```
In [1]: from matplotlib import pyplot as plt
import numpy as np
import pandas as pd
import seaborn as sns
In [2]: # SET THE RANDOM STATE
RANDOM_STATE = 0
```

## 1) Load the Data

```
In [3]: data = pd.read_csv("data/cs-training.csv", index_col=0)

# Take a peak at the data
print(f"Num Entries: {data.shape[0]}")
print(f"Num Features: {data.shape[1]}")
display(data.head())
print(f"Features: {list(data.columns)}")
```

Num Entries: 150000 Num Features: 11

	Serious Dlqin 2 yrs	Revolving Utilization Of Unsecured Lines	age	Number Of Time 30- 59 Days Past Due Not Worse	DebtRatio	MonthlyInco
1	1	0.766127	45	2	0.802982	912
2	0	0.957151	40	0	0.121876	260
3	0	0.658180	38	1	0.085113	304
4	0	0.233810	30	0	0.036050	330
5	0	0.907239	49	1	0.024926	6358

Features: ['SeriousDlqin2yrs', 'RevolvingUtilizationOfUnsecuredLines', 'age', 'NumberOfT ime30-59DaysPastDueNotWorse', 'DebtRatio', 'MonthlyIncome', 'NumberOfOpenCreditLinesAndL oans', 'NumberOfTimes90DaysLate', 'NumberRealEstateLoansOrLines', 'NumberOfTime60-89Days PastDueNotWorse', 'NumberOfDependents']

## 2) Data Exploration

In [4]: data.describe()

Out[4]:

	Serious DIqin 2 yrs	$Revolving {\bf Utilization Of Unsecured Lines}$	age	Number Of Time 30-59 Days Past Due Not Worse	DebtR
count	150000.000000	150000.000000	150000.000000	150000.000000	150000.000
mean	0.066840	6.048438	52.295207	0.421033	353.00
std	0.249746	249.755371	14.771866	4.192781	2037.81
min	0.000000	0.000000	0.000000	0.000000	0.000
25%	0.000000	0.029867	41.000000	0.000000	0.17!
50%	0.000000	0.154181	52.000000	0.000000	0.360
75%	0.000000	0.559046	63.000000	0.000000	0.86
max	1.000000	50708.000000	109.000000	98.000000	329664.000

#### **Label Balance Check**

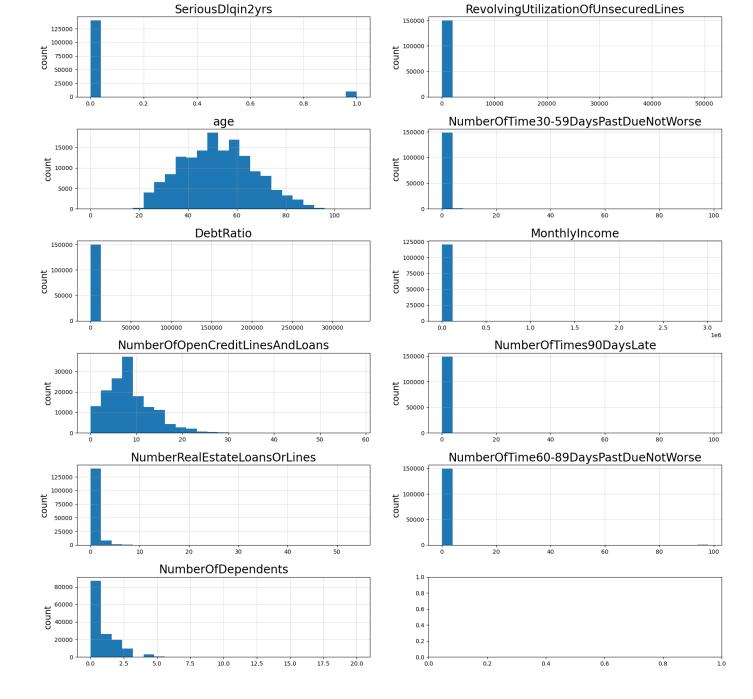
#### **Looking at Missing Values**

```
In [6]: # Print out the number of missing values for each feature
        data.isna().sum()
       SeriousDlqin2yrs
                                                   0
Out[6]:
       RevolvingUtilizationOfUnsecuredLines
                                                   0
       NumberOfTime30-59DaysPastDueNotWorse
       DebtRatio
       MonthlyIncome
                                              29731
       NumberOfOpenCreditLinesAndLoans
                                               0
       NumberOfTimes90DaysLate
       NumberRealEstateLoansOrLines
                                                  0
       NumberOfTime60-89DaysPastDueNotWorse
                                                 0
       NumberOfDependents
                                               3924
       dtype: int64
```

#### Visualizing the Distribution of Data for Each Feature

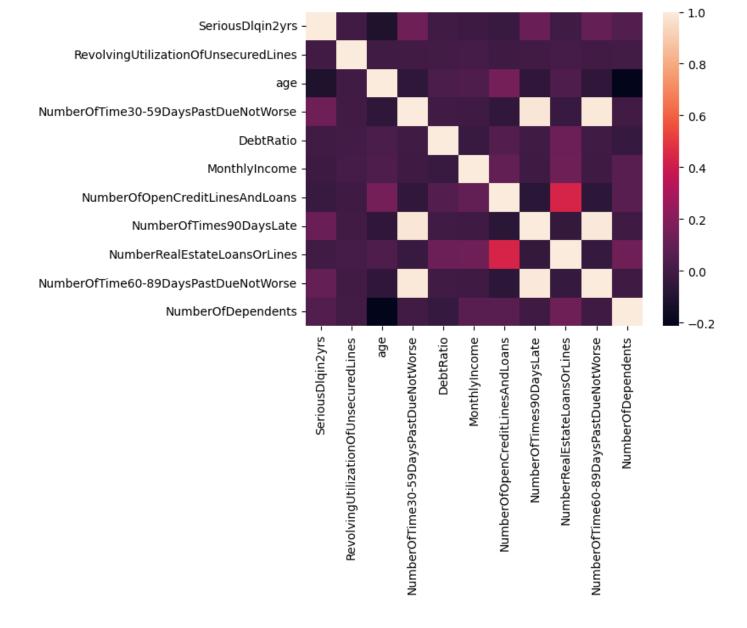
```
In [7]: # Generate an Image to Visualize the distribution
    fig, axs = plt.subplots(6, 2, figsize=(20,20))
    plt.subplots_adjust(hspace=0.4, wspace=0.2)

# Go through each feature and custom make a histogram for it
    for idx, col in enumerate(data.columns):
        fig_row, fig_col = idx//2, idx%2
        subplot = axs[fig_row, fig_col]
        subplot.hist(data[col], bins=25)
        subplot.set_title(col, fontsize=20)
        subplot.set_ylabel("count", fontsize=15)
        subplot.grid(alpha=0.4)
    plt.savefig("initial_hist.png")
```



