CS188 Project 2 - Binary Classification Comparative Methods

For this project we're going to attempt a binary classification of a dataset using multiple methods and compare results.

Our goals for this project will be to introduce you to several of the most common classification techniques, how to perform them and tweek parameters to optimize outcomes, how to produce and interpret results, and compare performance. You will be asked to analyze your findings and provide explanations for observed performance.

Specifically you will be asked to classify whether a **patient is suffering from heart disease** based on a host of potential medical factors.

DEFINITIONS

Binary Classification: In this case a complex dataset has an added 'target' label with one of two options. Your learning algorithm will try to assign one of these labels to the data.

Supervised Learning: This data is fully supervised, which means it's been fully labeled and we can trust the veracity of the labeling.

Background: The Dataset

For this exercise we will be using a subset of the UCI Heart Disease dataset, leveraging the fourteen most commonly used attributes. All identifying information about the patient has been scrubbed.

The dataset includes 14 columns. The information provided by each column is as follows:

- age: Age in years
- sex: (1 = male; 0 = female)
- **cp:** Chest pain type (0 = asymptomatic; 1 = atypical angina; 2 = non-anginal pain; 3 = typical angina)
- trestbps: Resting blood pressure (in mm Hg on admission to the hospital)
- cholserum: Cholestoral in mg/dl
- fbs Fasting blood sugar > 120 mg/dl (1 = true; 0 = false)
- restecg: Resting electrocardiographic results (0= showing probable or definite left ventricular hypertrophy by Estes' criteria; 1 = normal; 2 = having ST-T wave abnormality (T wave inversions and/or ST elevation or depression of > 0.05 mV))
- · thalach: Maximum heart rate achieved
- exang: Exercise induced angina (1 = yes; 0 = no)
- oldpeakST: Depression induced by exercise relative to rest
- slope: The slope of the peak exercise ST segment (0 = downsloping; 1 = flat; 2 = upsloping)
- ca: Number of major vessels (0-3) colored by flourosopy
- thal: 1 = normal; 2 = fixed defect; 7 = reversable defect

• Sick: Indicates the presence of Heart disease (True = Disease; False = No disease)

Loading Essentials and Helper Functions

```
#Here are a set of libraries we imported to complete this assignment.
In [1]:
            #Feel free to use these or equivalent libraries for your implementation
            import numpy as np # linear algebra
            import pandas as pd # data processing, CSV file I/O (e.g. pd.read_csv)
            import matplotlib.pyplot as plt # this is used for the plot the graph
            import os
            import seaborn as sns # used for plot interactive graph.
            from sklearn.model selection import train_test_split, cross_val_score, Grid
            from sklearn import metrics
            from sklearn.svm import SVC
            from sklearn.linear model import LogisticRegression
            from sklearn.neighbors import KNeighborsClassifier
            from sklearn.tree import DecisionTreeClassifier
            from sklearn.cluster import KMeans
            from sklearn.metrics import confusion matrix
            import sklearn.metrics.cluster as smc
            from sklearn.model selection import KFold
            from matplotlib import pyplot
            import itertools
            %matplotlib inline
            import random
            random.seed(42)
```

```
In [2]: # Helper function allowing you to export a graph
def save_fig(fig_id, tight_layout=True, fig_extension="png", resolution=300
    path = os.path.join(fig_id + "." + fig_extension)
    print("Saving figure", fig_id)
    if tight_layout:
        plt.tight_layout()
    plt.savefig(path, format=fig_extension, dpi=resolution)
```

```
In [3]:
            # Helper function that allows you to draw nicely formatted confusion matric
            def draw confusion matrix(y, yhat, classes):
                    Draws a confusion matrix for the given target and predictions
                    Adapted from scikit-learn and discussion example.
                plt.cla()
                plt.clf()
                matrix = confusion matrix(y, yhat)
                plt.imshow(matrix, interpolation='nearest', cmap=plt.cm.Blues)
                plt.title("Confusion Matrix")
                plt.colorbar()
                num_classes = len(classes)
                plt.xticks(np.arange(num classes), classes, rotation=90)
                plt.yticks(np.arange(num classes), classes)
                fmt = 'd'
                thresh = matrix.max() / 2.
                for i, j in itertools.product(range(matrix.shape[0]), range(matrix.shap
                    plt.text(j, i, format(matrix[i, j], fmt),
                             horizontalalignment="center",
                             color="white" if matrix[i, j] > thresh else "black")
                plt.ylabel('True label')
                plt.xlabel('Predicted label')
                plt.tight layout()
                plt.show()
```

[20 Points] Part 1. Load the Data and Analyze

Let's first load our dataset so we'll be able to work with it. (correct the relative path if your notebook is in a different directory than the csv file.)

Out[4]:

	age	sex	ср	trestbps	chol	fbs	restecg	thalach	exang	oldpeak	slope	са	thal	sic
0	63	1	3	145	233	1	0	150	0	2.3	0	0	1	False
1	37	1	2	130	250	0	1	187	0	3.5	0	0	2	Fals€
2	41	0	1	130	204	0	0	172	0	1.4	2	0	2	False
3	56	1	1	120	236	0	1	178	0	8.0	2	0	2	False
4	57	0	0	120	354	0	1	163	1	0.6	2	0	2	False

Question 1.1 Now that our data is loaded, let's take a closer look at the dataset we're working with. Use the head method to display some of the rows so we can visualize the types of data fields we'll be working with, then use the describe method, along with any additional methods you'd like to call to better help you understand what you're working with and what issues you might face.

