# "Customer Personality Analysis & Predictive Segmentation: A Real-World Application"

This Customer Personality Analysis (CPA) project is an illustration of my real-time work experience. It encapsulates the process and approach I have utilized in professional settings, making it an important asset in my portfolio. The primary objective is to leverage data analysis and machine learning techniques to understand and segment a company's diverse customer base. The ultimate goal is to support the development of targeted marketing strategies and enhance customer engagement.

In the professional project, the workflow is organized into four stages:

- 1. Customer Sentiment Analysis
- 2. Customer Segmentation
- 3. Predictive Model Development for Future Segmentation
- **4.Product Recommendations**

This portfolio project will focus on the second(Customer Segmentation) and third stages(Predictive Model Development for Future Segmentation), demonstrating my proficiency in customer segmentation and developing predictive models for future data segmentation.

# **Introduction of the Project**

My Customer Personality Analysis (CPA) project represents a crucial aspect of modern data-driven decision-making within the business context. Drawing from real-life experiences, this project serves to address the complexities and diversity within a company's customer base. Through machine learning and data analysis techniques, we aim to categorize customers into distinct segments based on their unique behaviors and characteristics. The goal of this project is to facilitate targeted marketing strategies, enhancing customer engagement, and promoting effective product recommendations.

# **Project Stages**

The project will proceed in two major stages, mirroring real-life project execution:

**Customer Segmentation**: This stage involves the use of clustering algorithms to classify customers into different groups based on their distinct attributes and purchasing behaviors. The derived segments will serve as a foundation for tailored marketing strategies and decision-making processes.

**Model Development for Future Data**: Building on the customer segmentation, this stage focuses on developing a predictive model that can handle future data. This model will enable the company to anticipate changes in customer behavior and adapt their strategies accordingly.

# **Business Requirements**

- 1. The core business requirement is to provide an analytical foundation that enables marketing and strategy teams to identify and target specific customer segments. Key business needs include:
- 2.Enabling targeted marketing campaigns by identifying customer segments more likely to respond positively.
- 3.Streamlining product development by understanding the unique needs and preferences of different customer segments.

3.Enhancing customer engagement by delivering personalized experiences based on individual customer segments.

# **Proposed Solution**

My solution focuses on leveraging machine learning techniques to meet the defined business requirements:

**Data Analysis and Clustering**: Conduct extensive data preprocessing and exploratory analysis to identify key features. Use clustering algorithms to create distinct customer segments.

**Predictive Model Development**: Post-segmentation, develop a robust predictive model using machine learning algorithms to anticipate future customer behaviors and preferences.

This solution aims to equip marketing teams with actionable insights that enhance customer satisfaction and drive business performance. By identifying customer segments and predicting future behaviors, companies can create more targeted and effective marketing strategies.

# **Approach**

Our approach to this project follows a structured, step-by-step methodology grounded in data science best practices. Each stage is thoughtfully designed to build upon the previous, ensuring a cohesive and comprehensive solution.

**Uderstanding the Data**: The first step involves a thorough understanding of the dataset, its variables, and its structure. This step is crucial for shaping the subsequent stages of the project.

**Data Preprocessing**: After understanding the dataset, we clean and preprocess the data. This involves handling missing values, potential outliers, and categorical variables, ensuring the data is ready for analysis.

**Exploratory Data Analysis (EDA)**: This stage involves unearthing patterns, spotting anomalies, testing hypotheses, and checking assumptions through visual and quantitative methods. It provides an in-depth understanding of the variables and their interrelationships, which aids in feature selection.

**Feature Selection**: Based on the insights from EDA, relevant features are selected for building the machine learning model. Feature selection is critical to improve the model's performance by eliminating irrelevant or redundant information.

**Customer Segmentation**: The preprocessed data is then fed into a clustering algorithm to group customers into distinct segments based on their attributes and behavior. This segmentation enables targeted marketing and personalized customer engagement.

**Model Development**: Once we have our customer segments, we develop a predictive model using a suitable machine learning algorithm. This model is trained on the current data and then validated using a separate test set.

**Model Evaluation and Optimization**: The model's performance is evaluated using appropriate metrics. If necessary, the model is fine-tuned and optimized to ensure the best possible performance.

**Prediction on Future Data**: The final step involves utilizing the trained model to make predictions on future data. This will allow the business to anticipate changes in customer behavior and adapt their strategies accordingly.

This approach ensures a systematic and thorough analysis of the customer data, leading to robust and reliable customer segments and predictions. It aims to provide a foundation upon which strategic business decisions can be made and future customer trends can be anticipated.

Project Work FLow-Chart

```
import pandas as pd
In [3]:
        import requests
        import numpy as np
        import seaborn as sns
        import matplotlib.pyplot as plt
        import plotly.express as px
        from scipy import stats
        import ydata profiling as pp
        from datetime import datetime
        import warnings
        warnings.filterwarnings('ignore')
```

# **Data Preprocessing**

#### 1.Data Collection

The data collection process in my project involved using the "requests" library to retrieve a CSV file from a specific URL. The content was decoded using UTF-8 encoding, replacing semicolons with commas. The decoded content was saved to a local file. This process ensured the successful acquisition of the necessary data for further analysis and modeling.

```
In [2]: | url = "https://raw.githubusercontent.com/amankharwal/Website-data/master/marketing campa
        filename = "marketing campaign.csv"
        response = requests.get(url)
        response.raise for status() # Check for any errors
        with open(filename, "wb") as file:
             file.write(response.content)
        print("CSV file downloaded successfully!")
        CSV file downloaded successfully!
        df = pd.read csv('marketing campaign.csv')
In [3]:
        df.head()
Out[3]:
           ID; Year_Birth; Education; Marital_Status; Income; Kidhome; Teenhome; Dt_Customer; Recency; MntWines; MntFruits; MntMea
```

```
0
1
2
3
```

4

Data is not separated my (,) comma, it separated by ;(colin)

```
In [4]: # updated code
url = "https://raw.githubusercontent.com/amankharwal/Website-data/master/marketing_campa
filename = "marketing_campaign_updated.csv"

response = requests.get(url)
response.raise_for_status() # Check for any errors

# Decode the content using UTF-8 and replace semicolons with commas
content = response.content.decode("utf-8")

with open(filename, "wb") as file:
    file.write(response.content)

print("CSV file downloaded successfully!")
```

CSV file downloaded successfully!

```
In [5]: #load the data
df = pd.read_csv('marketing_campaign_updated.csv', sep=';')
df.head()
```

Out[5]:		ID	Year_Birth	Education	Marital_Status	Income	Kidhome	Teenhome	Dt_Customer	Recency	MntWines
	0	5524	1957	Graduation	Single	58138.0	0	0	2012-09-04	58	635
	1	2174	1954	Graduation	Single	46344.0	1	1	2014-03-08	38	11
	2	4141	1965	Graduation	Together	71613.0	0	0	2013-08-21	26	426
	3	6182	1984	Graduation	Together	26646.0	1	0	2014-02-10	26	11
	4	5324	1981	PhD	Married	58293.0	1	0	2014-01-19	94	173

5 rows × 29 columns

```
In [6]: #download the data file in repo
    df.to_csv('final_data.csv', index=False)
```

# 2.Data Understanding (Attributes)

#### Attributes:

- 1. ID: Customer's unique identifier.
- 2. Year\_Birth: Customer's birth year.
- 3. Education: Customer's education level, ranging from basic education to a master's degree or higher.
- 4. Marital\_Status: Customer's marital status, including categories such as single, married, divorced, or widowed.
- 5. Income: Customer's yearly household income, representing the total income of all members in the household.
- 6. Kidhome: Number of children in the customer's household.
- 7. Teenhome: Number of teenagers in the customer's household.
- 8. Dt\_Customer: Date of customer's enrollment with the company, indicating when the customer became a registered member.

- 9. Recency: Number of days since the customer's last purchase, providing a measure of the customer's engagement and recent activity.
- 10. Complain: Indicates whether the customer has made a complaint in the last 2 years. (1 if the customer has complained, 0 otherwise)

#### Products:

- 1. MntWines: Amount spent on wine in the last 2 years, reflecting the customer's expenditure on wine products.
- 2. MntFruits: Amount spent on fruits in the last 2 years, representing the customer's expenditure on fruit products.
- 3. MntMeatProducts: Amount spent on meat in the last 2 years, indicating the customer's expenditure on meat products.
- 4. MntFishProducts: Amount spent on fish in the last 2 years, representing the customer's expenditure on fish products.
- 5. MntSweetProducts: Amount spent on sweets in the last 2 years, reflecting the customer's expenditure on sweet products.
- 6. MntGoldProds: Amount spent on gold products in the last 2 years, indicating the customer's expenditure on gold items.

#### Promotion:

- 1. NumDealsPurchases: Number of purchases made with a discount, indicating the customer's engagement with promotional deals.
- 2. AcceptedCmp1: Indicates whether the customer accepted the offer in the 1st campaign. (1 if the customer accepted, 0 otherwise)
- 3. AcceptedCmp2: Indicates whether the customer accepted the offer in the 2nd campaign. (1 if the customer accepted, 0 otherwise)
- 4. AcceptedCmp3: Indicates whether the customer accepted the offer in the 3rd campaign. (1 if the customer accepted, 0 otherwise)
- 5. AcceptedCmp4: Indicates whether the customer accepted the offer in the 4th campaign. (1 if the customer accepted, 0 otherwise)
- 6. AcceptedCmp5: Indicates whether the customer accepted the offer in the 5th campaign. (1 if the customer accepted, 0 otherwise)
- 7. Response: Indicates whether the customer accepted the offer in the last campaign. (1 if the customer accepted, 0 otherwise)

#### Place:

1. NumWebPurchases: Number of purchases made through the company's website, indicating the customer's online purchasing behavior.

- 2. NumCatalogPurchases: Number of purchases made using a catalogue, reflecting the customer's engagement with catalog-based purchases.
- 3. NumStorePurchases: Number of purchases made directly in stores, indicating the customer's offline purchasing behavior.
- 4. NumWebVisitsMonth: Number of visits to the company's website in the last month, representing the customer's online engagement and browsing activity.

Acknowledgement The dataset for this project is provided by Dr. Omar Romero-Hernandez.

Source: https://raw.githubusercontent.com/amankharwal/Website-data/master/marketing\_campaign.csv

```
In [7]: # display max columns

pd.set_option('display.max_columns', 50)

df.head()
```

Out[7]:		ID	Year_Birth	Education	Marital_Status	Income	Kidhome	Teenhome	Dt_Customer	Recency	MntWines
	0	5524	1957	Graduation	Single	58138.0	0	0	2012-09-04	58	635
	1	2174	1954	Graduation	Single	46344.0	1	1	2014-03-08	38	11
	2	4141	1965	Graduation	Together	71613.0	0	0	2013-08-21	26	426
	3	6182	1984	Graduation	Together	26646.0	1	0	2014-02-10	26	11
	4	5324	1981	PhD	Married	58293.0	1	0	2014-01-19	94	173

# 3.Data Inspection

A pandas profiling analysis was conducted on the dataset using the ProfileReport function from the pandas profiling library. The resulting profile report provided a comprehensive summary of the data's structure, statistics, and distributions. The report was saved as an HTML file for reference and further analysis. The profiling analysis facilitated data understanding by revealing patterns, identifying missing values, and offering insights for subsequent data preprocessing and analysis tasks.

```
In []: # ydata profiling is best way to summary of whole data

data_profile = pp.ProfileReport(df)

data_profile.to_file('data_profile.html')

data_profile
```

# 4. Data Cleaning

- 1. Dropped duplicate rows from the DataFrame.
- 2. Dropped rows with missing values (NaN) from the DataFrame.
- 3. Converted the "Dt\_Customer" column to a datetime format.
- 4. Removed the columns "Z\_CostContact" and "Z\_Revenue" as they contained only a single unique value.

5. Identified and counted the number of outliers for each numeric column using the interquartile range (IQR) method.

```
# drop duplicate rows
In [8]:
        df.drop duplicates (inplace=True)
In [9]: df.isnull().sum()
                               0
Out[9]: Year_Birth
                              0
       Education
       Marital_Status
Income
                             0
                            24
       Kidhome
                             0
       Teenhome
                              0
       Dt_Customer
                             0
                             0
       Recency
                             0
       MntWines
       MntFruits
                             0
       MntMeatProducts
MntFishProducts
       MntSweetProducts
       MntGoldProds
       NumDealsPurchases
                             0
       NumWebPurchases
       NumCatalogPurchases0NumStorePurchases0NumWebVisitsMonth0
                             0
       AcceptedCmp3
                             0
       AcceptedCmp4
       AcceptedCmp5
                             0
                             0
       AcceptedCmp1
       AcceptedCmp2
                             0
       Complain
                             0
       Z CostContact
       Z Revenue
                              0
       Response
       dtype: int64
```

To address missing values in the "income" attribute, it is decided to drop the null values to avoid potential distortions in clustering analysis, as income is a crucial feature for grouping the data accurately.

```
9 MntWines 2216 non-null int64
10 MntFruits 2216 non-null int64
           11 MntMeatProducts
                                       2216 non-null int64
           12 MntFishProducts
                                       2216 non-null int64
           13 MntSweetProducts 2216 non-null int64
14 MntGoldProds 2216 non-null int64
           15 NumDealsPurchases 2216 non-null int64
16 NumWebPurchases 2216 non-null int64
           17 NumCatalogPurchases 2216 non-null int64
           18 NumStorePurchases 2216 non-null int64
           19 NumWebVisitsMonth 2216 non-null int64
20 AcceptedCmp3 2216 non-null int64
21 AcceptedCmp4 2216 non-null int64
22 AcceptedCmp5 2216 non-null int64
23 AcceptedCmp1 2216 non-null int64
24 AcceptedCmp2 2216 non-null int64
25 Complain 2216 non-null int64
           26 Z_CostContact
27 Z_Revenue
28 Response
                                       2216 non-null int64
                                       2216 non-null int64
                                       2216 non-null int64
          dtypes: float64(1), int64(25), object(3)
          memory usage: 519.4+ KB
In [12]: # change dt Customer to datetime format
          df['Dt Customer'] = pd.to datetime(df['Dt Customer'])
In [13]: # check unique values in each feature
          df.nunique()
Out[13]: Year_Birth
                                     2216
                                      59
          Education
                                        5
          Marital Status
                                       8
                                    1974
          Income
          Kidhome
                                       3
          Teenhome
                                       3
          Dt Customer
                                     662
                                     100
          Recency
                                    776
          MntWines
                                    158
          MntFruits
          MntMeatProducts
                                    554
                                  182
176
          MntFishProducts
          MntSweetProducts
                                    212
          MntGoldProds
          NumDealsPurchases
                                     15
                                      15
          NumWebPurchases
          NumCatalogPurchases
                                      14
          NumStorePurchases
                                      14
          NumWebVisitsMonth
                                      16
          AcceptedCmp3
                                       2
          AcceptedCmp4
          AcceptedCmp5
          AcceptedCmp1
                                        2
          AcceptedCmp2
          Complain
                                         2
          Z CostContact
                                        1
          Z Revenue
                                         1
          Response
                                         2
          dtype: int64
```

After examining the unique values in each feature of the DataFrame df using the df.nunique() function, it was observed that the columns "Z\_CostContact" and "Z\_Revenue" contain only a single unique value. As

these columns do not contribute to the analysis and model development, it is decided to drop them from the dataset.

```
In [14]: # drop features
df.drop(columns=['Z_CostContact','Z_Revenue'], axis=1, inplace=True)
```

### **Outliers**

Outliers can have a significant impact on clustering models. They can distort the overall distribution and lead to biased cluster assignments. Outliers tend to pull cluster centroids towards them, resulting in less accurate cluster representation and potentially affecting the cluster boundaries. It is important to handle outliers carefully by either removing them or using robust clustering algorithms that are less sensitive to outliers to ensure more accurate and reliable clustering results.

{'Income': 8, 'Kidhome': 0, 'Teenhome': 0, 'Recency': 0, 'MntWines': 35, 'MntFruits': 24 6, 'MntMeatProducts': 174, 'MntFishProducts': 222, 'MntSweetProducts': 246, 'MntGoldProd s': 205, 'NumDealsPurchases': 84, 'NumWebPurchases': 3, 'NumCatalogPurchases': 23, 'NumS torePurchases': 0, 'NumWebVisitsMonth': 8}

The identification of outliers in the dataset prompts further analysis to understand their context and relationship to the data. These outliers may provide insights into unique patterns, market segments, or exceptional customer behavior. It is crucial to investigate their validity and consider their influence carefully during modeling by using robust clustering algorithms or outlier-resistant techniques.

