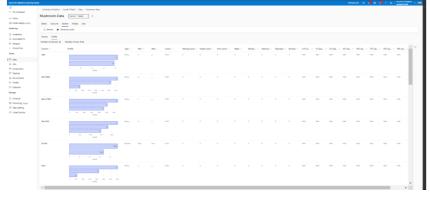
To create a dataset, select data from the Assets menu. After giving the name to the dataset, import the dataset(mushroom.csv) from local files. Destination storage type shall be **workspaceblobstore**. Upload the csv file into Azure. After uploading review the schema and click on create.



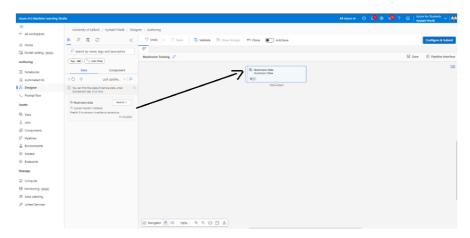
The dataset uploaded can be viewed under explore of data page.



The occurances of target attribute for each feature column can be visualized under profile.

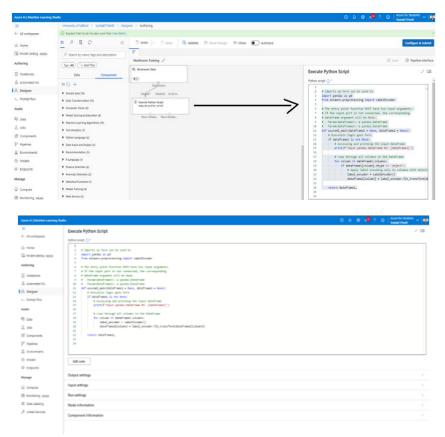


In designer page, load the dataset to canvas by dragging the dataset from data tab to the canvas

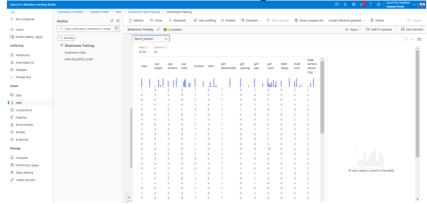


Add Transformations:

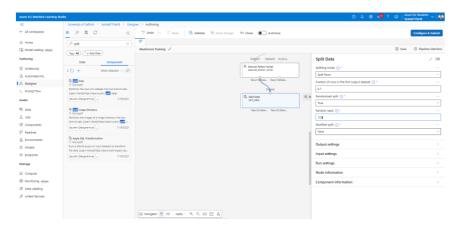
1. <u>Convert to Indicator Values</u>: Used to convert categorical elements to numerical. We execute the python script using the execute python script component to carry out label encoding.



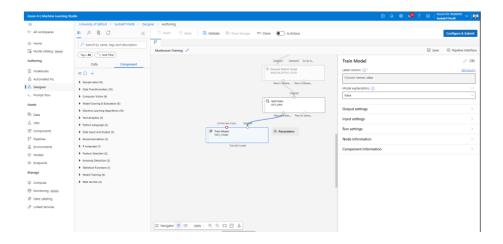
Below is the resultant dataset. All columns are having the encoded numerical characters.



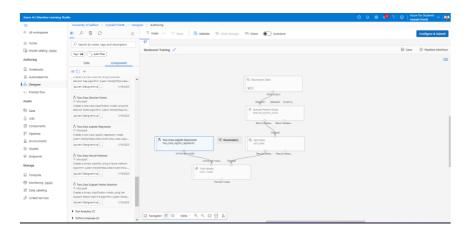
2. <u>Training the model:</u> Firstly, we need to split data into train and test data. We Use the split data transformation to achieve this. 70% will be given for test data.



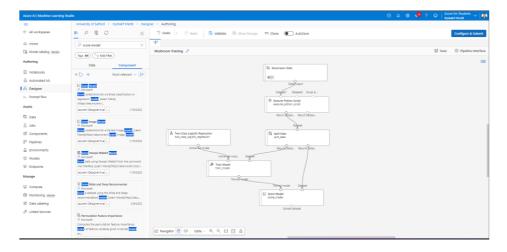
3. Next, we add the train model transformation. Connect split data resultant dataset to train model dataset input(right node). Provide the target variable in the train model.



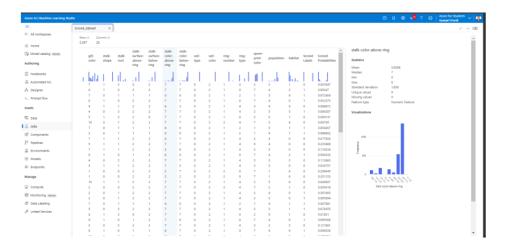
4. <u>Two Class Logistic Regression</u>: Add the Two Class Logistic Regression from the machine learning algorithm components. Connect the output of this to the left node of Train model.



5. **Scored Model:** Connect the output of train model to the trained model node input of score model and then connect the resultant dataset of split data to dataset node of scored model.

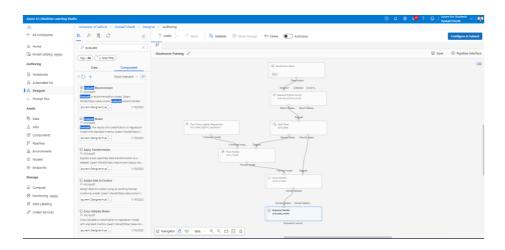


6. <u>Run the training pipeline:</u> Click on configure and submit. In the jobs page select the latest pipeline. Right click on **Scored model** and select Preview data to analyze the result.



The result will have scored labels and scored probability which displays what is the probability of the predicted result.

7. <u>Evaluate the model:</u> Add Evaluate model from components to the canvas. And then click again on configure and submit to run the pipeline.

