

## PREDICTIVE ANALYTICS MSBA 6420

Homework #3

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(put your name above)

Total grade: \_\_\_\_\_ out of \_\_\_\_100\_\_\_ points

There are 2 numbered questions.	Please answer the	m all and submit your	· assignment as a sii	ngle PDF file by
uploading it to the course website.				

Quation 1. (50 points) Use numeric prediction techniques to build a predictive model for the HW3.xlsx dataset. This dataset is provided on the course website and contains data about whether or not different consumers made a purchase in response to a test mailing of a certain catalog and, in case of a purchase, how much money each consumer spent. The data file has a brief description of all the attributes in a separate worksheet. Note that this dataset has two possible outcome variables: Purchase (0/1 value: whether or not the purchase was made) and Spending (numeric value: amount spent).

## Your tasks:

- (a) (20 points) Build numeric prediction models that predict Spending based on the other available customer information (obviously, not including the Purchase attribute among the inputs!). Use linear regression, k-NN, regression tree, SVM regression and Neural Network and ensembling models. Briefly discuss your explorations and present the best result (best predictive model) for each of these techniques. Compare the techniques; which of them provides the best predictive performance? Please make sure you use best practices for predictive modeling. (I.e., do you need to set which hyper-parameter? Normalize?)
- (b) (20 points) As a variation on this exercise, create a separate "restricted" dataset (i.e., a subset of the original dataset), which includes only purchase records (i.e., where Purchase = 1). Build numeric prediction models to predict Spending for this restricted dataset. All the same requirements as for task (a) apply.
- (c) (10 points) For each predictive modeling technique, discuss the predictive performance differences between the models built for task (a) vs. task (b): which models exhibit better predictive performance? Why do you think that is?