In [5]: ▶ data.describe()

Out[5]:

	age	sex	ср	trestbps	chol	fbs	restecg	
count	303.000000	303.000000	303.000000	303.000000	303.000000	303.000000	303.000000	3(
mean	54.366337	0.683168	0.966997	131.623762	246.264026	0.148515	0.528053	14
std	9.082101	0.466011	1.032052	17.538143	51.830751	0.356198	0.525860	1
min	29.000000	0.000000	0.000000	94.000000	126.000000	0.000000	0.000000	-
25%	47.500000	0.000000	0.000000	120.000000	211.000000	0.000000	0.000000	10
50%	55.000000	1.000000	1.000000	130.000000	240.000000	0.000000	1.000000	1
75%	61.000000	1.000000	2.000000	140.000000	274.500000	0.000000	1.000000	16
max	77.000000	1.000000	3.000000	200.000000	564.000000	1.000000	2.000000	2(

```
In [6]: ► data.info()
```

```
<class 'pandas.core.frame.DataFrame'>
RangeIndex: 303 entries, 0 to 302
Data columns (total 14 columns):
           303 non-null int64
age
           303 non-null int64
sex
ср
           303 non-null int64
trestbps
           303 non-null int64
           303 non-null int64
chol
fbs
           303 non-null int64
           303 non-null int64
restecg
thalach
           303 non-null int64
            303 non-null int64
exang
oldpeak
            303 non-null float64
            303 non-null int64
slope
            303 non-null int64
ca
thal
            303 non-null int64
sick
            303 non-null bool
dtypes: bool(1), float64(1), int64(12)
memory usage: 31.2 KB
```

Question 1.2 Discuss your data preprocessing strategy. Are their any datafield types that are problemmatic and why? Will there be any null values you will have to impute and how do you intend to do so? Finally,

for your numeric and categorical features, what if any, additional preprocessing steps will you take on those data elements?

[Use this area to discuss your data processing strategy]

```
In [7]: # The bool type, sick, will need to be changed to an int. The categorical v # will be scaled in the pipeline. The encoding will allow the categorical v # will normalize the numerical categories, allowing for more accurate and f # that have different ranges and average magnitudes will be judged similarl
```

Question 1.3 Before we begin our analysis we need to fix the field(s) that will be problematic. Specifically convert our boolean sick variable into a binary numeric target variable (values of either '0' or '1'), and then drop the original sick datafield from the dataframe.

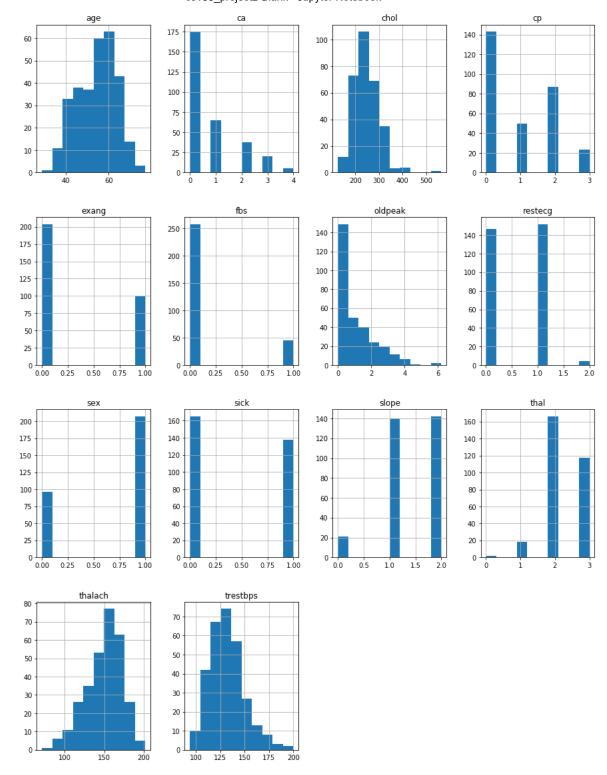
Out[8]:

	age	sex	ср	trestbps	chol	fbs	restecg	thalach	exang	oldpeak	slope	са	thal	sick
0	63	1	3	145	233	1	0	150	0	2.3	0	0	1	0
1	37	1	2	130	250	0	1	187	0	3.5	0	0	2	0
2	41	0	1	130	204	0	0	172	0	1.4	2	0	2	0
3	56	1	1	120	236	0	1	178	0	0.8	2	0	2	0
4	57	0	0	120	354	0	1	163	1	0.6	2	0	2	0