# 5. Feature Engineering & Label Encoding

Feature engineering is a critical step in the process of building a machine learning model, including clustering and analysis. It involves creating new features or modifying existing ones to improve model performance. Good features can capture important aspects of the data that might not be immediately apparent, and can help the model make better predictions or identify more meaningful clusters. In clustering, well-crafted features can help distinguish between different groups in the data more effectively, leading to more accurate and insightful segmentation of customers.

df['Total spend'] = df['MntWines'] + df['MntFishProducts'] + df['MntFruits'] + df['MntMe

```
# total purchase
         df['Total purchase'] = df['NumDealsPurchases'] + df['NumCatalogPurchases'] + df['NumWeb
         # total children in family
         df['Total children'] = df['Kidhome'] + df['Teenhome']
         #total accepted campaigns
         df['accepted camp'] = df['AcceptedCmp1'] +df['AcceptedCmp2'] + df['AcceptedCmp3'] + df['
In [18]: df['Marital_Status'].value counts()
Out[18]: Together
                   857
                   573
         Single
                   471
        Divorced 232
                    76
        Widow
         Alone
                      3
                      2
        Absurd
        YOLO
        Name: Marital Status, dtype: int64
In [19]: # change marital status into two cat Partner and single
         marital status mapping = {
            "Married": "Couple",
             "Together": "Couple",
             "Single": "Single",
             "Divorced": "Single",
             "Widow": "Single",
             "Alone": "Single",
             "Absurd": "Single",
             "YOLO": "Single"
         }
         df['Marital Status'] = df['Marital Status'].map(marital status mapping)
         df['Marital Status'].value counts()
         Couple
                1430
Out[19]:
                   786
         Single
         Name: Marital Status, dtype: int64
In [20]: # create new Feature
         total adults mapping = {
            "Couple": 2,
            "Single": 1
         df['Total adults'] = df['Marital Status'].map(total adults mapping)
         #total family size
         df['Family size'] = df['Total adults'] + df['Total children']
In [21]: # Get the current year
         current year = datetime.now().year
         # Extract the year from 'Dt Customer' and subtract from the current year
         df['Customer Since Years'] = current year - pd.to datetime(df['Dt Customer']).dt.year
In [22]: df['Education'].value counts()
        Graduation
                    1116
Out[22]:
```

#### Summary of above steps

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PhD

- Age Calculation: The age of each customer was calculated by subtracting the year of birth from the current year. This new feature, 'Age', provides information about the age of the customer, which could be a significant factor in their purchasing behavior.
- Total Spend: A new feature 'Total\_spend' was created by adding up the amount spent on different categories of products. This gives a holistic view of the customer's spending habits.
- Total Purchase: The total number of purchases made by each customer across different platforms was calculated and stored in the 'Total\_purchase' feature. This helps in understanding the overall purchasing activity of the customer.
- Family Size: The total number of adults and children in the family was calculated and stored in 'Total\_adults' and 'Total\_children' respectively. A 'Family\_size' feature was also created that sums up these two, providing a complete picture of the family size.
- Marital Status: The 'Marital\_Status' feature was simplified by categorizing the customers into 'Single' and 'Couple'. This simplification can make the data easier to analyze and interpret.
- Customer Since Years: The number of years a customer has been with the company was calculated by subtracting the year of joining from the current year. This 'Customer\_Since\_Years' feature can provide insights into customer loyalty and retention.
- Accepted Campaigns: A new feature 'accepted\_camp' was created that sums up the campaigns
  accepted by each customer. This can provide insights into the customer's responsiveness to marketing
  campaigns.
- Education Encoding: '2n Cycle' was replaced with 'Master' in the 'Education' feature for clarity. The 'Education' feature was then encoded into numerical values, which can be easier for machine learning algorithms to process.

# **EDA**

```
In [24]:
           df.describe()
                                                              Kidhome
                                                                         Teenhome
                                                                                                   MntWines
Out[24]:
                                 Year Birth
                                                   Income
                                                                                        Recency
                                                                                                                MntFruit:
                   2216.000000 2216.000000
                                              2216.000000 2216.000000 2216.000000 2216.000000
                                                                                                  2216.000000 2216.000000
           count
                   5588.353339 1968.820397
                                             52247.251354
                                                              0.441787
                                                                           0.505415
                                                                                       49.012635
                                                                                                  305.091606
                                                                                                                26.356047
           mean
```

std	3249.376275	11.985554	25173.076661	0.536896	0.544181	28.948352	337.327920	39.793917
min	0.000000	1893.000000	1730.000000	0.000000	0.000000	0.000000	0.000000	0.000000
25%	2814.750000	1959.000000	35303.000000	0.000000	0.000000	24.000000	24.000000	2.000000
50%	5458.500000	1970.000000	51381.500000	0.000000	0.000000	49.000000	174.500000	8.000000
75%	8421.750000	1977.000000	68522.000000	1.000000	1.000000	74.000000	505.000000	33.000000
max	11191.000000	1996.000000	666666.000000	2.000000	2.000000	99.000000	1493.000000	199.000000

Upon reviewing the descriptive statistics of the dataset, it's observed that there are significant variances in certain features, notably 'Income' and 'Age'.

The 'Age' feature has a maximum value of 130, which is unusually high and likely indicates the presence of outliers. Similarly, the 'Income' feature shows a substantial difference between the mean and the maximum value, suggesting potential outliers in this feature as well.

To ensure the accuracy of further analysis and modeling, it's crucial to address these outliers. This can be done by either removing the outliers or replacing them with more representative values, such as the median or mean of the feature.

The next step would be to conduct an outlier detection analysis to identify these extreme values and decide on the most appropriate method to handle them. This will help to improve the quality of the dataset and the reliability of subsequent insights derived from the data.

```
In [25]: # visualize outliers

# Box plot for 'Income'
fig1 = px.box(df, x='Income')
fig1.show()

# Box Plot for 'Age'
fig2 = px.box(df, x='Age')
fig2.show()
```

Upon reviewing the box plots for 'Income' and 'Age', it's observed that there are only a few outliers present in the data. Given their minimal presence, these outliers are unlikely to provide meaningful insights and may potentially skew the results of the analysis. Therefore, to maintain the integrity of the dataset and ensure more accurate results, the decision has been made to remove these outliers from the dataset.

```
In [26]: Q1 = df[['Income', 'Age']].quantile(0.25)
    Q3 = df[['Income', 'Age']].quantile(0.75)
    IQR = Q3 - Q1

    df = df[~((df[['Income', 'Age']] < (Q1 - 1.5 * IQR)) | (df[['Income', 'Age']] > (Q3 + 1.)

In [27]: # visualize outliers

# Box plot for 'Income'
fig1 = px.box(df, x='Income')
fig1.show()

# Box Plot for 'Age'
fig2 = px.box(df, x='Age')
fig2.show()
```

# **Dimensionality Reduction**

### **PCA**

Principal Component Analysis (PCA) is a statistical procedure that uses an orthogonal transformation to convert a set of observations of possibly correlated variables into a set of values of linearly uncorrelated variables called principal components.

PCA identifies the axes in the feature space along which the data varies the most. It does this by performing a covariance analysis between factors. In simple terms, it calculates an 'importance' score for each feature of your data, and then it orders these features by their score, giving you the components in order of significance.

This can help to mitigate the curse of dimensionality, improve computational efficiency, and make it easier to visualize the data. The transformed data (principal components) retain most of the variance in the data with fewer dimensions, which can lead to more meaningful and efficient clustering.

The filtered variables were selected based on expert advice from domain knowledge and expertise.

```
In [31]: #Before performing PCA, it is important to standardize the data to bring the values onto
    from sklearn.preprocessing import StandardScaler
    scaler = StandardScaler()

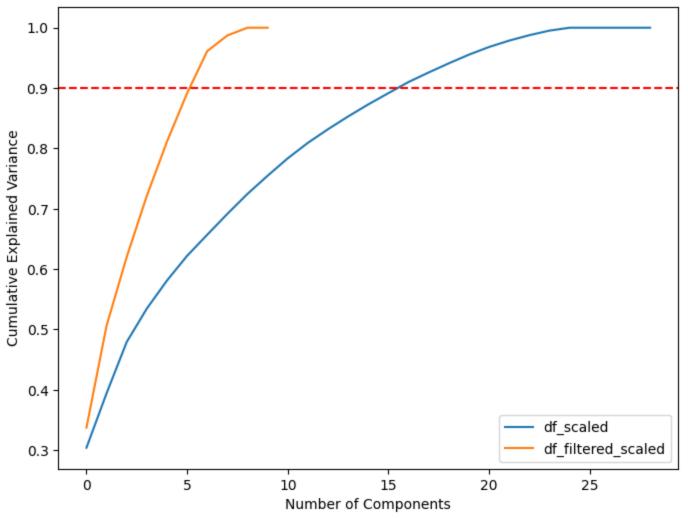
# Standardize the df features
    df_scaled = scaler.fit_transform(df)

# standardize the df_filtered features
    df_filtered_scaled = scaler.fit_transform(df_filtered)
```

# Check best n components for PCA Scree Plot

```
In [33]: # Create two PCA objects
    from sklearn.decomposition import PCA
    pca1 = PCA()
```

```
pca2 = PCA()
# Fit PCA on your data
pca1.fit(df scaled)
pca2.fit(df filtered scaled)
# Calculate cumulative explained variance for df scaled
cumulative explained variance scaled = np.cumsum(pcal.explained variance ratio )
# Calculate cumulative explained variance for df filtered scaled
cumulative explained variance filtered scaled = np.cumsum(pca2.explained variance ratio
# Plot the cumulative explained variance for df scaled
plt.figure(figsize=(8,6))
plt.plot(range(len(cumulative explained variance scaled)), cumulative explained variance
# Plot the cumulative explained variance for df filtered scaled
plt.plot(range(len(cumulative explained variance filtered scaled)), cumulative explained
# Add a horizontal line at 90% cumulative explained variance
plt.axhline(y=0.9, color='r', linestyle='--')
plt.xlabel('Number of Components')
plt.ylabel('Cumulative Explained Variance')
plt.legend()
plt.show()
```



Based on the scree plot for df\_scaled, a variance of 80% to 90% is achieved with 4 or 5 principal components. For df\_filtered\_scaled, a variance of 70% to 80% is achieved with 8 or 9 principal components. In PCA, it is desirable to retain a variance between 70% and 90% to capture a significant portion of the data's information while reducing dimensionality.

```
In [34]: | # perform PCA
         from sklearn.decomposition import PCA
         # Create two PCA objects
         pca1 = PCA(n components=8)
         pca2 = PCA(n components=5)
         # pca for df scaled
         df scaled pca = pcal.fit transform(df scaled)
         #pca for df filtered scaled
         df filtered scaled pca = pca2.fit transform(df filtered scaled)
         # Access the principal components for df scaled
         principal components scaled = pcal.components
         # Access the explained variance ratio for df scaled
         explained variance ratio scaled = pcal.explained variance ratio
         print("Explained Variance Ratio for df scaled:")
         print(explained_variance ratio scaled)
         print()
         # Access the principal components for df filtered scaled
         principal components filtered scaled = pca2.components
         # Access the explained variance ratio for df filtered scaled
         explained variance ratio filtered scaled = pca2.explained variance ratio
         print("Explained Variance Ratio for df filtered scaled:")
         print(explained variance ratio filtered scaled)
        Explained Variance Ratio for df scaled:
         [0.30409483 \ 0.09027176 \ 0.08496384 \ 0.05532585 \ 0.04647201 \ 0.04091918
         0.03495604 0.03451675]
        Explained Variance Ratio for df filtered scaled:
         [0.3373341 0.169114 0.11378058 0.10137992 0.08972968]
```

Looking at the explained variance ratio (EVR) for the df\_scaled dataset, the first component explains 30.4% of the variance, while the remaining seven components account for 39.2%, 8.5%, 5.5%, 4.6%, 4.1%, 3.5%, and 3.5% respectively. This suggests that the first component is contributing a substantial amount of information, but there is still a considerable amount of information distributed across the remaining components.

On the other hand, in the df\_filtered\_scaled dataset, the first component explains a significantly larger proportion of the variance, at 33.7%. The next four components account for 16.9%, 11.4%, 10.1%, and 8.9% respectively. This represents a more substantial proportion of the total variance explained by the first few components, indicating that this dataset may have a lower effective dimensionality.

Based on the above analysis, it would seem that the df\_filtered\_scaled dataset may be a better choice for further analysis or modeling. The reason for this is that a smaller number of components explain a larger proportion of the variance, meaning that we retain more information while reducing dimensionality. In other words, df\_filtered\_scaled is likely a more efficiently compressed representation of our data.

```
pca1 df = pd.DataFrame(data=df scaled pca, columns=[f"PC{i+1}" for i in range(df scaled
          # Create DataFrame for pca2 results
          pca2 df = pd.DataFrame(data=df filtered scaled pca, columns=[f"PC{i+1}" for i in range(d
In [36]:
          pcal df.head()
                 PC1
                           PC2
                                     PC3
                                                        PC5
                                                                  PC6
Out[36]:
                                              PC4
                                                                            PC7
                                                                                     PC8
             4.778575 -0.431022 -0.233758
                                          3.299100 -0.422712 -0.095591
                                                                        0.152742
                                                                                 0.249067
          1 -2.945829
                      0.206513 -0.741823 -0.939268 -0.662957 -0.797031
                                                                       0.577378
                                                                                 0.123228
             1.776005 -1.490874 -0.587736 -0.031005 -0.019317
                                                              0.114025 -0.932182
                                                                                 0.484745
            -2.915123
                      0.291962
                               -1.019391 -0.501065
                                                    0.898932
                                                              0.202037
                                                                      -0.951736
                                                                                 0.237706
            -0.375708 -1.210172
                                0.597005 -1.059922 -0.009606 -0.098461
                                                                       0.544075 -0.898794
In [37]:
          pca2 df.head()
                 PC1
                           PC2
                                     PC3
                                              PC4
                                                        PC5
Out[37]:
             3.109039
                     -0.958371 -0.196529 -1.800031 -0.853496
            -1.810221 -0.257530
                                1.794192 -0.027004 -1.575052
```

0.445606 -1.107744

1.687533 -0.179043

1.099736

1.548891

#### We got four dataframes

-2.013692 -0.876450

1.379175 0.009814 -0.661939

0.735736

1. df scaled

-0.307499

- 2. df\_filtered\_scaled
- 3.  $pca1_df$  ( PCA of df\_scaled data with n = 8)
- 4.  $pca2_df$  (PCA of  $df_filtered_scaled$  data with n = 5)

-0.649741

0.636703

# Stage1

# Clustering

We now have four dataframes: two original dataframes (df\_scaled and df\_filtered\_scaled) and two dimensionality-reduced dataframes using PCA (pca1\_df and pca2\_df). Our next step is to segment customers by implementing various clustering algorithms. Specifically, we will utilize K-Means++, Hierarchical Clustering, and DBSCAN.

Here's a brief overview of these three algorithms:

1. K-Means++: This is an extension of the K-Means clustering algorithm. In standard K-Means, the initial centers are randomly chosen which may lead to sub-optimal solutions. K-Means++ improves upon this by choosing initial centers in a way that accelerates convergence. This means that you get to the optimal solution faster.

- 2. Hierarchical Clustering: This is a type of clustering algorithm where the objective is to create a hierarchy of clusters. It starts with each object in its own cluster and then merges these atomic clusters into larger and larger clusters, until all objects are in a single cluster or a termination condition is met.
- 3. DBSCAN (Density-Based Spatial Clustering of Applications with Noise): This is a density-based clustering algorithm, which groups together points in high-density areas and identifies points in low-density areas as noise. It's different from K-Means and Hierarchical Clustering as it doesn't require the user to specify the number of clusters.

#### K-Means++

In the next step of our analysis, we will be using the Elbow Method and Within-Cluster-Sum-of-Squares (WCSS) to determine the optimal number of clusters for the K-Means++ algorithm.

Here's a brief explanation of these concepts:

- Elbow Method: This method involves running the K-Means algorithm multiple times over a loop, with an increasing number of cluster choice and then plotting a clustering score as a function of the number of clusters. The score could be within-cluster variance, average silhouette, or any other internal clustering validation indices. The optimal number of clusters is usually where the change in the clustering score begins to diminish, often called the "elbow" in the plot.
- Within-Cluster-Sum-of-Squares (WCSS): WCSS is a measure of the compactness of the clusters. It calculates the total distance of data points from their respective cluster centroids. The goal is to minimise this distance, as it means that the data points are closer to their centroids and hence, better clustered. However, as the number of clusters increases, the WCSS will always decrease. That's where the Elbow Method comes in handy, as it helps identify the point at which the rate of decrease sharply shifts, which is a reasonable balance between maximum compression of the data using minimal clusters and the precision of the data distribution.

By using the Elbow method with WCSS, we can determine an appropriate number of clusters for our K-Means++ algorithm without excessive computation. The optimal number of clusters is where we start to get diminishing returns in terms of reducing the WCSS, which is visually represented as an 'elbow' in the plot.