In [59]: import pandas as pd from matplotlib import pyplot as plt %matplotlib inline import numpy as np np.random.seed(13) import tensorflow as tf import pandas as pd from tensorflow.keras import backend as K from tensorflow.keras.models import Sequential from tensorflow.keras.datasets import mnist from tensorflow.keras.layers import Dense, Activation, Dropout from tensorflow.keras.utils import to categorical from tensorflow.keras.wrappers.scikit\_learn import KerasClassifier import matplotlib.pyplot as plt import scikitplot as skplt from sklearn import neighbors, datasets, tree, linear model, metrics, svm from sklearn.model\_selection import cross val score, train test split, KFold import itertools from itertools import permutations from sklearn.metrics import recall score, confusion matrix, mean squared error from sklearn.model\_selection import GridSearchCV 1. Import data In [60]: df=pd.read excel('HW3.xlsx') df=df.drop(columns=['sequence number', 'Purchase']) 2. Formating the data In [61]: c=['US', 'source a', 'source c', 'source b', 'source d', 'source e', 'source m', 'source o', 'source h', 'source r', 'source s', 'source t', 'source u', 'source p', 'source x', 'source w', 'Web order', 'Gender=male', 'Ad dress is res'] df[c] = df[c].astype(str)In [62]: df.dtypes Out[62]: US object source a object object source c source b object source d object source e object source m object source o object object source h object source r object source\_s source t object object source u object source p source\_x object object source w int64 last update days ago int64 1st\_update\_days\_ago int64 Web order object Gender=male object Address is res object Spending float64 dtype: object (a) In [63]: #Split Train and Test Data x\_train, x\_test, y\_train, y\_test = train\_test\_split(df.drop(columns=['Spending']), df['Spending'], test\_size =0.2, random state=9) #Standardization from sklearn.preprocessing import MinMaxScaler mms = MinMaxScaler() mms.fit(x train[['Freq','last update days ago','1st update days ago']]) x\_train[['Freq','last\_update\_days\_ago','1st\_update\_days\_ago']]=mms.transform(x\_train[['Freq','last\_upda te days ago', '1st update days ago']]) x test[['Freq','last update days ago','1st update days ago']]=mms.transform(x test[['Freq','last update days ago','1st update days ago']]) In [65]: #Elastic Net Regression from sklearn.linear\_model import ElasticNet grid={ 'alpha': [1,2,3,4,5,6,7,8,9,10], 'l1\_ratio':[0.3,0.5,0.7]} EN linear=ElasticNet(random state=99) model = GridSearchCV(estimator=EN linear, param grid=grid, cv=5,scoring='neg mean squared error') model.fit(x\_train,y\_train) print("With CV grid search, I found the best hyperparameter is alpha={} and L1 ratio={}.".format(model. best params ['alpha'], model.best params ['ll ratio'])) print("MSE on Test Data: {}".format(round((metrics.mean\_squared\_error(y\_test, model.predict(x\_test))),2 ) ) ) With CV grid search, I found the best hyperparameter is alpha=1 and L1 ratio=0.7. MSE on Test Data: 31795.78 In [66]: #KNN regression from sklearn.neighbors import KNeighborsRegressor grid={'n neighbors':[5,10,15,20,50,100,500], 'weights':['uniform','distance']} KNN=KNeighborsRegressor() model = GridSearchCV(estimator=KNN, param grid=grid, cv=5,scoring='neg mean squared error') model.fit(x train, y train) print("With CV grid search, I found the best hyperparameter is n\_neighbors={} and weights={}.".format(m odel.best params ['n neighbors'], model.best params ['weights'])) print("MSE on Test Data: {}".format(round((metrics.mean\_squared\_error(y\_test, model.predict(x\_test))),2 ))) With CV grid search, I found the best hyperparameter is n neighbors=5 and weights=distance. MSE on Test Data: 33890.18 In [67]: #Regression Tree from sklearn.tree import DecisionTreeRegressor grid={ 'max depth': [2,3,4,5,6,7,8,9,10,20], 'min\_samples\_split':[2,5,10,15,20]} Tree=DecisionTreeRegressor(random\_state=99) model = GridSearchCV(estimator=Tree, param\_grid=grid, cv=5,scoring='neg\_mean\_squared\_error') model.fit(x train,y train) print("With CV grid search, I found the best hyperparameter is max depth={} and min sample split={}.".f ormat(model.best\_params\_['max\_depth'],model.best\_params\_['min\_samples\_split'])) print("MSE on Test Data: {}".format(round((metrics.mean\_squared\_error(y\_test, model.predict(x\_test))),2 ))) With CV grid search, I found the best hyperparameter is max depth=7 and min sample split=15. MSE on Test Data: 23458.39 In [68]: | #SVM Regression from sklearn.svm import SVR grid={'kernel':['linear','rbf'], 'C': [1,2,3,4,5,6,7,8,9,10]} s=SVR()model = GridSearchCV(estimator=s, param\_grid=grid, cv=5,scoring='neg\_mean\_squared\_error') model.fit(x train, y train) print("With CV grid search, I found the best hyperparameter is kernel={} and C={}.".format(model.best