ECE 457B Assignment 4

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```
import numpy as np
import pandas as pd
from sklearn.model_selection import train_test_split
from sklearn.preprocessing import OneHotEncoder, MinMaxScaler
from sklearn.svm import SVC
from sklearn import metrics
from sklearn.metrics import confusion_matrix, classification_report, ConfusionMatrixDisplay
from sklearn.preprocessing import MinMaxScaler
```

```
1.a Data Preprocessing
        # data from white wine and red wine is already merged into a single winequality datasheet
        df = pd.read_csv("winequality.csv")
        print(df.head())
        print(df.shape)
           fixed acidity volatile acidity citric acid residual sugar chlorides \
        a
                    7.4
                                     0.70
                                                  0.00
                                                                  1.9
                                                                           0.076
                    7.8
                                     0.88
                                                  0.00
                                                                  2.6
                                                                           0.098
        2
                    7.8
                                     0.76
                                                  0.04
                                                                  2.3
                                                                           0.092
        3
                    11.2
                                     0.28
                                                  0.56
                                                                  1.9
                                                                           0.075
                                     0.70
                                                  0.00
                                                                  1.9
                                                                           0.076
                    7.4
           free sulfur dioxide total sulfur dioxide density
                                                             pH sulphates \
        0
                         11.0
                                               34.0
                                                      0.9978 3.51
                                                                         0.56
        1
                          25.0
                                               67.0
                                                      0.9968 3.20
                                                                         0.68
        2
                         15.0
                                               54.0 0.9970 3.26
                                                                         0.65
        3
                          17.0
                                               60.0
                                                     0.9980 3.16
                                                                         0.58
        4
                         11.0
                                               34.0 0.9978 3.51
                                                                         0.56
           alcohol quality type
               9.4
        0
                         5
                         5
        1
               9.8
        2
               9.8
                        5
                         6
        3
               9.8
                               0
        4
               9.4
                         5
        (6497, 13)
In [ ]: # column 1 -> 11 = features, column 12 = quality, column 13 = wine category
        (x, y) = df.iloc[:, 0:10], df.iloc[:, 11]
        # normalization
        n_x = MinMaxScaler().fit_transform(x)
        # 80% for training, 20% for testing
        x_train, x_test, y_train, y_test = train_test_split(n_x, y, test_size=0.20)
        print('Training set shape: {}'.format(x_train.shape))
        print('Training labels shape: {}'.format(y_train.shape))
        print('Test set shape: {}'.format(x_test.shape))
        print('Test label shape: {}'.format(y_test.shape))
        Training set shape: (5197, 10)
```

Training set shape: (5197, 10)
Training labels shape: (5197,)
Test set shape: (1300, 10)
Test label shape: (1300,)

Support Vector Machine

```
In [ ]: def SVM(C, kernal, g="scale"):
            svm = SVC(C=C, kernel=kernal, gamma=g)
            svm.fit(x_train, y_train)
            y_pred = svm.predict(x_test)
            acc svm = metrics.accuracy score(y test, y pred)
            print('SVM Accuracy for {} at C = {} and gamma = {}:'.format(kernal, C, g), acc_svm)
            # cm_svm = confusion_matrix(y_test, y_pred)
            # class names = ['0', '1', '2', '3', '4', '5', '6']
            # print(cm svm)
            # print(classification_report(y_test, y_pred, target_names=class_names))
            # disp1 = ConfusionMatrixDisplay(confusion matrix=cm svm)
            # disp1.plot()
            return acc_svm
In []: C1 = [1, 10, 50, 100]
        C2 = [1, 10, 20, 30]
        result = np.empty((3, 4), dtype=float)
        # RBF Kernal
        for idc, C in enumerate(C1):
            acc_svm = SVM(C, 'rbf')
            result[0, idc] = acc svm
        # Polv Kernal
        for idc, C in enumerate(C1):
            acc_svm = SVM(C, 'poly')
            result[1, idc] = acc svm
        # Linear Kernal
        for idc, C in enumerate(C2):
            acc_svm = SVM(C, 'linear')
            result[2, idc] = acc_svm
        SVM Accuracy for rbf at C = 1 and gamma = scale: 0.546923076923077
        SVM Accuracy for rbf at C = 10 and gamma = scale: 0.5930769230769231
        SVM Accuracy for rbf at C = 50 and gamma = scale: 0.5838461538461538
        SVM Accuracy for rbf at C = 100 and gamma = scale: 0.5892307692307692
        SVM Accuracy for poly at C = 1 and gamma = scale: 0.5592307692307692
        SVM Accuracy for poly at C = 10 and gamma = scale: 0.5738461538461539
        SVM Accuracy for poly at C = 50 and gamma = scale: 0.573076923076923
        SVM Accuracy for poly at C = 100 and gamma = scale: 0.5746153846153846
        SVM Accuracy for linear at C = 1 and gamma = scale: 0.5069230769230769
        SVM Accuracy for linear at C = 10 and gamma = scale: 0.5269230769230769
        SVM Accuracy for linear at C = 20 and gamma = scale: 0.5238461538461539
        SVM Accuracy for linear at C = 30 and gamma = scale: 0.5261538461538462
```

Accuracy result in table, the column represents the different regularization parameters used and the rows represents the different kernal used

