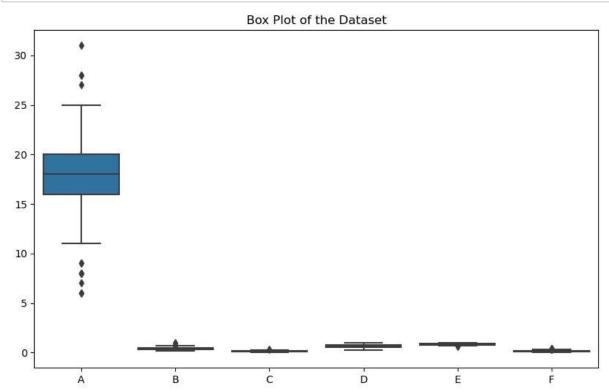
```
In [28]: import pandas as pd
import seaborn as sns
import matplotlib.pyplot as plt
from sklearn.ensemble import RandomForestRegressor
```

```
In [4]: # Read the CSV file into a DataFrame
df = pd.read_csv("Alumni Giving Regression (Edited).csv")

# Create a box plot
plt.figure(figsize=(10, 6))
sns.boxplot(data=df)
plt.title("Box Plot of the Dataset")
plt.show()
```



```
In [5]: # Function to identify and count outliers using IQR
def identify_outliers(column):
    Q1 = df[column].quantile(0.25)
    Q3 = df[column].quantile(0.75)
    IQR = Q3 - Q1
    lower_bound = Q1 - 1.5 * IQR
    upper_bound = Q3 + 1.5 * IQR
    outliers = df[(df[column] < lower_bound) | (df[column] > upper_bound)]
    return outliers

# Count and print outliers for each column
for column in df.columns:
    outliers = identify_outliers(column)
    print(f"Column '{column}': {len(outliers)} outliers")
```

```
Column 'A': 13 outliers
Column 'B': 7 outliers
Column 'C': 1 outliers
Column 'D': 0 outliers
Column 'E': 1 outliers
Column 'F': 7 outliers
```

#### Column 'A': 13 outliers

This suggests that there are 13 data points in Column 'A' that fall significantly outside the expected range based on the IQR. These points may be considered unusual or atypical compared to the majority of the data. Column 'B': 7 outliers

Similar to Column 'A', Column 'B' has 7 outliers. These values deviate from the typical range observed in the majority of the data. Column 'C': 1 outlier

Column 'C' has only 1 outlier, suggesting that most of the data in this column is within the expected range, with just a single data point being significantly different. Column 'D': 0 outliers

There are no outliers in Column 'D', indicating that the values in this column are relatively consistent and do not deviate significantly from the majority of the data. Column 'E': 1 outlier

Similar to Column 'C', Column 'E' has only 1 outlier, implying that the majority of the data in this column falls within the expected range. Column 'F': 7 outliers

Column 'F' has 7 outliers, suggesting that certain data points in this column significantly differ from the majority. These values may be considered unusual or anomalous.

#### Quantile method 1:

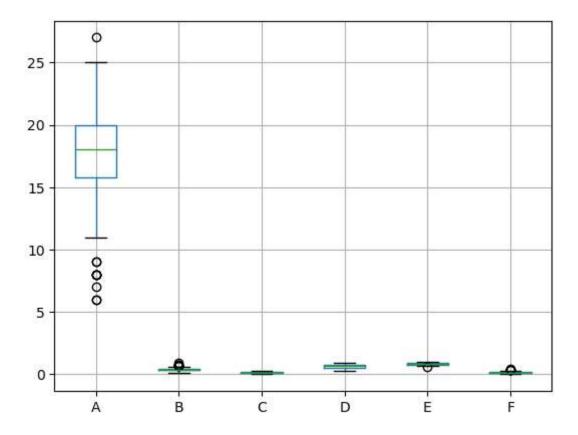
Calculates the 99th percentile (quantile) of the first column (iloc[:, 0]) and stores it in quantile99. Creates a new DataFrame df1 by excluding rows where the values in the first column are greater than the calculated 99th percentile. Plots a boxplot for the DataFrame df1.

```
In [15]: # Calculate the 99th percentile without missing values
   quantile99 = df.iloc[:, 0].dropna().quantile(0.99)

# Create a new DataFrame by excluding rows with missing values and values above
   df1 = df[df.iloc[:, 0].dropna() < quantile99]

# Plot a boxplot for the cleaned DataFrame
   df1.boxplot()</pre>
```

