### Homework 5

## **Question 1:**

a)

I used twenty percent of each class (0 and 1) of 'Whether' for testing and the remaining data for training.

### Code:

```
import numpy as np
import pandas as pd
dataframe = pd.read csv("transfusion.csv")
dataset for 0 = (dataframe.loc[dataframe['whether'] == 0])
dataset for 1 = (dataframe.loc[dataframe['whether'] == 1])
predictor count = 4
training set X for 0 = dataset for 0.values[(np.int(len(dataset for 0.values) *
0.20)):, :predictor count]
training set Y for \overline{0} = dataset for 0.values[:(np.int(len(dataset for 0.values)
* 0.20)), predictor count:]
training set X for \overline{1} = dataset for 1.values[(np.int(len(dataset for 1.values) *
0.20)):, :predictor count]
training_set_Y_for_1 = dataset_for_1.values[:(np.int(len(dataset_for_1.values)
* 0.20)), predictor count:]
testing set X for 0 = dataset for 0.values[:(np.int(len(dataset for 0.values) *
0.20)), :predictor count]
testing set Y for \overline{0} = dataset for 0.values[:(np.int(len(dataset for 0.values) *
0.20)), predictor count:]
testing set X for 1 = dataset for 1.values[:(np.int(len(dataset for 1.values) *
0.20)), :predictor count]
testing set Y for \overline{1} = dataset for 1.values[:(np.int(len(dataset for 1.values) *
0.20)), predictor count:]
training_set_X = np.vstack((training_set_X_for_0, training_set_X_for_1))
training_set_Y = np.vstack((training_set_Y_for_0, training_set_Y_for_1))
testing_set_X = np.vstack((testing_set_X_for_0, testing_set_X_for_1))
testing_set_Y = np.vstack((testing_set_Y_for_0, testing_set_Y_for_1))
```

# b) Supervised Learning:

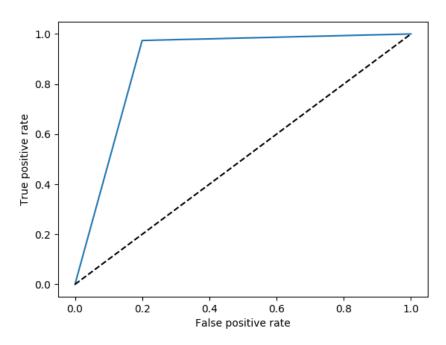
For normalization of data, I have used sklearn's preprocessing package and for imbalance reduction I have used SMOTE so that the minority class can be over-sampled. Then, I have used five fold cross-validation to find the best value for L1 penalty parameter and then, to get the best model for supervised classification, I have used GridSearchCV.

#### **Results:**

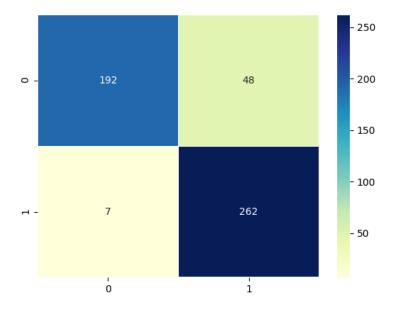
#### **Training Set:**

Best SVM Penalty Param = C = 0.966039211809 Best Score using C = 0.86740471823411

ROC Curve with AUC = 0.896977847584

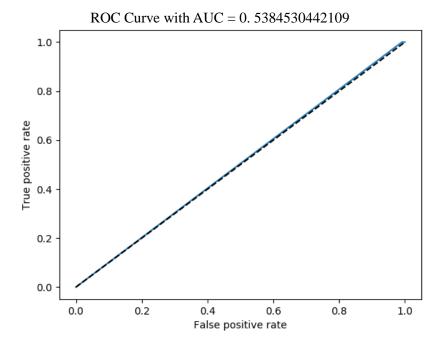


# **Confusion Matrix**

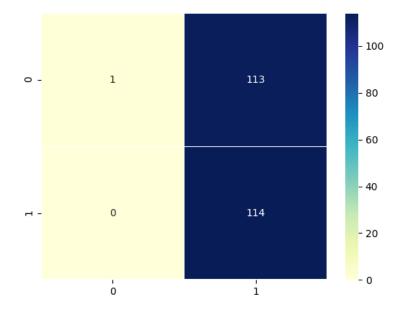


**Testing Set:** 

Prediction Accuracy of Testing Dataset = 0.5384530442109 Area under the Curve of Testing Dataset = 0.5384530442109







