This is the k-nearest neighbors workbook for ECE C147/C247 Assignment #2

Please follow the notebook linearly to implement k-nearest neighbors.

Please print out the workbook entirely when completed.

The goal of this workbook is to give you experience with the data, training and evaluating a simple classifier, k-fold cross validation, and as a Python refresher.

Import the appropriate libraries

```
In [69]: import numpy as np # for doing most of our calculations
import matplotlib.pyplot as plt# for plotting
from utils.data_utils import load_CIFAR10 # function to load the CIFAR-10 dataset.

# Load matplotlib images inline
%matplotlib inline

# These are important for reloading any code you write in external .py files.
# see http://stackoverflow.com/questions/1907993/autoreload-of-modules-in-ipython
%load_ext autoreload
%autoreload 2
```

The autoreload extension is already loaded. To reload it, use: %reload_ext autoreload

```
In [70]: # Set the path to the CIFAR-10 data
    cifar10_dir = './cifar-10-batches-py' # You need to update this Line
    X_train, y_train, X_test, y_test = load_CIFAR10(cifar10_dir)

# As a sanity check, we print out the size of the training and test data.
    print('Training data shape: ', X_train.shape)
    print('Training labels shape: ', y_train.shape)
    print('Test data shape: ', X_test.shape)
    print('Test labels shape: ', y_test.shape)
```

Training data shape: (50000, 32, 32, 3)
Training labels shape: (50000,)
Test data shape: (10000, 32, 32, 3)
Test labels shape: (10000,)



```
In [72]: # Subsample the data for more efficient code execution in this exercise
    num_training = 5000
    mask = list(range(num_training))
    X_train = X_train[mask]
    y_train = y_train[mask]

    num_test = 500
    mask = list(range(num_test))
    X_test = X_test[mask]
    y_test = y_test[mask]

# Reshape the image data into rows
    X_train = np.reshape(X_train, (X_train.shape[0], -1))
    X_test = np.reshape(X_test, (X_test.shape[0], -1))
    print(X_train.shape, X_test.shape)

(5000, 3072) (500, 3072)
```

K-nearest neighbors

In the following cells, you will build a KNN classifier and choose hyperparameters via k-fold cross-validation.

```
In [73]: # Import the KNN class
from nndl import KNN

In [74]: # Declare an instance of the knn class.
knn = KNN()

# Train the classifier.
# We have implemented the training of the KNN classifier.
# Look at the train function in the KNN class to see what this does.
knn.train(X=X_train, y=y_train)
```

Questions

- (1) Describe what is going on in the function knn.train().
- (2) What are the pros and cons of this training step?

Answers

- (1) The funciton knn.train() simply remembers the data points (images) and their labels.
- (2) The pros are that there is no training time and it is simple and fast; the cons are that it is memory intensive as we must store all the input data. Additionally, predicting the test data is expensive since we must get the distance to all points.

KNN prediction

In the following sections, you will implement the functions to calculate the distances of test points to training points, and from this information, predict the class of the KNN.

```
In [75]: # Implement the function compute_distances() in the KNN class.
# Do not worry about the input 'norm' for now; use the default definition of the norm
# in the code, which is the 2-norm.
# You should only have to fill out the clearly marked sections.

import time
time_start =time.time()

dists_L2 = knn.compute_distances(X=X_test)

print('Time to run code: {}'.format(time.time()-time_start))
print('Frobenius norm of L2 distances: {}'.format(np.linalg.norm(dists_L2, 'fro')))
```

Time to run code: 19.62835693359375 Frobenius norm of L2 distances: 7906696.077040902

Really slow code

Note: This probably took a while. This is because we use two for loops. We could increase the speed via vectorization, removing the for loops.

If you implemented this correctly, evaluating np.linalg.norm(dists_L2, 'fro') should return: ~7906696

KNN vectorization

The above code took far too long to run. If we wanted to optimize hyperparameters, it would be time-expensive. Thus, we will speed up the code by vectorizing it, removing the for loops.

```
In [76]: # Implement the function compute_L2_distances_vectorized() in the KNN class.
# In this function, you ought to achieve the same L2 distance but WITHOUT any for Loops.
# Note, this is SPECIFIC for the L2 norm.

time_start =time.time()
dists_L2_vectorized = knn.compute_L2_distances_vectorized(X=X_test)
print('Time to run code: {}'.format(time.time()-time_start))
print('Difference in L2 distances between your KNN implementations (should be 0): {}'.format(np.linalg.norm(dists_L2 - dists_L2_vectorized))
```

Time to run code: 0.11509299278259277 Difference in L2 distances between your KNN implementations (should be 0): 0.0

Speedup

Depending on your computer speed, you should see a 10-100x speed up from vectorization. On our computer, the vectorized form took 0.36 seconds while the naive implementation took 38.3 seconds.