## **Looking at Correlation Between Features**

Out[8]:



## 3) Data Wrangling

#### Extract the Features and the labels from the dataset

```
In [9]: X, y = data.iloc[:, 1:], data.iloc[:, 0]

print(f"Shape of X: {X.shape}")
 print(F"Shape of y: {y.shape}")

Shape of X: (150000, 10)
    Shape of y: (150000,)
```

#### Perform a statified shuffle split for splitting the taining and testing data

```
In [10]: from sklearn.model_selection import StratifiedShuffleSplit

# Create an instance of Stratified Shuffle Split
stratified_split = StratifiedShuffleSplit(n_splits=1, test_size=0.2, random_state=RANDOM)

# Perform the stratified shuffle split
for train_index, test_index in stratified_split.split(X, y):
    X_train, X_test = X.iloc[train_index], X.iloc[test_index]
```

```
y_train, y_test = y.iloc[train_index], y.iloc[test_index]

# Check the number of each type of label in the traning and test datasets
print(F"Delinquent Train: {y_train.sum()}")
print(F"Delinquent Test: {y_test.sum()}")
print()
print(f"Non-Delinquent Train: {(y_train==0).sum()}")
print(F"Non_Delinquent Test: {(y_test==0).sum()}")

Delinquent Train: 8021
Delinquent Train: 111979
Non_Delinquent Test: 27995
```

#### Use an imputer to fill the missing values

```
In [11]: from sklearn.impute import SimpleImputer
         # Create Instance of Imputer using median strategy
         imputer = SimpleImputer(strategy="median")
         # Fit to the training data
         imputer.fit(X train)
         # Save the index and columns before transformation
         index = X train.index
         cols = X train.columns
         # Perform the Imputer Tansformation
         X train = imputer.transform(X train)
         X train = pd.DataFrame(X train, columns=cols,
                                   index=index)
         # Check that the missing values were filled
        X train.isna().sum()
        RevolvingUtilizationOfUnsecuredLines
                                                 0
Out[11]:
                                                 0
        NumberOfTime30-59DaysPastDueNotWorse
        DebtRatio
        MonthlyIncome
        NumberOfOpenCreditLinesAndLoans
                                                 0
        NumberOfTimes90DaysLate
        NumberRealEstateLoansOrLines
        NumberOfTime60-89DaysPastDueNotWorse
                                                 0
        NumberOfDependents
        dtype: int64
```

#### Balance the Labels by undersampling the Major Class

```
In [12]: from sklearn.utils import resample
    from sklearn.utils import shuffle

# Need to recombine the X_train and y_train for the down_sampling
    train = X_train.copy()
    train["SeriousDlqin2yrs"] = y_train

# Get the counts for the minor class
    df_minor = train[train.SeriousDlqin2yrs == 1]
    df_major = train[train.SeriousDlqin2yrs == 0]

# Let's downsample the major class
    df_major_downsample = resample(df_major,
```

```
replace=False,
                               n samples=df minor.shape[0],
                               random state=RANDOM STATE)
# Combine the two classes back together
df downsample = pd.concat([df major downsample, df minor])
# Check that the two labels are equal
print("Check for if the labels are balanced: ")
print(df downsample.SeriousDlqin2yrs.value counts())
# Shuffle the dataframe before splitting
shuffled = shuffle(df downsample, random state=RANDOM STATE)
# Split the Downsampled data back into features & label
X train bal, y train bal = shuffled.iloc[:, :-1], shuffled["SeriousDlqin2yrs"]
# Check for proper splitting
print()
print(f"Shape of X: {X train bal.shape}")
print(f"Shape of y: {y train bal.shape}")
Check for if the labels are balanced:
0 8021
   8021
Name: SeriousDlqin2yrs, dtype: int64
Shape of X: (16042, 10)
Shape of y: (16042,)
```

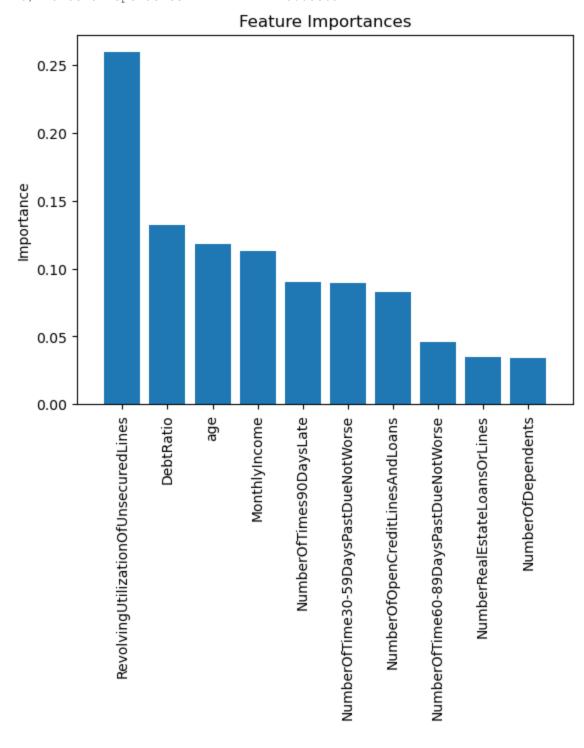
#### **Looking at Feature Importances**

```
In [13]:
        from sklearn.ensemble import RandomForestClassifier
         from sklearn.preprocessing import StandardScaler
         # Perform Standard Scaling
         scaler = StandardScaler()
         scaler.fit(X train bal)
         X scale = scaler.transform(X train bal)
         # Fit a random forest model to find feature importances
         feat forest = RandomForestClassifier(n estimators=500,
                                               random state=RANDOM STATE)
         feat forest.fit(X_scale, y_train_bal)
         importances = feat forest.feature importances
         # Sort features in descending order of importance
         # their indices are what should be sorted
         indices = np.argsort(importances)[::-1]
         # Define the features
         features = X_train_bal.columns
         # Print out feature importances
         for f in range(X scale.shape[1]):
            print("%2d) %-*s %f" % (f + 1, 30,
                                     features[indices[f]],
                                     importances[indices[f]]))
         # Make a figure showing the importances
         fig, axs = plt.subplots()
         axs.bar(range(X scale.shape[1]),
                importances[indices],
```