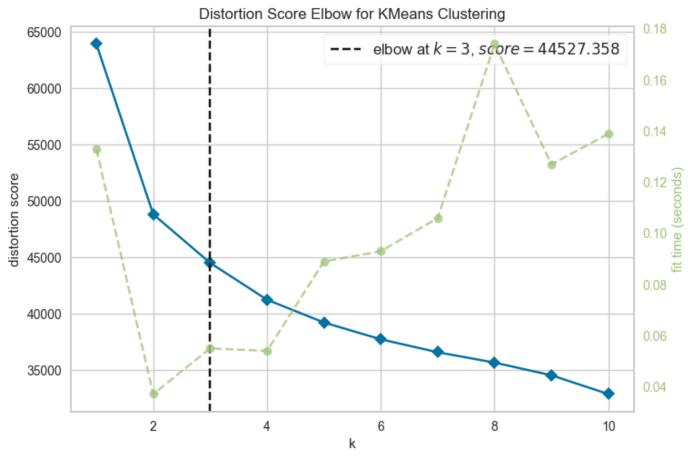
```
In [38]: from sklearn.cluster import KMeans
    from yellowbrick.cluster import KElbowVisualizer

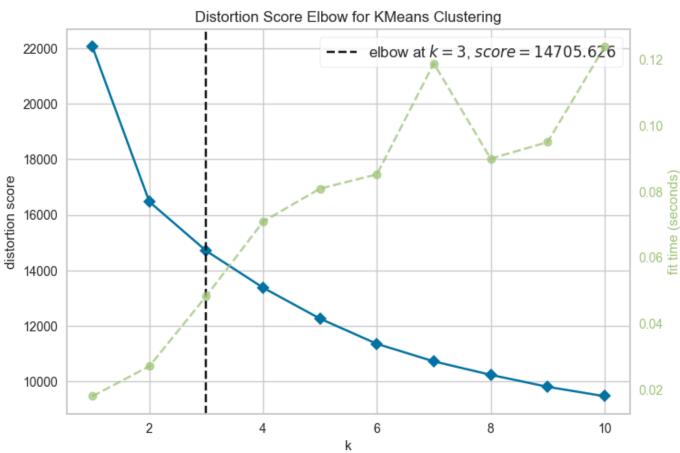
# Create a dictionary with your dataframes
dataframes = {
        'df_scaled': df_scaled,
        'df_filtered_scaled': df_filtered_scaled,
        'pcal_df': pcal_df,
        'pca2_df': pca2_df,
}

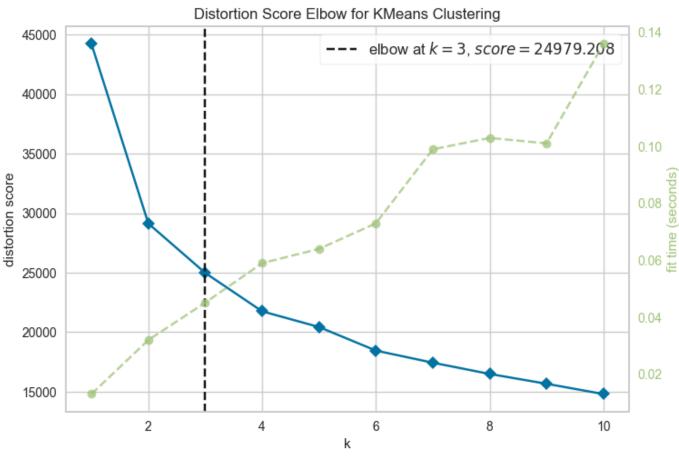
for name, df in dataframes.items():
        model = KMeans(init='k-means++', max_iter=300, n_init=10, random_state=0)
        visualizer = KElbowVisualizer(model, k=(1,11))

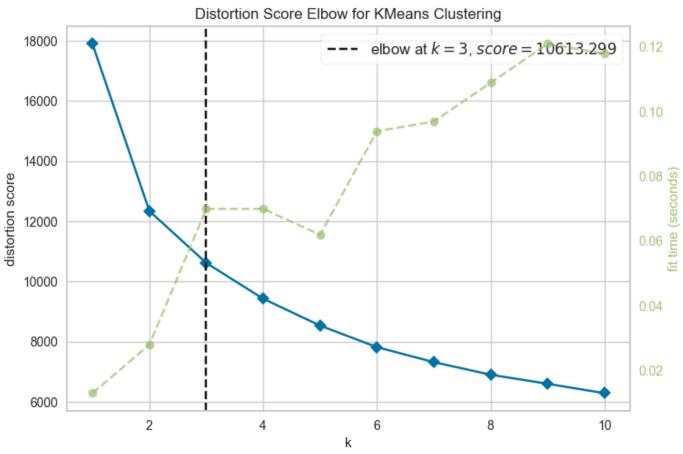
        visualizer.fit(df) # Fit the data to the visualizer

        visualizer.show() # Finalize and render the figure
```









#### Got K 3 for all dataframes

```
In [39]: df_scaled = pd.DataFrame(df_scaled)
    df_filtered_scaled = pd.DataFrame(df_filtered_scaled)
```

In [40]: import plotly.graph\_objects as go

```
from sklearn.cluster import KMeans
dataframes = {
    'df scaled': df scaled,
    'df filtered scaled': df filtered scaled,
    'pcal df': pcal df,
    'pca2 df': pca2 df,
for name, df in dataframes.items():
    model = KMeans(n clusters=3, init='k-means++', max iter=300, n init=10, random state
    y means = model.fit predict(df)
    fig = go.Figure()
    fig.add trace(go.Scatter(x=df.loc[y means == 0, df.columns[0]], y=df.loc[y means ==
                             mode='markers', name='Cluster 1'))
    fig.add trace(go.Scatter(x=df.loc[y means == 1, df.columns[0]], y=df.loc[y means ==
                             mode='markers', name='Cluster 2'))
    fig.add trace(go.Scatter(x=df.loc[y means == 2, df.columns[0]], y=df.loc[y means ==
                             mode='markers', name='Cluster 3'))
    fig.add trace(go.Scatter(x=model.cluster centers [:, 0], y=model.cluster centers [:,
                             mode='markers', name='Centroids', marker=dict(size=12, line
    fig.update layout(title=f'Clusters for {name}',
                      xaxis=dict(title='Feature 1' if 'pca' not in name else 'Principal
                      yaxis=dict(title='Feature 2' if 'pca' not in name else 'Principal
    fig.show()
```

The scatter plots of PCA1 and PCA2 show that the data points are well-distributed into distinct groups, indicating that the PCA transformation effectively captured the underlying structure of the data. In contrast, the scatter plots of the scaled DF and DF\_filtered reveal more noise and lack clear clustering patterns.

Moving forward, it is advisable to focus the analysis on PCA1 and PCA2 to gain further insights and perform subsequent clustering or classification tasks. These transformed components contain the most meaningful information from the original data while reducing noise and dimensionality. This approach allows for a more concise and accurate representation of the data, facilitating more effective analysis and decision-making.

# Validation K-Means++ using Silhouette Score

The silhouette score measures how similar an object is to its own cluster compared to other clusters. The silhouette score ranges from -1 to 1. If it is high, the object is well-matched to its own cluster and poorly matched to neighboring clusters.

```
In [42]: from sklearn.metrics import silhouette_samples
    from sklearn.metrics import silhouette_score

from matplotlib import cm

pca_dfs = {
        'pcal_df': pcal_df,
        'pca2_df': pca2_df,
}

# Define the range of clusters to check
clusters_range = range(2, 5)

for name, df in pca_dfs.items():
```

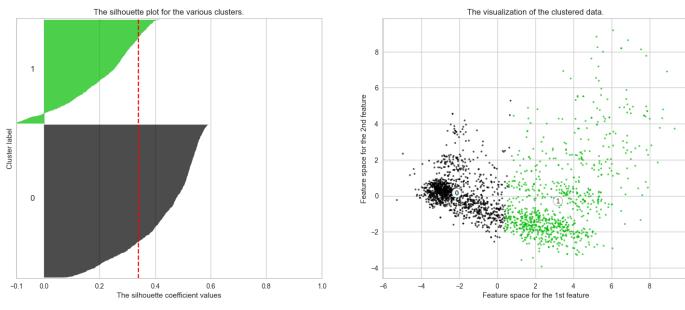
```
for k in clusters range:
   fig, (ax1, ax2) = plt.subplots(1, 2)
   fig.set size inches(18, 7)
    # The 1st subplot is the silhouette plot
   ax1.set xlim([-0.1, 1])
   ax1.set ylim([0, len(df) + (k + 1) * 10])
   model = KMeans(n clusters=k, init='k-means++', max iter=300, n init=10, random s
   labels = model.fit predict(df)
   silhouette avg = silhouette score(df, labels)
   print(f"For {name}, n clusters = {k}, the average silhouette score is : {silhoue
    # Compute the silhouette scores for each sample
   sample silhouette values = silhouette samples(df, labels)
   y lower = 10
   for i in range(k):
       ith cluster silhouette values = \
           sample silhouette values[labels == i]
       ith cluster silhouette values.sort()
        size cluster i = ith cluster silhouette values.shape[0]
       y upper = y lower + size cluster i
       color = cm.nipy spectral(float(i) / k)
       ax1.fill betweenx(np.arange(y lower, y upper),
                          0, ith cluster silhouette values,
                          facecolor=color, edgecolor=color, alpha=0.7)
       ax1.text(-0.05, y lower + 0.5 * size cluster i, str(i))
       y lower = y upper + 10
   ax1.set title("The silhouette plot for the various clusters.")
   ax1.set xlabel("The silhouette coefficient values")
   ax1.set ylabel("Cluster label")
   ax1.axvline(x=silhouette avg, color="red", linestyle="--")
   ax1.set yticks([]) # Clear the yaxis labels / ticks
   ax1.set xticks([-0.1, 0, 0.2, 0.4, 0.6, 0.8, 1])
    # 2nd Plot showing the actual clusters formed
   colors = cm.nipy spectral(labels.astype(float) / k)
   ax2.scatter(df.iloc[:, 0], df.iloc[:, 1], marker='.', s=30, lw=0, alpha=0.7,
                c=colors, edgecolor='k')
    # Draw white circles at cluster centers
   centers = model.cluster centers
   ax2.scatter(centers[:, 0], centers[:, 1], marker='o',
                c="white", alpha=1, s=200, edgecolor='k')
   for i, c in enumerate(centers):
       ax2.scatter(c[0], c[1], marker='$%d$' % i, alpha=1,
                   s=50, edgecolor='k')
   ax2.set title("The visualization of the clustered data.")
   ax2.set xlabel("Feature space for the 1st feature")
   ax2.set ylabel("Feature space for the 2nd feature")
   plt.suptitle(("Silhouette analysis for KMeans clustering on sample data "
                  "with n clusters = %d" % k),
                 fontsize=14, fontweight='bold')
plt.show()
```

For pcal\_df, n\_clusters = 2, the average silhouette\_score is : 0.3390445077860978

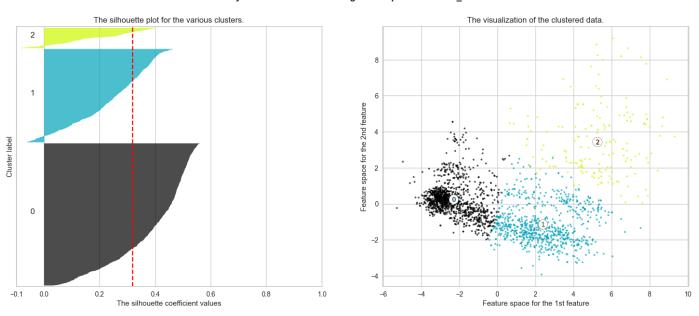
For pcal\_df, n\_clusters = 3, the average silhouette\_score is : 0.31803962197673913

For pcal\_df, n\_clusters = 4, the average silhouette\_score is : 0.2625098126276704

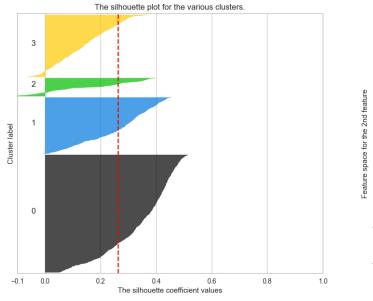
#### Silhouette analysis for KMeans clustering on sample data with n\_clusters = 2

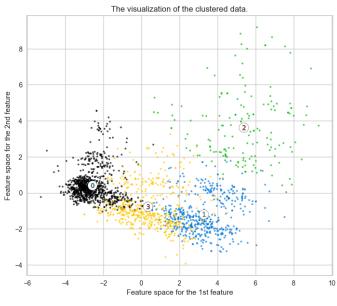


#### Silhouette analysis for KMeans clustering on sample data with n\_clusters = 3



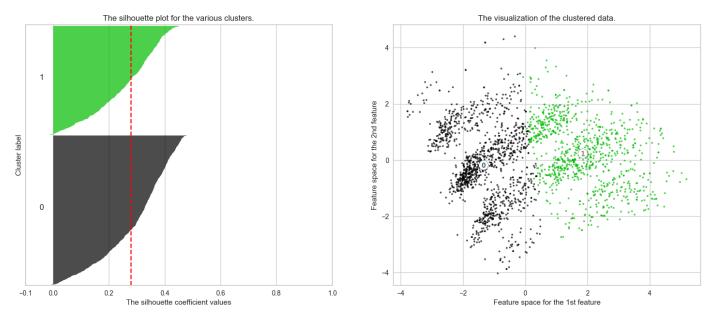
#### Silhouette analysis for KMeans clustering on sample data with n\_clusters = 4



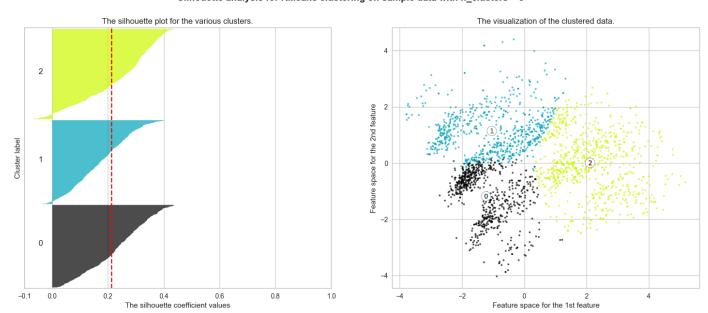


For pca2\_df, n\_clusters = 2, the average silhouette\_score is : 0.27939838848327153
For pca2\_df, n\_clusters = 3, the average silhouette\_score is : 0.21310144685275684
For pca2\_df, n\_clusters = 4, the average silhouette\_score is : 0.1898441675229436

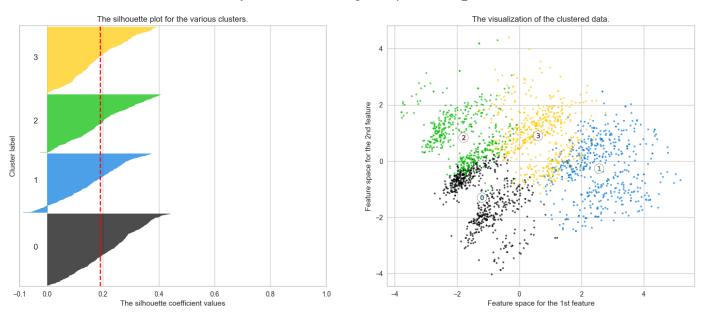
#### Silhouette analysis for KMeans clustering on sample data with n\_clusters = 2



Silhouette analysis for KMeans clustering on sample data with n\_clusters = 3



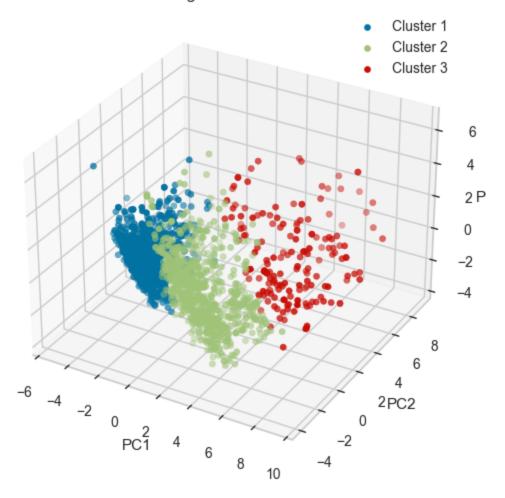
Silhouette analysis for KMeans clustering on sample data with n\_clusters = 4



The clustering analysis using pca1\_df with 3 clusters achieved a moderately good silhouette score of 0.315. This indicates a reasonable level of separation and coherence among the identified clusters. Compared to the alternative pca2\_df clustering solution, which obtained a lower silhouette score of 0.213, pca1\_df demonstrates a stronger clustering structure and better differentiation between the groups. However, it's essential to consider the specific context and domain knowledge when evaluating the quality of the clustering results. Further validation and exploration of the clusters using additional techniques are recommended to ensure the robustness and reliability of the chosen solution.

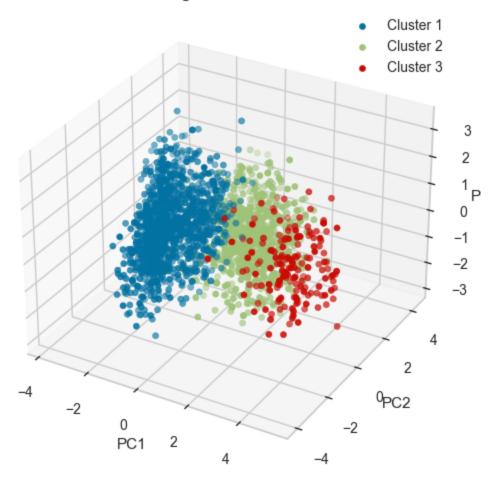
```
In [43]: # final pcal df 3d visual
         from mpl toolkits.mplot3d import Axes3D
         # Performing k-means clustering
         model = KMeans(n clusters=3, init='k-means++', max iter=300, n init=10, random state=0)
         pca1 df['Cluster'] = model.fit predict(pca1 df)
         # Create a 3D plot
         fig = plt.figure(figsize=(8, 6))
         ax = fig.add subplot(111, projection='3d')
         # Scatter plot for each cluster
         for cluster in range(3):
             cluster data = pca1 df[pca1 df['Cluster'] == cluster]
             ax.scatter(cluster data['PC1'], cluster data['PC2'], cluster data['PC3'], label=f'C1
         # Set labels and title
         ax.set xlabel('PC1')
         ax.set ylabel('PC2')
         ax.set zlabel('PC3')
         ax.set title('Clustering Visualization in 3D')
         # Add legend
         ax.legend()
         # Show the plot
         plt.show()
```

### Clustering Visualization in 3D



```
In [44]: # final pca2 df 3d visual
         # Performing k-means clustering
         model = KMeans(n clusters=3, init='k-means++', max iter=300, n init=10, random state=0)
         pca2 df['Cluster'] = model.fit predict(pca2 df)
         # Create a 3D plot
         fig = plt.figure(figsize=(8, 6))
         ax = fig.add subplot(111, projection='3d')
         # Scatter plot for each cluster
         for cluster in range(3):
             cluster data = pca2 df[pca1 df['Cluster'] == cluster]
             ax.scatter(cluster data['PC1'], cluster data['PC2'], cluster data['PC3'], label=f'C1
         # Set labels and title
         ax.set xlabel('PC1')
         ax.set ylabel('PC2')
         ax.set_zlabel('PC3')
         ax.set title('Clustering Visualization in 3D')
         # Add legend
         ax.legend()
         # Show the plot
         plt.show()
```

#### Clustering Visualization in 3D



```
In [45]:
        from plotly.subplots import make subplots
         # Counting the number of customers in each cluster for pcal df
        pca1 df counts = pca1 df['Cluster'].value counts().reset index()
        pca1 df counts.columns = ['Cluster', 'Count']
        pcal df counts = pcal df counts.sort values('Cluster')
         # Counting the number of customers in each cluster for pca2 df
        pca2 df counts = pca2 df['Cluster'].value counts().reset index()
        pca2 df counts.columns = ['Cluster', 'Count']
        pca2 df counts = pca2 df counts.sort values('Cluster')
         # Creating subplot figure
         fig = make subplots(rows=1, cols=2, subplot titles=("pca1 df", "pca2 df"))
         # Adding bar plot for pcal df
         fig.add trace(go.Bar(x=pca1 df counts['Cluster'], y=pca1 df counts['Count'], name="Count
                              textposition='auto'), row=1, col=1)
         # Adding bar plot for pca2 df
         fig.add trace(go.Bar(x=pca2 df counts['Cluster'], y=pca2 df counts['Count'], name="Count
                              textposition='auto'), row=1, col=2)
         # Update layout and axis labels
         fig.update layout(title="Distribution of Customers in Clusters",
                           xaxis=dict(title="Cluster"),
                           yaxis=dict(title="Count"),
                           showlegend=False)
         # Show the plot
         fig.show()
```

# **Hierarchical Clustering**

# Print the cluster labels

print(cluster labels1)

print("Cluster labels for pca1 df:")

Hierarchical clustering is a clustering algorithm that seeks to build a hierarchy of clusters. It does not require a predefined number of clusters but instead forms clusters by recursively merging or splitting them based on the similarity between data points.