Question 1.4 Now that we have a feel for the data-types for each of the variables, plot histograms of each field and attempt to ascertain how each variable performs (is it a binary, or limited selection, or does it follow a gradient? (Note: No need to describe each variable, but pick out a few you wish to highlight)

```
In [9]: ▶ data.hist(figsize=(15,20))
```

```
Out[9]:
         array([[<matplotlib.axes. subplots.AxesSubplot object at 0x000001C61775A308
         >,
                  <matplotlib.axes. subplots.AxesSubplot object at 0x000001C617A360C8</pre>
         >,
                  <matplotlib.axes. subplots.AxesSubplot object at 0x000001C617A69748</pre>
         >,
                  <matplotlib.axes. subplots.AxesSubplot object at 0x000001C617A9E4C8</pre>
         >],
                 [<matplotlib.axes. subplots.AxesSubplot object at 0x000001C617AD4E08</pre>
         >,
                  <matplotlib.axes. subplots.AxesSubplot object at 0x000001C617B0EF48</pre>
         >,
                  <matplotlib.axes. subplots.AxesSubplot object at 0x000001C617B4A0C8</pre>
         >,
                  <matplotlib.axes. subplots.AxesSubplot object at 0x000001C617B831C8</pre>
         >],
                 (<matplotlib.axes. subplots.AxesSubplot object at 0x000001C617B88D88</pre>
         >,
                  <matplotlib.axes. subplots.AxesSubplot object at 0x000001C617BC1F88</pre>
         >,
                  <matplotlib.axes._subplots.AxesSubplot object at 0x000001C617C28448</pre>
         >,
                  <matplotlib.axes. subplots.AxesSubplot object at 0x000001C617C61548</pre>
         >],
                 (<matplotlib.axes. subplots.AxesSubplot object at 0x000001C617CA0E88</pre>
         >,
                  <matplotlib.axes._subplots.AxesSubplot object at 0x000001C617CD2788</pre>
         >,
                  <matplotlib.axes. subplots.AxesSubplot object at 0x000001C617D0C848</pre>
         >,
                  <matplotlib.axes. subplots.AxesSubplot object at 0x000001C617D44A48</pre>
         >]],
                dtype=object)
```



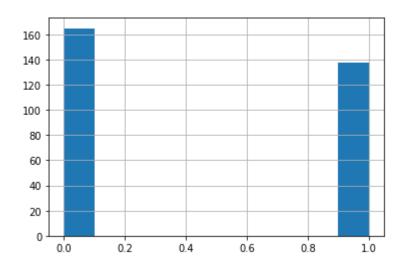
dataset. In this case, we want to confirm whether or not we have an equitable number of sick and healthy individuals to ensure that our classifier will have a sufficiently balanced dataset to adequately classify the two. Plot a histogram specifically of the sick target, and conduct a count of the number of sick and healthy individuals and report on the results:

```
In [10]: M data["sick"].hist()
data.groupby("sick").describe()
```

Out[10]:

	age								sex			са	
	count	mean	std	min	25%	50%	75%	max	count	mean		75%	max
sick													
0	165.0	52.496970	9.550651	29.0	44.0	52.0	59.0	76.0	165.0	0.563636		0.0	4.0
1	138.0	56.601449	7.962082	35.0	52.0	58.0	62.0	77.0	138.0	0.826087		2.0	4.0

2 rows × 104 columns



[Include description of findings here]

```
In [11]: ▶ # There are 165 "healthy" individuals and 138 "sick" individuals within the
```

Question 1.6 Balanced datasets are important to ensure that classifiers train adequately and don't overfit, however arbitrary balancing of a dataset might introduce its own issues. Discuss some of the problems that might arise by artificially balancing a dataset.

[Discuss problem here]

```
In [12]: # It could potentially introduce some additional biases into the model's be # not require any additional augmentations.
```

Question 1.9 Now that we have our dataframe prepared let's start analyzing our data. For this next question let's look at the correlations of our variables to our target value. First, map out the correlations between the values, and then discuss the relationships you observe. Do some research on the variables to understand why they may relate to the observed corellations. Intuitively, why do you think some variables correlate more highly than others (hint: one possible approach you can use the sns heatmap function to map the corr() method)?

```
In [13]:
             matrix = data.corr()
             matrix["sick"].sort values(ascending=False)
    Out[13]: sick
                          1.000000
             exang
                          0.436757
             oldpeak
                          0.430696
                          0.391724
             ca
                          0.344029
             thal
                          0.280937
             sex
             age
                          0.225439
             trestbps
                          0.144931
             chol
                          0.085239
             fbs
                          0.028046
                         -0.137230
             restecg
             slope
                         -0.345877
             thalach
                         -0.421741
                         -0.433798
             ср
             Name: sick, dtype: float64
```

[Discuss correlations here]

```
In [14]: # The factors that are the most correlated with sickness are exercise induc # and number of arteries colored by fluoroscopy. This makes sense, as angin # Depression induced by exercise, meaning the patients do not like to exerc # if they are blocked. Some negative correlations are chest pain, thalach, # possibly affecting its correlation with heart disease in the negative dir
```

[30 Points] Part 2. Prepare the Data

Before running our various learning methods, we need to do some additional prep to finalize our data. Specifically you'll have to cut the classification target from the data that will be used to classify, and then you'll have to divide the dataset into training and testing cohorts.

Specifically, we're going to ask you to prepare 2 batches of data: 1. Will simply be the raw numeric data that hasn't gone through any additional pre-processing. The other, will be data that you pipeline using your own selected methods. We will then feed both of these datasets into a classifier to showcase just how important this step can be!