1) RevolvingUtilizationOfUnsecuredLines 0.259576
2) DebtRatio 0.132057
3) age 0.118252
4) MonthlyIncome 0.113137
5) NumberOfTimes90DaysLate 0.090206
6) NumberOfTime30-59DaysPastDueNotWorse 0.089476
7) NumberOfOpenCreditLinesAndLoans 0.082814
8) NumberOfTime60-89DaysPastDueNotWorse 0.045476

9) NumberRealEstateLoansOrLines 0.035114

10) NumberOfDependents 0.033892



#### **Peform Dimensionality Reduction**

#### A) Principal Component Analysis

```
In [14]:
         from sklearn.decomposition import PCA
         from sklearn.pipeline import make pipeline
         # Scale the balanced label training set
         scaler = StandardScaler()
         scaler.fit(X train bal)
         X bal scaled = scaler.transform(X train bal)
         # Perform Principal Component Analysis
         pca = PCA()
         pca.fit(X bal scaled, y train bal)
         # Calculate cumulative sum of EVR
         cum sum = np.cumsum(pca.explained variance ratio )
         # Plot the Explain Variance by # Principal Components
         print(f"Explained Variance Ratio: \n{pca.explained_variance_ratio_}")
         print(F"Cumulative Sum of EVR: \n {cum sum}")
         plt.bar(range(1, 11), pca.explained variance ratio , alpha=0.5, align='center')
         plt.step(range(1, 11), cum sum, where='mid')
         plt.ylabel('Explained variance ratio')
         plt.xlabel('Principal components')
         plt.show()
         Explained Variance Ratio:
         [0.30587879 \ 0.16867738 \ 0.11931453 \ 0.10115368 \ 0.09935303 \ 0.07923475
          0.07456405 0.05026824 0.00102548 0.00053007]
         Cumulative Sum of EVR:
          \lceil 0.30587879 \ 0.47455617 \ 0.5938707 \ 0.69502438 \ 0.79437741 \ 0.87361216 
          0.94817621 0.99844445 0.99946993 1.
            1.0
            0.8
         Explained variance ratio
            0.6
            0.4
            0.2
            0.0
```

```
In [15]: # Set the Explained Variance Threshold required for
    # determining number of principal components to use
    THRESHOLD EVR = 0.9
```

Principal components

6

10

8

2

```
# Calculate num_components
n_components = np.argmax(cum_sum >= THRESHOLD_EVR) + 1
print(f"Number of Pricipal Components for PCA: {n_components}")

# Principal Component Analysis
pca = PCA(n_components=n_components)
X_train_bal_pca = pca.fit_transform(X_bal_scaled)
print(F"New shape of the data: {X_train_bal_pca.shape}")

Number of Pricipal Components for PCA: 7
New shape of the data: (16042, 7)
```

#### **B) Linear Discriminant Analysis**

```
In [16]: from sklearn.discriminant_analysis import LinearDiscriminantAnalysis as LDA

# Scale the balanced label training set
scaler = StandardScaler()
scaler.fit(X_train_bal)
X_bal_scaled = scaler.transform(X_train_bal)

# Perform Linear Discriminant Analysis
lda = LDA(n_components=1)
X_train_bal_lda = lda.fit_transform(X_bal_scaled, y_train_bal)
```

## 4) Building and Testing Models

```
In [17]: # Define the number of trials to run
# when performing nested cross validation
NUM_TRIALS = 10
```

## Nested Cross Validation Code to Assess Each Model Type to Determine which to move forward with

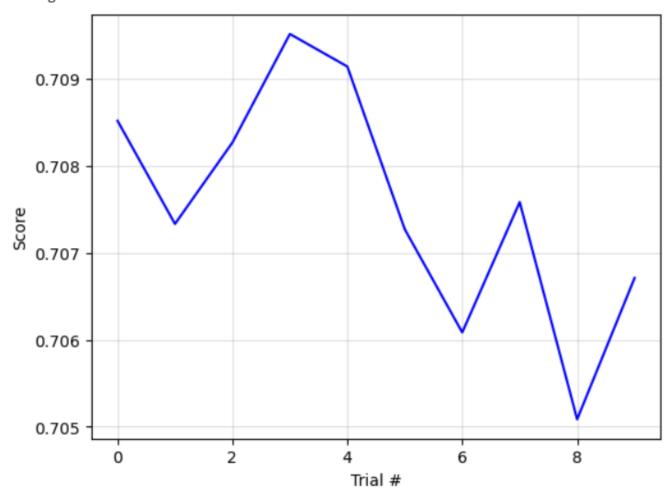
```
In [18]: def perform_nested_CV(num_trials, model, p grid, X, y):
             Returns a list of the scores and plots the nested
             cross-validation scors for each trial.
             # Make an array to store the socres
             nested scores = np.zeros(num trials)
             # Loop for each trial
             for i in range(num trials):
                # Choose cross-validation techniques for the inner and outer loops,
                 # independently of the dataset.
                 # E.g "GroupKFold", "LeaveOneOut", "LeaveOneGroupOut", etc.
                 inner cv = KFold(n splits=4, shuffle=True, random state=i)
                 outer cv = KFold(n splits=4, shuffle=True, random state=i)
                 # Perform parameter search for optimal model
                 # on the inner CV
                 clf = GridSearchCV(estimator=model,
                                    param grid=p grid,
                                    cv=inner cv)
                 # Perform outer CV to score the optimal model
                 # found on the inner CV
```

```
nested score = cross val score(clf, X=X, y=y, cv=outer cv)
       nested scores[i] = nested score.mean()
    return nested scores
def plot CV scores(nested scores):
   Generates a figure that plots the nested scores
   of the cross validations performed.
   # Print the average across all of the nested scores
   print(f"Average CV Score: {np.mean(nested scores)}")
   # Plot the scores for each of the nested CV
   # in nested scores file
   fig, axs = plt.subplots()
   axs.plot(nested scores, color="b")
   axs.set ylabel("Score")
   axs.set xlabel("Trial #")
   axs.grid(alpha=0.4)
   return fig
```