p arams ['kernel'], model.best params ['C'])) print("MSE on Test Data: {}".format(round((metrics.mean\_squared\_error(y\_test, model.predict(x\_test))),2 With CV grid search, I found the best hyperparameter is kernel=linear and C=10. MSE on Test Data: 27265.73 In [69]: # Random Search With Neural Network from tensorflow import keras from tensorflow.keras import layers from kerastuner.tuners import RandomSearch In [70]: def build\_model(hp): model = keras.Sequential() model.add(layers.Dense(units=hp.Int('units', min value=10, max\_value=100, step=10), activation='relu',input dim=22)) model.add(layers.Dense(units=hp.Int('units', min value=10, max value=100, step=10), activation='relu')) model.add(layers.Dense(1,activation='linear')) model.compile(optimizer=keras.optimizers.Adam(hp.Choice('learning rate', values=[0.01,0.001,0.0001])), loss='mse', metrics=['mse']) return model tuner=RandomSearch (build model, objective='mse', max trials=3, overwrite=True, executions per trial=3) #Keras cannot input object data type, so no matter the column is boolean or numeric we need to transfor m them to float32 x train float = np.asarray(x train).astype(np.float32) y\_train\_float = np.asarray(y\_train).astype(np.float32) x\_test\_float = np.asarray(x\_test).astype(np.float32) y\_test\_float = np.asarray(y\_test).astype(np.float32)  $tuner.search (x=x\_train\_float, y=y\_train\_float, epochs=200, batch\_size=32, validation\_data=(x\_test\_float, y\_test\_float, y\_test$ est\_float)) Trial 3 Complete [00h 00m 40s] mse: 18517.412109375 Best mse So Far: 5621.048177083333 Total elapsed time: 00h 01m 58s INFO:tensorflow:Oracle triggered exit In [72]: result=tuner.get\_best\_hyperparameters()[0].values print('The best 3 layers NN parameters would be {} neurons and {} learning rate.'.format(result['units' ],result['learning\_rate'])) print('----') print("The best model's mse on test data = 5621") print('----') print(tuner.results\_summary()) The best 3 layers NN parameters would be 80 neurons and 0.01 learning rate. The best model's mse on test data = 5621 \_\_\_\_\_\_ Results summary Results in ./untitled\_project Showing 10 best trials Objective (name='mse', direction='min') Trial summary Hyperparameters: units: 80 learning rate: 0.01 Score: 5621.048177083333 Trial summary Hyperparameters: units: 70 learning rate: 0.0001 Score: 18517.412109375 Trial summary Hyperparameters: units: 60 learning\_rate: 0.0001 Score: 20536.832682291668 None In [73]: #Random Forest Regression from sklearn.ensemble import RandomForestRegressor grid={'n estimators':[100,200,300,400,500], 'max depth': [1,2,3,4,5,6,7,8,9,10]} rf=RandomForestRegressor() model = GridSearchCV(estimator=rf, param\_grid=grid, cv=5,scoring='neg mean squared error') model.fit(x train,y train) print("With CV grid search, I found the best hyperparameter is n estimators={} and max depth={}.".forma t (model.best\_params\_['n\_estimators'], model.best\_params\_['max\_depth'])) print("MSE on Test Data: {}".format(round((metrics.mean\_squared\_error(y\_test, model.predict(x\_test))),2 ) ) ) With CV grid search, I found the best hyperparameter is n\_estimators=100 and max\_depth=8. MSE on Test Data: 23006.77 **Conclusion:** The 3 layers Neural Network with 80 Neurons in each layer and learning rate = 0.01 has the best predictive ability. The MSE on test data is around 5484. (b) In [76]: | df=pd.read excel('HW3.xlsx') #Include only the purchase data df=df[df['Purchase']==1] df=df.drop(columns=['sequence\_number','Purchase']) #Formating c=['US', 'source a', 'source c', 'source b', 'source d', 'source e', 'source m', 'source o', 'source h', 'source\_r', 'source s', 'source t', 'source u', 'source p', 'source x', 'source w', 'Web order', 'Gender=male', 'Ad dress is res'] df[c] = df[c].astype(str)#Split Train and Test Data x train, x test, y train, y test = train test split(df.drop(columns=['Spending']), df['Spending'], test size =0.2, random state=9) #Standardization from sklearn.preprocessing import MinMaxScaler mms = MinMaxScaler() mms.fit(x train[['Freq','last update days ago','1st update days ago']]) x\_train[['Freq','last\_update\_days\_ago','1st\_update\_days\_ago']]=mms.transform(x\_train[['Freq','last\_upda te days ago','1st update days ago']]) x test[['Freq','last update days ago','1st update days ago']]=mms.transform(x test[['Freq','last update \_days\_ago','1st\_update\_days\_ago']]) In [77]: #Elastic Net Regression from sklearn.linear\_model import ElasticNet grid={ 'alpha': [1,2,3,4,5,6,7,8,9,10], 'll ratio':[0.3,0.5,0.7]} EN linear=ElasticNet(random state=99) model = GridSearchCV(estimator=EN\_linear, param\_grid=grid, cv=5,scoring='neg\_mean\_squared\_error') model.fit(x\_train,y\_train) print("With CV grid search, I found the best hyperparameter is alpha={} and L1 