Question 2.1 Save the target column as a separate array and then drop it from the dataframe.

```
In [15]: N labels = data["sick"]
data = data.drop("sick", axis=1)
```

Question 2.2 First Create your 'Raw' unprocessed training data by dividing your dataframe into training and testing cohorts, with your training cohort consisting of 70% of your total dataframe (hint: use the train_test_split method) Output the resulting shapes of your training and testing samples to confirm that your split was successful.

Question 2.3 Now create a pipeline to conduct any additional preparation of the data you would like. Output the resulting array to ensure it was processed correctly.

```
In [17]:
             from sklearn.impute import SimpleImputer
             from sklearn.compose import ColumnTransformer
             from sklearn.pipeline import Pipeline
             from sklearn.preprocessing import StandardScaler
             from sklearn.preprocessing import OneHotEncoder
             from sklearn.base import BaseEstimator, TransformerMixin
             n pipeline = Pipeline([
                 ('std_scaler', StandardScaler())
                  1)
             numerical_features = ["age", "trestbps", "chol", "thalach", "oldpeak"]
             categorical_features = ["sex", "cp", "fbs", "restecg", "exang", "slope", "c
             f pipeline = ColumnTransformer([
                 ("num", n_pipeline, numerical_features),
                 ("cat", OneHotEncoder(handle unknown = 'ignore'), categorical features)
             1)
             data prepared = f pipeline.fit transform(data)
             print(data prepared)
             [[ 0.9521966
                            0.76395577 -0.25633371 ... 1.
                                                                     0.
                0.
              [-1.91531289 -0.09273778 0.07219949 ... 0.
                                                                     1.
                          1
              [-1.47415758 -0.09273778 -0.81677269 ... 0.
                                                                     1.
                0.
              [ 1.50364073  0.70684287 -1.029353  ...  0.
                                                                     0.
              [ 0.29046364 -0.09273778 -2.2275329 ... 0.
                                                                     0.
              [ 0.29046364 -0.09273778 -0.19835726 ... 0.
                                                                     1.
```

Question 2.4 Now create a separate, processed training data set by dividing your processed dataframe into training and testing cohorts, using the same settings as Q2.2 (REMEMBER TO USE DIFFERENT TRAINING AND TESTING VARIABLES SO AS NOT TO OVERWRITE YOUR PREVIOUS DATA). Output the resulting shapes of your training and testing samples to confirm that your split was successful, and describe what differences there are between your two training datasets.

[What differences are there between these two datasets?]

```
In []: # Once the data is prepared through the pipeline, there is a column created # which is why there are 30 entries rather than 13 for the x tuples. The pr # encoded.
```

[50 Points] Part 3. Learning Methods

We're finally ready to actually begin classifying our data. To do so we'll employ multiple learning methods and compare result.

Linear Decision Boundary Methods

SVM (Support Vector Machine)

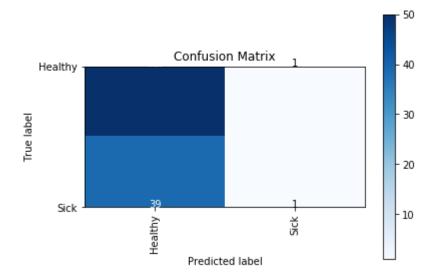
A Support Vector Machine (SVM) is a discriminative classifier formally defined by a separating hyperplane. In other words, given labeled training data (supervised learning), the algorithm outputs an optimal hyperplane which categorizes new examples. In two dimentional space this hyperplane is a line dividing a plane in two parts where in each class lay in either side.

Question 3.1.1 Implement a Support Vector Machine classifier on your RAW dataset. Review the <u>SVM Documentation (https://scikit-learn.org/stable/modules/generated/sklearn.svm.SVC.html)</u> for how to implement a model. For this implementation you can simply use the default settings, but set probability = True.

```
In [19]:  # SVM
svm_raw = SVC(probability=True, gamma='auto')
svm_raw.fit(x_train, y_train)
svm_raw_predict = svm_raw.predict(x_test)
```

Question 3.1.2 Report the accuracy, precision, recall, F1 Score, and confusion matrix of the resulting model.

Accuracy: 0.560440 Precision: 0.500000 Recall: 0.025000 F1 Score: 0.047619



Question 3.1.3 Discuss what each measure is reporting, why they are different, and why are each of these measures is significant. Explore why we might choose to evaluate the performance of differing models differently based on these factors. Try to give some specific examples of scenarios in which you might value one of these measures over the others.

[Provide explanation for each measure here]

```
In []:  # Accuracy is the number of successfully predicted instances over the total

# Precision measures the number of successfully predicted instances divided

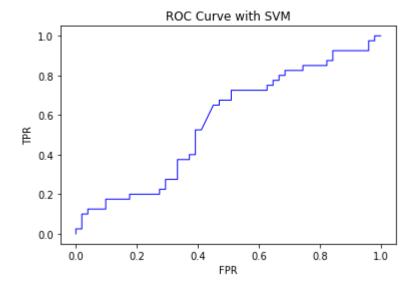
# predicted instances plus the false positives. The higher precision is, th

# Recall is the same as precision, but looks at the number of false negativ

# F1 Score is a weighted average between precision and recall, a wholistic

# false positives and false negatives.
```

Question 3.1.4 Plot a Receiver Operating Characteristic curve, or ROC curve, and describe what it is and what the results indicate