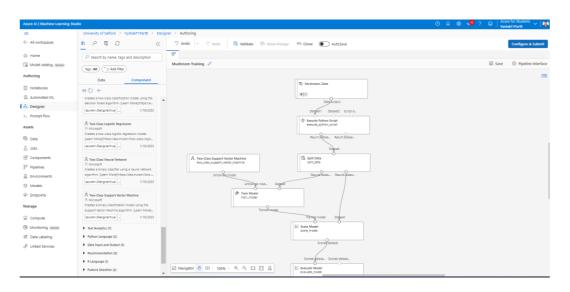


Once the run is completed go to job and right click on evaluate model to analyze the results.



The two-model logistic regression predicts with an accuracy of 94.8%. True positive is 1098. True Negative is 1212.

Two Class Support Vector Machine:

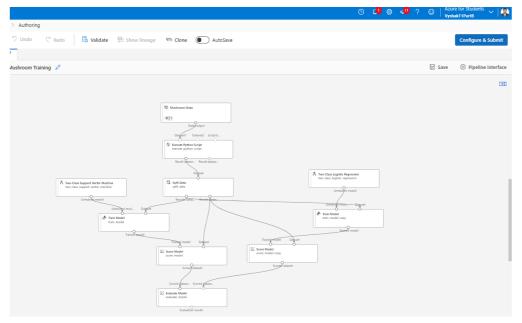


Once added, click on configure and submit. Once the pipeline run is completed, go to the jobs page and right click on Evaluate model to analyze the result.



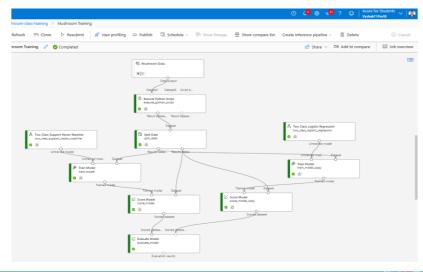
Model predicted at an accuracy of 94.1%. True positives are 1081. True Negatives are 1212.

<u>Training Model together</u>: Below figure shows running a pipeline with two models parallelly. We created two train models and score models. Then the machine learning algorithms were connected as explained previously. Then we ran the pipeline.



Once the pipeline run is completed, we will go to the jobs page to evaluate the model.

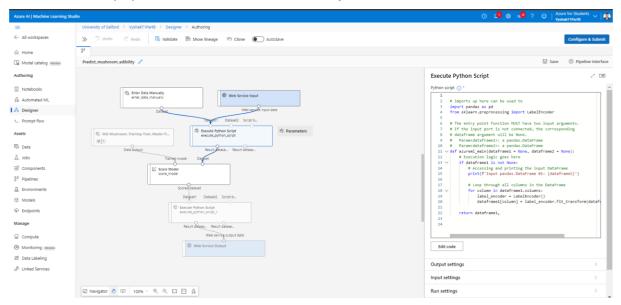
The evaluate model will now have two separate tabs with results of both machine learning models.



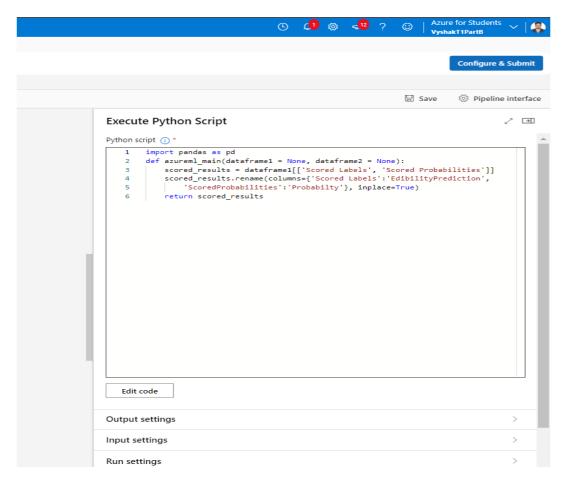


Inference Pipeline:

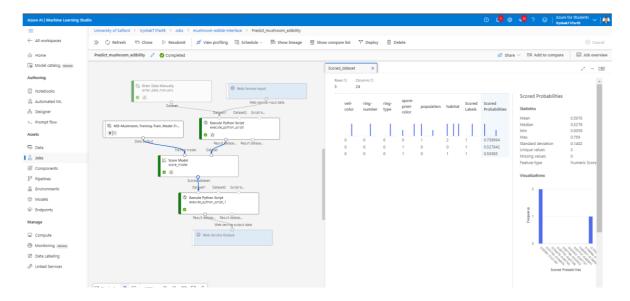
On jobs page, click on the mushroom training pipeline and create an inference pipeline. The canvas will be populated with few items already in it.



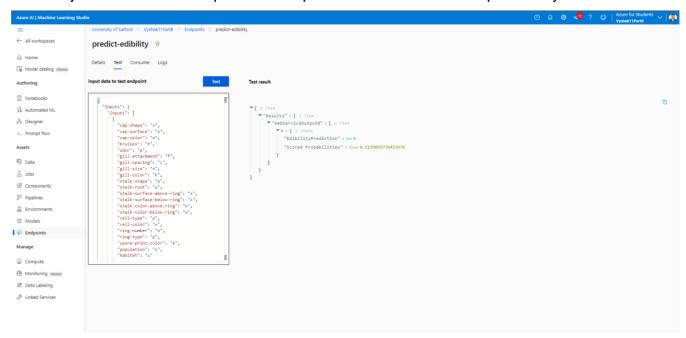
The first execute python script label will have label encoder script in it. In the second Execute python script, we will have to edit the following:



Run the inference pipeline and once it is completed. Analyze the scored labels and scored probabilities.



Next we will deploy our inference pipeline to test. Once the deployment is done, a test data is fed in json format. The output will be a prediction and the scored probability.



8. Evaluation of Results:

Python: From the below figure, we can confirm that the best performing model is Decision Tree Classifier which predicted with an accuracy of 94%.