Here are the steps involved in performing hierarchical clustering:

- Calculate the pairwise distance or similarity matrix between data points.
- Each data point is initially considered as a separate cluster.
- Merge the two closest clusters based on a distance metric (e.g., Euclidean distance, Manhattan distance) to form a new cluster.
- Update the similarity matrix with the distances between the new cluster and the remaining clusters.
- Repeat steps 3 and 4 until all data points are merged into a single cluster or until a desired number of clusters is reached.
- The result is a dendrogram, which is a tree-like structure representing the hierarchical clustering.

```
pcal df.head()
In [46]:
                 PC1
                                   PC3
                                                      PC5
                                                                        PC7
Out[46]:
                          PC2
                                            PC4
                                                               PC6
                                                                                 PC8
                                                                                     Cluster
             4.778575 -0.431022 -0.233758
                                         3.299100 -0.422712 -0.095591
                                                                    0.152742
                                                                             0.249067
                                                                                           1
            0.577378
                                                                              0.123228
             1.776005 -1.490874 -0.587736 -0.031005 -0.019317
                                                           0.114025 -0.932182
                                                                             0.484745
         3 -2.915123 0.291962 -1.019391 -0.501065
                                                  0.898932
                                                           0.202037 -0.951736
                                                                              0.237706
           -0.375708 -1.210172 0.597005 -1.059922 -0.009606 -0.098461
                                                                                           0
                                                                    0.544075 -0.898794
In [47]:
         pca2 df.head()
Out[47]:
                 PC1
                          PC2
                                   PC3
                                            PC4
                                                      PC5 Cluster
             3.109039 -0.958371 -0.196529 -1.800031 -0.853496
                                                               2
         1 -1.810221 -0.257530
                               1.794192 -0.027004 -1.575052
            1.379175 0.009814 -0.661939
                                        0.445606 -1.107744
                                                               2
           -2.013692 -0.876450 -0.649741
                                         1.687533 -0.179043
           -0.307499 0.735736 0.636703
                                                               1
                                         1.548891
                                                  1.099736
In [48]:
          #drop cluster feature from pcal df and pcal df
```

```
pca1_df.drop(columns=['Cluster'], inplace=True)
pca2_df.drop(columns=['Cluster'], inplace=True)

In [49]: from sklearn.cluster import AgglomerativeClustering

# Perform hierarchical clustering on pca1_df
model1 = AgglomerativeClustering(n_clusters=3)
cluster_labels1 = model1.fit_predict(pca1_df)
```

```
# Perform hierarchical clustering on pca2 df
        model2 = AgglomerativeClustering(n clusters=3)
         cluster labels2 = model2.fit predict(pca2 df)
         # Print the cluster labels
        print("Cluster labels for pca2 df:")
        print(cluster labels2)
        Cluster labels for pcal df:
        [0 1 0 ... 0 2 2]
        Cluster labels for pca2 df:
         [2 0 1 ... 2 1 0]
In [50]: dataframes = {
            'pcal df': pcal df,
             'pca2 df': pca2 df,
        for name, df in dataframes.items():
            model = AgglomerativeClustering(n clusters=3)
            cluster labels = model.fit predict(df)
            fig = go.Figure()
             fig.add trace(go.Scatter(x=df.iloc[cluster labels == 0, 0], y=df.iloc[cluster labels
                                     mode='markers', name='Cluster 1'))
            fig.add trace(go.Scatter(x=df.iloc[cluster labels == 1, 0], y=df.iloc[cluster labels
                                     mode='markers', name='Cluster 2'))
             fig.add trace(go.Scatter(x=df.iloc[cluster labels == 2, 0], y=df.iloc[cluster labels
                                     mode='markers', name='Cluster 3'))
             fig.update layout(title=f'Clusters for {name}',
                              xaxis=dict(title='Principal Component 1'),
                               yaxis=dict(title='Principal Component 2'))
             fig.show()
```

Based on the observation of the scatter plots, it appears that the clusters created by hierarchical clustering showed significant overlap. This suggests that the algorithm was not able to generate well-separated and distinct clusters for this particular dataset. Consequently, the effectiveness of hierarchical clustering in creating meaningful clusters for this dataset is questionable. Further analysis and alternative clustering algorithms may be necessary to achieve better cluster separation and differentiation.

# DBSCAN (Density-Based Spatial Clustering of Applications with Noise)

- DBSCAN is a density-based clustering algorithm that groups together data points based on their density in the feature space.
- It does not require specifying the number of clusters in advance and can discover clusters of arbitrary shape.
- DBSCAN defines clusters as dense regions separated by sparser regions of low-density data points.
- It works by defining a neighborhood around each data point and expanding the clusters by connecting neighboring points that have a sufficient density.
- Outliers or noise points that do not belong to any cluster are identified as points with low-density neighborhoods.

• DBSCAN's performance is influenced by two key parameters: epsilon (ε), which defines the radius of the neighborhood, and min\_samples, which sets the minimum number of points required to form a dense region.

```
In [51]: from sklearn.cluster import DBSCAN
         # Perform DBSCAN clustering on pcal df
         dbscan1 = DBSCAN(eps=0.5, min samples=5)
         cluster labels1 dbscan = dbscan1.fit predict(pca1 df)
         # Print the cluster labels
         print("DBSCAN cluster labels for pca1 df:")
         print(cluster labels1 dbscan)
         # Perform DBSCAN clustering on pca2 df
         dbscan2 = DBSCAN(eps=0.5, min samples=5)
         cluster labels2 dbscan = dbscan2.fit predict(pca2 df)
         # Print the cluster labels
         print("DBSCAN cluster labels for pca2 df:")
         print(cluster labels2 dbscan)
         DBSCAN cluster labels for pcal df:
         [-1 \ -1 \ -1 \ \dots \ -1 \ -1 \ -1]
         DBSCAN cluster labels for pca2 df:
         [-1 \ -1 \ 0 \ \dots \ -1 \ -1 \ -1]
In [52]: from sklearn.cluster import DBSCAN
         # Create an instance of DBSCAN to create non-spherical clusters based on data density
         # For pcal df
         db1 = DBSCAN(eps=0.726, min samples=26)
         # Fit to the dimensionality reduced dataset
         db1.fit(pca1 df)
         # Identify the clusters
         clusters1 = db1.labels
         # Display metrics/sample
         n clusters1 = len(set(clusters1)) - (1 if -1 in clusters1 else 0)
         n noise1 = list(clusters1).count(-1)
         print('Cluster Predictions for pca1 df')
         print('----')
         print("Number of clusters: %d" % n clusters1 )
         print("Number of noise points: %d" % n noise1)
         print('Number of points per cluster:')
         for i in range(n clusters1):
             print('Cluster', i, ':', len(clusters1[clusters1==i]))
         # Repeat the process for pca2 df
         db2 = DBSCAN(eps=0.726, min samples=26)
         # Fit to the dimensionality reduced dataset
         db2.fit(pca2 df)
         # Identify the clusters
         clusters2 = db2.labels
         # Display metrics/sample
         n_clusters2_ = len(set(clusters2)) - (1 if -1 in clusters2 else 0)
         n noise2 = list(clusters2).count(-1)
```

```
print('\nCluster Predictions for pca2 df')
print('----')
print("Number of clusters: %d" % n clusters2 )
print("Number of noise points: %d" % n noise2 )
print('Number of points per cluster:')
for i in range(n clusters2):
   print('Cluster', i, ':', len(clusters2[clusters2==i]))
Cluster Predictions for pcal df
_____
Number of clusters: 0
Number of noise points: 2205
Number of points per cluster:
Cluster Predictions for pca2 df
_____
Number of clusters: 5
Number of noise points: 1904
Number of points per cluster:
Cluster 0 : 92
Cluster 1: 54
Cluster 2 : 38
Cluster 3:81
Cluster 4: 36
```

#### **Hyperparamater Tunning**

#### Check Best eps and min\_sample

Best min samples: 20

```
In [53]: from sklearn.metrics import silhouette score
         # Define the parameter space
         eps values = [0.5, 0.6, 0.7, 0.8, 0.9, 1.0]
         min samples values = [5, 10, 15, 20, 25, 30]
         # Initialize variables to store the best parameters
        best eps = None
         best min samples = None
        best silhouette = -1
         # Perform the grid search
         for eps in eps values:
             for min samples in min samples values:
                 db = DBSCAN(eps=eps, min samples=min samples)
                 clusters = db.fit predict(pca1 df)
                 # Check if more than one cluster is created
                 if len(set(clusters)) > 1:
                     silhouette = silhouette score(pca1 df, clusters)
                     if silhouette > best silhouette:
                         best eps = eps
                         best min samples = min samples
                         best silhouette = silhouette
         print("Best eps:", best eps)
         print("Best min samples:", best min samples)
        Best eps: 1.0
```

```
In [54]: # Define the parameter space
    eps_values = [0.5, 0.6, 0.7, 0.8, 0.9, 1.0]
    min_samples_values = [5, 10, 15, 20, 25, 30]
```

```
# Initialize variables to store the best parameters
best eps = None
best min samples = None
best silhouette = -1
# Perform the grid search
for eps in eps values:
    for min samples in min samples values:
        db = DBSCAN(eps=eps, min samples=min samples)
        clusters = db.fit predict(pca2 df)
        # Check if more than one cluster is created
        if len(set(clusters)) > 1:
            silhouette = silhouette score(pca2 df, clusters)
            if silhouette > best silhouette:
               best eps = eps
                best min samples = min samples
                best silhouette = silhouette
print("Best eps:", best eps)
print("Best min samples:", best min samples)
Best eps: 1.0
Best min samples: 5
Remodel according to above values
```

```
In [55]: # Create an instance of DBSCAN to create non-spherical clusters based on data density
        # For pcal df
        db1 = DBSCAN(eps=1, min samples=20)
        # Fit to the dimensionality reduced dataset
        db1.fit(pca1 df)
        # Identify the clusters
        clusters1 = db1.labels
        # Display metrics/sample
        n clusters1 = len(set(clusters1)) - (1 if -1 in clusters1 else 0)
        n noise1 = list(clusters1).count(-1)
        print('Cluster Predictions for pcal df')
        print('----')
        print("Number of clusters: %d" % n clusters1 )
        print("Number of noise points: %d" % n noise1)
        print('Number of points per cluster:')
        for i in range(n clusters1):
            print('Cluster', i, ':', len(clusters1[clusters1==i]))
        Cluster Predictions for pcal df
        _____
        Number of clusters: 3
        Number of noise points: 1814
        Number of points per cluster:
        Cluster 0: 244
        Cluster 1: 76
        Cluster 2 : 71
In [ ]: # Create an instance of DBSCAN to create non-spherical clusters based on data density
        # For pcal df
```

db1 = DBSCAN(eps=1, min samples=5)

db1.fit(pca2 df)

# Fit to the dimensionality reduced dataset

```
# Identify the clusters
clusters1 = db1.labels_

# Display metrics/sample
n_clusters1_ = len(set(clusters1)) - (1 if -1 in clusters1 else 0)
n_noise1_ = list(clusters1).count(-1)

print('Cluster Predictions for pcal_df')
print('------')
print("Number of clusters: %d" % n_clusters1_)
print("Number of noise points: %d" % n_noise1_)
print('Number of points per cluster:')
for i in range(n_clusters1_):
    print('Cluster', i, ':', len(clusters1[clusters1==i]))
```

Based on these results, using DBSCAN with the current parameter settings is not recommended for these datasets. The parameters, especially epsilon and the minimum sample count, should be fine-tuned based on the characteristics of the data. If tuning the parameters doesn't improve the results, it might be more appropriate to use a different clustering algorithm that is more suited to the data's structure. Alternative algorithms to consider could include K-means, Hierarchical Clustering, or Gaussian Mixture Models.

# **Conclusive Report on Clustering Analysis**

In the course of our exploratory data analysis, we have applied various clustering algorithms, specifically K-means++, Hierarchical Clustering, and DBSCAN, to our datasets: pca1\_df and pca2\_df. These dataframes represent two distinct sets of principal components derived from our original dataset.

After an exhaustive and iterative process of model tuning and evaluation, we found that the K-means++ algorithm performed the most effectively in revealing the underlying structure of our data. The clusters formed by the K-means++ algorithm were more cohesive and better separated than those generated by the Hierarchical Clustering and DBSCAN algorithms.

To further explore and validate our clustering results from the K-means++ model, we have conducted an indepth exploratory data analysis (EDA) on the formed clusters for both pca1\_df and pca2\_df. This process included assessing the distributions, central tendencies, and dispersions of our data within and across the identified clusters.

We also employed visual analytics, creating an interactive dashboard using Power BI. This step greatly facilitated the presentation of our clustering results, enabling us to interactively explore each cluster and their respective properties. This dashboard offered a visually intuitive understanding of the cluster formations and their characteristics, providing significant insights into the structure of our data.

Based on the insights from our EDA and Power BI visualization, we were able to evaluate and compare the clustering results between pca1\_df and pca2\_df. These evaluations will allow us to ascertain which set of principal components, and consequently, which feature space better represents the structure of our original data.

In conclusion, the application of K-means++ has yielded valuable insights into the hidden structure of our datasets. This, combined with the subsequent EDA and Power BI dashboard creation, has provided us with a comprehensive understanding of our data and will be instrumental in guiding our future data-driven decision-making processes.

# **Feature Engineering**

As a part of our feature engineering strategy, we have enriched the original datasets df1 and df2 by incorporating the cluster assignments from our K-means++ algorithm applied to the respective PCA-transformed dataframes, pca1\_df and pca2\_df. This additional 'Cluster' feature serves as an indicator of the underlying group structure detected in our high-dimensional datasets, thus providing a valuable source of information for further exploratory data analysis.

Subsequently, these enhanced datasets will be visualized in a Power BI dashboard, enabling us to interactively investigate the patterns and relationships within and between the identified clusters. Additionally, these enriched datasets will be employed to develop predictive models, leveraging the newly added 'Cluster' feature to enhance the model's ability to recognize and adapt to the inherent groupings in future data.