As it can be observed in the results, ROC and confusion matric that the classifier could not classify class 0 successfully which resulted in a large number of misclassifications.

```
import numpy as np
import pandas as pd
from sklearn.preprocessing import normalize
from sklearn.svm import LinearSVC
```

```
from sklearn.model selection import KFold
from sklearn.model selection import GridSearchCV
import matplotlib.pyplot as plt
from sklearn.metrics import roc curve
from sklearn.utils import shuffle
from imblearn.combine import SMOTEENN, SMOTETomek
from imblearn.over sampling import SMOTE
from sklearn.metrics import confusion matrix
from sklearn.metrics import auc
from sklearn.metrics import mean squared error
from scipy.sparse import coo matrix, hstack
import seaborn as sns
dataframe = pd.read csv("transfusion.csv")
dataset for 0 = (dataframe.loc[dataframe['whether'] == 0])
dataset for 1 = (dataframe.loc[dataframe['whether'] == 1])
predictor count = 4
training set X for 0 = dataset for 0.values[(np.int(len(dataset for 0.values) * 0.20));,
:predictor count]
training_set_Y_for_0 = dataset_for_0.values[:(np.int(len(dataset_for_0.values) * 0.20)),
predictor count:]
training set X for 1 = dataset for 1.values[(np.int(len(dataset for 1.values) * 0.20));,
:predictor count]
training set Y for 1 = dataset for 1.values[:(np.int(len(dataset for 1.values) * 0.20)),
predictor count:]
testing set X for 0 = dataset for 0.values[:(np.int(len(dataset for 0.values) * 0.20)),
:predictor count]
testing set Y for 0 = dataset for 0.values[:(np.int(len(dataset for 0.values) * 0.20)),
predictor count:]
testing_set_X_for_1 = dataset_for_1.values[:(np.int(len(dataset_for 1.values) * 0.20)),
:predictor count]
testing set Y for 1 = dataset for 1.values[:(np.int(len(dataset for 1.values) * 0.20)),
predictor_count:]
training X = normalize(np.vstack((training set X for 0, training set X for 1)),
norm='12')
training Y = np.vstack((training set Y for 0, training set Y for 1))
testing X = \text{normalize}(\text{np.vstack}((\text{testing set } X \text{ for } 0, \text{ testing set } X \text{ for } 1)), \text{ norm='}12')
testing_Y = np.vstack((testing_set_Y_for_0, testing_set_Y_for_1))
training X new = shuffle(hstack([coo matrix(training X),
coo matrix(training Y)]).toarray(), random state=15)
testing X new = shuffle(hstack([coo matrix(testing X),
coo matrix(testing Y)]).toarray(), random state=36)
training_X = training_X_new[:, :predictor_count]
training_Y = training_X_new[:, predictor_count:]
testing_X = testing_X_new[:, :predictor_count]
testing_Y = testing_X_new[:, predictor_count:]
sampling = SMOTEENN(random state=5, kind smote='svm')
training smote X, training smote Y = sampling.fit sample(training X, training Y.ravel())
testing_smote_X, testing_smote_Y = SMOTE().fit_sample(testing_X, testing_Y.ravel())
testing_smote_Y=testing_smote_Y.reshape(len(testing_smote_Y), 1)
testing X new = shuffle(hstack([coo matrix(testing smote X),
coo matrix(testing smote Y)]).toarray(), random state=65)
testing X = testing X new[:, :predictor_count]
testing Y = testing X new[:, predictor count:]
classifier = GridSearchCV(LinearSVC(penalty='11', dual=False, tol=0.001), [{'C':
np.linspace(0.0001, 1, 250)}], cv=KFold(5), refit=True, n_jobs=4)
```

```
classifier.fit(training_smote_X, training_smote_Y.ravel())
predictions = classifier.predict(training smote X)
accuracy = 1 - mean squared error(training smote Y, predictions)
best SVM param = classifier.best params
best score = classifier.best score
test set predictions = classifier.predict(testing X)
accuracy = 1 - mean squared error(testing Y, test set predictions)
fpr_rf_lm, tpr_rf_lm, _ = roc_curve(training_smote_Y, predictions, pos_label=1)
roc auc = auc(fpr rf lm, tpr rf lm)
plt.plot(fpr_rf_lm, tpr_rf_lm, label=str(roc auc))
plt.xlabel('False positive rate')
plt.ylabel('True positive rate')
plt.plot([0, 1], [0, 1], 'k--')
plt.show()
fpr rf lm, tpr rf lm, = roc curve(testing Y, test set predictions, pos label=1)
roc auc = auc(fpr rf lm, tpr rf lm)
plt.plot(fpr_rf_lm, tpr_rf_lm, label=str(roc_auc))
plt.title('ROC curve- (Test-Set), AUC: '+str(roc auc))
plt.xlabel('False positive rate')
plt.ylabel('True positive rate')
plt.plot([0, 1], [0, 1], 'k--')
plt.show()
confusion matrix=confusion matrix(training_smote_Y, predictions)
sns.heatmap(confusion matrix, cmap="Blues", annot=True, linewidths=0.5, fmt='d')
plt.show()
confusion matrix = confusion matrix(testing Y, test set predictions)
sns.heatmap(confusion matrix, cmap="Blues", annot=True, linewidths=0.5, fmt='d')
plt.show()
```

#### c) Semi-Supervised Learning

- i) For semi-supervised learning, I have used Linear SVM with L1 penalty and the labeled training data was used to fit the model. Then, I used GridSearchCV to perform five fold cross-validation on this model.
- ii) Then, to find the unlabeled data point closet to the decision boundary I used the decision function of LinearSVC.
- iii) Then, I added this data point labeled by the model to the labeled set of data and removed it from the unlabeled set of data.
- iv) Then, I calculated the misclassification rate and accuracy of the predictions made by the model on test dataset.

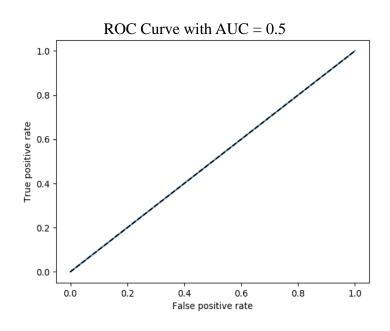
#### **Results:**

### **Training:**

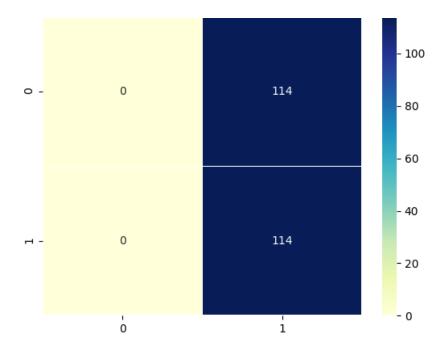
Best SVM Penalty Param for index 0 = C = 8.9891282324Best Score using C for index 0 = 0.932497317Best SVM Penalty Param for index 216 = C = 6.293474182 Best Score using C for index 216 = 0.8413402377