Comparing the results



Azure: Two model logistic regression gave good accuracy with 94.8%. the precision and

recall metrics—95.6% and 93.4%, respectively—are incredibly high. This indicates that 95.6% of the time the model properly identifies edible mushrooms and 93.4% of the time it correctly identifies poisonous mushrooms.

In Conclusion, both the Python Decision Tree model and the Microsoft Azure logistic regression model showed remarkable accuracy when assessing the classification performance for the mushroom dataset. The Decision Tree demonstrated its strong capacity to distinguish between edible and deadly mushrooms with an accuracy of 94%, along with good precision and recall scores. In contrast, the logistic regression model on Azure performed better, demonstrating its effectiveness in appropriately classifying edible and hazardous mushrooms with an accuracy of 94.8%. With precision and recall metrics of 95.6% and 93.4%, respectively, these numbers demonstrate the model's remarkable categorization accuracy. These findings support the effectiveness of both models in handling the task of predicting the edibility of mushrooms, offering trustworthy and precise insights into their classification.

Part Two: Clustering

1. Introduction

Imagine you have a vast library of user reviews spanning a wide spectrum of establishments, from cozy cafes and bustling marketplaces to serene temples and vibrant art galleries. This is essentially the Google review dataset, a treasure trove of insights into people's preferences and experiences across various venues.

The Google review dataset is a collection of user reviews on a variety of establishments, including churches and resorts, as well as restaurants and beauty spas. Each record in the dataset is distinguished by a unique user ID and average ratings across all sorts of venues, revealing vital information about user preferences and opinions. The dataset provides a rich supply of data for clustering analysis, which aims to group comparable individuals based on their average ratings across many attributes.

Each entry in this dataset is like a unique voice, express ing their opinions and ratings for different types of establishments. Just as a book reviewer might have distinct tastes in genres, a user in this dataset may prefer lively cafes and adventurous restaurants while another might gravitate towards serene temples and intimate art galleries.

This information acts as a social compass, directing us to people who share our tastes and preferences, allowing us to discover new experiences and connections throughout the vast geography of human preferences.

2. Research Objective

"Using clustering methods, investigate patterns in user preferences in the Google review dataset."

The major goal of this clustering study is to find hidden patterns and structures in the Google review dataset. We intend to group users with similar reviewing patterns and preferences using two clustering techniques, potentially revealing distinct subsets of the user community. Finally, we need to decide on which clustering technique can be utilized for future applications.

3. Overview of dataset

The source of this dataset is https://archive.ics.uci.edu/dataset/485/tarvel+review+ratings
User ratings from Google reviews are used to create this data set. Attraction reviews from 24 categories from around Europe are evaluated. Google user ratings vary from 1 to 5, with the average user rating derived for each category. The dataset have 5456 instances and 26 features. Below table describes the independent attributes.

| Attributes | Description |
|-------------|--|
| User | Unique user id |
| Category 2 | Average ratings on churches |
| Category 3 | Average ratings on beaches |
| Category 4 | bruises=t,no=f |
| Category 5 | Average ratings on parks |
| Category 6 | Average ratings on theatres |
| Category 7 | Average ratings on museums |
| Category 8 | Average ratings on malls |
| Category 9 | Average ratings on zoo |
| Category 10 | Average ratings on restaurants |
| Category 11 | Average ratings on pubs/bars |
| Category 12 | Average ratings on local services |
| Category 13 | Average ratings on burger/pizza shops |
| Category 14 | Average ratings on hotels/other lodgings |
| Category 15 | Average ratings on juice bars |
| Category 16 | Average ratings on art galleries |
| Category 17 | Average ratings on dance clubs |
| Category 18 | Average ratings on swimming pools |
| Category 19 | Average ratings on gyms |
| Category 20 | Average ratings on bakeries |
| Category 21 | Average ratings on beauty & spas |
| Category 22 | Average ratings on cafes |
| Category 23 | Average ratings on view points |
| Category 24 | Average ratings on monuments |

4. Exploratory Data Analysis

Importing necessary libraries: Below figure shows the necessary python libraries that we require to perform our clustering task. We will make use some of the popular python libraries like pandas, numpy, matplotlib and seaborn. Apart from these we will also require some libraries for our clustering which we will see in detail in the later sections. Here we set OMP_NUM_THREADS to 1 to limit the parallelism used by the KMeans algorithm.

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

#To run befoire importing KMeans
import os
os.environ["OMP_NUM_THREADS"] = '1'
```

<u>Loading the dataset:</u> using pandas' read_csv function, we load the dataset of csv format for analysis.

```
#Import the google review ratings dataset
review_dataset = pd.read_csv("google_review_ratings.csv")
```

<u>Dimensions of dataset:</u> shape is used to check the dimensions of dataset. This returns a tuple with the number of rows and columns that the dataset has. Our current dataset has 5456 rows and 26 columns.