This method illustrates a cohesive blend of unsupervised learning (clustering) and dimensionality reduction (PCA) techniques that not only assists in revealing the intrinsic structure of our data, but also adds a new dimension for data exploration and modeling.

```
# Performing k-means clustering
In [57]:
         model = KMeans(n clusters=3, init='k-means++', max iter=300, n init=10, random state=0)
         pca1 df['Cluster'] = model.fit predict(pca1 df)
         # Store the cluster labels
         cluster labels1 = pca1 df['Cluster'].values
         # Add cluster labels to the original dataframe df1
         df1['Cluster'] = cluster labels1
         # Repeat the same for pca2 df and df2
         model = KMeans(n clusters=3, init='k-means++', max iter=300, n init=10, random state=0)
         pca2 df['Cluster'] = model.fit predict(pca2 df)
         # Store the cluster labels
         cluster labels2 = pca2 df['Cluster'].values
         # Add cluster labels to the original dataframe df2
         df2['Cluster'] = cluster labels2
In [58]: # save in csv file
         df1.to csv('group1 pca1.csv', index label=False)
         df2.to csv('group2 pca2.csv', index label=False)
```

# **Evaluation (EDA)**

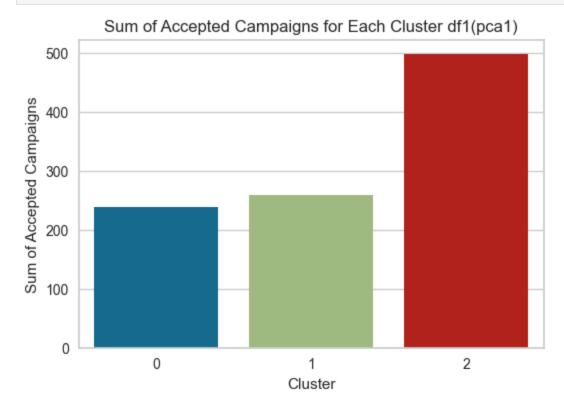
#### Sum of accepted campaigns by groups

```
In [59]: #df1(pcal group)

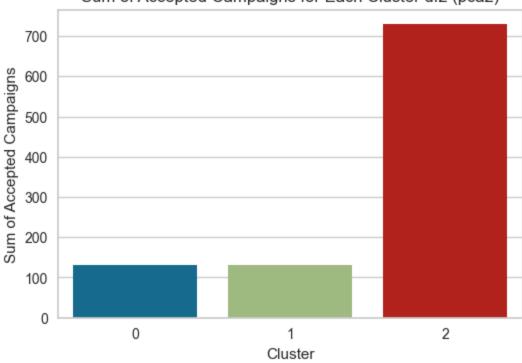
plt.figure(figsize=(6, 4))
    sns.barplot(x="Cluster", y="accepted_camp", data=df1, estimator=sum, ci=None)
    plt.title('Sum of Accepted Campaigns for Each Cluster df1(pca1)')
    plt.xlabel('Cluster')
    plt.ylabel('Sum of Accepted Campaigns')
    plt.show()

#df2(pca2 group)

plt.figure(figsize=(6, 4))
    sns.barplot(x="Cluster", y="accepted_camp", data=df2, estimator=sum, ci=None)
    plt.title('Sum of Accepted Campaigns for Each Cluster df2 (pca2)')
    plt.xlabel('Cluster')
    plt.ylabel('Sum of Accepted Campaigns')
    plt.show()
```



### Sum of Accepted Campaigns for Each Cluster df2 (pca2)



### **Income VS Spent by Clusters**

```
color_discrete_sequence=px.colors.qualitative.D3) # Use a discrete colo
fig.show()
#Create a scatter plot df2
fig = px.scatter(df2,
                 x="Income",
                 y="Total spend",
                 color='Cluster', # Different clusters will have different colors
                 hover data=['Cluster'], # This will show cluster number on hover
                 title="Income vs Spending by Cluster Group df2(pca2)",
                 labels={
                     "Income": "Income",
                     "Total spend": "Total Spending",
                     "Cluster": "Cluster Group"
                 },
                 color discrete sequence=px.colors.qualitative.D3) # Use a discrete colo
fig.show()
```

### Cluster groups by Total Spend

```
# Boxplot for 'Total spend' across different clusters in df1
In [61]:
         fig = px.box(df1,
                      x='Cluster',
                      y='Total spend',
                      color='Cluster',
                      labels={'Total spend': 'Total Spending', 'Cluster': 'Cluster Group'},
                      title='Distribution of Total Spending by Cluster Group df1(pca1)',
                      color discrete sequence=px.colors.qualitative.D3) # Use a discrete color se
         fig.show()
         # Boxplot for 'Total spend' across different clusters in df2
         fig = px.box(df2,
                      x='Cluster',
                     y='Total spend',
                      color='Cluster',
                      labels={'Total_spend': 'Total Spending', 'Cluster': 'Cluster Group'},
                      title='Distribution of Total Spending by Cluster Group df2(pca2)',
                      color discrete sequence=px.colors.qualitative.D3) # Use a discrete color se
         fig.show()
```

# Family Size, Education, Total children vs Group

```
In [62]: fig, axs = plt.subplots(nrows=3, figsize=(10, 10))

# Plot 'Education' vs 'Cluster'
df1.groupby(['Education', 'Cluster']).size().unstack().plot(kind='bar', ax=axs[0])
axs[0].set_title('Education by Cluster Group pcal')

# Plot 'Total_children' vs 'Cluster'
df1.groupby(['Total_children', 'Cluster']).size().unstack().plot(kind='bar', ax=axs[1])
axs[1].set_title('Total Children by Cluster Group pcal')

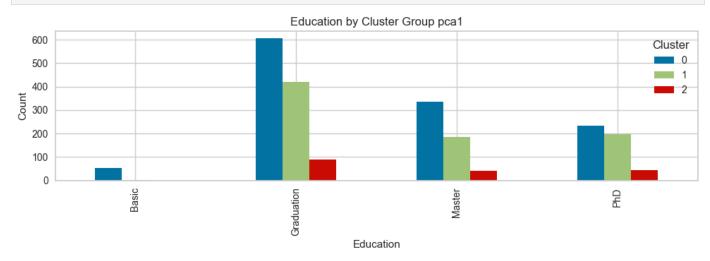
# Plot 'Family_size' vs 'Cluster'
df1.groupby(['Family_size', 'Cluster']).size().unstack().plot(kind='bar', ax=axs[2])
axs[2].set_title('Family size by Cluster Group pcal')
```

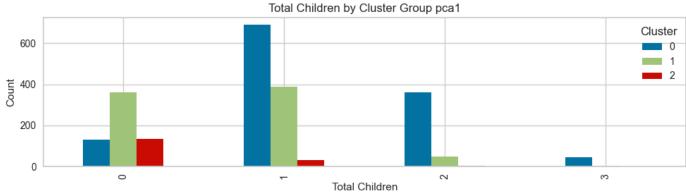
```
# Set x-axis labels
axs[0].set_xlabel('Education')
axs[1].set_xlabel('Total Children')
axs[2].set_xlabel('Family Size')

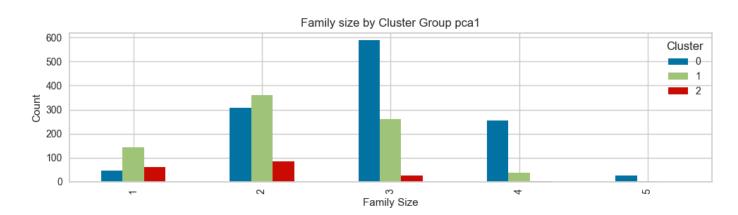
# Set y-axis labels
axs[0].set_ylabel('Count')
axs[1].set_ylabel('Count')
axs[2].set_ylabel('Count')

# Adjust spacing between subplots
plt.tight_layout()

# Show the plot
plt.show()
```



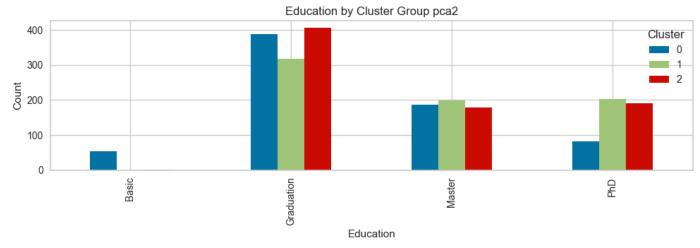


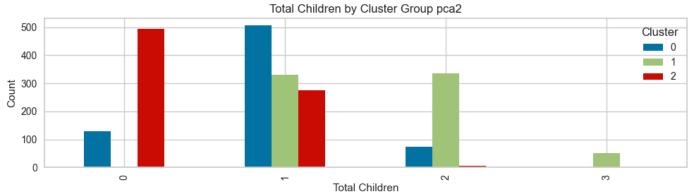


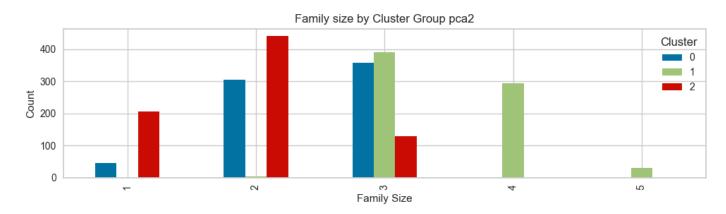
```
In [63]: fig, axs = plt.subplots(nrows=3, figsize=(10, 10))

# Plot 'Education' vs 'Cluster'
df2.groupby(['Education', 'Cluster']).size().unstack().plot(kind='bar', ax=axs[0])
axs[0].set_title('Education by Cluster Group pca2')
```

```
# Plot 'Total children' vs 'Cluster'
df2.groupby(['Total children', 'Cluster']).size().unstack().plot(kind='bar', ax=axs[1])
axs[1].set title('Total Children by Cluster Group pca2')
# Plot 'Family size' vs 'Cluster'
df2.groupby(['Family size', 'Cluster']).size().unstack().plot(kind='bar', ax=axs[2])
axs[2].set title('Family size by Cluster Group pca2')
# Set x-axis labels
axs[0].set xlabel('Education')
axs[1].set xlabel('Total Children')
axs[2].set xlabel('Family Size')
# Set y-axis labels
axs[0].set_ylabel('Count')
axs[1].set ylabel('Count')
axs[2].set ylabel('Count')
# Adjust spacing between subplots
plt.tight layout()
# Show the plot
plt.show()
```

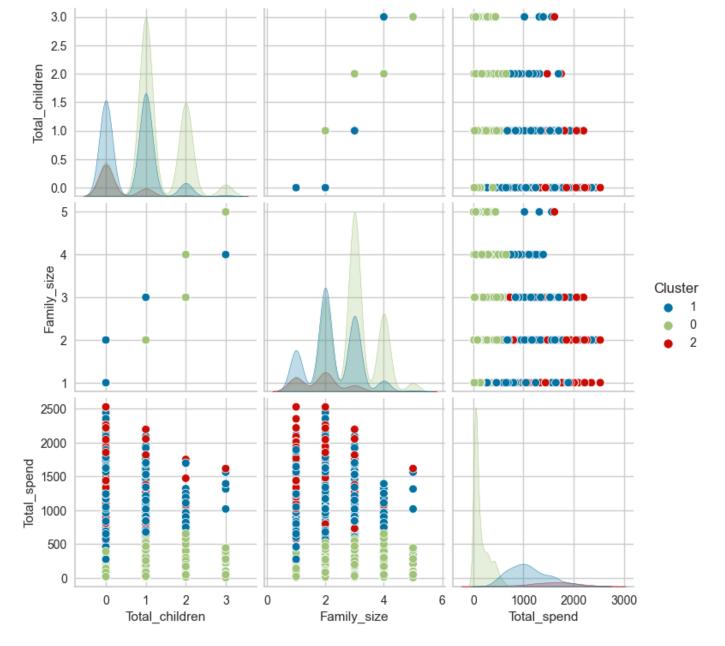






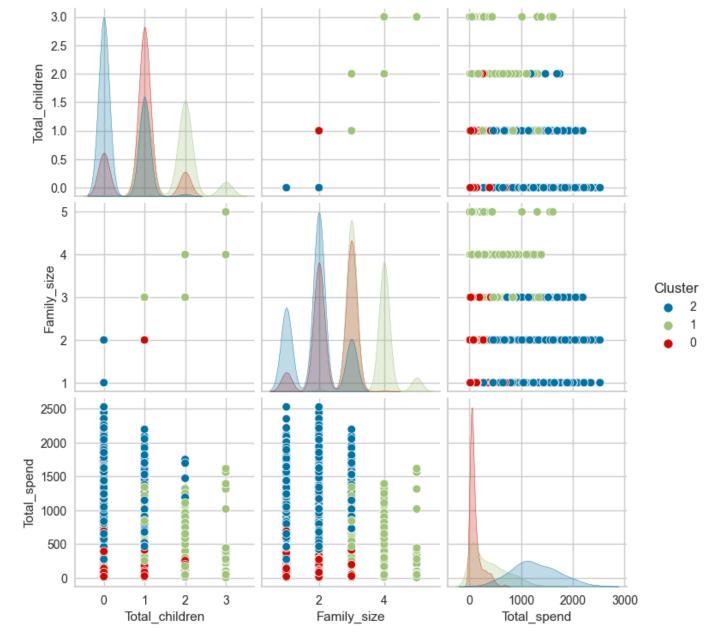
```
In [64]: # Subset of features df1
df_subset = df1[['Education', 'Total_children', 'Family_size', 'Total_spend', 'Cluster']
# Convert 'Cluster' to string type so that pairplot treats it as a categorical variable
df_subset['Cluster'] = df_subset['Cluster'].astype(str)

# Pairplot
sns.pairplot(df_subset, hue='Cluster')
plt.show()
```



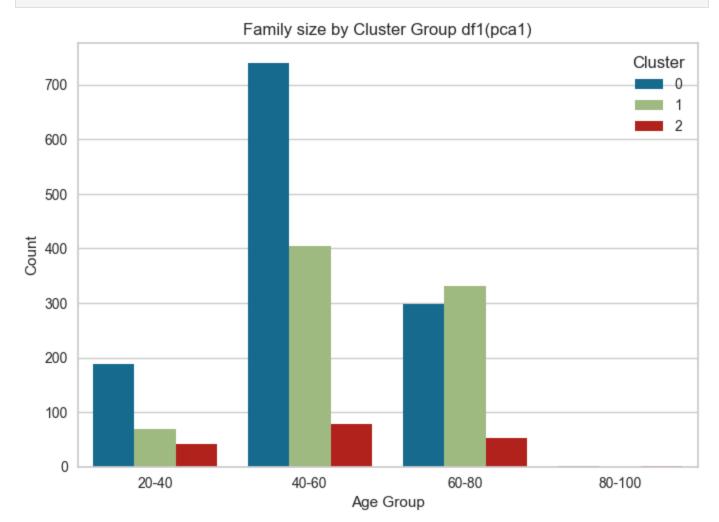
```
In [65]: # Subset of features df2
df_subset = df2[['Education', 'Total_children', 'Family_size', 'Total_spend', 'Cluster']
# Convert 'Cluster' to string type so that pairplot treats it as a categorical variable
df_subset['Cluster'] = df_subset['Cluster'].astype(str)

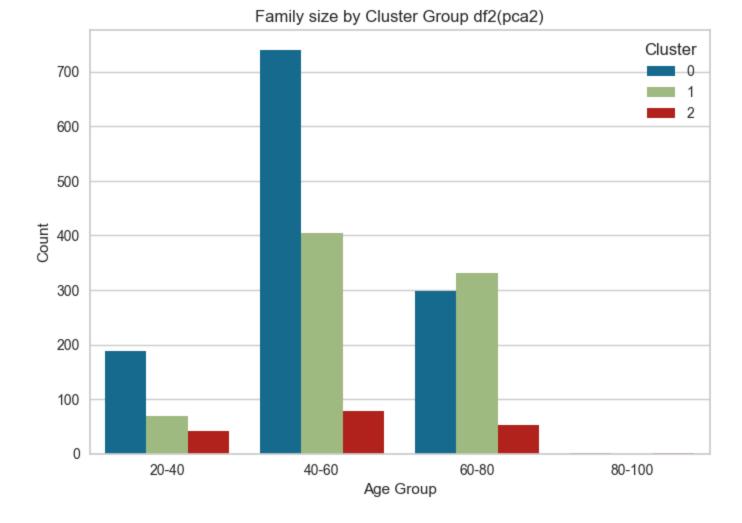
# Pairplot
sns.pairplot(df_subset, hue='Cluster')
plt.show()
```



```
# Define the bin edges and labels for age bins
In [66]:
         age bins = [20, 40, 60, 80, 100]
         age labels = ['20-40', '40-60', '60-80', '80-100']
         # Bin the 'Age' variable using pd.cut()
         df1['Age bins'] = pd.cut(df1['Age'], bins=age bins, labels=age labels)
         # Plot 'Age bins' vs 'Cluster' using countplot
         sns.countplot(data=df1, x='Age bins', hue='Cluster')
         plt.title('Family size by Cluster Group df1(pca1)')
         plt.xlabel('Age Group')
        plt.ylabel('Count')
         plt.show()
         # Define the bin edges and labels for age bins df2
         age bins = [20, 40, 60, 80, 100]
         age_labels = ['20-40', '40-60', '60-80', '80-100']
         # Bin the 'Age' variable using pd.cut()
         df2['Age bins'] = pd.cut(df1['Age'], bins=age bins, labels=age labels)
         # Plot 'Age bins' vs 'Cluster' using countplot
         sns.countplot(data=df1, x='Age bins', hue='Cluster')
         plt.title('Family size by Cluster Group df2(pca2)')
         plt.xlabel('Age Group')
```

plt.ylabel('Count')
plt.show()





After rigorous exploratory data analysis and insightful visualizations through a Power BI dashboard, the stakeholders and marketing strategy teams have decided to move forward with the PCA1 cluster group. The decision was based on the cluster group's superior ability to highlight underlying patterns and segregate the data into meaningful segments.

The process uncovered several interlinked relationships between features that present valuable insights for targeted marketing and customer engagement strategies. These insights, derived from the PCA1 cluster group analysis, will significantly influence the teams' future course of action and help design personalized, data-driven marketing campaigns.

### **Cluster 0: High Spending Customers**

- This group typically has higher income, usually exceeding 80K.
- They exhibit a higher spending range of 1300 to 2000, indicating strong purchasing power.
- Predominantly made up of single individuals, and most of them do not have children or have only one child.
- Show high engagement with campaigns, with a majority having accepted campaign offerings in the past.
- Considering the high income and spending, targeted marketing for premium and exclusive products/services could resonate well with this group.

#### **Cluster 1: Low-Income Family Group**

- This group consists of lower income individuals, typically earning between 0 to 50K.
- Their spending capacity is comparably lower, ranging between 10 to 250.

- The group is largely made up of couples, most of whom have at least 1 or 2 children.
- A high proportion of individuals in this group are educated.
- The family size is typically larger, ranging from 3 to 5 with no singles in this group.
- The age range is quite mature, from 40 to 60 years old.
- Discounts, family-centric offerings, and educational resources might be well-received by this group.

### **Cluster 2: Middle-Income Singles with Children**

- This cluster comprises a middle-income group, with earnings between 50K and 80K.
- They spend an average amount on purchases, with the range falling between 1000 to 1600.
- Most are single and have only one child, indicating a small family size not exceeding 3 members.
- Given the balance between income and spending, this group may respond well to value-for-money offerings.
- Targeted marketing strategies may include single parent-focused campaigns or promotions for moderate to high-priced items.

### **Cluster 0: Low-Income Family Group**

- This group consists of lower income individuals, typically earning between 0 to 50K.
- Their spending capacity is comparably lower, ranging between 10 to 250.
- The group is largely made up of couples, most of whom have at least 1 or 2 children.
- A high proportion of individuals in this group are educated.
- The family size is typically larger, ranging from 3 to 5 with no singles in this group.
- The age range is quite mature, from 40 to 60 years old.
- Discounts, family-centric offerings, and educational resources might be well-received by this group.