[Describe what an ROC Curve is and what the results mean here]

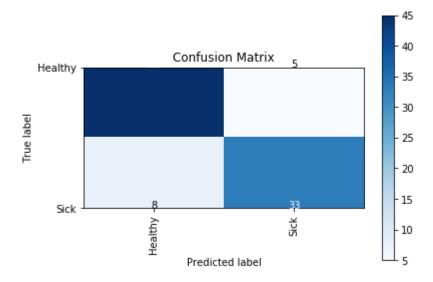
```
In []: # A ROC Curve plots the True Positive Rate (Precision) against the False Po # It can be used to measure the effectiveness of a particular model on a da # The more separability the model provides, the less the curves will inters # The closer the AUC is to 1, the more effective the model is.
```

Question 3.1.5 Rerun, using the exact same settings, only this time use your processed data as inputs.

Question 3.1.6 Report the accuracy, precision, recall, F1 Score, confusion matrix, and plot the ROC Curve of the resulting model.

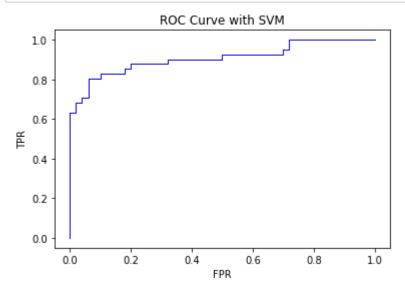
```
In [23]: # Print scores for processed SVM
print("%-12s %f" % ('Accuracy:', metrics.accuracy_score(y_test2, svm_proces
print("%-12s %f" % ('Precision:', metrics.precision_score(y_test2, svm_procested)
print("%-12s %f" % ('Recall:', metrics.recall_score(y_test2, svm_processed_print("%-12s %f" % ('F1 Score:', metrics.f1_score(y_test2, svm_processed_print("%-12s %f" % ('y_test2, svm_processed_print(", 'Sick'))
```

Accuracy: 0.857143 Precision: 0.868421 Recall: 0.804878 F1 Score: 0.835443



```
In [24]: # ROC Curve for processed data
svm_processed_score = svm_processed.decision_function(x_test2)
fpr_svm_processed, tpr_svm_processed, thresholds_processed = metrics.roc_cu

pyplot.figure()
pyplot.plot(fpr_svm_processed, tpr_svm_processed, color='blue', lw=1)
pyplot.title("ROC Curve with SVM")
pyplot.xlabel('FPR')
pyplot.ylabel('TPR')
pyplot.show()
auc_score2 = metrics.roc_auc_score(y_test2, svm_processed_predict)
print("The AUC Score for the SVM is: ", auc_score2)
```



The AUC Score for the SVM is: 0.852439024390244

Question 3.1.7 Hopefully you've noticed a dramatic change in performance. Discuss why you think your new data has had such a dramatic impact.

[Provide explanation here]

```
In []: # Because the numerical features in this data set have different ranges and # properly separate the data if it is not scaled. Additionally, the model d # the categorical features without onehotencoding them. Once the data is pr # much more effectively.
```

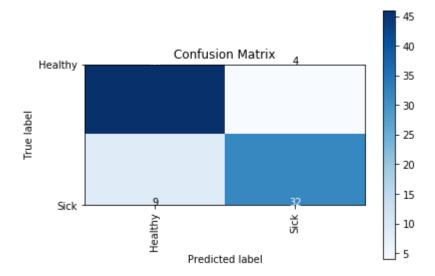
Question 3.1.8 Rerun your SVM, but now modify your model parameter kernel to equal 'linear'. Again report your Accuracy, Precision, Recall,

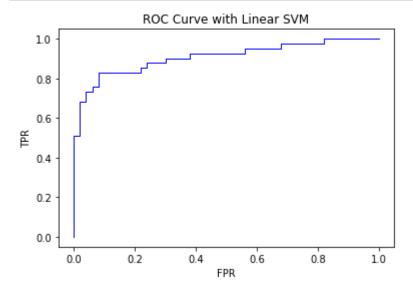
F1 scores, and Confusion matrix and plot the new ROC curve.

```
In [25]: # SVM
svm_processed2 = SVC(kernel='linear', probability=True, gamma='auto')
svm_processed2.fit(x_train2, y_train2)
svm_processed_predict2 = svm_processed2.predict(x_test2)
```

In [26]: # Print scores for Linear SVM print("%-12s %f" % ('Accuracy:', metrics.accuracy_score(y_test2, svm_proces print("%-12s %f" % ('Precision:', metrics.precision_score(y_test2, svm_proce) print("%-12s %f" % ('Recall:', metrics.recall_score(y_test2, svm_processed_print("%-12s %f" % ('F1 Score:', metrics.f1_score(y_test2, svm_processed_prediate) print("%-12s %f" % ('F1 Score:', metrics.f1_score(y_test2, svm_processed_prediate))

Accuracy: 0.857143 Precision: 0.888889 Recall: 0.780488 F1 Score: 0.831169





The AUC Score for the Linear SVM is: 0.8502439024390245

Question 3.1.9 Explain the what the new results you've achieved mean. Read the documentation to understand what you've changed about your model and explain why changing that input parameter might impact the results in the manner you've observed.