```
# print the dataframe
print(df)
```

```
C[0] C[1] C[2] C[3] RBF 0.546923 0.593077 0.583846 0.589231 Poly 0.559231 0.573846 0.573077 0.574615 Linear 0.506923 0.526923 0.523846 0.526154
```

1.b

Comparing the accuracy result from the table above, we can see that RBF kernal performed the best overall with an average accuracy of 57.826925% compared to 57.019225% for poly kernal and only 52.12115% for linear kernal. This make sense since guassian kernal is generally the perfered function in svm. It is suitable for non-linear data and helps to make proper separation when there is no prior knowledge of data. On the otherhand the linear kernal is the most basic kernal and is mostly preferred for text-classification and linear kernal is just a more generalized representation of the linear kernal. Furthermore, we can see that increasing the regularization parameter c sees improvement in training and test accuracy as well, since a higher value of the regularization parameter will penalize the model more for misclassifying training examples and lead to a smaller margin, while a lower value of the regularization parameter will allow more margin violations and lead to a larger margin.

1.c Improving the model

```
In [ ]: C = [1, 10, 50, 100]
        gamma = [10**-1, 10**0, 10*1, 10**2]
        result = np.empty((len(C), len(gamma)), dtype=float)
        for idc, c in enumerate(C):
            for idg, g in enumerate(gamma):
                acc_svm = SVM(c, 'rbf', g)
                result[idc, idg] = acc_svm
        SVM Accuracy for rbf at C = 1 and gamma = 0.1: 0.49538461538461537
        SVM Accuracy for rbf at C = 1 and gamma = 1: 0.5169230769230769
        SVM Accuracy for rbf at C = 1 and gamma = 10: 0.5776923076923077
        SVM Accuracy for rbf at C = 1 and gamma = 100: 0.6215384615384615
        SVM Accuracy for rbf at C = 10 and gamma = 0.1: 0.5246153846153846
        SVM Accuracy for rbf at C = 10 and gamma = 1: 0.5284615384615384
        SVM Accuracy for rbf at C = 10 and gamma = 10: 0.5884615384615385
        SVM Accuracy for rbf at C = 10 and gamma = 100: 0.6376923076923077
        SVM Accuracy for rbf at C = 50 and gamma = 0.1: 0.5223076923076924
        SVM Accuracy for rbf at C = 50 and gamma = 1: 0.5453846153846154
        SVM Accuracy for rbf at C = 50 and gamma = 10: 0.5953846153846154
        SVM Accuracy for rbf at C = 50 and gamma = 100: 0.64
        SVM Accuracy for rbf at C = 100 and gamma = 0.1: 0.5269230769230769
        SVM Accuracy for rbf at C = 100 and gamma = 1: 0.5592307692307692
        SVM Accuracy for rbf at C = 100 and gamma = 10: 0.6046153846153847
        SVM Accuracy for rbf at C = 100 and gamma = 100: 0.64
```

Accuracy result for rbf kernal in table, the column represents the different regularization parameters used and the rows represents the different gamma parameters used.

print the dataframe print(df)

	0.1	1.0	10.0	100.0
1	0.495385	0.516923	0.577692	0.621538
10	0.524615	0.528462	0.588462	0.637692
50	0.522308	0.545385	0.595385	0.640000
100	0.526923	0.559231	0.604615	0.640000

Here we see that not only does a larger regularization parameter kernal improves the model, but also does a larger gamma parameter. In general a high gamma value may lead to overfitting, where the model captures noise in the training data and performs poorly on new, unseen data. A low gamma value may lead to underfitting, where the model is too simple to capture the underlying patterns in the data and performs poorly on both the training and testing data. Here since the number of data in the wine-dataset is very limited and consist of a low number of feature, a higher gamma is more suitable in order to prevent underfitting. Here we reached the optimal accuracy of 0.64 using C = 100 and gamma = 100