# **Testing:**

Area Under the Curve = 0.5Accuracy = 0.5



### **Confusion Matrix**



Results show that semi-supervised learning fails at predicting the test dataset correctly. The number of misclassifications is huge—an entire class is misclassified.

```
import numpy as np
import pandas as pd
from sklearn.preprocessing import normalize
from sklearn.svm import LinearSVC
from sklearn.model selection import KFold
from sklearn.model selection import GridSearchCV
import matplotlib.pyplot as plt
from sklearn.metrics import roc curve
from sklearn.utils import shuffle
from imblearn.combine import SMOTEENN, SMOTETomek
from imblearn.over sampling import SMOTE
from sklearn.metrics import confusion matrix
from sklearn.metrics import auc
from sklearn.metrics import mean squared error
from scipy.sparse import coo matrix, hstack
import seaborn as sns
dataframe = pd.read csv("transfusion.csv")
dataset for 0 = (dataframe.loc[dataframe['whether'] == 0])
dataset for 1 = (dataframe.loc[dataframe['whether'] == 1])
predictor count = 4
percentage = 0.20
training set X for 0 = dataset for 0.values[(np.int(len(dataset for 0.values) *
percentage)):, :predictor count]
training_set_Y_for_0 = dataset_for_0.values[:(np.int(len(dataset for 0.values) *
percentage)), predictor_count:]
training_set_X_for_1 = dataset_for_1.values[(np.int(len(dataset for 1.values) *
percentage)):, :predictor count]
training set Y for 1 = dataset for 1.values[:(np.int(len(dataset for 1.values) *
percentage)), predictor count:]
testing_set_X_for_0 = dataset_for_0.values[:(np.int(len(dataset for 0.values) *
percentage)), :predictor_count]
testing_set_Y_for_0 = dataset_for_0.values[:(np.int(len(dataset_for_0.values) *
percentage)), predictor_count:]
testing_set_X_for_1 = dataset_for 1.values[:(np.int(len(dataset for 1.values) *
percentage)), :predictor_count]
testing_set_Y_for_1 = dataset_for_1.values[:(np.int(len(dataset for 1.values) *
percentage)), predictor_count:]
testing_set_for_Y = np.vstack((testing_set_Y_for_0, testing_set_Y_for_1))
testing_set_for_X = normalize(np.vstack((testing_set_X_for_0, testing_set_X_for_1)),
norm='12')
training set X for 0 half = int(len(training set X for 0) * 0.50)
training set X for 1 half = int(len(training set X for 1) * 0.50)
labelled_training_set_X_for_0 = training_set_X_for_0[training_set_X_for_0_half:, :]
labelled_training_set_for_Y_for_0 = training_set_Y_for_0[training_set_X_for_0_half:, :]
unlabelled_training_set_X_for_0 = training_set_X_for_0[:training_set_X_for_0 half, :]
unlabelled_training_set_Y_for_0 = training_set_Y_for_0[:training_set_X_for_0_half, :]
labelled training set X for 1 = training set X for 1[training set X for 1 half:, :]
labelled_training_set Y for 1 = training_set Y for 1[training_set X for 1 half:, :]
unlabelled training set X for 1 = training set X for 1[:training set X for 1 half, :]
unlabelled_training_set_Y_for_1 = training_set_Y_for_1[:training_set_X_for_1_half, :]
labelled training set for Y = np.vstack((labelled training set for Y for 0,
labelled training set Y for 1))
unlabelled training set for Y = np.vstack((unlabelled training set Y for 0,
unlabelled_training_set_Y_for_1))
unlabelled training set for X = normalize(np.vstack((unlabelled training set X for 0,
```

```
unlabelled training set X for 1)), norm='12')
X combined train unlabelled = shuffle(hstack([coo matrix(unlabelled training set for X),
coo_matrix(unlabelled_training_set_for_Y)]).toarray(), random_state=10)
unlabelled training set for X = X combined train unlabelled[:, :predictor count]
unlabelled training set for Y = X combined train unlabelled[:, predictor count:]
sampling=SMOTEENN(random state=5, kind smote='svm')
X_train_unlabelled_smote, Y_train_unlabelled_smote =
sampling.fit_sample(unlabelled_training_set_for_X,
unlabelled training set for Y.ravel())
X train labelled=normalize(np.vstack((labelled training set X for 0,
labelled training set X for 1)), norm='12')
index to use = hstack([coo matrix(X train labelled),
coo matrix(labelled training set for Y)]).toarray()
X combined train labelled = shuffle(index to use, random state=10)
X train labelled=X combined train labelled[:, :predictor count]
labelled training set for Y= X combined train labelled[:, predictor count:]
sampling=SMOTEENN(random_state=5, kind_smote='svm')
X_train_labelled_smote, Y_train_labelled_smote = sampling.fit_sample(X_train_labelled,
labelled_training_set_for_Y.ravel())
X combined test = shuffle(hstack([coo matrix(testing set for X),
coo matrix(testing set for Y)]).toarray(), random state=35)
testing set for X= X combined test[:, :predictor count]
testing set for Y= X combined test[:, predictor count:]
X test smote, Y test smote = SMOTE().fit sample(testing set for X,
testing_set_for_Y.ravel())
Y test smote=Y test smote.reshape(len(Y test smote),1)
X combined test = shuffle(hstack([coo_matrix(X_test_smote),
coo matrix(Y test smote)]).toarray(), random state=35)
testing set for X= X combined test[:, :predictor count]
testing set for Y= X combined test[:, predictor count:]
training_iterations = len(X_train_unlabelled_smote)
for i in range(1, training iterations):
    classifier = GridSearchCV(LinearSVC(penalty='11', dual=False, tol=0.001), [{'C':
np.linspace(0.0001, 10, 50)}], cv=KFold(5), refit=True, n jobs=4)
    classifier.fit(X train labelled smote, Y train labelled smote.ravel())
    scores = classifier.cv_results_['mean_test_score']
    scores std = classifier.cv results ['std test score']
    best SVM param = classifier.best params
    best score = classifier.best_score_
    X margin = np.abs(classifier.decision function(X train unlabelled smote))
    dataset = pd.DataFrame(data={'index': np.arange(0, len(X_margin), 1),
'distance of margin': X margin})
    dataset = dataset.sort_values(by='distance_of_margin')
    check arr = X train labelled smote
    check Y arr = Y train labelled smote
    to use indices = []
    top_data = pd.DataFrame()
    top_data = dataset.iloc[:1]
```