```
#determine the number of rows and column in the dataset
review_dataset.shape
(5456, 26)
```

There are 5456 instances and 26 features in this dataset

<u>Displaying the first few rows of dataset:</u> This is done by using the head() function which print the first 5 rows in the dataset.

| | User | Category 1 | Category 2 | Category 3 | Category 4 | Category 5 | Category 6 | Category 7 | Category 8 | | Category 16 | Category 17 | Category 18 | Category 19 | Category 20 |
|---|-----------|---------------|------------|------------|---------------|---------------|---------------|---------------|---------------|--------|----------------|----------------|----------------|----------------|----------------|
| 0 | User 1 | 0.0 | 0.0 | 3.63 | 3.65 | 5.0 | 2.92 | 5.0 | 2.35 | 2.33 . | 0.59 | 0.5 | 0.0 | 0.5 | 0.0 |
| 1 | User 2 | 0.0 | 0.0 | 3.63 | 3.65 | 5.0 | 2.92 | 5.0 | 2.64 | 2.33 | 0.59 | 0.5 | 0.0 | 0.5 | 0.0 |
| 2 | User 3 | 0.0 | 0.0 | 3.63 | 3.63 | 5.0 | 2.92 | 5.0 | 2.64 | 2.33 | 0.59 | 0.5 | 0.0 | 0.5 | 0.0 |
| 3 | User 4 | 0.0 | 0.5 | 3.63 | 3.63 | 5.0 | 2.92 | 5.0 | 2.35 | 2.33 . | 0.59 | 0.5 | 0.0 | 0.5 | 0.0 |
| 4 | User 5 | 0.0 | 0.0 | 3.63 | 3.63 | 5.0 | 2.92 | 5.0 | 2.64 | 2.33 . | 0.59 | 0.5 | 0.0 | 0.5 | 0.0 |

<u>Summary of the dataframe:</u> The info() method is used to print the concise summary of the dataframe. From the below figure we can interpret the following

 All the features in the dataset consists of categorical values except category 11 which we will convert to numeric in next step.

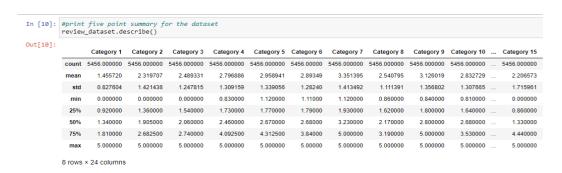
```
cclass 'pandas.core.frame.DataFrame'>
RangeIndex: 5456 entries, 0 to 5455
Data columns (total 26 columns):
    # Column    Non-Null Count Dty
                                                                                                                                                                                                                                                                                             5456 non-null
                                                                          User
Category 1
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Category 3
Category 4
Category 5
Category 7
Category 7
Category 8
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Category 10
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                                                                                                                                                                                                                                                                                               5456
                                                                                 Category
                                                                                                                                                                                                                               14
   dtypes: float
memory usage:
```

Figure 1

<u>Five Point Summary:</u> Figure below gives the five-point summary of the dataset. Below are some of the observations from that

- The mean ratings across all categories generally range from approximately 1.45 to 3.35. The highest average rating is observed in Category 7, while the lowest mean rating is in Category 18.
- Category22 has the highest SD hence, it has more dispersed ratings. Whereas category1 has the lowest meaning the ratings are consistent.

- A higher median often indicates that a considerable portion of users rated that
 category significantly higher, indicating a more positive overall mood or experience.
 A lower median indicates that a significant number of people rated it lower, implying
 a less pleasant emotion or experience on average.
- Looking at the five point summary, we can assume that there wont be significant outliers.



<u>Changing datatype of feature column:</u> One of the columns is a categorical column but has elements of numerical datatype. Hence, we will change the datatype from object to numerical. This is shown in Figure below

<u>Check for missing values:</u> isnull() is used to check if there is any null values in each of our column. Sum() counts the number of occurrences of null value in each of these column. From the below figure column 26 have plenty of null values and removing this wont affect much of our analysis. Hence we will remove that in the next step.

Removing unwanted columns: Column 26 has plenty null values. As discussed above we will remove this column using the drop function.

```
#drop column 26 since it has more null values
review_dataset = review_dataset.drop('Unnamed: 25', axis = 1)
```

Renaming the existing columns: For better readability we are renaming the existing columns as shown in the figure below

| ev | 'Cate 'Cate 'Se the | | paRating', '(pardens' pod p(columns=ne | 'Category 1 Category 21 | 7': 'Swimm ': 'CafeRa | ingPools','C ting','Categ | ategory 18': | 'GymRatin | g','Catego | ory 19': 'Bakery 23': 'Monuments | у', | tGalleries | , |
|--------|---------------------------|------------------|--|----------------------------|--------------------------|------------------------------|----------------------|----------------|-------------------|-------------------------------------|-----|--------------|---|
| ev | | | | | | | | | | | | | |
| 21 | UserID | ChurchRating | ResortRating | BeachRating | ParkRating | TheatreRating | MuseumRating | MallRating | ZooRating | RestaurantRating | | ArtGalleries | D |
| | UserID User 1 | ChurchRating 0.0 | ResortRating 0.0 | BeachRating 3.63 | ParkRating 3.65 | TheatreRating 5.0 | MuseumRating 2.92 | MallRating 5.0 | ZooRating 2.35 | RestaurantRating 2.33 | | ArtGalleries | D |
| 0 | | | | | | | | | | | | | D |
| 0 | User 1 | 0.0 | 0.0 | 3.63 | 3.65 | 5.0 | 2.92 | 5.0 | 2.35 | 2.33 | | 1.74 | |
|) I | User 1 User 2 | 0.0 | 0.0 | 3.63 3.63 | 3.65 3.65 | 5.0 5.0 | 2.92 | 5.0 5.0 | 2.35 2.64 | 2.33 2.33 | | 1.74 1.74 | |

5. Clustering K-Mean Clustering technique

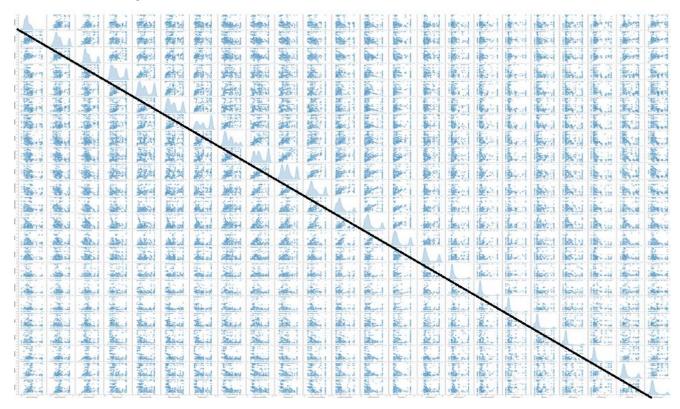
Normalization:

Normalizing the data before to applying K-Means clustering is a best practice for ensuring fair and effective clustering results, which provide a more accurate picture of the underlying patterns in the data. Even though there is no much scale differences, it is always a best practice to do normalization before applying KMeans for the following reasons:

- K-Means is sensitive to feature scale. Normalizing the data guarantees that each feature is assigned equal weight in the clustering process.
- Normalizing the data can improve the K-Means algorithm's convergence speed and efficiency.

Below figure shows the normalizing of numerical columns using z score normalization which is also known as standardization. We use the **standardscaler** class from scikit-learn to achieve that. In fit_transform(), the mean and standard deviation of each feature are calculated, and the data is then standardized by subtracting the mean and dividing by the standard deviation. Resultant features will have a mean of 0 and SD of 1.

<u>Scatter Plot:</u> To get an initial guess on the number of clusters we will plot a scatter plot that shows the pairwise relationship between the variables. The KDE plot along diagonal will provide insights into the distribution which will further give us an initial guess on the number of clusters by analyzing their peaks. From figure below we can interpret most of the variables have a **peak** of 3. So, we will first run our K Means with **3 clusters**.



<u>Fitting K-Means to dataset</u>: The below figure shows the code to fit the K-Means clustering algorithm to the scaled dataset. For this we create an instance of KMeans class and pass the number of clusters as 3. K-means++ is a smart initialization method for the centroids. The fit_predict() is used to fit the KMeans model to standardized data and result is stored in y_kmeans_3. Cluster label for each data point is assigned to labels variable.