```
In []: from __future__ import division
    from __future__ import print_function
    import pandas as pd
    import numpy as np
    import matplotlib.pyplot as plt
    from sklearn.model_selection import train_test_split
    import tensorflow.keras as keras
    from tensorflow.keras.models import Sequential
    from tensorflow.keras.layers import Dense, Conv2D, MaxPool2D, Flatten, Dropout
    from tensorflow.keras.datasets import cifar10
    from tensorflow.keras.utils import to_categorical
```

Data Preprocessing

```
In []: (train_images, train_labels), (test_images, test_labels) = cifar10.load_data()
    train_labels = to_categorical(train_labels, num_classes=10)
    test_labels = to_categorical(test_labels, num_classes=10)
    print('train images shape:', train_images.shape)
    print('train labels shape:', train_labels.shape)
    print('test images shape:', test_images.shape)
    print('test labels shape:', test_labels.shape)

train images shape: (50000, 32, 32, 3)
    train labels shape: (50000, 10)
    test images shape: (10000, 32, 32, 3)
    test labels shape: (10000, 10)
```

Data Normalization

```
In []: train_images = train_images.astype(np.float32)
    mean = np.mean(train_images)
    std = np.std(train_images)
    train_images = (train_images - mean) / std
    test_images = (test_images - mean) / std

In []: if len(train_labels.shape)==1:
        train_labels = keras.utils.to_categorical(train_labels, num_classes=10)
        test_labels = keras.utils.to_categorical(test_labels, num_classes=10)
    # print(train_labels.shape, test_labels.shape)
    # print(train_labels[0:3])
# print('Old labels: {}'.format(np.argmax(train_labels[:3], axis=1)))

result = np.empty((3, 4), dtype=float)
```

Multilayer Perceptron

```
In []: # define the model type (still empty)
    mlp = keras.models.Sequential()
    # add a Layer that just flattens the input (no weights here)
    mlp.add(Flatten(input_shape=(32, 32, 3)))
# add the first hidden Layer with 64 neurons, an activation function of sigmoid, and an input
    mlp.add(Dense(512, activation='sigmoid', input_shape=(3072,)))
# add the second hidden Layer with 64 neurons and sigmoid activation function
```

```
# add the output layer with 10 units and Softmax activation function
      mlp.add(Dense(10, activation='softmax'))
In [ ]: mlp.compile(loss='categorical crossentropy', optimizer='adam', metrics=['accuracy'])
      # Randomly sample 20% of the training set
      x train, x test, y train, y test = train test split(train images, train labels, train size=0.
      mlp h = mlp.fit(x train, y train, epochs=5, batch size=32, validation data=(test images, test
      result[0, 0] = mlp h.history['accuracy'][-1]
      result[0, 1] = mlp_h.history['val_accuracy'][-1]
      # evalute test accuracy
      mlp test = mlp.evaluate(test images, test labels, verbose=1)
      result[0, 2] = mlp test[0]
      result[0, 3] = mlp test[1]
      Epoch 1/5
      al_loss: 1.8159 - val_accuracy: 0.3667
      Epoch 2/5
      al_loss: 1.7204 - val_accuracy: 0.3932
      Epoch 3/5
      al loss: 1.6940 - val accuracy: 0.4004
      al_loss: 1.6626 - val_accuracy: 0.4166
      Epoch 5/5
      al_loss: 1.6722 - val_accuracy: 0.4098
      CNN1
In [ ]: cnn = Sequential()
      cnn.add(Conv2D(filters=64, kernel_size=(3, 3), padding='same', activation='relu', input_shape
      cnn.add(Conv2D(64, kernel_size=(3, 3), padding='same', activation='relu'))
      cnn.add(Flatten())
      cnn.add(Dense(512, activation='sigmoid'))
      cnn.add(Dense(512, activation='sigmoid'))
      cnn.add(Dense(10, activation='softmax'))
In [ ]: cnn.compile(loss='categorical crossentropy', optimizer='adam', metrics=['accuracy'])
      # Randomly sample 20% of the training set
      x_train, x_test, y_train, y_test = train_test_split(train_images, train_labels, train_size=0.
      cnn h = cnn.fit(x train.reshape(-1, 32, 32, 3),
            y train,
            batch size=32,
            epochs=5,
            validation_data=(test_images.reshape(-1, 32, 32, 3), test_labels),
            verbose=1)
      result[1, 0] = cnn_h.history['accuracy'][-1]
      result[1, 1] = cnn_h.history['val_accuracy'][-1]
      # evalute test accuracy
      cnn_test = cnn.evaluate(test_images.reshape(-1, 32, 32, 3), test_labels)
      result[1, 2] = cnn_test[0]
      result[1, 3] = cnn_test[1]
```

mlp.add(Dense(512, activation='sigmoid'))