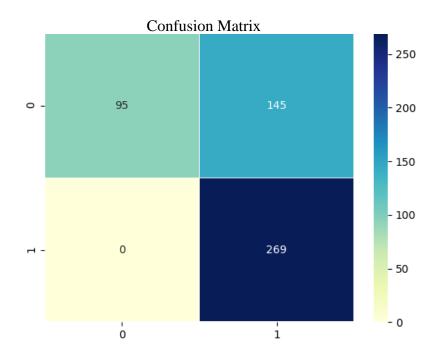
```
for j in range (0, 1):
        index to use = np.int(top data.values[j, 1])
       to use indices.append(index to use)
       check arr = np.vstack((check arr,
np.array(X train unlabelled smote[index to use].reshape(1, 4))))
        Y train unlabelled smote[index to use] =
classifier.predict(X train unlabelled smote[index to use:index to use + 1, :])
        check_Y_arr = np.vstack((check_Y_arr.reshape(len(check_Y_arr),1),
(Y train unlabelled smote[index to use].reshape(1, 1))))
    X train unlabelled smote = np.delete(X train unlabelled smote, to use indices,
axis=0)
    Y train unlabelled smote = np.delete(Y train unlabelled smote, to use indices,
axis=0)
    X train labelled smote = check arr
    Y train labelled smote = check Y arr
    if i == training iterations - 1:
        test set predictions = classifier.predict(testing set for X)
        accuracy = 1 - mean_squared_error(testing_set_for_Y, test_set_predictions)
        fpr_rf_lm, tpr_rf_lm, _ = roc_curve(testing_set_for_Y, test_set_predictions,
pos_label=1)
        roc auc = auc(fpr rf lm, tpr rf lm)
        plt.plot(fpr rf lm, tpr rf lm, label=str(roc auc))
        plt.xlabel('False Positive Rate')
       plt.ylabel('True Positive Rate')
        plt.plot([0, 1], [0, 1], 'k--')
        plt.show()
        confusion matrix = confusion matrix(testing set for Y, test set predictions)
        sns.heatmap(confusion matrix, cmap="Blues", annot=True, linewidths=0.5, fmt='d')
        plt.show()
```

#### d) Unsupervised Learning:

- i) To avoid getting into a local minima, I used hyper parameter *n\_init* to run the k-means algorithm hundred times and then, out of all the different values of number of clusters best value is chosen which has the least distortion.
- ii) The parameter **Kmeans.cluster\_centers\_** gives the co-ordinates of the center for each of the clusters. For the Euclidean distance of each of the data points from the center of each cluster, I used *kmeans.transform()* which transformed the dataset into [Number\_of\_samples X Number\_of\_clusters] matrix. Each data point falls into the cluster it has minimum Euclidean distance from its center. To get thirty data points closet to each cluster, I sorted this matrix. Then, to label each cluster with either 0 or 1, I used majority polling.
- iii) To calculate the misclassification rate and accuracy of the model, I used the k-mean model to predict the labels of the data points and compared them to their true labels.

#### **Results:**

Centroid of first cluster = [0.06910311 0.00396899 0.99224684 0.08623242] Centroid of second cluster = [0.00615781 0.00399807 0.99951781 0.02551618] Accuracy = 0.7151277013752455



```
import numpy as np
import pandas as pd
from sklearn.preprocessing import normalize
from sklearn.cluster import KMeans
import matplotlib.pyplot as plt
from sklearn.utils import shuffle
from imblearn.combine import SMOTEENN
from imblearn.over sampling import SMOTE
from sklearn.metrics import confusion matrix
from scipy.sparse import coo matrix, hstack
import seaborn as sns
dataframe = pd.read csv("transfusion.csv")
dataset for 0 = (dataframe.loc[dataframe['whether'] == 0])
dataset for 1 = (dataframe.loc[dataframe['whether'] == 1])
predictor count = 4
percentage = 0.20
training set X for 0 = dataset for 0.values[(np.int(len(dataset for 0.values) *
percentage)):, :predictor count]
training set Y for 0 = dataset for 0.values[:(np.int(len(dataset for 0.values) *
percentage)), predictor count:]
training set X for 1 = dataset for 1.values[(np.int(len(dataset for 1.values) *
percentage)):, :predictor count]
training_set_Y_for_1 = dataset_for_1.values[:(np.int(len(dataset_for_1.values) *
percentage)), predictor count:]
testing_set_X_for_0 = dataset_for_0.values[:(np.int(len(dataset for 0.values) *
percentage)), :predictor_count]
testing_set_Y_for_0 = dataset_for_0.values[:(np.int(len(dataset for 0.values) *
percentage)), predictor count:]
testing set X for 1 = dataset for 1.values[:(np.int(len(dataset for 1.values) *
percentage)), :predictor count]
testing_set_Y_for_1 = dataset_for_1.values[:(np.int(len(dataset_for_1.values) *
percentage)), predictor count:]
```

```
training data X = normalize(np.vstack((training set X for 0, training set X for 1)),
norm='12')
training data Y = np.vstack((training set Y for 0, training set Y for 1))
testing data X = normalize(np.vstack((testing set X for 0, testing set X for 1)),
testing data Y = np.vstack((testing set Y for 0, testing set Y for 1))
training data X all = shuffle(hstack([coo matrix(training data X),
coo matrix(training data Y)]).toarray(), random state=1)
training_data_X= training_data_X_all[:, :predictor_count]
training_data_Y= training_data_X_all[:, predictor_count:]
training_data_X, training_data_Y = SMOTEENN(random_state = 5,
kind smote='svm').fit sample(training data X, training data Y.ravel())
training data Y=training data Y.reshape(len(training data Y), 1)
testing data X all = shuffle(hstack([coo matrix(testing data X),
coo_matrix(testing_data_Y)]).toarray(), random_state=5)
testing data X= testing data X all[:, :predictor count]
testing data Y= testing data X all[:, predictor count:]
X test smote, Y test smote = SMOTE().fit sample(testing data X,
testing_data_Y.ravel())
Y_test_smote=Y_test_smote.reshape(len(Y_test_smote),1)
testing_data_X_all = shuffle(hstack([coo_matrix(X_test_smote),
coo_matrix(Y_test_smote)]).toarray(), random_state=5)
testing data X = testing data X all[:, :predictor count]
testing data Y = testing data X all[:, predictor count:]
model = KMeans(n clusters=2, n init=20, random state=5)
model.fit(training data X)
inertia = model.inertia
centroids = model.cluster centers
labels = model.labels
labels = labels.reshape(len(labels), 1)
training_data_X_new = model.transform(training_data_X)
data = {'index': np.arange(0, len(training_data_X_new), 1),
'distance_of_margin_for_first_cluster': training_data_X_new[:, :1].reshape(-1),
'distance of margin for second cluster': training data X new[:, 1:2].reshape(-1)}
data frame = pd.DataFrame(data=data)
top \overline{30} for first cluster =
(data frame.sort values(by='distance of margin for first cluster')).iloc[:30]
top_30_for_second_cluster =
(data_frame.sort_values(by='distance_of_margin_for_second_cluster')).iloc[:30]
zero counter for first cluster = 0
one counter for first cluster = 0
zero counter for second cluster = 0
one_counter_for_second_cluster = 0
for i in range (0, 30):
    label = int(training data Y[int(top 30 for first cluster.values[i, 2]), 0])
    if label == 1:
        one counter for first cluster = one counter for first cluster+ 1
    else:
        zero_counter_for_first_cluster = zero_counter_for_first_cluster + 1
    label = int(training_data_Y[int(top_30_for_second_cluster.values[i, 2]), 0])
    if label == 1:
        one counter for second cluster = one counter for second cluster + 1
    else:
        zero counter for second cluster = zero counter for second cluster + 1
if one counter for first cluster >= zero counter for first cluster:
    class_of_first_cluster = 1
```

```
else:
    class_of_first_cluster = 0

if one_counter_for_second_cluster >= zero_counter_for_second_cluster:
    class_of_second_cluster = 1
else:
    class_of_second_cluster = 0

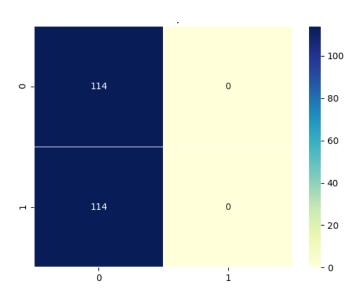
no_of_missclassifications = 0
for i in range(0, len(labels)):
    if int(labels[i, 0]) != int(training_data_Y[i, 0]):
        no_of_missclassifications = no_of_missclassifications + 1

error = no_of_missclassifications / len(labels)
accuracy = 1 - error
confusion_grid=confusion_matrix(training_data_Y, labels)
sns.heatmap(confusion_grid,cmap="Blues",annot=True, linewidths=0.6,fmt='d')
plt.show()
```