```
#Fitting K-Means to the dataset
from sklearn.cluster import KMeans
kmeans_3 = KMeans(n_clusters = 3, init = 'k-means++', random_state = 15)
y_kmeans_3 = kmeans_3.fit_predict(scaled_review)
labels = kmeans_3.labels_
```

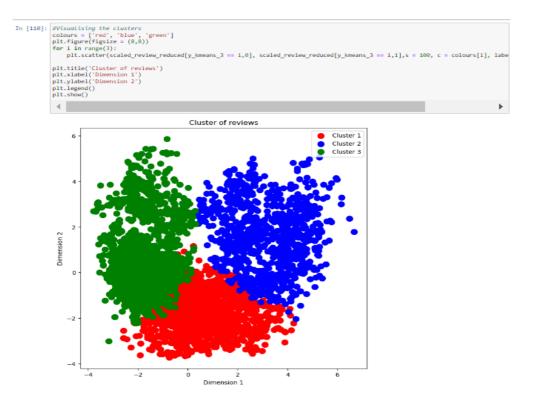
<u>Silhouette Analysis:</u> The ideal number of clusters can be determined by silhouette analysis. The ideal number of clusters is sometimes thought to be the one that optimizes the average silhouette coefficient. Greater isolation and cohesiveness between clusters are indicated by higher silhouette coefficients. Below code shows the Silhoette Coefficient for 3 clusters. **The coefficient score is 0.145**. We will do further analysis with more clusters and determine the best by analysing the silhouette score received.

```
In [17]: #Calculate the Silhouette coefficient for 3 clusters
    from sklearn.metrics import silhouette_samples, silhouette_score
    silhouette_score(scaled_review,labels)
Out[17]: 0.1457822651039671
```

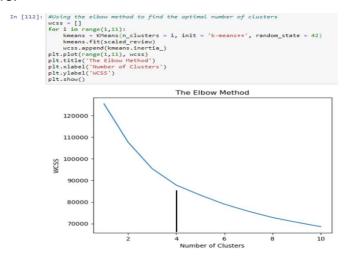
<u>Principal Component Analysis:</u> Since our dataset has 24 components we cannot visualize our clusters in 24 Dimensions, hence we will use PCA technique to reduce the dimensionality. In PCA we assign a value for number of components inorder to retain more variance in our visualization which is set to 10. Then for the purpose of visualizing in 2D space we use a subset of these components. In the below code we could get a cumulative variance of 74.3%. The interesting fact was with 2 components I could only get 34.4% variance.

Visualising the clusters:

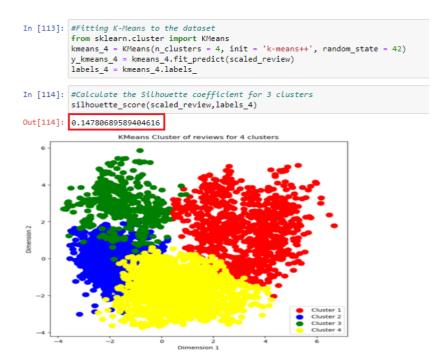
Below figure shows the code for visualizing the clusters on a 2D space. We will be visualizing the 3 clusters below. In the next step we will use a elbow curve to check if we could get more clarity by increasing the clusters.



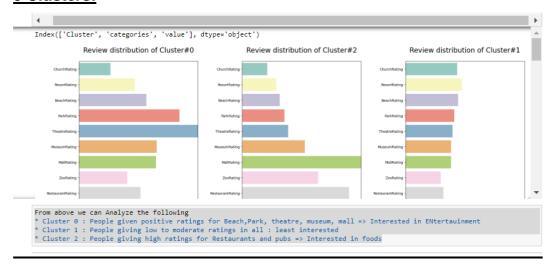
Elbow Curve: Figure below shows the implementation and output of the elbow method that we are using to determine the number of clusters from the elbow curve. WCSS in Y axis is calculated for each value of k and this decreases as k value increases forming a sharp decrease which is considered as an elbow and we take the value of k at this point. In the output below we can interpret that there is an elbow at k=4. This means we can perform KMeans with 4 clusters.



KMeans with 4 clusters : Below figure shows the Silhouette score and scatter plot with 4 clusters. The silhouette score has improved very minimal compared to the one with 3 clusters. However we will visualize and check what features are present in each cluster by taking 4 clusters.



KMeans Cluster Profiling - Visualizing the features assigned to each cluster: 3 Clusters:



4 Clusters:



From above we can Analyze the following

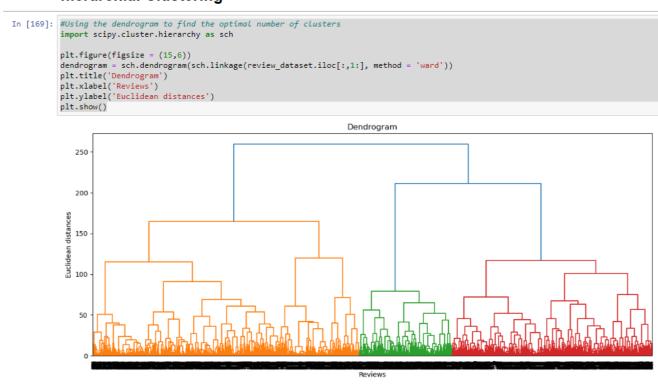
- Cluster 1 : People giving low to moderate ratings in all
- Cluster 2 : People giving high ratings for mall restaurants and pubs
- Cluster 3: People giving high ratings for Burgers, Hotels, Juice Bar, Art Galleries
 Cluster 4: People giving high ratings for Beach, Park, Museum, Theatre

From the above two figures we can interpret that having 4 clusters help us group more precisely. Hence for K Means 4 cluster will be a good choice.

Hierachial Clustering

The below code plots a dendogram which helps us visualize the hierarchial clustering. From the output, we can interpret that we can take the number of clusters as 3 if we imagine a vertical line is drawn through the tallest vertical line.

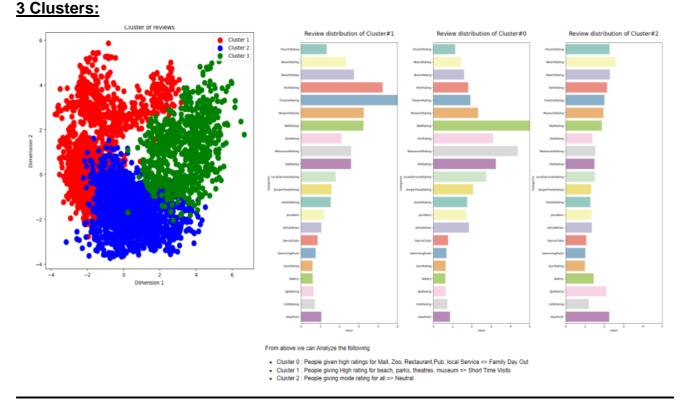
hierarchial Clustering



<u>Performing hierarchical clustering with 3 clusters:</u> The below code uses scikit-learn's AgglomerativeClustering algorithm to perfrom clustering with 3 clusters. As we can observe the Silhouette score turned out to be 0.165.