## Out[15]: <Axes: >

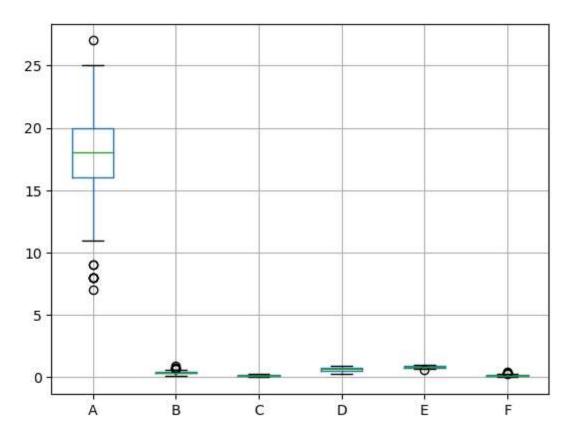


quantile method 2: Calculates the 1st percentile (quantile1) and the 99th percentile (quantile99) of the first column. Creates a new DataFrame df2 by including rows where the values in the first column are greater than the calculated 1st percentile and less than the 99th percentile. Plots a boxplot for the DataFrame df2.

```
In [16]: # Calculate the 1st and 99th percentiles without missing values
    quantile1 = df.iloc[:, 0].dropna().quantile(0.01)
    quantile99 = df.iloc[:, 0].dropna().quantile(0.99)

# Create a new DataFrame by including rows without missing values and values wordf2 = df[(df.iloc[:, 0].dropna() > quantile1) & (df.iloc[:, 0].dropna() < quantile1) & (df.iloc[:, 0].dropna() < quantile2.boxplot for the cleaned DataFrame
    df2.boxplot()</pre>
```

## Out[16]: <Axes: >



Differences: Quantile Calculation:

Method 1 calculates only the 99th percentile. Method 2 calculates both the 1st and 99th percentiles.

In [17]: pip install scikit-learn

Requirement already satisfied: scikit-learn in c:\users\anas\anaconda3\lib\si te-packages (1.3.0)
Requirement already satisfied: numpy>=1.17.3 in c:\users\anas\anaconda3\lib\s ite-packages (from scikit-learn) (1.24.3)
Requirement already satisfied: scipy>=1.5.0 in c:\users\anas\anaconda3\lib\si te-packages (from scikit-learn) (1.10.1)
Requirement already satisfied: joblib>=1.1.1 in c:\users\anas\anaconda3\lib\s ite-packages (from scikit-learn) (1.2.0)
Requirement already satisfied: threadpoolctl>=2.0.0 in c:\users\anas\anaconda3\lib\site-packages (from scikit-learn) (2.2.0)

Note: you may need to restart the kernel to use updated packages.

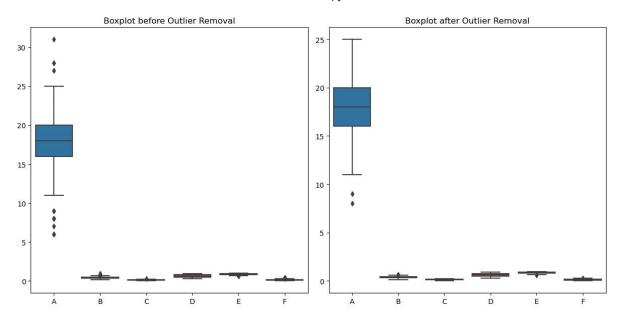
# **Outlier**

```
In [23]: import pandas as pd
         # Drop "NA" values from the DataFrame
         df cleaned = df.dropna()
         # Set the quantile threshold
         quantile_threshold = 0.99
        # Identify the upper bound for each column based on the quantile
         upper bound quantile = df cleaned.quantile(quantile threshold)
        # Calculate the lower and upper bounds for each column based on IQR
        Q1 = df cleaned.quantile(0.25)
        Q3 = df cleaned.quantile(0.75)
        IQR = Q3 - Q1
         lower bound = Q1 - 1.5 * IQR
         upper bound = Q3 + 1.5 * IQR
        # Identify outliers using the upper bound
        outliers = (df_cleaned > upper_bound).any(axis=1)
        # Display rows with outliers
        print("Rows with outliers:")
         print(df_cleaned[outliers])
         Rows with outliers:
              Α
                    В
                          C
                               D
                                     Ε
                                           F
         2
             18
                 0.24
                       0.17 0.66 0.87
                                        0.31
         3
                      0.00 0.81 0.88 0.11
                 0.74
         4
                 0.95
                      0.00 0.86 0.92 0.28
              8
         15
             19
                 0.46 0.16 0.67 0.85 0.34
         23
                 0.71 0.06 0.94 0.97 0.36
              8
         27
             28
                 0.23
                      0.22 0.46 0.81 0.06
         65
              7
                 0.75 0.06 0.94 0.97 0.30
         68
             27 0.33 0.16 0.50 0.78 0.09
         72
             12 0.55 0.10 0.96 0.98 0.41
         82
              6 0.69 0.08 0.92 0.97 0.32
         87
                 0.70 0.13 0.95 0.98 0.34
              6
         91
             28 0.23 0.16 0.51 0.86 0.15
         103 19
                 0.34 0.31 0.45 0.68 0.05
         106 31
                 0.28 0.23 0.63 0.86 0.16
         108
              9 0.63 0.11 0.89 0.97 0.39
         110 11 0.70 0.01 0.65 0.88 0.22
```