### iii. Following results were obtained after predicting the testing dataset on this model:

### Accuracy = 0.5

#### Confusion Matrix



This approach has misclassified class 1, similar to semi-supervised learning technique, therefore, it's performance is unsatisfactory.

```
import numpy as np
import pandas as pd
from sklearn.preprocessing import normalize
from sklearn.cluster import KMeans
import matplotlib.pyplot as plt
from sklearn.utils import shuffle
from imblearn.combine import SMOTEENN
from imblearn.over_sampling import SMOTE
from sklearn.metrics import confusion_matrix
from scipy.sparse import coo_matrix, hstack
import seaborn as sns
```

```
dataframe = pd.read csv("transfusion.csv")
dataset for 0 = (dataframe.loc[dataframe['whether'] == 0])
dataset for 1 = (dataframe.loc[dataframe['whether'] == 1])
predictor count = 4
percentage = 0.20
training set X for 0 = dataset for 0.values[(np.int(len(dataset for 0.values) *
percentage)):, :predictor count]
training set Y for 0 = dataset for 0.values[:(np.int(len(dataset for 0.values) *
percentage)), predictor count:]
training set X for 1 = dataset for 1.values[(np.int(len(dataset for 1.values) *
percentage)):, :predictor count]
training set Y for 1 = dataset for 1.values[:(np.int(len(dataset for 1.values) *
percentage)), predictor count:]
testing set X for 0 = dataset for 0.values[:(np.int(len(dataset for 0.values) *
percentage)), :predictor count]
testing_set_Y_for_0 = dataset_for_0.values[:(np.int(len(dataset_for_0.values) *
percentage)), predictor_count:]
testing_set_X_for_1 = dataset_for_1.values[:(np.int(len(dataset_for_1.values) *
percentage)), :predictor_count]
testing_set_Y_for_1 = dataset_for_1.values[:(np.int(len(dataset_for_1.values) *
percentage)), predictor count:]
training_data_X = normalize(np.vstack((training_set_X_for_0, training_set_X_for_1)),
norm='12')
training data Y = np.vstack((training set Y for 0, training set Y for 1))
testing_data_X = normalize(np.vstack((testing_set_X_for_0, testing_set_X_for_1)),
norm='12')
testing data Y = np.vstack((testing set Y for 0, testing set Y for 1))
training data X all = shuffle(hstack([coo matrix(training data X),
coo matrix(training data Y)]).toarray(), random state=1)
training_data_X= training_data_X_all[:, :predictor_count]
training_data_Y= training_data_X_all[:, predictor_count:]
training_data_X, training_data_Y = SMOTEENN(random_state = 5,
kind smote='svm').fit sample(training data X, training data Y.ravel())
training data Y=training data Y.reshape(len(training data Y), 1)
testing data X all = shuffle(hstack([coo matrix(testing data X),
coo_matrix(testing_data_Y)]).toarray(), random_state=5)
testing data X= testing data X all[:, :predictor count]
testing data Y= testing data X all[:, predictor count:]
X test smote, Y test smote = SMOTE().fit sample(testing data X, testing data Y.ravel())
Y_test_smote=Y_test_smote.reshape(len(Y_test_smote),1)
testing data X all = shuffle(hstack([coo matrix(X test smote),
coo_matrix(Y_test_smote)]).toarray(), random_state=5)
testing data X = testing data X all[:, :predictor count]
testing data Y = testing data X all[:, predictor count:]
model = KMeans(n clusters=2, n init=20, random state=5)
model.fit(training_data_X)
inertia = model.inertia
centroids = model.cluster centers
predictions=model.predict(testing data X)
predictions=predictions.reshape(len(predictions), 1)
no of missclassifications = 0
for i in range(0, len(predictions)):
```

### e) Comparison:

The results show that the as compared to unsupervised learning and semi-supervised learning, supervised learning has performed best in classifying this dataset because the size of dataset is small. Even the semi-supervised learning technique works better than the unsupervised learning.

This can be concluded from the result that on small datasets supervised learning works best whereas unsupervised learning should be used on large datasets.

On our given dataset, supervised learning gives the best results.