```
In [170]: # Fitting Hierarchial Clustering to the dataset
    from sklearn.cluster import AgglomerativeClustering
    hc_3 = AgglomerativeClustering(n_clusters=3, affinity = 'euclidean', linkage = 'ward')
    y_hc_3 = hc_3.fit_predict(review_dataset.iloc[:,1:])
    labels_hc_3 = hc_3.labels_
In [171]: #Calculate the Silhouette coefficient for 3 clusters
    silhouette_score(review_dataset.iloc[:,1:],labels_hc_3)
Out[171]: 0.1654648378380093
```

<u>Hierarchial Cluster Profiling - Visualizing the features assigned to each cluster:</u>



6. Evaluation of Results

- Analyzing the Silhouette score, 3 clusters generated by hierarchial Clustering would be recommended for further Model Building process.
- We were able to categorize the clusters to 3 different categories: Food and Entertainment, Short Time Visits, Neutral.
- These grouping can help to build any recommending application according to the interests of the users.

Finally, for further model building, the Silhouette score strongly suggests using three clusters in hierarchical clustering. The three found clusters (Food and Entertainment, Short Time Visits, and Neutral) provide useful information for developing applications that provide tailored recommendations. These clusters show different user preferences, ranging from a neutral interest group to a desire for short visits and interests in food and entertainment. The outcomes offer a tactical basis for augmenting user involvement, permitting customized offerings, and guiding focused promotional endeavors.

Part Three: Text Classification and Sentiment Analysis

1. Introduction

Envision discovering an immense collection of film critiques, every one conveying its distinct viewpoint and emotion. This is the core of the IMDB dataset, an extensive compilation of movie reviews that provides an insight into the realm of public opinion and cinema criticism. We will use the capabilities of natural language processing (NLP), namely the Multinomial Naive Bayes (MultinomialNB) approach, to explore this linguistic terrain. Like a proficient language interpreter, our algorithm will assist us in identifying the positive and negative sentiments contained in these reviews.

We aim to evaluate MultinomialNB's performance in this assignment by determining how well it can discern the underlying emotion of movie reviews. Understanding how it performs will help us recognize its advantages and disadvantages in the context of text-based analysis.

2. Research Objective

"To apply Multinomial Naive Bayes for text classification and perform sentiment analysis on the IMDB dataset to discern sentiment patterns and evaluate classification performance."

The goal of this study is to delve into the complexities of sentiment analysis by investigating how Multinomial Naive Bayes, a commonly used text classification algorithm, performs on the IMDB dataset. The project intends to use this classification approach to reveal the underlying sentiment structure in movie reviews, distinguishing between positive and negative thoughts.

This study is significant because it has the potential to contribute to a better understanding of sentiment analysis applications in the context of movie reviews. Analyzing the sentiment distribution across various reviews provides insights into audience perspectives, perhaps assisting filmmakers, critics, and platforms in measuring movie response. Furthermore, assessing the accuracy and performance metrics of the classification model provides a quantitative layer, confirming the dependability of the chosen approach for sentiment analysis.

3. Overview of dataset

IMDB dataset (Maas, A. L., Daly, R. E., Pham, P. T., Huang, D., Ng, A. Y., & Potts, C. 2011, June 1) having 50K movie reviews for natural language processing or Text analytics. This is a dataset for binary sentiment classification containing substantially more data than previous benchmark datasets. We provide a set of 25,000 highly polar movie reviews for training and 25,000 for testing. So, predict the number of positive and negative reviews

using either classification or deep learning algorithms.

For this task we will be using the unsupervised dataset of imdb review dataset that has 50000 instances in it. To download the dataset, we will follow some of the below steps:

Step1: Installing the dataset via the Python Package Index (PyPi)