### **Cluster 1: Middle-Income Singles with Children**

- This cluster comprises a middle-income group, with earnings between 50K and 80K.
- They spend an average amount on purchases, with the range falling between 1000 to 1600.
- Most are single and have only one child, indicating a small family size not exceeding 3 members.
- Given the balance between income and spending, this group may respond well to value-for-money offerings.
- Targeted marketing strategies may include single parent-focused campaigns or promotions for moderate to high-priced items.

### **Cluster 2: High Spending Customers**

- This group typically has higher income, usually exceeding 80K.
- They exhibit a higher spending range of 1300 to 2000, indicating strong purchasing power.
- Predominantly made up of single individuals, and most of them do not have children or have only one child.
- Show high engagement with campaigns, with a majority having accepted campaign offerings in the past.
- Considering the high income and spending, targeted marketing for premium and exclusive products/services could resonate well with this group.

### **Business Recommendations and Marketing Strategies:**

1. **Personalized Campaigns:** Develop personalized marketing campaigns for each cluster group, considering their unique characteristics and preferences.

- 2. **Premium Offerings for Cluster 2:** Focus on promoting premium products and exclusive services to this group, given their high income and spending power.
- 3. **Value Deals for Cluster 0:** Offer value deals, discounts, or bundle deals to cater to the lower income and spending power of this group.
- 4. **Family Packages for Cluster 0:** Create family packages or child-friendly products/services to attract and cater to the larger families in Cluster 0.
- 5. **Balanced Approach for Cluster 1:** Offer a balanced approach with value for money products and emphasize quality and durability.
- 6. **Education Focused Marketing:** Provide detailed information and content marketing to the highly educated Cluster 1.
- 7. **Single Parent Campaigns for Cluster 1:** Tailor campaigns to cater to the unique challenges and needs of single-parent households in Cluster 1.
- 8. **Lifecycle Marketing:** Track customers through different life stages and adjust marketing strategies accordingly.
- 9. **Customer Feedback:** Seek regular feedback from each cluster group to understand their evolving needs and make necessary adjustments.
- 10. **Predictive Analytics:** Utilize predictive analytics to anticipate future buying behavior and tailor marketing efforts proactively.

## STAGE 2

## Predictive Model Development for Future Data

"Developing a predictive model using cluster groups and the target variable to classify future data and assign customers to their respective cluster groups."

```
In [4]: #load the data
model_df = pd.read_csv(r"group1_pca1.csv")
model_df.head(5)
```

Out[4]:		ID	Year_Birth	Education	Marital_Status	Income	Kidhome	Teenhome	Dt_Customer	Recency	MntWines
	0	5524	1957	Graduation	Single	58138.0	0	0	2012-09-04	58	635
	1	2174	1954	Graduation	Single	46344.0	1	1	2014-03-08	38	11
	2	4141	1965	Graduation	Couple	71613.0	0	0	2013-08-21	26	426
	3	6182	1984	Graduation	Couple	26646.0	1	0	2014-02-10	26	11
	4	5324	1981	PhD	Couple	58293.0	1	0	2014-01-19	94	173

5 rows × 37 columns

```
model_df2 = model_df.copy()
model_df3 = model_df.copy()
model_df4 = model_df.copy()
```

### **Experiment 1 (SMOTE)**

```
In [6]:
        # drop unimportant features
        model_df.drop(columns=['ID', 'Year_Birth', 'Education','Marital Status','Kidhome',
               'Teenhome', 'Dt Customer', 'MntWines', 'MntFruits',
               'MntMeatProducts', 'MntFishProducts', 'MntSweetProducts',
               'MntGoldProds', 'AcceptedCmp3', 'AcceptedCmp4', 'AcceptedCmp5', 'AcceptedCmp1',
               'AcceptedCmp2','NumDealsPurchases', 'NumWebPurchases','NumStorePurchases','NumCat
In [7]: # train test split the data
        from sklearn.model selection import train test split
        X = model df.drop(columns=['Cluster'], axis=1)
        y = model df['Cluster']
        X train, X test, y train, y test = train test split(X,y, test size=0.2, random state=9)
        y train.value counts()
In [8]:
             975
Out[8]:
             653
             136
        Name: Cluster, dtype: int64
```

To address the class imbalance in the target variable, consider using the SMOTE (Synthetic Minority Over-sampling Technique) algorithm.

SMOTE (Synthetic Minority Over-sampling Technique) is a technique used to address class imbalance problems in machine learning. It creates synthetic samples of the minority class to balance out the number of samples between classes, thereby helping to improve model performance and robustness. SMOTE works by considering the k-nearest neighbors of a data point in the feature space and generating new instances along the lines joining the neighbors. While it's a powerful technique to counter class imbalance, it should be applied thoughtfully as it can sometimes introduce noise by generating synthetic instances without considering the instances from other classes.

```
instances without considering the instances from other classes.
In [9]: from imblearn.over sampling import SMOTE
         # Create an instance of SMOTE
         sm = SMOTE(random state=42)
         # Fit SMOTE to the training data
         X train, y train = sm.fit resample(X train, y train)
        y train.value counts()
In [10]:
              975
Out[10]:
              975
              975
        Name: Cluster, dtype: int64
         from sklearn.metrics import accuracy score, recall score, f1 score, precision score
In [11]:
         from sklearn.metrics import classification report
         from sklearn.linear model import LogisticRegression
         from sklearn.naive bayes import GaussianNB
         from sklearn.tree import DecisionTreeClassifier
         from sklearn.ensemble import RandomForestClassifier, AdaBoostClassifier
         from sklearn.svm import SVC
         from sklearn.neighbors import KNeighborsClassifier
```

```
from xgboost import XGBClassifier
from sklearn.neural network import MLPClassifier
from sklearn.model selection import cross val score
# Defining the models
classifiers = {
   "Logistic Regression": LogisticRegression(),
   "Naive Bayes": GaussianNB(),
   "Decision Tree": DecisionTreeClassifier(),
   "Random Forest": RandomForestClassifier(),
    "Support Vector Machines (SVM)": SVC(),
    "K-Nearest Neighbors (KNN)": KNeighborsClassifier(),
   "Gradient Boosting": XGBClassifier(eval metric='mlogloss'),
   "Neural Networks (MLP)": MLPClassifier(),
    "AdaBoost": AdaBoostClassifier()
# Dataframe to store results
experiment 1 = pd.DataFrame(columns=["Model", "Training Accuracy", "Testing Accuracy", "
# Loop over classifiers
for classifier name, classifier in classifiers.items():
    # Train the classifier
   classifier.fit(X train, y train)
    # Make predictions on training set
   y train pred = classifier.predict(X train)
    # Make predictions on testing set
   y test pred = classifier.predict(X test)
    # Calculate metrics
   training_accuracy = accuracy_score(y_train, y_train_pred)
   testing accuracy = accuracy_score(y_test, y_test_pred)
   cv score = cross val score(classifier, X, y, cv=5).mean()
   precision = precision score(y test, y test pred, average='weighted')
   recall = recall score(y test, y test pred, average='weighted')
   f1 = f1 score(y test, y test pred, average='weighted')
    # Append results to dataframe
    experiment 1 = experiment 1.append({
       "Model": classifier name,
       "Training Accuracy": training accuracy,
       "Testing Accuracy": testing accuracy,
       "CV Score": cv score,
       "Precision": precision,
       "Recall": recall,
       "F1 Score": f1
    }, ignore index=True)
# Display results
experiment 1.head(10)
```

Out[11]:		Model	<b>Training Accuracy</b>	Accuracy Testing Accuracy		Precision	Recall	F1 Score
	0	Logistic Regression	0.787009	0.845805	0.874376	0.883256	0.845805	0.855174
	1	Naive Bayes	0.917607	0.929705	0.936961	0.937507	0.929705	0.931688
	2	Decision Tree	1.000000	0.968254	0.965986	0.968254	0.968254	0.968254
	3	Random Forest	1.000000	0.979592	0.974603	0.979844	0.979592	0.979637
	4	Support Vector Machines (SVM)	0.745299	0.829932	0.829025	0.850708	0.829932	0.836761
	5	K-Nearest Neighbors (KNN)	0.914872	0.825397	0.868934	0.858275	0.825397	0.836054

6	Gradient Boosting	1.000000	0.972789 0.973696 0.972872 0.972789 0.972814	.972789 0.973696 0.9
7	Neural Networks (MLP)	0.778803	0.852608	.852608 0.654422 0.8
8	AdaBoost	0.969915	0.970522 0.964626 0.970562 0.970522 0.970523	.970522 0.964626 0.9

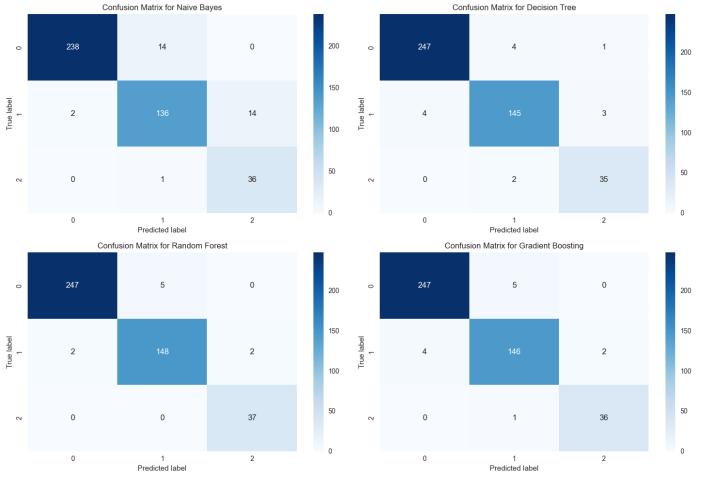
### **Evaluation**

Among these models, Naive Bayes, Decision Tree, Random Forest, and Gradient Boosting performed well in terms of accuracy, precision, recall, and F1 score. Further evaluation was conducted using visualizations such as confusion matrix, learning curve, and class prediction error.

- Confusion matrix: A visualization that shows the performance of a classification model by comparing
  actual and predicted class labels. It helps in understanding the distribution of true positives, true
  negatives, false positives, and false negatives.
- Learning curve: A plot that depicts the performance of a model as training data size increases. It helps in assessing model performance and identifying issues such as overfitting or underfitting.
- Class prediction error: A plot that shows the difference between the true class distribution and the predicted class distribution. It helps in identifying which classes are misclassified more frequently.

These visualizations were used to gain further insights into the models' performance and identify areas for improvement.

```
from sklearn.metrics import confusion matrix
In [12]:
         # Instantiate the classifiers
         classifiers = {
            "Naive Bayes": GaussianNB(),
             "Decision Tree": DecisionTreeClassifier(),
             "Random Forest": RandomForestClassifier(),
             "Gradient Boosting": XGBClassifier(eval metric='mlogloss')
         fig, axs = plt.subplots(2, 2, figsize=(15, 10)) # Create a 2 by 2 subplot
         for (classifier name, classifier), ax in zip(classifiers.items(), axs.ravel()):
             # Fit the classifier
            classifier.fit(X train, y train)
             # Predict the test set results
             y pred = classifier.predict(X test)
             # Create a confusion matrix
             cf matrix = confusion matrix(y test, y pred)
             # Plot the confusion matrix
             sns.heatmap(cf matrix, annot=True, fmt='d', cmap='Blues', ax=ax)
             ax.set title(f'Confusion Matrix for {classifier name}')
             ax.set ylabel('True label')
             ax.set xlabel('Predicted label')
         plt.tight layout()
         plt.show()
```

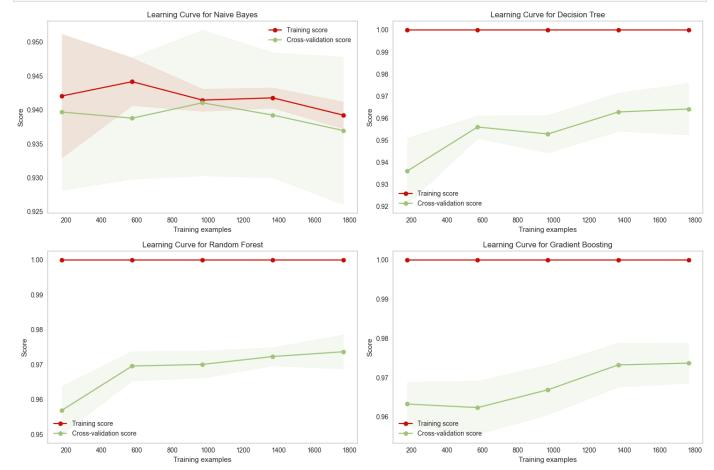


In [13]: from sklearn.model selection import learning curve def plot learning curve (estimator, title, X, y, axes, cv=None, n jobs=None, train sizes= axes.set title(title) axes.set xlabel("Training examples") axes.set ylabel("Score") train sizes, train scores, test scores, fit times, learning curve(estimator, X, y, cv=cv, n jobs=n jobs, train sizes=train sizes, return times=True) train scores mean = np.mean(train scores, axis=1) train scores std = np.std(train scores, axis=1) test scores mean = np.mean(test scores, axis=1) test scores std = np.std(test\_scores, axis=1) axes.grid() axes.fill between(train sizes, train scores mean - train scores std, train scores mean + train scores std, alpha=0.1, color="r") axes.fill between(train sizes, test scores mean - test scores std, test scores mean + test scores std, alpha=0.1, color="g") axes.plot(train sizes, train scores mean, 'o-', color="r", label="Training score") axes.plot(train sizes, test scores mean, 'o-', color="g", label="Cross-validation score") axes.legend(loc="best") fig, axs = plt.subplots(2, 2, figsize=(15, 10)) # Create a 2 by 2 subplot # Make sure your 'selected classifiers' is a list or ordered dict, not a normal dict. # Normal dict doesn't preserve order before Python 3.7

```
selected_classifiers = [
    ("Naive Bayes", GaussianNB()),
    ("Decision Tree", DecisionTreeClassifier()),
    ("Random Forest", RandomForestClassifier()),
    ("Gradient Boosting", XGBClassifier(eval_metric='mlogloss')),
]

for (name, clf), ax in zip(selected_classifiers, axs.ravel()):
    plot_learning_curve(clf, f'Learning Curve for {name}', X, y, ax, cv=5)

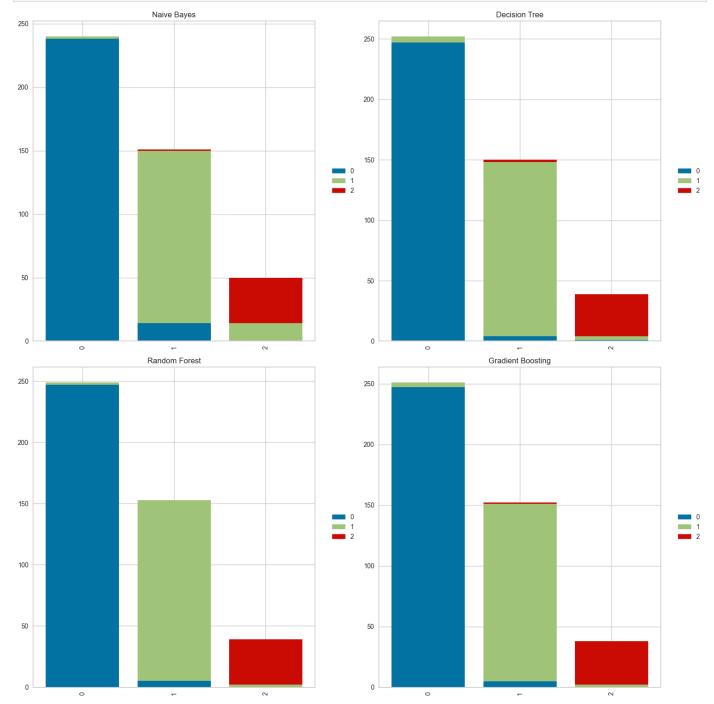
plt.tight_layout()
plt.show()
```



```
from yellowbrick.classifier import ClassPredictionError
In [14]:
         fig, axes = plt.subplots(nrows=2, ncols=2, figsize=(15,15))
         # Define the dictionary of classifiers
         selected classifiers = {
             "Naive Bayes": GaussianNB(),
             "Decision Tree": DecisionTreeClassifier(),
             "Random Forest": RandomForestClassifier(),
             "Gradient Boosting": XGBClassifier(eval metric='mlogloss')
         for (classifier name, classifier), ax in zip(selected classifiers.items(), axes.flatten(
             # Instantiate the classification model and visualizer
             visualizer = ClassPredictionError(classifier, classes=[0,1,2], ax=ax)
             # Fit the training data to the visualizer
             visualizer.fit(X train, y train)
             # Evaluate the model on the test data
             visualizer.score(X test, y test)
```

```
# Set the title for the subplot
ax.set_title(classifier_name)

# Display the figure
plt.tight_layout()
plt.show()
```



## **Experiment 2 (Without SMOTE)**

```
In [15]: # train test split the data

X = model_df.drop(columns=['Cluster'], axis=1)
y = model_df['Cluster']

X_train, X_test, y_train, y_test = train_test_split(X,y, test_size=0.2, random_state=9)
```

In [16]: y\_train.value\_counts()

```
1
             653
             136
        Name: Cluster, dtype: int64
In [17]: # Defining the models
         classifiers = {
             "Logistic Regression": LogisticRegression(),
             "Naive Bayes": GaussianNB(),
             "Decision Tree": DecisionTreeClassifier(),
             "Random Forest": RandomForestClassifier(),
             "Support Vector Machines (SVM)": SVC(),
             "K-Nearest Neighbors (KNN)": KNeighborsClassifier(),
             "Gradient Boosting": XGBClassifier(eval metric='mlogloss'),
             "Neural Networks (MLP)": MLPClassifier(),
             "AdaBoost": AdaBoostClassifier()
         # Dataframe to store results
         experiment 2 = pd.DataFrame(columns=["Model", "Training Accuracy", "Testing Accuracy", "
         # Loop over classifiers
         for classifier name, classifier in classifiers.items():
             # Train the classifier
             classifier.fit(X train, y train)
             # Make predictions on training set
             y train pred = classifier.predict(X train)
             # Make predictions on testing set
             y test pred = classifier.predict(X test)
             # Calculate metrics
             training accuracy = accuracy score(y train, y train pred)
             testing accuracy = accuracy score(y test, y test pred)
             cv score = cross val score(classifier, X, y, cv=5).mean()
             precision = precision score(y test, y test pred, average='weighted')
             recall = recall score(y test, y test pred, average='weighted')
             f1 = f1 score(y test, y test pred, average='weighted')
             # Append results to dataframe
             experiment 2 = experiment 2.append({
                 "Model": classifier name,
                 "Training Accuracy": training accuracy,
                 "Testing Accuracy": testing accuracy,
                 "CV Score": cv score,
                 "Precision": precision,
                 "Recall": recall,
                 "F1 Score": f1
             }, ignore index=True)
         # Display results
         experiment 2.head(10)
Out[17]:
```