# **Question 2:**

### a) Optimal value of k:

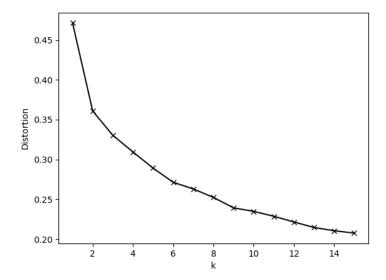
I have used Elbow method to find the optimal number of clusters to made for the given dataset, which determines the best value of *k* by examining the variance percentage.

In this method, we run k-means clustering on the dataset for a range of values of k and for each value of k calculate the sum of squared errors (SSE). The optimal value of k is selected when an increment of another cluster doesn't give a significant improvement in our model, but if adding another cluster reduces the variance significantly, it will be the optimal value of k (if a bigger value doesn't give a significantly better result by reducing the variance). Therefore, when the improvement significantly decreases, we get a pointy graph resulting in an elbow shape.

On our given dataset, using a range of different values (as shown below), the optimal value of k is 2 because the most drastic improvement/reduction in variance or distortion is observed with k = 2, hence elbow can be seen when number of clusters is 2.

#### **Results:**

Distortions for k=1 to 14:[0.6723847012398, 0.520943856709, 0.4823057134, 0.4523059701247, 0.414097234191, 0.39092375093124, 0.381098723509, 0.369071324097178, 0.3478152340987, 0.3423094709255, 0.328913489619, 0.32203847209, 0.3180732508676, 0.30896132419029]



```
import pandas as pd
import numpy as np
import matplotlib.pyplot as plt
from sklearn.utils import shuffle
from scipy.spatial.distance import cdist
from sklearn.cluster import KMeans
df = pd.read csv("Frogs_MFCCs.csv")
df.replace('Bufonidae', int(1), inplace=True)
df.replace('Dendrobatidae', int(2), inplace=True)
df.replace('Hylidae', int(3), inplace=True)
df.replace('Leptodactylidae', int(4), inplace=True)
df.replace('Adenomera', int(5), inplace=True)
df.replace('Ameerega', int(6), inplace=True)
df.replace('Dendropsophus', int(7), inplace=True)
df.replace('Hypsiboas', int(8), inplace=True)
df.replace('Leptodactylus', int(9), inplace=True)
df.replace('Osteocephalus', int(10), inplace=True)
df.replace('Rhinella', int(11), inplace=True)
df.replace('Scinax', int(12), inplace=True)
df.replace('AdenomeraAndre', int(13), inplace=True)
df.replace('AdenomeraHylaedactylus', int(14), inplace=True)
df.replace('Ameeregatrivittata', int(15), inplace=True)
df.replace('HylaMinuta', int(16), inplace=True)
df.replace('HypsiboasCinerascens', int(17), inplace=True)
df.replace('HypsiboasCordobae', int(18), inplace=True)
df.replace('LeptodactylusFuscus', int(19), inplace=True)
df.replace('OsteocephalusOophagus', int(20), inplace=True)
df.replace('Rhinellagranulosa', int(21), inplace=True)
df.replace('ScinaxRuber', int(22), inplace=True)
df.drop('RecordID', inplace=True, axis=1)
df.to csv('Frogs MFCCs new.csv', index=False)
shuffled data = shuffle(df.values, random state=0)
predictors = shuffled data[:, :22]
class_label = shuffled_data[:, 22:]
```

```
number of clusters = range(1, 15)
variances = []
for k in number of clusters:
    model = KMeans(n clusters=k)
    model.fit(predictors)
    variance = sum(np.min(cdist(predictors, model.cluster_centers_, 'euclidean'),
axis=1)) / predictors.shape[0]
    variances.append(variance)
plt.plot(number_of_clusters, variances, 'bx-')
plt.xlabel('K')
plt.ylabel('Distortion')
plt.show()
diff = []
biggest_diff = 0
for i in range(0, len(variances)-1):
    difference in variances = np.abs(variances[i + 1] - variances[i])
    if difference in variances > biggest diff:
        biggest diff = difference in variances
best value of k = k+2
```

### b. K-mean Clustering:

I have clustered the data into two clusters using k-mean clustering because the optimal value of k and determined the species and genus of each of these clusters by majority polling after determining the family of each cluster.

### **Results:**

```
First Cluster:
    Family = Hylidae
    Genus = Hypsiboas
    Species = HypsiboasCordobae

Second Cluster:
    Family = Leptodactylidae
    Genus = Adenomera
    Species = AdenomeraHylaedactylus
```