## 1) Logistic Regression Model - Nested CV - (Supervised)

Code is commented out because it takes a significant amount of time to run. The results can be visualized in the cell below the code.

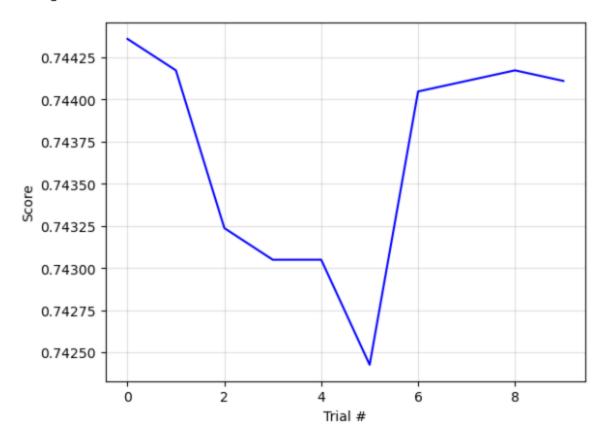
Average CV Score: 0.70754917897229



## 2) Support Vector Machine (SVM) - Nested CV - (Supervised)

Code is commented out because it takes a significant amount of time to run. The results can be visualized in the cell below the code.

Average CV Score: 0.7436730257378245



## 3) K-Means Clustering - (Unsupervised)

#### A) Perform PCA Dimensionality Reduction to 2-D

```
In [21]: # Principal Component Analysis for 2 Principal Components
         pca 2 = PCA(n components=2)
         X train bal pca 2 = pca 2.fit transform(X bal scaled)
         print(F"New shape of the data: {X train bal pca 2.shape}")
         print()
         # Determine contribution of each feature to each principal component
         dataset pca 2 = pd.DataFrame(pca 2.components , columns=cols, index=["PC 1", "PC 2"])
         # Most Important Features for PC 1
         print("Most Important Features for PC 1: ")
         print(dataset pca 2[abs(dataset pca 2) > 0.3].iloc[0].dropna())
        print()
         # Most Important Features for PC 2
         print("Most Important Features for PC 2: ")
        print(dataset pca 2[abs(dataset pca 2) > 0.3].iloc[1].dropna())
         print()
        New shape of the data: (16042, 2)
        Most Important Features for PC 1:
        NumberOfTime30-59DaysPastDueNotWorse 0.560500
        NumberOfTimes90DaysLate
                                                0.563166
        NumberOfTime60-89DaysPastDueNotWorse 0.562465
        Name: PC 1, dtype: float64
        Most Important Features for PC 2:
        MonthlyIncome
                                            0.443187
```

NumberOfOpenCreditLinesAndLoans 0.563457 NumberRealEstateLoansOrLines 0.592006 Name: PC 2, dtype: float64

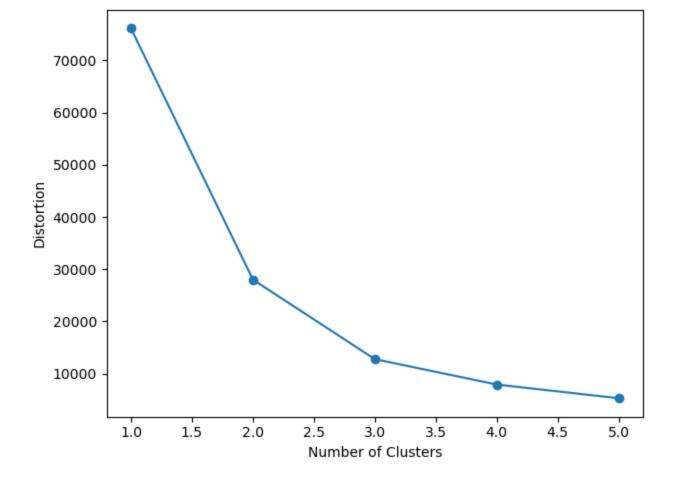
#### B) Elbow Plot

```
In [22]: def elbow_plot(data):
             # Get the distortions for each cluster number
             distortions = []
             for i in range (1, 6):
                 km = KMeans(n clusters=i,
                         init='k-means++',
                         n init=10,
                         max iter=300,
                         random state=0)
                 km.fit(data)
                 distortions.append(km.inertia )
             # Make the figure
             fig, axs = plt.subplots()
             axs.plot(range(1, 6), distortions, marker="o")
             axs.set xlabel("Number of Clusters")
             axs.set ylabel("Distortion")
             plt.tight layout()
             return fig
```

```
In [23]: from sklearn.cluster import KMeans

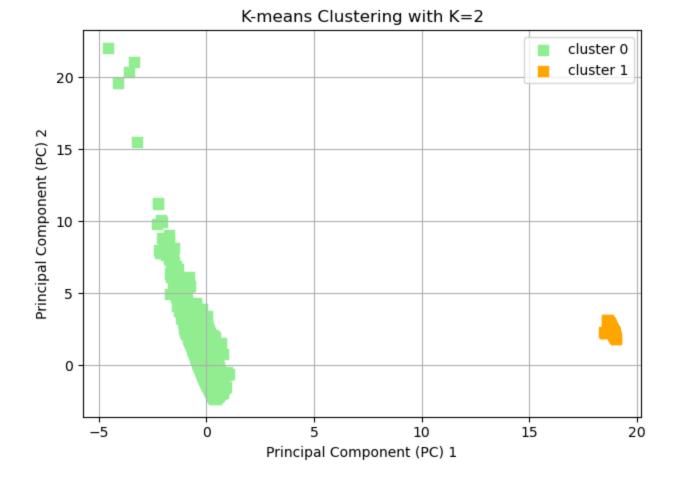
# Make an elbow plot
plt.show(elbow_plot(X_train_bal_pca_2))

# Based on the elbow plot, I would say that the
# ideal number of clusters is 2. After two clusters
# the distortion does not drastically decrease per cluster added.
```