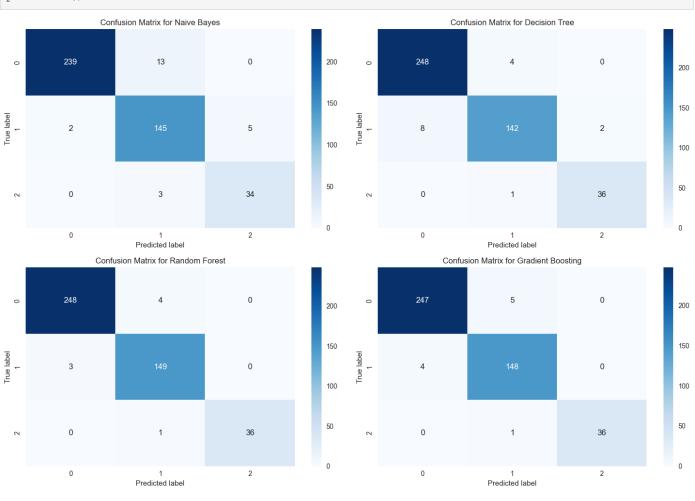
Out[16]: 0

975

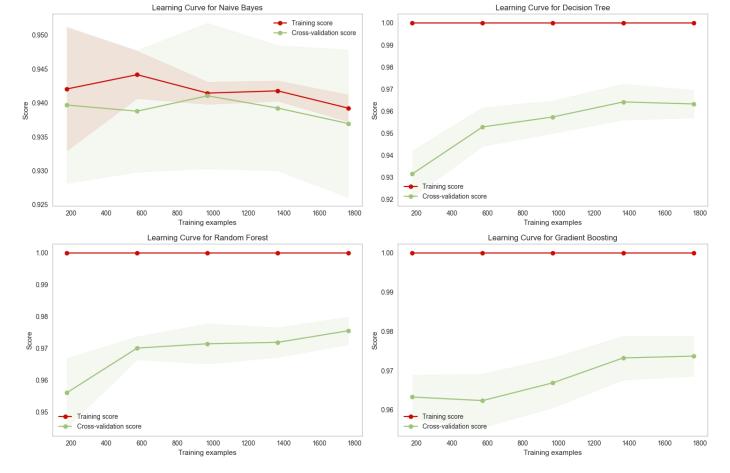
:		Model	<b>Training Accuracy</b>	<b>Testing Accuracy</b>	CV Score	Precision	Recall	F1 Score
	0	Logistic Regression	0.880952	0.868481	0.874376	0.864575	0.868481	0.856618
	1	Naive Bayes	0.936508	0.947846	0.936961	0.950248	0.947846	0.948455
	2	Decision Tree	1.000000	0.977324	0.965986	0.977363	0.977324	0.977294
	3	Random Forest	1.000000	0.981859	0.975964	0.981980	0.981859	0.981887
	4	Support Vector Machines (SVM)	0.826531	0.843537	0.829025	0.785101	0.843537	0.810306
	5	K-Nearest Neighbors (KNN)	0.907029	0.879819	0.868934	0.875134	0.879819	0.874393

6	Gradient Boosting	1.000000	0.977324 0.973696 0.977465 0.977324 0.977363
7	Neural Networks (MLP)	0.824830	0.814059 0.695692 0.795318 0.814059 0.786857
8	AdaBoost	0.963152	0.977324 0.964626 0.977363 0.977324 0.977294

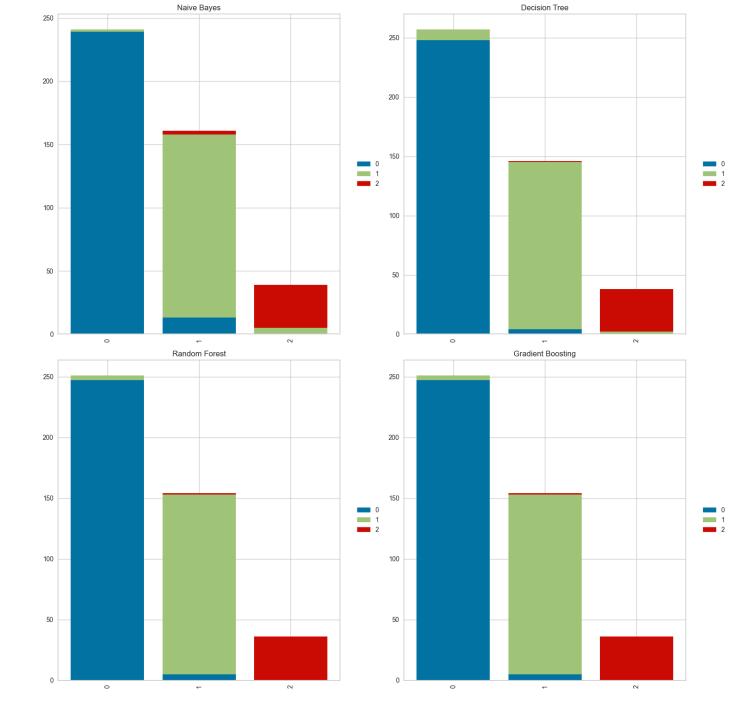
```
selected classifiers = {
In [18]:
             "Naive Bayes": GaussianNB(),
             "Decision Tree": DecisionTreeClassifier(),
             "Random Forest": RandomForestClassifier(),
             "Gradient Boosting": XGBClassifier(eval metric='mlogloss'),
         fig, axs = plt.subplots(2, 2, figsize=(15, 10)) # Create a 2 by 2 subplot
         for (classifier name, classifier), ax in zip(selected classifiers.items(), axs.ravel()):
             # Fit the classifier
             classifier.fit(X train, y train)
             # Predict the test set results
             y pred = classifier.predict(X test)
             # Create a confusion matrix
             cf matrix = confusion matrix(y test, y pred)
             # Plot the confusion matrix
             sns.heatmap(cf matrix, annot=True, fmt='d', cmap='Blues', ax=ax)
             ax.set title(f'Confusion Matrix for {classifier name}')
             ax.set ylabel('True label')
             ax.set xlabel('Predicted label')
        plt.tight layout()
         plt.show()
```



```
In [19]: def plot_learning_curve(estimator, title, X, y, axes, cv=None, n jobs=None, train sizes=
            axes.set title(title)
             axes.set xlabel("Training examples")
            axes.set ylabel("Score")
             train sizes, train scores, test scores, fit times, = \
                 learning curve(estimator, X, y, cv=cv, n_jobs=n_jobs,
                                train sizes=train sizes,
                                return times=True)
             train scores mean = np.mean(train scores, axis=1)
             train scores std = np.std(train scores, axis=1)
             test scores mean = np.mean(test scores, axis=1)
             test scores std = np.std(test scores, axis=1)
             axes.grid()
             axes.fill between(train sizes, train scores mean - train scores std,
                             train scores mean + train scores std, alpha=0.1,
                              color="r")
            axes.fill between (train sizes, test scores mean - test scores std,
                             test scores mean + test scores std, alpha=0.1,
                              color="q")
             axes.plot(train sizes, train scores mean, 'o-', color="r",
                     label="Training score")
             axes.plot(train sizes, test scores mean, 'o-', color="g",
                     label="Cross-validation score")
             axes.legend(loc="best")
         fig, axs = plt.subplots(2, 2, figsize=(15, 10)) # Create a 2 by 2 subplot
        selected classifiers = [
            ("Naive Bayes", GaussianNB()),
             ("Decision Tree", DecisionTreeClassifier()),
             ("Random Forest", RandomForestClassifier()),
             ("Gradient Boosting", XGBClassifier(eval metric='mlogloss')),
         1
        for (name, clf), ax in zip(selected classifiers, axs.ravel()):
            plot learning curve(clf, f'Learning Curve for {name}', X, y, ax, cv=5)
        plt.tight layout()
        plt.show()
```



```
In [20]:
         fig, axes = plt.subplots(nrows=2, ncols=2, figsize=(15,15))
         # Define the dictionary of classifiers
         selected classifiers = {
             "Naive Bayes": GaussianNB(),
             "Decision Tree": DecisionTreeClassifier(),
             "Random Forest": RandomForestClassifier(),
             "Gradient Boosting": XGBClassifier(eval metric='mlogloss')
         for (classifier name, classifier), ax in zip(selected classifiers.items(), axes.flatten(
             # Instantiate the classification model and visualizer
             visualizer = ClassPredictionError(classifier, classes=[0,1,2], ax=ax)
             # Fit the training data to the visualizer
             visualizer.fit(X train, y train)
             # Evaluate the model on the test data
             visualizer.score(X test, y test)
             # Set the title for the subplot
             ax.set title(classifier name)
         # Display the figure
        plt.tight layout()
         plt.show()
```



# **Experiment 3 (More Features in Dataframe)**

model\_df2.head() In [21]: Out[21]: Year\_Birth Marital\_Status Income Kidhome Teenhome Dt\_Customer Recency Education **MntWines** Single 58138.0 0 5524 1957 Graduation 0 0 2012-09-04 635 58 **1** 2174 1954 Graduation Single 46344.0 1 2014-03-08 38 11 **2** 4141 1965 Graduation Couple 71613.0 0 2013-08-21 26 426 **3** 6182 1984 Graduation Couple 26646.0 0 2014-02-10 26 11 **4** 5324 1981 PhD Couple 58293.0 2014-01-19 173

5 rows × 37 columns

```
In [22]: # drop unimportant features
        model df2.drop(columns=['ID', 'Year Birth', 'Education','Marital Status','Dt Customer'],
In [23]: | # train test split the data
        from sklearn.model selection import train_test_split
        X = model df2.drop(columns=['Cluster'], axis=1)
        y = model df2['Cluster']
        X train, X test, y train, y test = train test split(X,y, test size=0.2, random state=9)
        y train.value counts()
            975
Out[23]:
             653
             136
        Name: Cluster, dtype: int64
        # Create an instance of SMOTE
In [24]:
        sm = SMOTE(random state=42)
         # Fit SMOTE to the training data
        X train, y train = sm.fit resample(X train, y train)
        y train.value counts()
        1 975
Out[24]:
        0
            975
             975
        Name: Cluster, dtype: int64
In [25]: # Defining the models
         classifiers = {
            "Logistic Regression": LogisticRegression(),
            "Naive Bayes": GaussianNB(),
             "Decision Tree": DecisionTreeClassifier(),
             "Random Forest": RandomForestClassifier(),
             "Support Vector Machines (SVM)": SVC(),
             "K-Nearest Neighbors (KNN)": KNeighborsClassifier(),
             "Gradient Boosting": XGBClassifier(eval metric='mlogloss'),
             "Neural Networks (MLP)": MLPClassifier(),
             "AdaBoost": AdaBoostClassifier()
         # Dataframe to store results
         experiment 3 = pd.DataFrame(columns=["Model", "Training Accuracy", "Testing Accuracy", "
         # Loop over classifiers
         for classifier name, classifier in classifiers.items():
             # Train the classifier
            classifier.fit(X train, y train)
             # Make predictions on training set
             y train pred = classifier.predict(X train)
             # Make predictions on testing set
            y test pred = classifier.predict(X test)
             # Calculate metrics
             training accuracy = accuracy score(y train, y train pred)
             testing accuracy = accuracy score(y test, y test pred)
             cv score = cross val score(classifier, X, y, cv=10).mean()
            precision = precision score(y test, y test pred, average='weighted')
            recall = recall score(y test, y test pred, average='weighted')
             f1 = f1_score(y_test, y_test_pred, average='weighted')
             # Append results to dataframe
```

```
experiment_3 = experiment_3.append({
    "Model": classifier_name,
    "Training Accuracy": training_accuracy,
    "Testing Accuracy": testing_accuracy,
    "CV Score": cv_score,
    "Precision": precision,
    "Recall": recall,
    "F1 Score": f1
}, ignore_index=True)

# Display results
experiment_3.head(10)
```

#### Out[25]: Model Training Accuracy Testing Accuracy CV Score Precision Recall F1 Score 0 0.757949 0.782313 0.873030 0.834536 0.782313 0.797400 Logistic Regression 0.903590 0.931983 0.939648 0.929705 0.932684 Naive Bayes 0.929705 2 **Decision Tree** 1.000000 0.975057 0.970068 0.975207 0.975057 0.975037 3 Random Forest 1.000000 0.990930 0.979142 0.990930 0.990930 0.990930 Support Vector Machines (SVM) 0.747350 0.829932 0.830370 0.850708 0.829932 0.836761 5 K-Nearest Neighbors (KNN) 0.934017 0.852608 0.871199 0.873874 0.852608 0.859340 6 **Gradient Boosting** 1.000000 0.990930 0.983221 0.990930 0.990930 0.990930

0.816068

0.960684

0.823129

0.789605

0.864929 0.823129 0.835782

0.981859 0.957820 0.981931 0.981859 0.981823

7

8

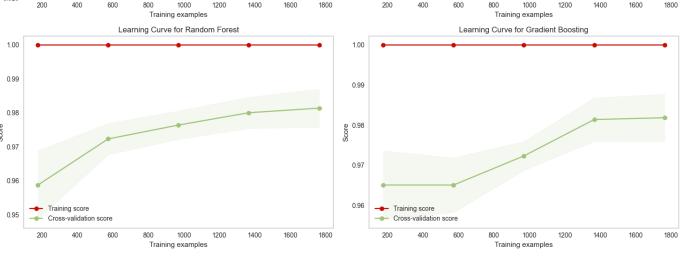
Neural Networks (MLP)

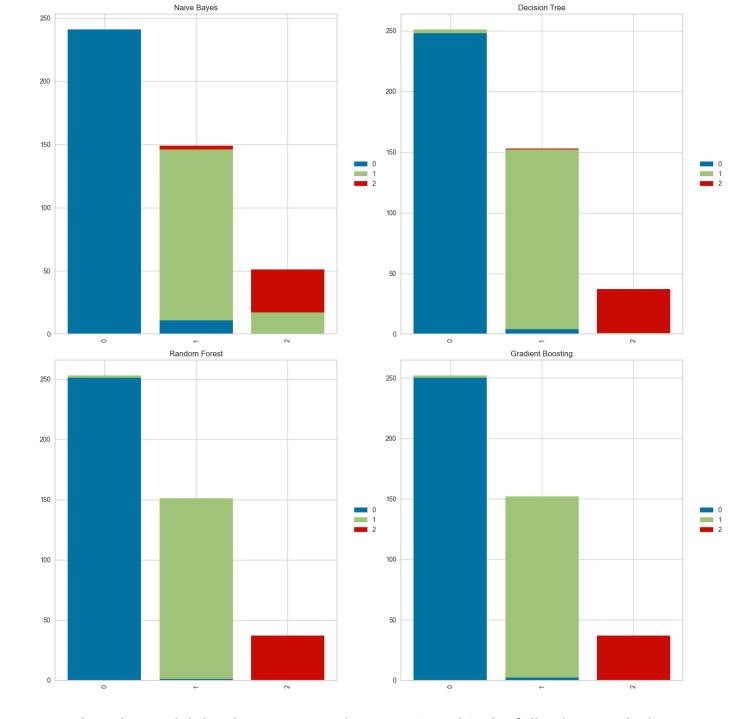
AdaBoost

```
# Instantiate the classifiers
In [26]:
         classifiers = {
             "Naive Bayes": GaussianNB(),
             "Decision Tree": DecisionTreeClassifier(),
             "Random Forest": RandomForestClassifier(),
             "Gradient Boosting": XGBClassifier(eval metric='mlogloss')
         fig, axs = plt.subplots(2, 2, figsize=(15, 10)) # Create a 2 by 2 subplot
         for (classifier name, classifier), ax in zip(classifiers.items(), axs.ravel()):
             # Fit the classifier
             classifier.fit(X train, y train)
             # Predict the test set results
             y pred = classifier.predict(X test)
             # Create a confusion matrix
             cf matrix = confusion matrix(y test, y pred)
             # Plot the confusion matrix
             sns.heatmap(cf matrix, annot=True, fmt='d', cmap='Blues', ax=ax)
             ax.set title(f'Confusion Matrix for {classifier_name}')
             ax.set ylabel('True label')
             ax.set xlabel('Predicted label')
         plt.tight layout()
         plt.show()
         def plot learning curve(estimator, title, X, y, axes, cv=None, n jobs=None, train sizes=
             axes.set title(title)
```

```
axes.set xlabel("Training examples")
    axes.set ylabel("Score")
    train sizes, train scores, test scores, fit times, = \
        learning curve (estimator, X, y, cv=cv, n jobs=n jobs,
                       train sizes=train sizes,
                       return times=True)
    train scores mean = np.mean(train scores, axis=1)
    train scores std = np.std(train scores, axis=1)
    test scores mean = np.mean(test_scores, axis=1)
    test scores std = np.std(test scores, axis=1)
    axes.grid()
    axes.fill between(train sizes, train scores mean - train scores std,
                     train scores mean + train scores std, alpha=0.1,
                     color="r")
    axes.fill between(train sizes, test scores mean - test scores std,
                     test scores mean + test scores std, alpha=0.1,
                     color="q")
    axes.plot(train sizes, train scores mean, 'o-', color="r",
             label="Training score")
    axes.plot(train_sizes, test_scores_mean, 'o-', color="g",
             label="Cross-validation score")
    axes.legend(loc="best")
fig, axs = plt.subplots(2, 2, figsize=(15, 10)) # Create a 2 by 2 subplot
selected classifiers = [
    ("Naive Bayes", GaussianNB()),
    ("Decision Tree", DecisionTreeClassifier()),
    ("Random Forest", RandomForestClassifier()),
    ("Gradient Boosting", XGBClassifier(eval metric='mlogloss')),
]
for (name, clf), ax in zip(selected classifiers, axs.ravel()):
    plot learning curve(clf, f'Learning Curve for {name}', X, y, ax, cv=5)
plt.tight layout()
plt.show()
fig, axes = plt.subplots(nrows=2, ncols=2, figsize=(15,15))
# Define the dictionary of classifiers
selected classifiers = {
    "Naive Bayes": GaussianNB(),
    "Decision Tree": DecisionTreeClassifier(),
    "Random Forest": RandomForestClassifier(),
    "Gradient Boosting": XGBClassifier(eval metric='mlogloss')
for (classifier name, classifier), ax in zip(selected classifiers.items(), axes.flatten(
    # Instantiate the classification model and visualizer
    visualizer = ClassPredictionError(classifier, classes=[0,1,2], ax=ax)
    # Fit the training data to the visualizer
    visualizer.fit(X train, y train)
    # Evaluate the model on the test data
   visualizer.score(X test, y test)
    # Set the title for the subplot
    ax.set title(classifier name)
# Display the figure
```

plt.tight\_layout() plt.show() Confusion Matrix for Naive Bayes Confusion Matrix for Decision Tree True label True label Predicted label Predicted label Confusion Matrix for Random Forest Confusion Matrix for Gradient Boosting True label Predicted label Predicted label Learning Curve for Naive Bayes Learning Curve for Decision Tree Training score 0.960 1.00 Cross-validation score 0.955 0.99 0.950 0.98 0.945 0.97 0.940 0.96 0.935 0.95 0.930 0.94 0.925 Training score Cross-validation score 0.93 0.920 Training examples Training examples Learning Curve for Gradient Boosting Learning Curve for Random Forest 1.00 1.00 0.99 0.99 0.98