[Provide explanation here]

```
In [28]: | # Linear is faster but less accurate than other mappings like rbf. As most # it is difficult to create the most accurate division using a linear kerne # However, after processing this data, it seems to be fairly separable, whi # boost performance. An rbf re-maps the data set to a space that has a more #Thus, the 'rbf' kernel has a slightly higher AUC score compared to the lin
```

Logistic Regression

Knowing that we're dealing with a linearly configured dataset, let's now try another classifier that's

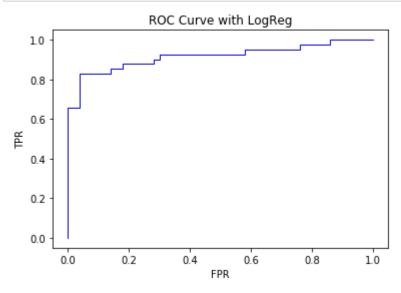
well known for handling linear models: Logistic Regression. Logistic regression is a statistical model that in its basic form uses a logistic function to model a binary dependent variable.

Question 3.2.1 Implement a Logistical Regression Classifier. Review the Logistical Regression Documentation (https://scikit-learn.org/stable/modules/generated/sklearn.linear_model.LogisticRegres for how to implement the model. For this initial model set the solver = 'sag' and max_iter= 10). Report on the same four metrics as the SVM and graph the resulting ROC curve.

```
# Logistic Regression
In [29]:
             log reg = LogisticRegression(solver='sag', max iter=10)
             log_reg.fit(x_train2, y_train2)
             log reg predict = log reg.predict(x test2)
             C:\Users\ethan\Anaconda3\lib\site-packages\sklearn\linear model\sag.py:337:
             ConvergenceWarning: The max iter was reached which means the coef did not
             converge
               "the coef did not converge", ConvergenceWarning)
             # Print scores for Logisitic Regression, max iter=10
In [30]:
             print("%-12s %f" % ('Accuracy:', metrics.accuracy_score(y_test2, log_reg_pr
             print("%-12s %f" % ('Precision:', metrics.precision_score(y_test2, log_reg_
             print("%-12s %f" % ('Recall:', metrics.recall score(y test2, log reg predic
             print("%-12s %f" % ('F1 Score:', metrics.f1 score(y test2, log reg predict)
                          0.890110
             Accuracy:
             Precision:
                          0.918919
             Recall:
                          0.829268
             F1 Score:
                          0.871795
```

```
In [31]: No log_reg_score = log_reg.decision_function(x_test2)
fpr_log, tpr_log, thresholds_log = metrics.roc_curve(y_test2, log_reg_score)

pyplot.figure()
pyplot.plot(fpr_log, tpr_log, color='blue', lw=1)
pyplot.title("ROC Curve with LogReg")
pyplot.xlabel('FPR')
pyplot.ylabel('TPR')
pyplot.show()
auc_score = metrics.roc_auc_score(y_test2, log_reg_predict)
print("The AUC Score for the Logistic Regression Model is: ", auc_score)
```



The AUC Score for the Logistic Regression Model is: 0.8846341463414633

Question 3.2.2 Did you notice that when you ran the previous model you got the following warning: "ConvergenceWarning: The max_iter was reached which means the coef_ did not converge". Check the documentation and see if you can implement a fix for this problem, and again report your results.

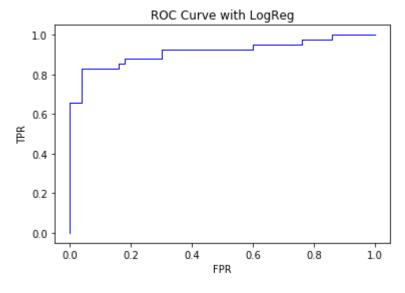
```
In [32]: # Logistic Regression
log_reg2 = LogisticRegression(solver='sag', max_iter=60)
log_reg2.fit(x_train2, y_train2)
log_reg_predict2 = log_reg2.predict(x_test2)
```

```
In [33]: # Print scores for Logisitic Regression, max_iter=100
print("%-12s %f" % ('Accuracy:', metrics.accuracy_score(y_test2, log_reg_pr
print("%-12s %f" % ('Precision:', metrics.precision_score(y_test2, log_reg_
print("%-12s %f" % ('Recall:', metrics.recall_score(y_test2, log_reg_predict)
print("%-12s %f" % ('F1 Score:', metrics.f1_score(y_test2, log_reg_predict)
Accuracy: 0.901099
```

Accuracy: 0.901099 Precision: 0.944444 Recall: 0.829268 F1 Score: 0.883117

```
In [34]: N log_reg_score2 = log_reg2.decision_function(x_test2)
fpr_log2, tpr_log2, thresholds_log2 = metrics.roc_curve(y_test2, log_reg_sc

pyplot.figure()
pyplot.plot(fpr_log2, tpr_log2, color='blue', lw=1)
pyplot.title("ROC Curve with LogReg")
pyplot.xlabel('FPR')
pyplot.ylabel('TPR')
pyplot.show()
auc_score = metrics.roc_auc_score(y_test2, log_reg_predict2)
print("The AUC Score for the Logistic Regression Model is: ", auc_score)
```



The AUC Score for the Logistic Regression Model is: 0.8946341463414633

Question 3.2.3 Explain what you changed, and why that produced an improved outcome.