#### C) Perform K-means Clustering

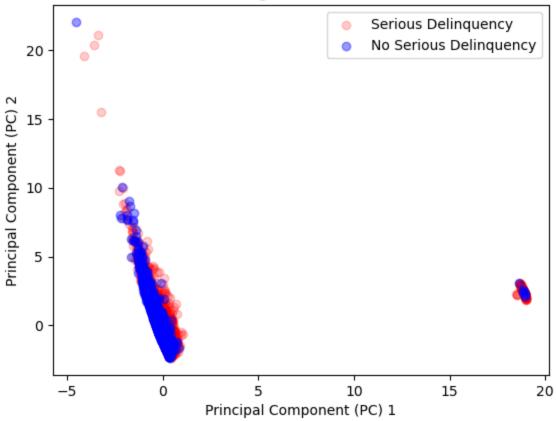
```
def print cluster(model, n clusters, X):
In [24]:
             y km = model.fit predict(X)
             color list = ['lightgreen', 'orange', 'lightblue', 'red', 'yellow', 'brown', 'cyan']
             for i in range(n clusters):
                 plt.scatter(X[y km == i, 0],
                 X[y km == i, 1],
                 s=50,
                 c=color list[i],
                 marker='s',
                 label='cluster ' + str(i))
             plt.xlabel("Principal Component (PC) 1")
             plt.ylabel("Principal Component (PC) 2")
             plt.title(f"K-means Clustering with K={n clusters}")
             plt.legend()
             plt.grid()
             plt.tight layout()
             plt.show()
         km = KMeans(n clusters=2,
                 init='k-means++',
                 n init=10,
                 max iter=300,
                 tol=1e-04,
                 random state=0)
        print cluster(km, 2, X train bal pca 2)
```



#### D) Analyze Labels in Clusters

```
In [25]: fig, axs = plt.subplots()
         axs.scatter(X train bal pca 2[y train bal == 1, 0],
                     X train bal pca 2[y train bal == 1, 1],
                     color="red",
                     alpha=0.2
                     label="Serious Delinquency")
         axs.scatter(X train bal pca 2[y train bal == 0, 0],
                     X train bal pca 2[y train bal == 0, 1],
                     color="blue",
                     alpha=0.4,
                    label="No Serious Delinquency")
         axs.legend()
         plt.xlabel("Principal Component (PC) 1")
         plt.ylabel("Principal Component (PC) 2")
         plt.title(f"Visualizing Labels of Clusters")
         plt.show()
```

#### Visualizing Labels of Clusters

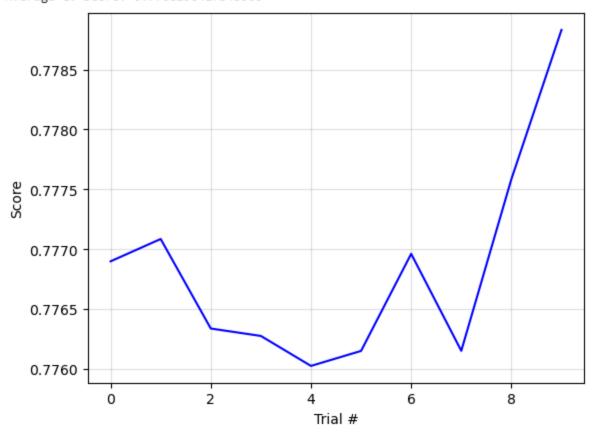


## 4) Random Forest - Nested CV - (Supervised)

Code is commented out because it takes a significant amount of time to run. The results can be visualized in the cell below the code.

```
In [26]:
         from sklearn.ensemble import RandomForestClassifier
         # # Set up possible values of parameters to optimize
         # p grid = {"max depth": [4, 6, 8],
                    "min samples leaf": [30, 100]}
         # # Create the random forest classifier
         # rfc = RandomForestClassifier(n estimators=10,
                                        criterion='gini',
                                        random state=RANDOM STATE)
         # # Define the features and label
         # # for nested CV
         # X = X train bal
         # y = y train bal
         # # Run the nested CV
         # nested CV scores = perform nested CV(10, rfc, p grid, X, y)
         # plt.show(plot CV scores(nested CV scores))
```





## 5) Random Forest Classifier Hyperparameter Tuning

## **Code for Testing Parameter Values**

```
In [27]:
         from sklearn.model selection import validation curve
         def plot scores(param name, train mean, train std, test mean, test std):
             Makes a plot of the parameter values and their
             corresponding training and testing accuracies.
             fig, axs = plt.subplots()
             # Plot the accuracy values
             axs.plot(param range, train mean,
                      color="blue",
                      marker="o",
                      markersize=5,
                      label="Training")
             axs.plot(param range,
                      test mean,
                      color="green",
                      linestyle="--",
                      marker="s",
                      markersize=5,
                      label="Validation")
             # Plot the standard deviation range
             axs.fill between (param range,
                              train mean + train std,
```