As it can be seen in the results that the family of first cluster is Hylindae, genus is Hypsiboas and species is HypsiboasCordobae, whereas the family of second cluster is Leptodactylidae, genus is Adenomera and species is AdenomeraHylaedactylus.

```
import pandas as pd
import numpy as np
from sklearn.model selection import train test split
from sklearn.utils import shuffle
from sklearn.cluster import KMeans
import random
df = pd.read csv("Frogs MFCCs.csv")
df.replace('Bufonidae', int(1), inplace=True)
df.replace('Dendrobatidae', int(2), inplace=True)
df.replace('Hylidae', int(3), inplace=True)
df.replace('Leptodactylidae', int(4), inplace=True)
df.replace('Adenomera', int(5), inplace=True)
df.replace('Ameerega', int(6), inplace=True)
df.replace('Dendropsophus', int(7), inplace=True)
df.replace('Hypsiboas', int(8), inplace=True)
df.replace('Leptodactylus',int(9),inplace=True)
df.replace('Osteocephalus', int(10), inplace=True)
df.replace('Rhinella', int(11), inplace=True)
df.replace('Scinax', int(12), inplace=True)
df.replace('AdenomeraAndre', int(13), inplace=True)
df.replace('AdenomeraHylaedactylus', int(14), inplace=True)
df.replace('Ameeregatrivittata', int(15), inplace=True)
df.replace('HylaMinuta', int(16), inplace=True)
df.replace('HypsiboasCinerascens', int(17), inplace=True)
df.replace('HypsiboasCordobae', int(18), inplace=True)
df.replace('LeptodactylusFuscus', int(19), inplace=True)
df.replace('OsteocephalusOophagus',int(20),inplace=True)
df.replace('Rhinellagranulosa', int(21), inplace=True)
df.replace('ScinaxRuber', int(22), inplace=True)
df.drop('RecordID', inplace=True, axis=1)
df.to csv('Frogs MFCCs new.csv',index=False)
shuffled data = shuffle(df.values, random_state=5)
training x, testing x, training y, testing y =
train_test_split(shuffled_data[:,:22], shuffled_data[:,22:], train_size=0.7)
family dataset Y = shuffled data[:, 22:][:, :1]
genus dataset \overline{Y} = shuffled data[:, 22:][:, 1:2]
species dataset Y = shuffled data[:, 22:][:, 2:]
model = KMeans(n_clusters=2, n_init=20, random state=5)
model.fit(shuffled_data[:,:22])
labels = model.labels
print(model.cluster centers )
print(model.inertia_)
family count for first cluster = \{1:0, 2:0, 3:0, 4:0\}
family count for second cluster = \{1:0, 2:0, 3:0, 4:0\}
for i in range(1, len(labels)):
    if int(labels[i]) == 0:
        family_count_for_first_cluster[int(family_dataset_Y[i])] =
family count for first cluster[int(family dataset Y[i])] + 1
    else:
```

```
family count for second cluster[int(family dataset Y[i])] =
family_count_for_second_cluster[int(family_dataset_Y[i])] + 1
majority family for 1 index = int(max(family count for first cluster.keys(),
key=(lambda label: family count for first cluster[label])))
majority family for 2 index = int(max(family count for second cluster.keys(),
key=(lambda label: family count for second cluster[label])))
family labels = {1: 'Bufonidae', 2:
'Dendrobatidae', 3: 'Hylidae', 4: 'Leptodactylidae'}
family of first cluster = family labels[majority family for 1 index]
family of second cluster = family labels[majority family for 2 index]
genus count for first cluster = \{5:0, 6:0, 7:0, 8:0, 9:0, 10:0, 11:0, 12:0\}
genus count for second cluster = {5:0, 6:0, 7:0, 8:0, 9:0, 10:0, 11:0, 12:0}
for i in range(1, len(labels)):
    if int(labels[i]) == 0:
        genus count for first cluster[int(genus dataset Y[i])] =
genus count for first cluster[int(genus dataset Y[i])] + 1
        genus_count_for_second_cluster[int(genus_dataset_Y[i])] =
genus count for second cluster[int(genus dataset Y[i])] + 1
majority_genus_for_1_index = int(max(genus_count_for_first_cluster.keys(),
key=(lambda label: genus count for first cluster[label])))
majority genus for 2 index = int(max(genus count for second cluster.keys(),
key=(lambda label: genus count for second cluster[label])))
genus labels =
{5:'Adenomera', 6:'Ameerega', 7:'Dendropsophus', 8:'Hypsiboas', 9:'Leptodactylus', 10:'O
steocephalus',11:'Rhinella',12:'Scinax'}
genus of first cluster = genus labels[majority genus for 1 index]
genus of second cluster = genus labels[majority genus for 2 index]
species_count_for_first_cluster = {13:0, 14:0, 15:0, 16:0, 17:0, 18:0, 19:0, 20:0,
21:0, 2\overline{2}:0
species count for second cluster = {13:0, 14:0, 15:0, 16:0, 17:0, 18:0, 19:0, 20:0,
21:0, 22:0}
for i in range(1, len(labels)):
    if int(labels[i]) == 1:
        species_count_for_first_cluster[int(species_dataset_Y[i])] =
species_count_for_first_cluster[int(species_dataset_Y[i])] + 1
        species count for second cluster[int(species dataset Y[i])] =
species count for second cluster[int(species dataset Y[i])] + 1
majority species for 1 index = int(max(species count for first cluster.keys(),
key=(lambda label: species count for first cluster[label])))
majority_species_for_2_index = int(max(species_count_for_second_cluster.keys(),
key=(lambda label: species count for second cluster[label])))
species labels =
{13:'AdenomeraAndre',14:'AdenomeraHylaedactylus',15:'Ameeregatrivittata',16:'HylaMi
nuta',17: 'HypsiboasCinerascens',18: 'HypsiboasCordobae',19: 'LeptodactylusFuscus',20:
'OsteocephalusOophagus',21:'Rhinellagranulosa',22:'ScinaxRuber'}
species of first cluster = species labels[majority species for 1 index]
species of second cluster = species labels[majority species for 2 index]
```

### c. Hamming Loss of the Majority Triplets:

I have calculated the hamming loss by comparing the true labels to the labels predicted by clustering and then, calculated the mean of these hamming losses to come up with an average hamming loss of the given dataset.

#### **Results:**

```
Family of first cluster = Hylidae
Family of second cluster = Leptodactylidae
Hamming Loss = 0.2309237535

Genus of first cluster = Hypsiboas
Genus of second cluster = Adenomera
Hamming Loss = 0.29915674786

Species of first cluster = HypsiboasCordobae
Species of second cluster = AdenomeraHylaedactylus
Hamming Loss = 0.372407918324
```