[Provide explanation here]

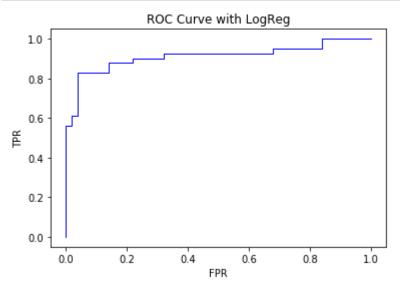
```
In []: # The max iteration for the Logistic Regression model was being reached bef # The solution to this is increase the maximum number of iterations, allowi # However, the performance was only improved marginally (0.01) according to # Logistic Regression can achieve a high level of performance fairly quickl
```

Question 3.2.4 Rerun your logistic classifier, but modify the penalty = 'none', solver='sag' and again report the results.

```
In [35]:
          # Logistic Regression
             log_reg3 = LogisticRegression(penalty='none', solver='sag', max_iter=500)
             log reg3.fit(x train2, y train2)
             log reg predict3 = log reg3.predict(x test2)
             C:\Users\ethan\Anaconda3\lib\site-packages\sklearn\linear model\sag.py:337:
             ConvergenceWarning: The max iter was reached which means the coef did not
             converge
               "the coef_ did not converge", ConvergenceWarning)
In [36]:
          # Print scores for Logisitic Regression, penalty=none
             print("%-12s %f" % ('Accuracy:', metrics.accuracy_score(y_test2, log_reg_pr
             print("%-12s %f" % ('Precision:', metrics.precision_score(y_test2, log_reg_
             print("%-12s %f" % ('Recall:', metrics.recall_score(y_test2, log_reg_predic
             print("%-12s %f" % ('F1 Score:', metrics.f1_score(y_test2, log_reg_predict3
             Accuracy:
                          0.901099
             Precision:
                          0.944444
             Recall:
                          0.829268
```

F1 Score:

0.883117



The AUC Score for the Logistic Regression Model is: 0.8946341463414633

Question 3.2.5 Explain what what the penalty parameter is doing in this function, what the solver method is, and why this combination likely produced a more optimal outcome.

[Provide explanation here]

```
In [38]: # It has no change in performance.

# Like stochastic gradient (SG) methods, the SAG method's iteration cost is

#However, by incorporating a memory of previous gradient values the SAG met

# black-box SG methods. No regularization is applied to the data set, beca

# This is possibly because the Logistic Regression model can perform well e
```

Question 3.2.6 Both logistic regression and linear SVM are trying to classify data points using a linear decision boundary, then what's the difference between their ways to find this boundary?

[Provide Answer here:]

```
In [39]: | # SVM tries to find the widest possible margin, while LogReg tries to maxim # the correct classification. SVM minizes hinge loss, while LogReg minimiz # to outliers.
```

Clustering Approaches

Let us now try a different approach to classification using a clustering algorithm. Specifically, we're going to be using K-Nearest Neighbor, one of the most popular clustering approaches.

K-Nearest Neighbor

Question 3.3.1 Implement a K-Nearest Neighbor algorithm on our data and report the results. For this initial implementation simply use the default settings. Refer to the KNN Documentation (https://scikit-learn.org/stable/modules/generated/sklearn.neighbors.KNeighborsClass for details on implementation. Report on the accuracy of the resulting model.

```
In [45]: # k-Nearest Neighbors algorithm
    neighbors = KNeighborsClassifier()
    neighbors.fit(x_train2, y_train2)
    neighbors_predict = neighbors.predict(x_test2)

ascore = metrics.accuracy_score(y_test2, neighbors_predict)
    print("Accuracy Score: ", ascore)
```

Accuracy Score: 0.8461538461538461

Question 3.3.2 For clustering algorithms, we use different measures to determine the effectiveness of the model. Specifically here, we're interested in the Homogeneity Score, Completeness Score, V-Measure, Adjusted Rand Score, and Adjusted Mutual Information. Calculate each score (hint review the SKlearn Metrics Clustering documentation for how to implement).

Homogeneity Score: 0.37620925388444065 Completeness Score: 0.3791512192913599

V-Measure: 0.3776745074289339

Adjusted Rand Score: 0.4735103958323643 Adjusted Mutual Info: 0.37110042222474654

C:\Users\ethan\Anaconda3\lib\site-packages\sklearn\metrics\cluster\supervis
ed.py:746: FutureWarning: The behavior of AMI will change in version 0.22.
To match the behavior of 'v_measure_score', AMI will use average_method='ar
ithmetic' by default.
FutureWarning)

Question 3.3.3 Explain what each score means and interpret the results for this particular model.

[Input answer here]

```
In []: # Homogeneity checks to see if each cluster only has entries of a single cl # perfectly homogenous. This 5-neighbor model is not very homogenous.

# Completeness is the inverse of homogeneity. It checks to see if each entr # cluster. As completeness is related to homogeneity, this 5-neighbor model # V-Measure is the harmonic mean between homogeneity and completeness. As b # Adjusted Rand Score checks the similarity between clusters.

# Adjusted Mutual Info is the measure of mutual dependence between x and y. # that larger clusters could have a higher degree of mutual information.
```

As we're beginning to see, the input parameters for your model can dramatically impact the performance of the model. How do you know which settings to choose? Studying the models and studying your datasets are critical as they can help you anticipate which models and settings are likely to produce optimal results. However sometimes that isn't enough, and a brute force method is necessary to determine which parameters to use. For this next question we'll attempt to optimize a parameter using a brute force approach.