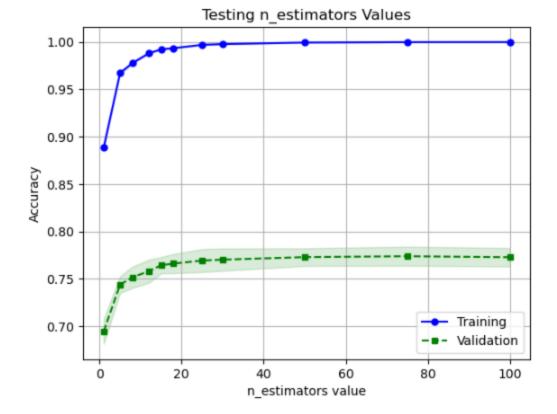
```
train mean - train std,
                     alpha=0.15,
                     color="blue")
    axs.fill between (param range,
                    test mean + test std,
                     test mean - test std,
                     alpha=0.15,
                     color="green")
    # Make the formatting better
   axs.set title(f"Testing {param name} Values")
    axs.set ylabel("Accuracy")
    axs.set xlabel(f"{param name} value")
   plt.grid(0.4)
   plt.legend()
   return fig
def run param test(model, param name, param range, X train, y train):
   Perform a validation curve and plot the results.
   model: sklearn model to test
   param name: string matching parameter to test for model
   param 1st: possible values for testing given parameter
   X train: data to use to train and test the model
    y train: labels for the data
    # Run the validation curve
    train scores, test scores = validation curve(estimator=model,
                                                 X=X train,
                                                 y=y train,
                                                 param name=param name,
                                                 param range=param_range,
                                                  cv=10)
    # Calculate the mean and standard deivatins
    train mean = np.mean(train scores, axis=1)
    train std = np.std(train scores, axis=1)
    test mean = np.mean(test scores, axis=1)
    test std = np.std(test scores, axis=1)
    # Plot the results
   plt.show(plot scores(param name, train mean, train std, test mean, test std))
    # return dictionary
    return train scores, test scores
```

## A) Testing N\_Estimators

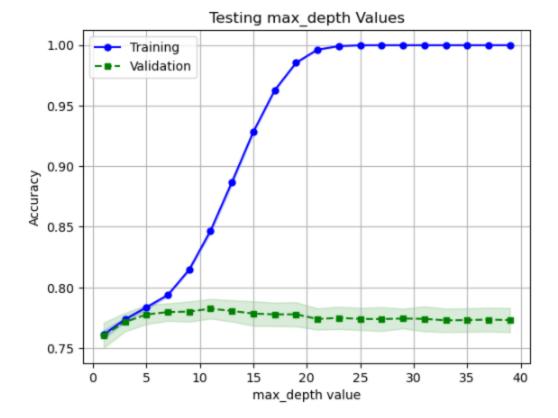
```
In [28]: # # Define the model for testing
# rfc = RandomForestClassifier(criterion="gini",
# random_state=RANDOM_STATE)

# # Define the parameters for testing
# param_name = 'n_estimators'
# param_range = [1, 5, 8, 12, 15, 18, 25, 30, 50, 75, 100]

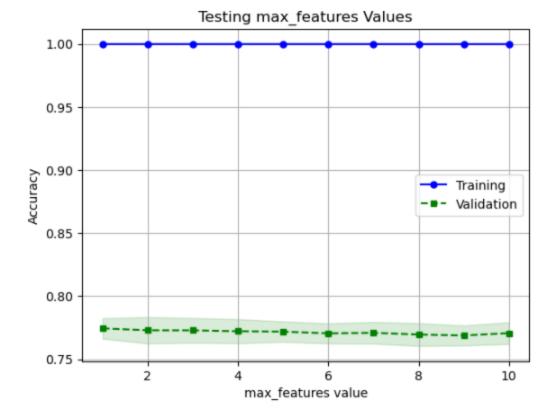
# # run the testing
# train_scores, test_scores = run_param_test(rfc, param_name, param_range, X_train_bal,
```



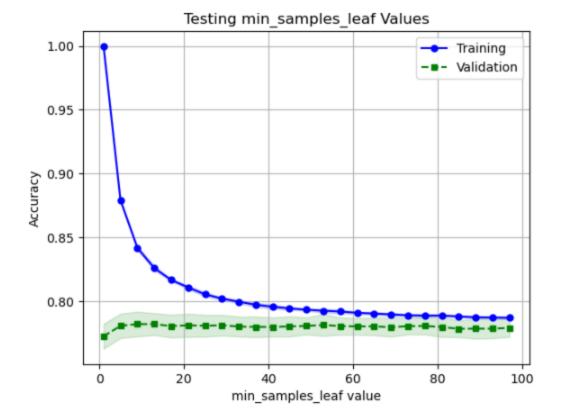
## B) Testing Max Depth



## C) Testing Max Features



## D) Testing Min Samples Leaf



## E) Perform Randomized CV Search

In [32]:

from sklearn.model selection import RandomizedSearchCV

```
from scipy.stats import randint
         from sklearn.ensemble import RandomForestClassifier
         # Make the parameters to search
         param distribs = {'n estimators': randint(low=15, high=25),
                           'max_features': randint(low=1, high=10),
                           'max depth': randint(low=3, high=8),
                           'min samples leaf': randint(low=45, high=100)
         # Define the model
         rfc = RandomForestClassifier(criterion="gini",
                                       random state=RANDOM STATE)
         # Perform the randomized search
         rnd search = RandomizedSearchCV(rfc,
                                          param distributions=param_distribs,
                                          n iter=50,
                                          cv=5,
                                          scoring='accuracy',
                                          random state=RANDOM STATE)
         rnd search.fit(X train bal, y train bal)
         RandomizedSearchCV(cv=5, estimator=RandomForestClassifier(random state=0),
Out[32]:
                            n iter=50,
                            param distributions={'max depth': <scipy.stats. distn infrastructure.</pre>
         rv discrete frozen object at 0x000002C37CF8E130>,
                                                  'max features': <scipy.stats. distn infrastructu
         re.rv discrete frozen object at 0x000002C37CF39EB0>,
                                                  'min_samples_leaf': <scipy.stats. distn infrastr</pre>
         ucture.rv discrete frozen object at 0x000002C30827DF70>,
                                                  'n estimators': <scipy.stats. distn infrastructu
         re.rv discrete frozen object at 0x000002C37CF39DF0>},
                            random state=0, scoring='accuracy')
In [33]: print(f"Best Model Accuracy: {rnd search.best score }")
```

```
Best Model Accuracy: 0.7811989968985296
        Best Model Parameters: {'max depth': 7, 'max features': 5, 'min samples leaf': 48, 'n es
        timators': 19}
In [34]: # Collect the results
         cvres = rnd search.cv results
         # Iterate through scores and params and add to results list
         results = []
         for mean score, params in zip(cvres["mean test score"], cvres["params"]):
            results.append((mean score, params))
         # Print score and parameters in decreasing order of accuracy
         results.sort(key=lambda x: -x[0])
         for result in results:
            print(*result)
        0.7811989968985296 {'max depth': 7, 'max features': 5, 'min samples leaf': 48, 'n estima
        tors': 19}
        0.780824776637403 {'max depth': 7, 'max features': 6, 'min samples leaf': 45, 'n estimat
        ors': 18}
        0.7803885036551655 {'max depth': 7, 'max features': 9, 'min samples leaf': 59, 'n estima
        tors': 19}
        0.7800769966638407 {'max depth': 7, 'max features': 9, 'min samples leaf': 69, 'n estima
        tors': 18}
        0.7797028346864219 {'max depth': 6, 'max features': 8, 'min samples leaf': 47, 'n estima
        tors': 24}
        0.7796406265420898 {'max depth': 7, 'max features': 7, 'min samples leaf': 66, 'n estima
        tors': 22}
        0.7792664645646713 {'max depth': 7, 'max features': 4, 'min samples leaf': 90, 'n estima
        0.7792040038576044 {'max depth': 6, 'max features': 6, 'min samples leaf': 95, 'n estima
        tors': 19}
        0.7788924774383765 {'max depth': 6, 'max features': 9, 'min samples leaf': 71, 'n estima
        tors': 17}
        0.778767730875367 {'max depth': 6, 'max features': 6, 'min samples leaf': 57, 'n estimat
        ors': 24}
        0.7785807178843169 {'max depth': 6, 'max features': 6, 'min samples leaf': 60, 'n estima
        tors': 15}
        0.7784561461724312 {'max depth': 6, 'max features': 4, 'min samples leaf': 84, 'n estima
        tors': 24}
        0.7784560101771125 {'max depth': 6, 'max features': 4, 'min samples leaf': 64, 'n estima
        tors': 22}
        0.7784557187585726 {'max depth': 5, 'max features': 4, 'min samples leaf': 87, 'n estima
        tors': 21}
        0.7783313996094214 {'max depth': 6, 'max features': 9, 'min samples leaf': 56, 'n estima
        tors': 17}
        0.7783312441862 {'max depth': 5, 'max features': 5, 'min samples leaf': 56, 'n estimator
        0.7776455169337485 {'max depth': 6, 'max features': 8, 'min samples leaf': 60, 'n estima
        tors': 20}
        0.7766482049783611 {'max depth': 5, 'max features': 4, 'min samples leaf': 65, 'n estima
        tors': 16}
        0.776585705415489 {'max depth': 5, 'max features': 4, 'min samples leaf': 75, 'n estimat
        ors': 18}
        0.7764613085547273 {'max depth': 6, 'max features': 2, 'min samples leaf': 85, 'n estima
        tors': 15}
        0.7763989061313683 {'max depth': 6, 'max features': 3, 'min samples leaf': 73, 'n estima
        tors': 17}
        0.7758377311628999 {'max depth': 5, 'max features': 4, 'min samples leaf': 69, 'n estima
        tors': 16}
        0.7752770613199006 {'max depth': 6, 'max features': 3, 'min samples leaf': 72, 'n estima
        tors': 15}
        0.7751521787615723 {'max depth': 5, 'max features': 9, 'min samples leaf': 99, 'n estima
```

print(f"Best Model Parameters: {rnd\_search.best\_params\_}")

```
tors': 21}
0.7750896597707972 {'max depth': 5, 'max features': 4, 'min samples leaf': 75, 'n estima
tors': 17}
0.7733448204045822 {'max depth': 6, 'max features': 1, 'min samples leaf': 76, 'n estima
tors': 20}
0.7727837620035297 {'max depth': 5, 'max features': 2, 'min samples leaf': 81, 'n estima
tors': 21}
0.7725967101566743 {'max depth': 4, 'max features': 9, 'min samples leaf': 46, 'n estima
tors': 16}
0.7724722744401074 {'max depth': 4, 'max features': 2, 'min samples leaf': 76, 'n estima
tors': 21}
0.7724721384447887 {'max depth': 5, 'max features': 1, 'min samples leaf': 84, 'n estima
tors': 20}
0.77228528087696 {'max depth': 6, 'max features': 1, 'min samples leaf': 77, 'n estimato
rs': 21}
0.7717240087689782 {'max depth': 3, 'max features': 2, 'min samples leaf': 54, 'n estima
tors': 24}
0.7717240087689782 {'max depth': 3, 'max features': 2, 'min samples leaf': 50, 'n estima
tors': 24}
0.7713499050752676 {'max depth': 4, 'max features': 9, 'min samples leaf': 61, 'n estima
tors': 23}
0.7710386312187746 {'max depth': 6, 'max features': 1, 'min samples leaf': 85, 'n estima
tors': 23}
0.7706642555344267 {'max depth': 3, 'max features': 2, 'min samples leaf': 82, 'n estima
tors': 24}
0.7700410861285553 {'max depth': 4, 'max features': 1, 'min samples leaf': 55, 'n estima
tors': 17}
0.7698538205747706 {'max depth': 4, 'max features': 4, 'min samples leaf': 83, 'n estima
tors': 22}
0.7683577555021763 {'max depth': 3, 'max features': 5, 'min samples leaf': 68, 'n estima
tors': 18}
0.7679838655153952 {'max depth': 3, 'max features': 5, 'min samples leaf': 48, 'n estima
tors': 15}
0.7679836712363685 {'max depth': 4, 'max features': 1, 'min samples leaf': 99, 'n estima
tors': 20}
0.7676721836729461 {'max depth': 3, 'max features': 5, 'min samples leaf': 62, 'n estima
tors': 19}
0.7676720476776274 {'max depth': 5, 'max features': 1, 'min samples leaf': 74, 'n estima
tors': 18}
0.7672983131140675 {'max depth': 3, 'max features': 5, 'min samples leaf': 96, 'n estima
tors': 15}
0.7661138327444089 {'max depth': 3, 'max features': 2, 'min samples leaf': 95, 'n estima
0.7657399621855302 {'max depth': 3, 'max features': 1, 'min samples leaf': 81, 'n estima
tors': 20}
0.7583215341204483 {'max depth': 3, 'max features': 9, 'min samples leaf': 57, 'n estima
tors': 16}
0.758134482273593 {'max depth': 3, 'max features': 9, 'min samples leaf': 60, 'n estimat
ors': 24}
0.7580721575618449 {'max depth': 3, 'max features': 9, 'min samples leaf': 58, 'n estima
tors': 23}
0.7577604368635905 {'max depth': 3, 'max features': 7, 'min samples leaf': 64, 'n estima
tors': 18}
```