```
In [10]: #To install dataset package for installing immdb dataset

#These are the datasets

Requirement already satisfied: datasets in c:\users\vysha\anaconda3\lib\site-packages (2.15.0)

Requirement already satisfied: dillo3.8,>=0.3.0 in c:\users\vysha\anaconda3\lib\site-packages (from datasets) (0.3.7)

Requirement already satisfied: aiohttp in c:\users\vysha\anaconda3\lib\site-packages (from datasets) (3.9.1)

Requirement already satisfied: numby=1.17 in c:\users\vysha\anaconda3\lib\site-packages (from datasets) (1.21.5)

Requirement already satisfied: xxhash in c:\users\vysha\anaconda3\lib\site-packages (from datasets) (1.2.1.5)

Requirement already satisfied: multiprocess in c:\users\vysha\anaconda3\lib\site-packages (from datasets) (1.0.0.1)

Requirement already satisfied: fsspec[http]<2023.10.0; nc:\users\vysha\anaconda3\lib\site-packages (from datasets) (0.70.15)

Requirement already satisfied: tqdm>=4.62.1 in c:\users\vysha\anaconda3\lib\site-packages (from datasets) (4.64.1)

Requirement already satisfied: huggingface-hub>=0.18.0 in c:\users\vysha\anaconda3\lib\site-packages (from datasets) (4.64.1)

Requirement already satisfied: pyarrow-hofix in c:\users\vysha\anaconda3\lib\site-packages (from datasets) (0.6)

Requirement already satisfied: pyarrow-hofix in c:\users\vysha\anaconda3\lib\site-packages (from datasets) (0.6)

Requirement already satisfied: pyaml>=5.1 in c:\users\vysha\anaconda3\lib\site-packages (from datasets) (21.3)

Requirement already satisfied: requests>=2.19.0 in c:\users\vysha\anaconda3\lib\site-packages (from datasets) (2.28.1)

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Requirement already satisfied: requests>=2.19.0 in c:\users\vysha\anaconda3\lib\site-packages (from datasets) (2.28.1)

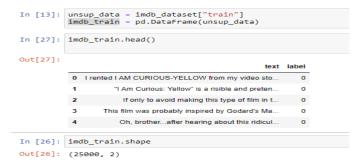
Requirement already satisfied: attrs>=1.1.1 in c:\users\vysha\anaconda3\lib\site-packages (from aiohttp->datasets) (2.0.0)

Requirement already satisfied: attrs>=17.3.0 in c:\users\vysha\anaconda3\lib\site-packages (fr
```

Step2: Loading the imdb dataset from the dataset library. Figure below shows that by running the load_dataset module of dataset library, it downloads all the files of imdb dataset. The resultant imdb_dataset is a dictionary with train, test and unsupervised datasets.



Step 3: Accessing the train data from imdb_dataset dictionary. Figure below shows that we are using our train data for our analysis. This data should have 25000 reviews.



Steps to load the dataset is from: https://huggingface.co/docs/datasets/load hub.

Information about above https://toolbox.google.com/datasetsearch the source: For information, through following more dataset please the link, go http://ai.stanford.edu/~amaas/data/sentiment/

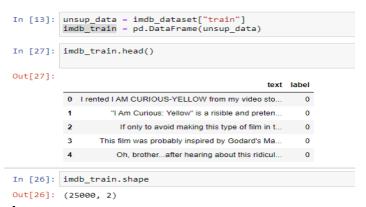
4. Text Classification

<u>Importing the required libraries</u>: Below figure shows the required libraries. Apart from the previous we will be making use of some nltk libraries like stopwords, wordnet for text processing and word cloud to display the categorized review. We will see this in detail in later sections.

```
In [12]: import numpy as np import pandas as pd import pandas as pd import machical population in the import washoring as ns import re from wordcloud import hitk nltk.download(['stopwords', punkt', 'omm-1.4', 'omm-1.4',
```

Loading the dataset and checking the dimensions

In the below screenshot, the train data is loaded and a dataframe is created using the dataframe module of the pandas library. The used dataset have 25000 rows and 2 columns.



Check for class imbalance

The figure below shows the result of the class imbalance. For this we use target column label and count the unique occurrences. We can observe that the class is clearly balanced with equal number of 0 and 1 results.

Text Preprocessing:

The preprocessing function incorporates various steps to clean and prepare the text data for analysis. The following key aspects are covered:

- Lowercasing: Consistent casing for every words to reduce dimensionality.
- Removing special characters: Special symbols, URLs and HTML tags donot contribute much to the sentiment analysis. Hence, we will remove those.
- Removing Punctuations.
- Removing Stopwords of English language.
- Lemitization to reduce words to their root form to achieve standardization.
- Removing emojis as they don't contribute enough for sentiment analysis.

Below figure shows the function defined to achieve this

```
In [69]: # preprocessing function

sw = nltk.corpus.stopwords.words('english') # call stopwords from nltk
lemmatizer = nltk.stem.wordMetlemmatizer() # call Lemmatisation from nltk