Question 3.3.4 Parameter Optimization. The KNN Algorithm includes an n_neighbors attribute that specifies how many neighbors to use when developing the cluster (The default value is 5, which is what your

previous model used.) Lets now try n values of: 1, 2, 3, 5, 10, 20, 50, and 100. Run your model for each value and report the 6 measures (5 clustering specific plus accuracy) for each. Report on which n value produces the best accuracy and V-Measure. (HINT leverage python's ability to loop to run through the array and generate results without needing to manually code each iteration).

Number of Neighbors: 1

Accuracy Score: 0.7142857142857143 Homogeneity Score: 0.13106249480889112 Completeness Score: 0.1335172792933158

V-Measure: 0.1322784992439942

Adjusted Rand Score: 0.17462753853419777 Adjusted Mutual Info: 0.12394330029337142

Number of Neighbors: 2

Accuracy Score: 0.7582417582417582 Homogeneity Score: 0.22100295999912523 Completeness Score: 0.2587248857389366

V-Measure: 0.23838084898228307

Adjusted Rand Score: 0.2587837596136649 Adjusted Mutual Info: 0.21458661619903746

Number of Neighbors: 3

Accuracy Score: 0.8351648351648352 Homogeneity Score: 0.3500777415855479 Completeness Score: 0.35131561039893566

V-Measure: 0.3506955836528328

Adjusted Rand Score: 0.4432259334132422 Adjusted Mutual Info: 0.3447556589975192

Number of Neighbors: 5

Accuracy Score: 0.8461538461538461 Homogeneity Score: 0.37620925388444065 Completeness Score: 0.3791512192913599

V-Measure: 0.3776745074289339

Adjusted Rand Score: 0.4735103958323643 Adjusted Mutual Info: 0.37110042222474654

Number of Neighbors: 10

Accuracy Score: 0.8681318681318682 Homogeneity Score: 0.44235681956589074 Completeness Score: 0.4569431960541842

V-Measure: 0.4495317144844716

Adjusted Rand Score: 0.5370204950536377 Adjusted Mutual Info: 0.43778584925774844 Number of Neighbors: 20

Accuracy Score: 0.8901098901098901 Homogeneity Score: 0.5108996695970632 Completeness Score: 0.5277461939386586

V-Measure: 0.5191863089821956

Adjusted Rand Score: 0.6044162457736778 Adjusted Mutual Info: 0.5068905412396896

Number of Neighbors: 50

Accuracy Score: 0.8791208791208791 Homogeneity Score: 0.4709144471353802 Completeness Score: 0.4828835034650588

V-Measure: 0.4768238764234958

Adjusted Rand Score: 0.5702248698545591 Adjusted Mutual Info: 0.4665786747127186

Number of Neighbors: 100

Accuracy Score: 0.8571428571428571 Homogeneity Score: 0.427411840711569 Completeness Score: 0.453636397941567

V-Measure: 0.4401338300258592

Adjusted Rand Score: 0.5048090052723155 Adjusted Mutual Info: 0.4227138044218406

C:\Users\ethan\Anaconda3\lib\site-packages\sklearn\metrics\cluster\sup ervised.py:746: FutureWarning: The behavior of AMI will change in vers ion 0.22. To match the behavior of 'v_measure_score', AMI will use ave rage_method='arithmetic' by default.

FutureWarning)

C:\Users\ethan\Anaconda3\lib\site-packages\sklearn\metrics\cluster\sup ervised.py:746: FutureWarning: The behavior of AMI will change in vers ion 0.22. To match the behavior of 'v_measure_score', AMI will use ave rage_method='arithmetic' by default.

FutureWarning)

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FutureWarning)

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FutureWarning)

C:\Users\ethan\Anaconda3\lib\site-packages\sklearn\metrics\cluster\sup
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FutureWarning)

C:\Users\ethan\Anaconda3\lib\site-packages\sklearn\metrics\cluster\sup ervised.py:746: FutureWarning: The behavior of AMI will change in vers ion 0.22. To match the behavior of 'v_measure_score', AMI will use ave rage_method='arithmetic' by default.

FutureWarning)

[Describe optimal results here]

```
In []: 

# 20 Neighbors has the best performance in terms of accuracy and v-measure.

# and completeness, meaning that the partitions for each cluster are the mo
```

Question 3.3.5 When are clustering algorithms most effective, and what do you think explains the comparative results we achieved?

[Input answer here]

```
In []: N # Clustering algorithms are useful when there are distinct classifications # With a high degree of inter-subject similarities, it is beneficial to clu # items that are virtually indistinguishable from each other. As there are # clustering will not produce the most optimal results. There are many nuan # health. As clustering evaluates the data entries based on their similarit # in something as nuanced as heart disease.

# Up to a certain threshold, there is a benefit to increasing the number of # partition, the model is more effectively able to separate the data. Howev # Thus, increasing the value of K to 50 or 100 (respectively around 1/4 and # actually decrease the model's effectiveness.
```