ratio={}.".format(model. best\_params\_['alpha'], model.best\_params\_['l1\_ratio'])) print("MSE on Test Data: {}".format(round((metrics.mean squared error(y test, model.predict(x test))),2 With CV grid search, I found the best hyperparameter is alpha=1 and L1 ratio=0.7. MSE on Test Data: 58009.48 In [78]: #KNN regression from sklearn.neighbors import KNeighborsRegressor grid={ 'n\_neighbors': [5,10,15,20,50,100,500], 'weights':['uniform','distance']} KNN=KNeighborsRegressor() model = GridSearchCV(estimator=KNN, param\_grid=grid, cv=5,scoring='neg\_mean\_squared\_error') model.fit(x\_train,y\_train) print("With CV grid search, I found the best hyperparameter is n neighbors={} and weights={}.".format(m odel.best params ['n neighbors'], model.best params ['weights'])) print("MSE on Test Data: {}".format(round((metrics.mean squared error(y test, model.predict(x test))),2 ))) With CV grid search, I found the best hyperparameter is n neighbors=5 and weights=distance. MSE on Test Data: 62928.14 In [79]: #Regression Tree from sklearn.tree import DecisionTreeRegressor grid={ 'max\_depth': [2,3,4,5,6,7,8,9,10,20], 'min\_samples\_split':[2,5,10,15,20]} Tree=DecisionTreeRegressor(random state=99) model = GridSearchCV(estimator=Tree, param grid=grid, cv=5,scoring='neg mean squared error') model.fit(x\_train,y\_train) print("With CV grid search, I found the best hyperparameter is max depth={} and min sample split={}.".f ormat(model.best params ['max depth'], model.best params ['min samples split'])) print("MSE on Test Data: {}".format(round((metrics.mean\_squared\_error(y\_test, model.predict(x\_test))),2 ) ) ) With CV grid search, I found the best hyperparameter is max depth=3 and min sample split=2. MSE on Test Data: 38713.35 In [80]: #SVM Regression from sklearn.svm import SVR grid={'kernel':['linear','rbf'], 'C':[1,2,3,4,5,6,7,8,9,10]} model = GridSearchCV(estimator=s, param grid=grid, cv=5,scoring='neg mean squared error') model.fit(x\_train,y\_train) print("With CV grid search, I found the best hyperparameter is kernel={} and C={}.".format(model.best p arams\_['kernel'], model.best\_params\_['C'])) print("MSE on Test Data: {}".format(round((metrics.mean\_squared\_error(y\_test, model.predict(x\_test))),2 ))) With CV grid search, I found the best hyperparameter is kernel=linear and C=10. MSE on Test Data: 51984.36 In [81]: # Random Search With Neural Network from tensorflow import keras from tensorflow.keras import layers from kerastuner.tuners import RandomSearch In [82]: def build model(hp): model = keras.Sequential() model.add(layers.Dense(units=hp.Int('units', min\_value=10, max value=100, step=10), activation='relu',input\_dim=22)) model.add(layers.Dense(units=hp.Int('units', min value=10, max value=100, step=10), activation='relu')) model.add(layers.Dense(1,activation='linear')) model.compile(optimizer=keras.optimizers.Adam(hp.Choice('learning rate', values=[0.01,0.001,0.0001])), loss='mse', metrics=['mse']) return model tuner=RandomSearch(build\_model, objective='mse', max trials=3, overwrite=True, executions\_per\_trial=3) #Keras cannot input object data type, so no matter the column is boolean or numeric we need to transfor m them to float32 x\_train\_float = np.asarray(x\_train).astype(np.float32) y\_train\_float = np.asarray(y\_train).astype(np.float32) x\_test\_float = np.asarray(x\_test).astype(np.float32) y\_test\_float = np.asarray(y\_test).astype(np.float32)  $tuner.search (x=x\_train\_float, y=y\_train\_float, epochs=200, batch\_size=32, validation\_data=(x\_test\_float, y\_test\_float, y\_test$ est\_float)) Trial 3 Complete [00h 00m 27s] mse: 37634.412760416664 Best mse So Far: 9512.133463541666 Total elapsed time: 00h 01m 20s INFO:tensorflow:Oracle triggered exit In [83]: result=tuner.get\_best\_hyperparameters()[0].values print('The best 3 layers NN parameters would be {} neurons and {} learning rate.'.format(result['units' ],result['learning\_rate'])) print('----') print("The best model's mse on test data = 9512") print(tuner.results\_summary()) The best 3 layers NN parameters would be 80 neurons and 0.01 learning rate. The best model's mse on test data = 9512 Results summary Results in ./untitled\_project Showing 10 best trials Objective(name='mse', direction='min') Trial summary Hyperparameters: units: 80 learning\_rate: 0.01 Score: 9512.133463541666 Trial summary Hyperparameters: units: 60 learning\_rate: 0.001 Score: 19139.3359375 Trial summary Hyperparameters: units: 90 learning rate: 0.0001 Score: 37634.412760416664 None In [84]: #Random Forest Regression from sklearn.ensemble import RandomForestRegressor grid={'n estimators':[100,200,300,400,500], 'max\_depth': [1,2,3,4,5,6,7,8,9,10]} rf=RandomForestRegressor() model = GridSearchCV(estimator=rf, param grid=grid, cv=5, scoring='neg mean squared error')