## 6) Training Final Model Based on the Randomized Search

```
In [35]: print(f"Best Model Parameters: {rnd_search.best_params_}")

Best Model Parameters: {'max_depth': 7, 'max_features': 5, 'min_samples_leaf': 48, 'n_es timators': 19}
```

In [36]: # Define the model

Out[36]: RandomForestClassifier(max\_depth=7, max\_features=5, min\_samples\_leaf=48, n estimators=19, random state=0)

# 7) Preparing Test Data And Testing Model for Accuracy

#### **Preparing Test Data**

```
In [37]: # Check missing values before using the imputer
         X train.isna().sum()
Out[37]: RevolvingUtilizationOfUnsecuredLines
                                                 \cap
        NumberOfTime30-59DaysPastDueNotWorse
         DebtRatio
        MonthlyIncome
         NumberOfOpenCreditLinesAndLoans
         NumberOfTimes90DaysLate
         NumberRealEstateLoansOrLines
        NumberOfTime60-89DaysPastDueNotWorse
                                                 0
        NumberOfDependents
         dtype: int64
In [38]: # Prepare the test data in the same manner as the training data
         # Save the index and columns before transformation
         index = X test.index
         cols = X test.columns
         # Fill the missing values with same values used in training set
         X test = imputer.transform(X test)
         X test = pd.DataFrame(X test,
                                columns=cols,
                                index=index)
         # Check that the missing values were filled
         X test.isna().sum()
        RevolvingUtilizationOfUnsecuredLines
                                                 0
Out[38]:
        NumberOfTime30-59DaysPastDueNotWorse
         DebtRatio
        MonthlyIncome
        NumberOfOpenCreditLinesAndLoans
        NumberOfTimes90DaysLate
         NumberRealEstateLoansOrLines
         NumberOfTime60-89DaysPastDueNotWorse
                                                 0
         NumberOfDependents
         dtype: int64
```

```
from sklearn.metrics import accuracy score
In [39]:
         from sklearn.metrics import f1 score
         # Make the predictions
         y pred = final rfc.predict(X test)
         # Calculate probabilities for ROC curve
         y probs = final rfc.predict proba(X test)[:, 1]
In [40]:
        TN = np.sum((y pred==0) & (y test==0))
         TP = np.sum((y pred==1) & (y test==1))
         FP = np.sum((y pred==1) & (y test==0))
         FN = np.sum((y pred==0) & (y test==1))
         print(F"Number of Samples: {len(y pred)}")
         print(F"Number Positive: {np.sum(y test==1)}")
         print(F"Number Negative: {np.sum(y test==0)}")
        print("----")
         print(F"True Positive (TN): {TP}")
         print(F"True Negative (TN): {TN}")
        print(F"False Positive (TN): {FP}")
         print(F"False Negative (TN): {FN}")
        Number of Samples: 30000
        Number Positive: 2005
        Number Negative: 27995
         _____
        True Positive (TN): 1579
        True Negative (TN): 21815
        False Positive (TN): 6180
        False Negative (TN): 426
In [41]: # Score the predictions
         accuracy = accuracy score(y test, y pred)
         precision = TP / (TP + FP)
         recall = TP / (TP + FN)
         f1 = (2*precision*recall) / (precision + recall)
         print(f"Model Accuracy: {np.round(accuracy, 4)}")
         print(F"Precision: {np.round(precision, 4)}")
        print(f"Recall: {np.round(recall, 4)}")
        print(f"F1 Score: {np.round(f1, 4)}")
        Model Accuracy: 0.7798
        Precision: 0.2035
        Recall: 0.7875
        F1 Score: 0.3234
In [42]: from sklearn.metrics import roc curve, roc auc score, auc
         # Calculate the Area Under Curve (AUC)
         fpr, tpr, thresholds = roc curve(y test, y probs)
         roc auc = auc(fpr, tpr)
         print("AUC:", roc auc)
         # Plot the ROC curve
         fig, axs = plt.subplots()
         axs.plot(fpr, tpr,
                  color="orange",
                  lw=2,
                  label=f"Final Model (Area Under Curve = {np.round(roc auc, 4)})")
         axs.plot([0, 1], [0, 1],
                  color="darkblue",
```

```
lw=2,
         linestyle='--',
         label="Random Classifier")
axs.plot([0, 1], [1, 1],
         color="green",
         lw=2,
         linestyle='--',
         label="Perfect Classifier")
axs.plot([0, 0], [0, 1],
         color="green",
         lw=2,
         linestyle='--')
axs.set xlim([-0.01, 1.0])
axs.set ylim([0.0, 1.05])
axs.set xlabel('False Positive Rate')
axs.set ylabel('True Positive Rate')
axs.set title('Receiver Operating Characteristic (ROC) Curve')
plt.legend(loc="lower right")
plt.show()
```

AUC: 0.8607407004902461