```
Epoch 1/5
val_loss: 1.4280 - val_accuracy: 0.4888
Epoch 2/5
val_loss: 1.3226 - val_accuracy: 0.5263
Epoch 3/5
val_loss: 1.3553 - val_accuracy: 0.5463
Epoch 4/5
val loss: 1.4616 - val accuracy: 0.5778
Epoch 5/5
val_loss: 1.5138 - val_accuracy: 0.5924
```

CNN₂

```
In [ ]: cnn2 = Sequential()
        cnn2.add(Conv2D(filters=64, kernel size=(3, 3), padding='same', activation='relu'
        , input_shape=(32, 32, 3)))
        cnn2.add(MaxPool2D())
        cnn2.add(Conv2D(64, kernel size=(3, 3), padding='same', activation='relu'))
        cnn2.add(MaxPool2D())
        cnn2.add(Flatten())
        cnn2.add(Dense(512, activation='sigmoid'))
        cnn2.add(Dropout(0.2))
        cnn2.add(Dense(512, activation='sigmoid'))
        cnn2.add(Dropout(0.2))
        cnn2.add(Dense(10, activation='softmax'))
In [ ]: cnn2.compile(loss='categorical crossentropy', optimizer='adam', metrics=['accuracy'])
        # Randomly sample 20% of the training set
        x_train, x_test, y_train, y_test = train_test_split(train_images, train_labels, train_size=0.
        cnn2_h = cnn2.fit(x_train.reshape(-1, 32, 32, 3),
                y_train,
                batch size=32,
                epochs=5,
                validation data=(test images.reshape(-1, 32, 32, 3), test labels),
                verbose=1)
        result[2, 0] = cnn2 h.history['accuracy'][-1]
        result[2, 1] = cnn2_h.history['val_accuracy'][-1]
        # evalute test accuracy
        cnn2_test = cnn2.evaluate(test_images.reshape(-1, 32, 32, 3),test_labels)
        result[2, 2] = cnn2_test[0]
        result[2, 3] = cnn2_test[1]
```

```
Epoch 1/5
val_loss: 1.4044 - val_accuracy: 0.4844
Epoch 2/5
val_loss: 1.2677 - val_accuracy: 0.5428
Epoch 3/5
val_loss: 1.1537 - val_accuracy: 0.5882
Epoch 4/5
val loss: 1.1154 - val accuracy: 0.5986
Epoch 5/5
val_loss: 1.1196 - val_accuracy: 0.6126
```

2.a

From the table below, we can see that in general MLP performs worse than CNN both in terms of the training accuracy and test accuracy. This makes sense because CNNs are primarily used for image recognition tasks by taking advantage of the spatial structure of an image, learning features such as edges, lines, and textures. On the other hand, MLPs are more general purposed for classification, regression, and other tasks. However, an advantage of MLP is the simple architecture of MLP compared to CNN, we can see that it trains significantly faster than CNN in general.

0.6126 1.119578

0.6126

2.b

CNN2

0.7481

As we can see from the figure below, the training time for the first CNN model is significantlylonger than then the second one. This is because the 2 additional max pooling layer reduces the size of the images by half. Between the two CNN we can see that although the first CNN shows signficantly higher training accuracy, it shows slightly lower validation accuracy and subsequently lower test accuracy, a clear sign of overfitting. This is because it lacks the dropout layer from the second CNN model which provides regularization to reduce a overfitting.

```
In []: # plot accuracy for the first CNN model
    plt.plot(cnn_h.history['accuracy'], label='cnn1 accuracy')
    plt.plot(cnn_h.history['val_accuracy'], label='cnn1 val_accuracy')

# plot accuracy for the second CNN model
    plt.plot(cnn2_h.history['accuracy'], label='cnn2 accuracy')
    plt.plot(cnn2_h.history['val_accuracy'], label='cnn2 val_accuracy')

plt.legend(loc='best')
    plt.show()
```