model.fit(x\_train,y\_train)

MSE on Test Data: 39281.85

Therefore, out models' performances drop dramatically.

)))

(c)

print("With CV grid search, I found the best hyperparameter is n\_estimators={} and max\_depth={}.".forma

print("MSE on Test Data: {}".format(round((metrics.mean\_squared\_error(y\_test, model.predict(x\_test))),2

For full data, including both purchased and non-purchased data, tree based models (Regression Tree, Ramdom Forest Regression) and Neural Network have the better performances. On the other hand, for purchased data only, tree based models (Regression Tree,

excluded, all models' performances droppped dramatically. I believe the reason is because those who do not purchase and spend 0 dollars have very similar traits. On the other hand, those who do purchase and spend more than 0 dollar have very diverse traits.

Random Forest Regression) and Neural Network also have the better performances. In general, after the purchased data is

With CV grid search, I found the best hyperparameter is n\_estimators=400 and max\_depth=4.

t (model.best\_params\_['n\_estimators'], model.best\_params\_['max\_depth']))

Quation 2. (50 points) Download the dataset on spam vs. non-spam emails from the following URL: <a href="http://archive.ics.uci.edu/ml/datasets/Spambase">http://archive.ics.uci.edu/ml/datasets/Spambase</a>. Specifically, (i) file "spambase.data" contains the actual data, and (ii) files "spambase.names" and "spambase.DOCUMENTATION" contain the description of the data. This dataset has 4601 records, each record representing a different email message. Each record is described with 58 attributes (indicated in the aforementioned .names file): attributes 1-57 represent various content-based characteristics already extracted from each email message (related to the frequency of certain words or certain punctuation symbols in a message as well as to the usage of capital letters in a message), and the last attribute represents the class label for each message (spam or non-spam).

Task: The general task for this assignment is to build two different models for detecting spam messages (based on the email characteristics that are given): (i) the best possible model that you can build in terms of the overall predictive accuracy (i.e., not taking any cost information into account), and (ii) the best cost-sensitive classification model that you can build in terms of the average misclassification cost.

Some specific instructions for your assignment/write-up:

- Start working on the assignment early.
- Make sure to explore multiple classification techniques (we have learned quite a few of them in the class by now).
  - o Also, make sure to explore different hyper-param for each technique (for example, try several different values of k for k-NN) to find which configurations work best for this application.
- Make sure to explore the impact of various data pre-processing techniques (e.g., normalization).
- When building cost-sensitive prediction models, use 10:1 cost ratio for different misclassification errors. (It should be pretty clear which of the two errors false positive or false negative is the costlier one in this scenario.)
- In general, use best practices when evaluating the models: nested CV, discuss the confusion matrix and some relevant performance metrics (accuracy, precision, recall, f-measure, AUC, average misclassification cost...), show some visual indications of model performance (ROC curves, lift charts).
- As a deliverable, produce a write-up describing your aforementioned explorations. Report the performances of different models that you tried. Discuss the best models in two different tasks in detail (which parameters worked best, what was the performance), provide some comparisons. (upload your code as well) Draw some conclusions from the assignment.
- Evaluation: 50 points:
  - O Performance: 30 points (based on the performance achieved by your best reported models).
  - Exploration/write-up: 20 points (based on the comprehensiveness of your exploration, i.e., when searching for the best performing model, did you evaluate and report just one or two techniques, or did you try a number of different variations, based on what you know from the class?).