# defining the preprocessing function
def preprocess(text):

text = text.lower() #to convert into lowercase

text = re.sub(r"[^a-zA-Z^.!,¿]+", " ", text) # replacing everything with space except (a-z, A-Z, ".", "?", "!", ",")

text = re.sub(r"http\$+", "", text) #Removing URLs

html=re.compile(r'<.*?>')

text = html.sub(r'',text) #Removing html tags

punctuations = '@#!?+&*[]-%.:/();$>><|{^\, ' + "\" + '-'
for p in punctuations:
 text = text.replace(p,'') #Removing punctuations

text = [word.lower() for word in text.split() if word.lower() not in sw] #removing stopwords

text = [lemmatizer.lemmatize(word) for word in text if lemmatizer.lemmatize(word) not in sw]

text = ".join(text) #Lemmatisation

emoji_pattern = re.compile("[" u"\U00001F64F" # emoticons
 u"\U00001F650>\U00001F6FF" # symbols | pictographs
 u"\U00001F650>\U0001F6FF" # symbols | pictographs
 u"\U00002762>\U00001F6FF" # flags
 u"\U000027
```

Next, as shown below we take all the elements in the text column which holds all the review and call the preprocess function to eliminate all those which are not required for our analysis. This will return a new column 'review' that will have the preprocessed text which will look something like this

<u>Text Vectorization:</u> We create a matrix of token counts from a series of text documents using scikit-learn's CountVectorizer. The maximum number of words is limited to 1000 by setting the max_feature option to 1000. This is necessary since, if the parameter is left unset, our dataset generates an enormous matrix of tokens and could result in a memory usage problem. Fit_transform tokenizes a document, each row representing a word, and returns a space matrix. This matrix is finally transformed into a dataframe called "X."

```
In [71]: from sklearn.feature_extraction.text import CountVectorizer
      vectorizer = CountVectorizer(max features=1000)
      X = vectorizer.fit_transform(imdb_train['review'].map(''.join))
      X = pd.DataFrame(X.toarray())
Out[71]:
        0 1 2 3 4 5 6 7 8 9 ... 990 991 992 993 994 995 996 997 998 999
       1 1 0 0 0 0 0 0 0 0 0 ... 0 0
       3 0 0 0 0 0 1 0 1 0 0
                                0
                                   0
                                      0
                                         0
                                            0
       4 0 0 0 0 0 0 0 0 1 ... 0 0 0 3 0 0 0
      5 rows × 1000 columns
```

Model Building: Below figure shows assigning of target variable and splitting the dataset to train and test data.

```
In [72]: from sklearn.model_selection import train_test_split

y = imdb_train['label']

X_train, X_test, y_train, y_test = train_test_split(X,y,train_size = 0.8, test_size = 0.2, random_state = 99)
```

In the below figure we make use of Multinomial Naive Bayes, a popular text classification task. Finally the model.fit() will fit the model to the training data.

Evaluation of result: The model predicts with an accuracy of 84% with 2105 TP and 2072 TN.

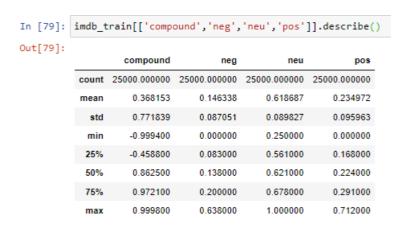
5. Sentiment Analysis

SentimentIntensityAnalyzer is a part of NLTK designed for sentiment analysis using VADER analysis tool. It returns a dictionary sentiment with values like neg, compound, pos and neu.

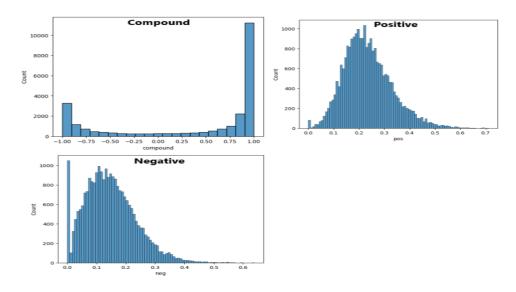
```
In [76]: from nltk.sentiment.vader import SentimentIntensityAnalyzer
sentiment = SentimentIntensityAnalyzer()
```

Below code runs a lambda function for every values and create additional columns in the existing dataframe with sentiment score for each value.

The below five point summary for the compound variable have a positive mean and median value suggesting an overall positive sentiment with less negative sentiment. The maximum value suggest that a very high scoring text have a positive sentiment.



Below graphs show the distribution of compound, positive and negative scores. Positive looks like a normal distribution and negative seems to be right skewed.



6. Evaluation of Result



In conclusion, we successfully distinguished between positive and negative reviews with an amazing 84% accuracy rate in our text classification and sentiment analysis on the IMDb review dataset using the Multinomial Naive Bayes (MNB) model. The model's strong performance is demonstrated by the confusion matrix, which has 2072 true negatives, 2105 true positives, 441 erroneous positives, and 382 false negatives. The model's accuracy in making predictions is supported by its balanced performance for both classes as demonstrated by measures like as precision, recall, and F1-score.

Additionally, the sentiment analysis word cloud provides a visual depiction of the most influential terms in the dataset and includes both positive and negative ratings. This qualitative investigation offers insightful information on the common opinions voiced by reviewers. terms like "Good, brilliant action, excellent, interesting, best, beautiful, funny" are indicative of positive sentiments, whereas terms like "mean, poor, unfortunately, waste time, awful, never, problem" are indicative of negative sentiments.

Part Four: References and Appendices

Mushroom Dataset: https://archive.ics.uci.edu/dataset/73/mushroom

Google Review Dataset: https://archive.ics.uci.edu/dataset/485/tarvel+review+ratings

IMDB Review Dataset: https://ai.stanford.edu/~amaas/data/sentiment/

IMDB Dataset: Maas, A. L., Daly, R. E., Pham, P. T., Huang, D., Ng, A. Y., & Potts, C. (2011, June 1). *Learning Word Vectors for Sentiment Analysis*. ACLWeb; Association for Computational Linguistics. http://www.aclweb.org/anthology/P11-1015