Based on the model development experiments 1, 2, and 3, the following conclusions can be made:

**Feature Selection**: The chosen experiment was Experiment 3, which involved the use of almost all features and the application of the Synthetic Minority Over-sampling Technique (SMOTE) to handle class imbalance. Despite differences in feature selection and handling of class imbalance among the experiments, all models performed similarly. This suggests that using a more comprehensive set of features, as in Experiment 3, provided a well-fitted input for the classification models without negatively impacting performance.

**SMOTE's Impact**: Experiment 3, which made use of SMOTE to balance the classes, was chosen over Experiment 2 which did not use SMOTE. This decision underscores the importance of handling class imbalance in the dataset. While the impact of SMOTE was not drastically apparent in the difference between the performances of the models, its utilization is critical in datasets with imbalanced classes to ensure that the minority class is not ignored.

**Model Performance**: Among all the classification algorithms tested, the Gradient Boosting method, specifically the XGBoost Classifier, showed the best performance with fewer errors in the Class Prediction

Error plot. It accurately identified the main target customers (class 1), which was the objective of this exercise.

Given these findings, the XGBoost Classifier was chosen for further hyperparameter tuning. This process of optimization helps enhance the performance of the model by adjusting the model parameters to their ideal values. Once this is achieved, the model will be saved and deployed. The deployment of the model allows it to be used in practical applications, providing predictions on new, unseen data.

In conclusion, the analysis underscores the importance of careful feature selection, handling of class imbalance, and selection of the right model in creating robust and accurate predictive models. The chosen experiment and model, Experiment 3 and the XGBoost Classifier respectively, were the best fit for this dataset and problem, providing reliable predictions and highlighting the main target customer group (class 1) accurately. It also indicates the importance of model optimization via hyperparameter tuning to maximize the performance of the chosen model.

## **Final Model Selected Development**

## Hyperparameter Tunning

### **GradianteBoost (XGB Classifier)**

To ensure an accurate evaluation, the dataset model\_df2 will be split into three subsets: training, validation, and testing. The training data will be used for model training, the validation data for hyperparameter optimization, and the testing data for final evaluation. This approach ensures a robust assessment of the model's performance.

### Train-Test-Valid

```
from sklearn.model selection import train test split
In [27]:
         from imblearn.over sampling import SMOTE
         from sklearn.metrics import accuracy score
         from sklearn.model selection import RandomizedSearchCV
         from xgboost import XGBClassifier
         # Define features (X) and target (y)
         X = model df2.drop('Cluster', axis=1)
         y = model df2['Cluster']
         # Split data into training, validation, and test sets (70% - 15% - 15%)
         X temp, X test, y temp, y test = train test split(X, y, test size=0.15, random state=42)
         X train, X val, y train, y val = train test split(X temp, y temp, test size=0.18, random
         # Apply SMOTE to handle class imbalance in the training set
         smote = SMOTE()
         X train smote, y train smote = smote.fit resample(X train, y train)
         #Define the hyperparameters for RandomizedSearch
         param distributions = {
             'learning rate': [0.01, 0.1, 0.2, 0.3],
             'max depth': [3, 5, 7, 10],
             'min child weight': [1, 3, 5],
             'subsample': [0.5, 0.7, 1],
             'colsample bytree': [0.5, 0.7, 1],
             'n estimators' : [100, 200, 500],
             'objective': ['binary:logistic']
         # Initialize the XGBClassifier
         xgb classifier = XGBClassifier(eval metric='mlogloss')
         # Initialize RandomizedSearchCV and fit the training data
```

```
random search.fit(X train smote, y train smote)
         # Get the best parameters
         best parameters = random search.best params
         print("Best parameters found: ", best parameters)
         # Train the model using the best parameters on the training data
        best classifier = XGBClassifier(**best parameters, eval metric='mlogloss')
         best classifier.fit(X train smote, y train smote)
         # Predict the validation set results
         y val pred = best classifier.predict(X val)
         # Calculate and print the validation accuracy
         val accuracy = accuracy score(y val, y val pred)
         print("Validation Accuracy: ", val accuracy)
         # Predict the test set results
         y test pred = best classifier.predict(X test)
         # Calculate and print the test accuracy
         test accuracy = accuracy score(y test, y test pred)
         print("Test Accuracy: ", test accuracy)
        Fitting 5 folds for each of 10 candidates, totalling 50 fits
        Best parameters found: {'subsample': 0.7, 'objective': 'binary:logistic', 'n estimator
        s': 100, 'min child weight': 5, 'max depth': 5, 'learning rate': 0.2, 'colsample bytre
        Validation Accuracy: 0.9822485207100592
        Test Accuracy: 0.9667673716012085
In [28]: from sklearn.metrics import confusion matrix
         from sklearn.model selection import learning curve
         from yellowbrick.classifier import ClassPredictionError
         # Function for plotting learning curve
         def plot learning curve (estimator, title, X, y, ax=None):
             train sizes, train scores, test scores = learning curve(estimator, X, y, cv=10)
             train scores mean = np.mean(train scores, axis=1)
             test scores mean = np.mean(test scores, axis=1)
            ax.grid()
            ax.plot(train sizes, train scores mean, 'o-', color="r", label="Training score")
            ax.plot(train sizes, test scores mean, 'o-', color="g", label="Cross-validation scor
            ax.set title(title)
            ax.set xlabel("Training examples")
            ax.set ylabel("Score")
            ax.legend(loc="best")
         # Function for plotting confusion matrix
         def plot confusion matrix(y true, y pred, title, ax=None):
            cf matrix = confusion_matrix(y_true, y_pred)
            sns.heatmap(cf matrix, annot=True, fmt='d', cmap='Blues', ax=ax)
            ax.set title(f'Confusion Matrix for {title}')
            ax.set ylabel('True label')
             ax.set xlabel('Predicted label')
         # Function for plotting class prediction error
         def plot class prediction error(estimator, X, y, classes, title, ax=None):
            visualizer = ClassPredictionError(estimator, classes=classes, ax=ax)
            visualizer.fit(X, y)
            visualizer.score(X, y)
            ax.set title(title)
```

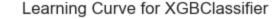
random search = RandomizedSearchCV(xgb classifier, param distributions, cv=5, scoring='a

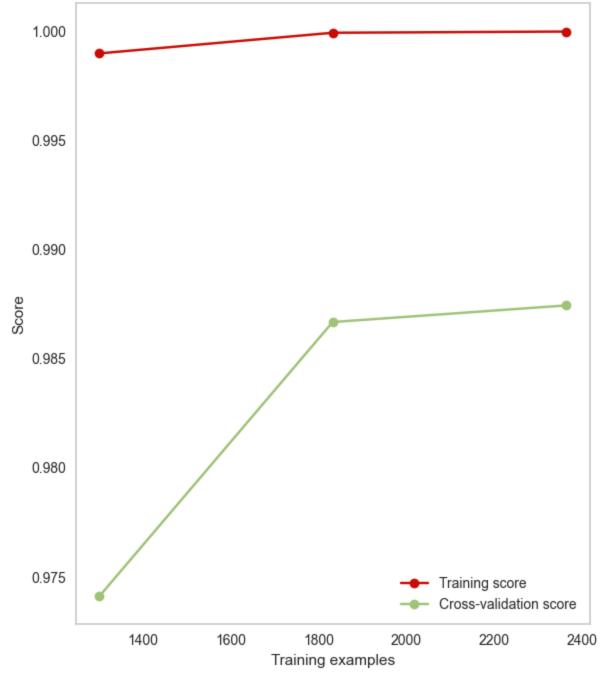
```
# Plotting
fig, axes = plt.subplots(nrows=3, ncols=1, figsize=(7, 21))

# Plot the learning curve
plot_learning_curve(best_classifier, 'Learning Curve for XGBClassifier', X_train_smote,

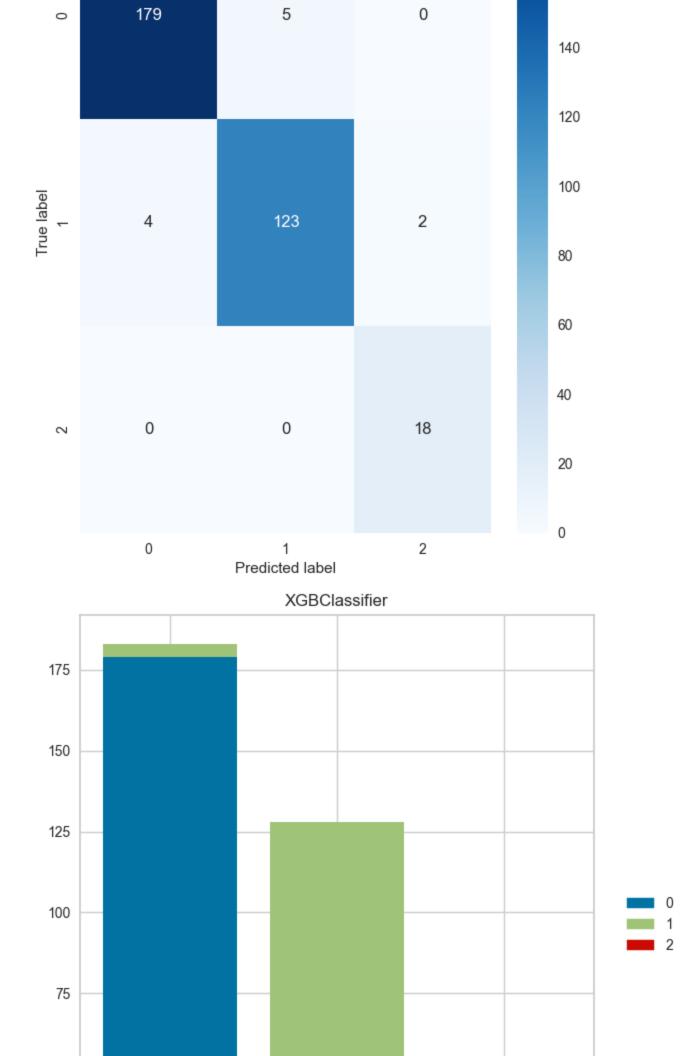
# Plot the confusion matrix
plot_confusion_matrix(y_test, y_test_pred, 'XGBClassifier', axes[1])

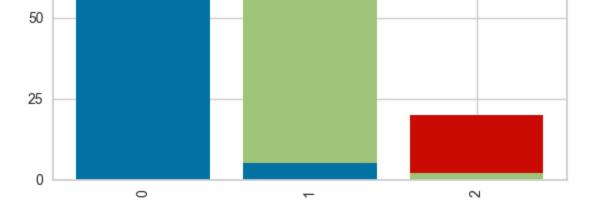
# Plot the class prediction error
plot_class_prediction_error(best_classifier, X_test, y_test, [0, 1, 2], 'XGBClassifier',
plt.tight_layout()
plt.show()
```





Confusion Matrix for XGBClassifier





### Conclusion

The purpose of the presented analysis was to develop an effective predictive model that can identify the target customers. Three separate experiments were conducted with varying features and data manipulation techniques, which included SMOTE for handling class imbalance.

The experiments compared a set of classification models including logistic regression, Naive Bayes, decision tree, random forest, SVM, K-Nearest Neighbors (KNN), gradient boosting (XGB Classifier), neural networks (MLP), and AdaBoost. Among these, the gradient boosting model, specifically XGBoost, consistently performed the best across all experiments.

Experiment 3, which incorporated nearly all features and applied SMOTE for oversampling, was the most successful in terms of performance metrics, implying that both a more comprehensive feature set and balanced data contribute positively to the model's performance.

Moreover, XGBoost's performance was verified by splitting the data into training, validation, and test sets. Hyperparameter optimization was also performed to ensure that the best parameters were selected for the final model. This was initially done with GridSearchCV, but due to computational constraints, RandomizedSearchCV was utilized as a more efficient alternative.

The accuracy scores from both validation and test data demonstrate that the XGBoost model generalized well, reducing the likelihood of overfitting. Confusion matrix, learning curve, and class prediction error plots further confirmed the model's good performance.

To conclude, the XGBoost model developed in this analysis has shown to be a strong predictor for the target customers. It has a good balance between bias and variance, which makes it a reliable tool for new, unseen data. Therefore, this model was chosen as the final model, and will be saved and deployed for further use.

### Save the Model

```
In [29]: import pickle

filename = 'final_model.sav'
pickle.dump(best_classifier, open(filename, 'wb'))

In [30]: # load the model from disk
loaded_model = pickle.load(open(filename, 'rb'))
```

### Load Model and evaluate the Model

```
In [31]: | # Load the model
         loaded model = pickle.load(open(filename, 'rb'))
         # New data
         data = {
             'Income': [24882.0, 22979.0, 27071.0, 36957.0, 70044.0],
             'Kidhome': [1, 1, 1, 1, 0],
             'Teenhome': [0, 0, 0, 1, 1],
             'Recency': [52, 29, 90, 43, 46],
             'MntWines': [1, 16, 8, 100, 1073],
             'MntFruits': [4, 17, 3, 2, 0],
             'MntMeatProducts': [10, 19, 19, 16, 250],
             'MntFishProducts': [29, 20, 0, 2, 153],
             'MntSweetProducts': [0, 21, 2, 1, 14],
             'MntGoldProds': [36, 22, 3, 31, 14],
             'NumDealsPurchases': [1, 3, 2, 4, 4],
             'NumWebPurchases': [1, 3, 2, 3, 7],
             'NumCatalogPurchases': [1, 2, 0, 2, 10],
             'NumStorePurchases': [2, 2, 3, 2, 5],
             'NumWebVisitsMonth': [6, 8, 6, 9, 5],
             'AcceptedCmp3': [1, 0, 0, 0, 0],
             'AcceptedCmp4': [0, 0, 0, 0, 0],
             'AcceptedCmp5': [0, 0, 0, 0, 0],
             'AcceptedCmp1': [0, 0, 0, 0, 0],
             'AcceptedCmp2': [0, 0, 0, 0, 0],
             'Complain': [0, 0, 0, 0, 0],
             'Response': [0, 1, 0, 1, 0],
             'Age': [45, 53, 55, 70, 69],
             'Total spend': [80, 115, 35, 152, 1504],
             'Total purchase': [5, 10, 7, 11, 26],
             'Total children': [1, 1, 1, 2, 1],
             'accepted camp': [1, 1, 0, 1, 0],
             'Total adults': [2, 2, 1, 1, 1],
             'Family size': [3, 3, 2, 3, 2],
             'Customer Since Years': [11, 11, 9, 11, 10],
             'Education encode': [1, 2, 3, 4, 2]
         # Convert to DataFrame
         new data = pd.DataFrame(data)
         # Make predictions using the loaded model
         predictions = loaded model.predict(new data)
         # Print the predictions
         print(predictions)
```

[0 0 0 0 1]

In [32]: X train.head(5)

ut[32]:		Income	Kidhome	Teenhome	Recency	MntWines	MntFruits	MntMeatProducts	MntFishProducts	MntSw
	190	24882.0	1	0	52	1	4	10	29	
	2002	22979.0	1	0	29	16	17	19	20	
	1842	27071.0	1	0	90	8	3	19	0	
	1409	36957.0	1	1	43	100	2	16	2	
	1735	70044.0	0	1	46	1073	0	250	153	

5 rows × 31 columns

```
In [33]: y_train.head(5)

Out[33]: 190      0
      2002      0
      1842      0
      1409      0
      1735      1
      Name: Cluster, dtype: int64
```

## Predicted values and and y\_train values matches

```
In [ ]:
```