4. Data Normalization In [74]: from sklearn.preprocessing import MinMaxScaler scaler=MinMaxScaler() scaler=scaler.fit(x train) x train[x train.columns] = scaler.transform(x train[x train.columns]) x\_test[x\_test.columns] = scaler.transform(x\_test[x\_test.columns]) 5. Nested Grid Search CV to find the best model, using accuracy In [75]: # Create the CV inner\_cv = KFold(n\_splits=5, shuffle=True, random\_state=9) outer cv = KFold(n splits=5, shuffle=True, random state=9) # Create the Classifier def create\_model(activation, nb\_hidden): model = Sequential() model.add(Dense(nb hidden, input dim=57, activation=activation)) model.add(Dense(nb\_hidden, activation=activation)) model.add(Dense(nb\_hidden, activation=activation)) model.add(Dense(1, activation='sigmoid')) model.compile(loss='binary crossentropy', optimizer='adam', metrics=['accuracy']) NN = KerasClassifier(build\_fn=create\_model, epochs=30, batch size=256, verbose=0) t=tree.DecisionTreeClassifier(random state=9) knn=neighbors.KNeighborsClassifier() s=svm.SVC(random\_state=9) # Set up the parameter grid NN grid = {'activation':['relu', 'sigmoid'], 'nb hidden':[10,20,30,40,50,60,70,80,90,100]} tree\_grid={'criterion':['gini','entropy'], 'max\_depth':list(range(30))} knn\_grid={'weights':['uniform','distance'], 'n neighbors':list(range(5,31))} svm\_grid={'C':[0.1,1,5,10,50,100], 'gamma': [1,5,10,15,20,25,30,50,100], 'kernel':['linear','rbf']} #Nested CV for Neural Network clf = GridSearchCV(estimator=NN, param\_grid=NN\_grid, cv=inner\_cv,scoring='accuracy') nested\_score = cross\_val\_score(clf, X=x\_train, y=y\_train, cv=outer\_cv,scoring='accuracy') NN\_result=nested\_score.mean() #Nested CV for SVM clf = GridSearchCV(estimator=s, param\_grid=svm\_grid, cv=inner\_cv,scoring='accuracy') nested\_score = cross\_val\_score(clf, X=x\_train, y=y\_train, cv=outer\_cv,scoring='accuracy') svm result=nested score.mean() #Nested CV for Decision Tree clf = GridSearchCV(estimator=t, param grid=tree grid, cv=inner cv,scoring='accuracy') nested\_score = cross\_val\_score(clf, X=x\_train, y=y\_train, cv=outer\_cv,scoring='accuracy') tree\_result=nested\_score.mean() #Nested CV for KNN clf = GridSearchCV(estimator=knn, param\_grid=knn\_grid, cv=inner\_cv,scoring='accuracy') nested score = cross val score(clf, X=x train, y=y train, cv=outer cv,scoring='accuracy') knn result=nested score.mean() print('Average Performance of Neural Network Classifier: {}%'.format(round(NN result\*100,2))) print('Average Performance of SVM Classifier: {}%'.format(round(svm\_result\*100,2))) print('Average Performance of Decision Tree Classifier: {}%'.format(round(tree\_result\*100,2))) print('Average Performance of KNN Classifier: {}%'.format(round(knn result\*100,2))) Average Performance of Neural Network Classifier: 94.64% Average Performance of SVM Classifier: 94.08% Average Performance of Decision Tree Classifier: 93.14% Average Performance of KNN Classifier: 91.94% 6-1. Use GridSearch to find best hyper-parameters for 4-layers NN model

## In [77]: #Grid Search Verison # Create the Classifier model = Sequential() return model #Result

model = keras.Sequential()

return model

m them to float32

est float))

In [79]: #Get the best model

Results summary

Trial summary Hyperparameters:

units: 100

units: 60

units: 50

units: 30

units: 90

tem=[]

prob=tem y pred=[]

In [83]:

else:

#create ROC curve

plt.plot(fpr,tpr)

plt.show()

True

Predict

1.0

0.6

0.2

0.0

In [84]:

0.0

#Grid Search Verison

return model

# Create the Classifier

model = Sequential()

def create model(activation, nb hidden):

#Grid Search CV for Neural Network

odel3.best params ['nb hidden']))

e/checkpoint#loading mechanics for details.

Recall positive on Test Data: 95.77%

y pred=NNmodel3.predict(x test)

1.5

50 340

In [85]: | #create the confucion matrix

516

(150+50) / 921 = 0.217.

True Predict

 $\Omega$ 

The best parameters activation=relu & nb hidden=100 Prediction Accuracy Score on Test Data: 92.94%

y\_pred=[ y\_pred[i][0] for i in range(len(x test))] tem=pd.DataFrame({'True':y\_test,'Predict':y\_pred})

print(tem.groupby(['Predict','True']).size().unstack())

class\_weight =  $\{0: 1., 1:10.\}$ 

3.predict(x test))\*100,2)))

test),pos\_label=1)\*100,2)))

0.2

True Positive Rate

plt.title('ROC Curve')

542

In [80]: #result

tuner=RandomSearch (build model,

Trial 5 Complete [00h 00m 08s] accuracy: 0.9723672270774841

Total elapsed time: 00h 00m 42s

],result['learning rate'])) print('-----

print(tuner.results summary())

Results in ./untitled\_project

Showing 10 best trials

learning rate: 0.01

learning\_rate: 0.01

learning\_rate: 0.01

learning rate: 0.01

learning\_rate: 0.0001 Score: 0.8956345677375793

In [81]: | prob=NNmodel2.predict(x\_test)

for i in range(len(prob)): tem.append(prob[i][0])

for i in range(len(prob)): **if** prob[i]>0.5:

y\_pred.append('1')

y\_pred.append('0')

plt.ylabel('True Positive Rate') plt.xlabel('False Positive Rate')

1

23

24 332

tem=pd.DataFrame({'True':y test, 'Predict':y pred})

ROC Curve

0.6

NN model with different class weights.

model.add(Dense(nb hidden, activation=activation)) model.add(Dense(nb hidden, activation=activation))