# **Outliers Removed**

In [ ]:

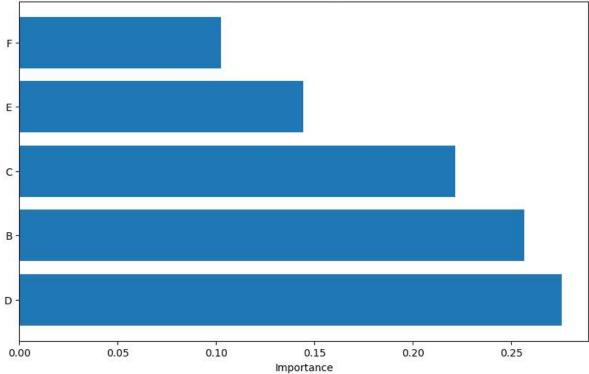
```
import pandas as pd
In [32]:
         import matplotlib.pyplot as plt
         import seaborn as sns
         # Drop "NA" values from the DataFrame
         df_cleaned = df.dropna()
         # Set the quantile threshold
         quantile_threshold = 0.99
         # Identify the upper bound for each column based on the quantile
         upper bound quantile = df cleaned.quantile(quantile threshold)
         # Calculate the lower and upper bounds for each column based on IQR
         Q1 = df_cleaned.quantile(0.25)
         Q3 = df cleaned.quantile(0.75)
         IOR = 03 - 01
         lower_bound = Q1 - 1.5 * IQR
         upper bound = Q3 + 1.5 * IQR
         # Identify outliers using the upper bound
         outliers = (df_cleaned > upper_bound).any(axis=1)
         # Remove rows with outliers
         df_no_outliers = df_cleaned[~outliers]
         # Plot boxplots before and after outlier removal
         plt.figure(figsize=(12, 6))
         plt.subplot(1, 2, 1)
         sns.boxplot(data=df_cleaned)
         plt.title('Boxplot before Outlier Removal')
         plt.subplot(1, 2, 2)
         sns.boxplot(data=df_no_outliers)
         plt.title('Boxplot after Outlier Removal')
         plt.tight_layout()
         plt.show()
```



```
import pandas as pd
In [26]:
         from sklearn.ensemble import RandomForestRegressor
         import matplotlib.pyplot as plt
         # Assuming df no outliers is your DataFrame without outliers
         # Separate the target variable (Column 'A') and features
         target = 'A'
         X_no_outliers = df_no_outliers.drop(target, axis=1)
         y no outliers = df no outliers[target]
         # Create a RandomForestRegressor model
         model no outliers = RandomForestRegressor()
         # Fit the model to the data without outliers
         model no outliers.fit(X no outliers, y no outliers)
         # Get feature importances from the model
         feature importances no outliers = model no outliers.feature importances
         # Create a DataFrame to display feature importances
         feature_ranking_no_outliers = pd.DataFrame({'Feature': X_no_outliers.columns,
         # Sort features by importance in descending order
         feature ranking no outliers = feature ranking no outliers.sort values(by='Impor
         # Display feature importances
         print("Feature Ranking without outliers:")
         print(feature_ranking_no_outliers)
         # Plot feature importances
         plt.figure(figsize=(10, 6))
         plt.barh(feature_ranking_no_outliers['Feature'], feature_ranking_no_outliers[']
         plt.xlabel('Importance')
         plt.title('Feature Importance Ranking without Outliers')
         plt.show()
```

Feature Ranking without outliers:





these importance values suggest the contribution of each feature to the predictive power of the model. Features with higher importance values are considered more influential in predicting the target variable.

```
import pandas as pd
In [27]:
         from sklearn.ensemble import RandomForestRegressor
         from sklearn.model_selection import train_test_split
         from sklearn.metrics import mean_squared_error
         import matplotlib.pyplot as plt
         # Assuming df no outliers is your DataFrame without outliers
         # Separate the target variable (Column 'A') and top three features
         target = 'A'
         top_three_features = ['D', 'B', 'C'] # Update with the top three features from
         X top three = df no outliers[top three features]
         y_top_three = df_no_outliers[target]
         # Split the data into training and testing sets
         X_train, X_test, y_train, y_test = train_test_split(X_top_three, y_top_three,
         # Create a RandomForestRegressor model
         model top three = RandomForestRegressor()
         # Fit the model to the training data
         model top three.fit(X train, y train)
         # Predict on the test set
         y pred = model top three.predict(X test)
         # Evaluate the model
         mse = mean_squared_error(y_test, y_pred)
         print(f"Mean Squared Error on the Test Set: {mse}")
         # Plot actual vs. predicted values
         plt.figure(figsize=(8, 6))
         plt.scatter(y_test, y_pred, alpha=0.5)
         plt.xlabel('Actual Values')
         plt.ylabel('Predicted Values')
         plt.title('Actual vs. Predicted Values for Top Three Features')
         plt.show()
```

Mean Squared Error on the Test Set: 7.237886363636364