Average Hamming Loss = 0.29233209406812

### **Majority Triplet of first cluster:**

Family = Hylidae, Genus = Hypsiboas, Species = HypsiboasCordobae

### **Majority Triplet of second cluster:**

Family = Leptodactylidae, Genus = Adenomera, Species = AdenomeraHylaedactylus

```
import pandas as pd
import numpy as np
from sklearn.model selection import train test split
from sklearn.utils import shuffle
from sklearn.cluster import KMeans
from sklearn.metrics import hamming loss
df = pd.read csv("Frogs_MFCCs.csv")
df.replace('Bufonidae', int(1), inplace=True)
df.replace('Dendrobatidae', int(2), inplace=True)
df.replace('Hylidae', int(3), inplace=True)
df.replace('Leptodactylidae', int(4), inplace=True)
df.replace('Adenomera', int(5), inplace=True)
df.replace('Ameerega', int(6), inplace=True)
df.replace('Dendropsophus', int(7), inplace=True)
df.replace('Hypsiboas', int(8), inplace=True)
df.replace('Leptodactylus', int(9), inplace=True)
df.replace('Osteocephalus', int(10), inplace=True)
df.replace('Rhinella', int(11), inplace=True)
df.replace('Scinax',int(12),inplace=True)
df.replace('AdenomeraAndre', int(13), inplace=True)
df.replace('AdenomeraHylaedactylus', int(14), inplace=True)
df.replace('Ameeregatrivittata', int(15), inplace=True)
df.replace('HylaMinuta', int(16), inplace=True)
df.replace('HypsiboasCinerascens', int(17), inplace=True)
```

```
df.replace('HypsiboasCordobae', int(18), inplace=True)
df.replace('LeptodactylusFuscus', int(19), inplace=True)
df.replace('OsteocephalusOophagus', int(20), inplace=True)
df.replace('Rhinellagranulosa', int(21), inplace=True)
df.replace('ScinaxRuber', int(22), inplace=True)
df.drop('RecordID', inplace=True, axis=1)
df.to csv('Frogs MFCCs new.csv',index=False)
shuffled data = shuffle(df.values, random state=5)
training x, testing x, training y, testing y = train test split(shuffled data[:,:22],
shuffled data[:,22:], train size=0.7)
family dataset Y = shuffled data[:, 22:][:, :1]
genus dataset \overline{Y} = shuffled data[:, 22:][:, 1:2]
species dataset Y = shuffled data[:, 22:][:, 2:]
model = KMeans(n clusters=2, n init=20, random state=5)
model.fit(shuffled data[:,:22])
labels = model.labels
print(model.cluster centers)
print(model.inertia_)
family_count_for_first_cluster = {1:0, 2:0, 3:0, 4:0}
family_count_for_second_cluster = \{1:0, 2:0, 3:0, 4:0\}
for i in range(1, len(labels)):
    if int(labels[i]) == 0:
        family_count_for_first_cluster[int(family_dataset_Y[i])] =
family_count_for_first_cluster[int(family_dataset_Y[i])] + 1
    else:
        family count for second cluster[int(family dataset Y[i])] =
family count for_second_cluster[int(family_dataset_Y[i])] + 1
majority family for 1 index = int(max(family count for first cluster.keys(),
key=(lambda label: family count for first cluster[label])))
majority family for 2 index = int(max(family count for second cluster.keys(),
key=(lambda label: family_count_for_second_cluster[label])))
family labels = {1: 'Bufonidae', 2: 'Dendrobatidae', 3: 'Hylidae', 4: 'Leptodactylidae'}
family of first cluster = family labels[majority family for 1 index]
family of second cluster = family labels [majority family for 2 index]
family predictions = np.arange(0, len(labels), 1)
for i in range(1, len(labels)):
    if labels[i] == 1:
        family predictions[i] = majority family for 1 index
    else:
        family predictions[i] = majority family for 2 index
family hamming loss = hamming loss(family dataset Y,
family predictions.reshape(len(family predictions), 1))
genus count for first cluster = {5:0, 6:0, 7:0, 8:0, 9:0, 10:0, 11:0, 12:0}
genus_count_for_second_cluster = {5:0, 6:0, 7:0, 8:0, 9:0, 10:0, 11:0, 12:0}
for i in range(1, len(labels)):
    if int(labels[i]) == 0:
        genus count for first cluster[int(genus dataset Y[i])] =
genus_count_for_first_cluster[int(genus_dataset_Y[i])] + 1
        genus count for second cluster[int(genus dataset Y[i])] =
genus count for second cluster[int(genus dataset Y[i])] + 1
majority_genus_for_1_index = int(max(genus_count_for_first_cluster.keys(),
key=(lambda label: genus count for first cluster[label])))
majority genus_for_2_index = int(max(genus_count_for_second_cluster.keys(),
```

```
key=(lambda label: genus count for second cluster[label])))
genus labels =
{5:'Adenomera', 6:'Ameerega', 7:'Dendropsophus', 8:'Hypsiboas', 9:'Leptodactylus', 10:'Ost
eocephalus',11:'Rhinella',12:'Scinax'}
genus of first cluster = genus labels[majority genus for 1 index]
genus of second cluster = genus labels[majority genus for 2 index]
genus predictions = np.arange(0, len(labels), 1)
for i in range(1, len(labels)):
    if labels[i] == 1:
        genus predictions[i] = majority genus for 1 index
        genus predictions[i] = majority genus for 2 index
genus hamming loss = hamming loss (genus dataset Y,
genus predictions.reshape(len(genus predictions), 1))
species count for first cluster = {13:0, 14:0, 15:0, 16:0, 17:0, 18:0, 19:0, 20:0,
21:0, 22:0}
species count for second cluster = {13:0, 14:0, 15:0, 16:0, 17:0, 18:0, 19:0, 20:0,
21:0, 22:0}
for i in range(1, len(labels)):
    if int(labels[i]) == 1:
        species count for first cluster[int(species dataset Y[i])] =
species count for first cluster[int(species dataset Y[i])] + 1
    else:
        species count for second cluster[int(species dataset Y[i])] =
species count for second cluster[int(species dataset Y[i])] + 1
majority species for 1 index = int(max(species count for first cluster.keys(),
key=(lambda label: species count for first cluster[label])))
majority species for 2 index = int(max(species count for second cluster.keys(),
key=(lambda label: species count for second cluster[label])))
species_labels =
{13:'AdenomeraAndre',14:'AdenomeraHylaedactylus',15:'Ameeregatrivittata',16:'HylaMinu
ta',17:'HypsiboasCinerascens',18:'HypsiboasCordobae',19:'LeptodactylusFuscus',20:'Ost
eocephalusOophagus',21:'Rhinellagranulosa',22:'ScinaxRuber'}
species of first cluster = species labels[majority species for 1 index]
species of second cluster = species labels[majority species for 2 index]
species predictions = np.arange(0, len(labels), 1)
for i in range(1, len(labels)):
    if labels[i] == 1:
        species predictions[i] = majority species for 1 index
    else:
        species_predictions[i] = majority_species for 2 index
species hamming loss = hamming loss(species dataset Y,
species predictions.reshape(len(species predictions), 1))
average hamming loss = np.mean(np.array([family hamming loss, genus hamming loss,
species hamming loss]))
```

# **Question 3:**