NNmodel3.fit(x\_train, y\_train,class\_weight=class\_weight)

model.add(Dense(1, activation='sigmoid'))

False Positive Rate

0.8

matrix and cost structure, the average misclassification rate = (230+24) / 921 = 0.2757

model.add(Dense(nb hidden, input dim=57, activation=activation))

1.0

According to the Lift Curve, I find the SVM model is really good at capturing spam cases. In addition according to the confustion

10. According to the cost model, the consequence of mis-classifying positive cases are more severe. The weight for class 1 is 10, and the weight for class 0 is 1. Hence, this time I try to find the best 4-layers

NN grid = { 'activation':['relu', 'sigmoid'], 'nb hidden':[10,20,30,40,50,60,70,80,90,100]}

model.compile(loss='binary crossentropy', optimizer='adam', metrics=['accuracy'])

print('The best parameters activation={} & nb hidden={}'.format(NNmodel3.best params ['activation'],NNm

print("Prediction Accuracy Score on Test Data: {}%".format(round(metrics.accuracy score(y test, NNmodel

print("Recall positive on Test Data: {}%".format(round(metrics.recall score(y test, NNmodel3.predict(x

WARNING: tensorflow: A checkpoint was restored (e.g. tf.train.Checkpoint.restore or tf.keras.Model.load weights) but not all checkpointed values were used. See above for specific issues. Use expect partia 1() on the load status object, e.g. tf.train.Checkpoint.restore(...).expect partial(), to silence the se warnings, or use assert consumed() to make the check explicit. See https://www.tensorflow.org/guid

According to the confustion matrix, after taking class wights into account, the average misclassification rate drop dramatically to

NN = KerasClassifier(build fn=create model, epochs=30, batch size=256, verbose=0)

NNmodel3 = GridSearchCV(estimator=NN, param grid=NN grid, cv=5,scoring='accuracy')

WARNING: tensorflow: Unresolved object in checkpoint: (root).optimizer.iter WARNING: tensorflow: Unresolved object in checkpoint: (root).optimizer.beta 1 WARNING:tensorflow:Unresolved object in checkpoint: (root).optimizer.beta\_2 WARNING: tensorflow: Unresolved object in checkpoint: (root).optimizer.decay

WARNING: tensorflow: Unresolved object in checkpoint: (root).optimizer.learning rate

print(tem.groupby(['Predict','True']).size().unstack())

fpr, tpr, \_ = metrics.roc\_curve(y\_test, prob,pos\_label=1)

#create the confucion matrix

Score: 0.9880287885665894

Score: 0.9828082680702209

Score: 0.9807830810546875

Score: 0.9723672270774841

Best accuracy So Far: 0.9880287885665894

result=tuner.get\_best\_hyperparameters()[0].values

The best model's accuracy on test data = 98%

Objective (name='accuracy', direction='max')

print("The best model's accuracy on test data = 98%") print('----')

7. Plot the ROC Curve & Confusion Matrix

The best 4-layers NN parameters would be 100 neurons and 0.01 learning rate.

INFO:tensorflow:Oracle triggered exit

NNmodel2=tuner.get\_best\_models()[0]

model.add(layers.Dense(units=hp.Int('units',

model.add(layers.Dense(units=hp.Int('units',

model.add(layers.Dense(units=hp.Int('units',

model.add(layers.Dense(1,activation='sigmoid'))

metrics=['accuracy'])

max trials=5, overwrite=True,

x\_train\_float = np.asarray(x\_train).astype(np.float32) y\_train\_float = np.asarray(y\_train).astype(np.float32) x\_test\_float = np.asarray(x\_test).astype(np.float32) y\_test\_float = np.asarray(y\_test).astype(np.float32)

seed=99,

loss='binary\_crossentropy',

objective='accuracy',

executions\_per\_trial=5)

min value=10, max\_value=100, step=10), activation='relu',input dim=57))

> min value=10, max value=100, step=10),

> min value=10, max value=100, step=10),

#Keras cannot input object data type, so no matter the column is boolean or numeric we need to transfor

tuner.search(x=x train float,y=y train float,epochs=30,batch size=512,validation data=(x test float,y t

print('The best 4-layers NN parameters would be {} neurons and {} learning rate.'.format(result['units'

values=[0.01,0.001,0.0001])),

activation='relu'))

activation='relu'))

model.compile(optimizer=keras.optimizers.Adam(hp.Choice('learning rate',

In [1]: import pandas as pd

%matplotlib inline

import numpy as np np.random.seed(13)

import tensorflow as tf import pandas as pd

from tensorflow import keras

from tensorflow.keras import layers

import matplotlib.pyplot as plt

from itertools import permutations

warnings.filterwarnings('ignore')

import scikitplot as skplt

import itertools

import warnings

1. Import Data

In [70]:

from kerastuner.tuners import RandomSearch

from matplotlib import pyplot as plt

from tensorflow.keras import backend as K from tensorflow.keras.models import Sequential from tensorflow.keras.datasets import mnist

from tensorflow.keras.utils import to categorical

from tensorflow.keras.layers import Dense, Activation, Dropout

from sklearn.metrics import recall score, confusion matrix

stuner` is deprecated, please use `import keras tuner`.

from sklearn.model\_selection import GridSearchCV

df=pd.read csv('spambase.data', header=None)

'word\_freq\_report','word\_freq\_addresses',

erage', 'capital run length longest',

df['spam'] = df['spam'].astype(str)

2. Train-Test Data Split

In [72]: from imblearn.over\_sampling import SMOTE

oversample = SMOTE()

In [73]: y train=y train.astype(int)

y\_test=y\_test.astype(int)

dom state=9)

req\_over','word\_freq\_remove','word freq internet',

'word\_freq\_your','word\_freq\_font','word freq 000',

eq\_lab','word\_freq\_labs','word\_freq\_telnet','word\_freq\_857',

'word\_freq\_table','word\_freq\_conference','char\_freq\_;',

'capital run length total','spam']

3. SMOTE to fix imbalance problem

x train, y train = oversample.fit resample(x train, y train)

d freq parts', 'word freq pm', 'word freq direct', 'word freq cs',

from tensorflow.keras.wrappers.scikit\_learn import KerasClassifier

from sklearn import neighbors, datasets, tree, linear model, metrics, svm from sklearn.model\_selection import cross val score, train test split, KFold

/opt/anaconda3/lib/python3.7/site-packages/ipykernel\_launcher.py:19: DeprecationWarning: `import kera

df.columns=['word\_freq\_make','word\_freq\_address','word\_freq\_all','word\_freq\_3d','word\_freq\_our','word\_f

In [71]: x\_train,x\_test,y\_train,y\_test = train\_test\_split(df.drop(columns=['spam']),df['spam'],test\_size=0.2,ran

'word freq order', 'word freq mail', 'word freq receive', 'word freq will', 'word freq people',

'word\_freq\_free','word\_freq\_business','word\_freq\_email','word\_freq\_you','word\_freq\_credit',

'word\_freq\_money','word\_freq\_hp','word\_freq\_hpl','word\_freq\_george','word\_freq\_650','word\_fr

'word\_freq\_data','word\_freq\_415','word\_freq\_85','word\_freq\_technology','word\_freq\_1999','wor

'word\_freq\_meeting','word\_freq\_original','word\_freq\_project','word\_freq\_re','word\_freq\_edu',

'char\_freq\_(','char\_freq\_!','char\_freq\_\$','char\_freq\_#','capital\_run\_length\_av

The best parameters activation=relu & nb hidden=100 Prediction Accuracy Score on Test Data: 94.46% Recall positive on Test Data: 93.8% 6-2. Use RandomSearch to find best hyper-parameters for 4-layers NN model In [78]: #Random Search Version def build model(hp):

NN grid = {'activation':['relu', 'sigmoid'], 'nb hidden':[10,20,30,40,50,60,70,80,90,100]} def create model(activation, nb hidden): model.add(Dense(nb\_hidden, input\_dim=57, activation=activation)) model.add(Dense(nb hidden, activation=activation)) model.add(Dense(nb hidden, activation=activation)) model.add(Dense(1, activation='sigmoid')) model.compile(loss='binary\_crossentropy', optimizer='adam', metrics=['accuracy']) NN = KerasClassifier(build fn=create model, epochs=30, batch size=256, verbose=0) #Grid Search CV for Neural Network NNmodel = GridSearchCV(estimator=NN, param grid=NN grid, cv=5,scoring='accuracy') NNmodel.fit(x\_train, y\_train) print('The best parameters activation={} & nb hidden={}'.format(NNmodel.best params ['activation'], NNmo del.best params ['nb hidden'])) print("Prediction Accuracy Score on Test Data: {}%".format(round(metrics.accuracy score(y test, NNmodel  $.predict(x_test))*100,2)))$ print("Recall positive on Test Data: {}%".format(round(metrics.recall score(y test, NNmodel.predict(x t est),pos label=1)\*100,2)))