Implementing the prediction

Now that we have functions to calculate the distances from a test point to given training points, we now implement the function that will predict the test point labels.

0.726

If you implemented this correctly, the error should be: 0.726.

This means that the k-nearest neighbors classifier is right 27.4% of the time, which is not great, considering that chance levels are 10%.

Optimizing KNN hyperparameters

In this section, we'll take the KNN classifier that you have constructed and perform cross-validation to choose a best value of k, as well as a best choice of norm.

Create training and validation folds

First, we will create the training and validation folds for use in k-fold cross validation.

```
In [80]: # Create the dataset folds for cross-valdiation.
        num_folds = 5
        X_train_folds = []
        y_train_folds = []
        np.random.seed(123)
        idx = np.random.permutation(num_training)
        print(idx, idx.shape)
            Split the training data into num_folds (i.e., 5) folds.
            X_train_folds is a list, where X_train_folds[i] contains the
               data points in fold i.
           y_train_folds is also a list, where y_train_folds[i] contains
              the corresponding labels for the data in X_train_folds[i]
        X_train_shuffle = X_train[idx]
        y_train_shuffle = y_train[idx]
        X_train_folds = np.array_split(X_train_shuffle, num_folds)
        y_train_folds = np.array_split(y_train_shuffle, num_folds)
        # END YOUR CODE HERE
        print(y_train_folds[0].shape)
        [2648 2456 4557 ... 1346 3454 3582] (5000,)
```

Optimizing the number of nearest neighbors hyperparameter.

In this section, we select different numbers of nearest neighbors and assess which one has the lowest k-fold cross validation error.

(1000,)

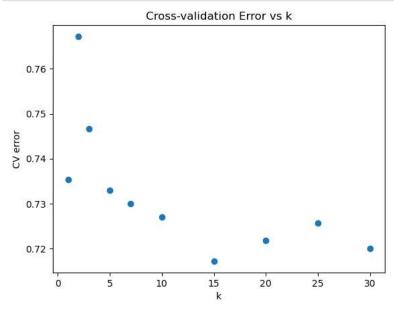
```
In [81]: time_start =time.time()
       ks = [1, 2, 3, 5, 7, 10, 15, 20, 25, 30]
        # ============ #
       # YOUR CODE HERE:
          Calculate the cross-validation error for each k in ks, testing
          the trained model on each of the 5 folds. Average these errors
          together and make a plot of k vs. cross-validation error. Since
          we are assuming L2 distance here, please use the vectorized code!
       # Otherwise, you might be waiting a long time.
       error_list = []
        for k in ks:
           error = 0
           for i in np.arange(num_folds):
              X_fold_train = np.concatenate(X_train_folds[:i] + X_train_folds[i + 1:], axis=0)
              y_fold_train = np.concatenate(y_train_folds[:i] + y_train_folds[i + 1:], axis=0)
              X_fold_val = X_train_folds[i]
              y_fold_val = y_train_folds[i]
              knn.train(X_fold_train, y_fold_train)
              cur_pred = knn.predict_labels(knn.compute_L2_distances_vectorized(X_fold_val), k)
              error += (1 - np.mean(np.equal(y_fold_val, cur_pred)))
           error_list.append(error / num_folds)
        # END YOUR CODE HERE
        print('Computation time: %.2f'%(time.time()-time_start))
```

Computation time: 19.94

```
In [82]: plt.scatter(ks, error_list)
plt.title('Cross-validation Error vs k')
plt.ylabel('CV error')
plt.xlabel('k')
plt.show()

for i in range(len(error_list)):
    print(str(ks[i]) + ': ' + str(error_list[i]))

print()
best_error = min(error_list)
k_idx = error_list.index(best_error)
best_k = ks[k_idx]
print("Best k, error: " + str(best_k) + ', ' + str(best_error))
```



```
1: 0.7354

2: 0.7672

3: 0.7465999999999999

5: 0.733

7: 0.73000000000000001

10: 0.727

15: 0.7172

20: 0.7218

25: 0.725599999999999

30: 0.72

Best k, error: 15, 0.7172
```

Questions:

- (1) What value of k is best amongst the tested k's?
- (2) What is the cross-validation error for this value of k?

Answers:

- (1) k=15
- (2) 0.7172

Optimizing the norm

Next, we test three different norms (the 1, 2, and infinity norms) and see which distance metric results in the best cross-validation performance.

```
In [83]: time_start =time.time()
        L1_norm = lambda x: np.linalg.norm(x, ord=1)
        L2_norm = lambda x: np.linalg.norm(x, ord=2)
        Linf_norm = lambda x: np.linalg.norm(x, ord= np.inf)
        norms = [L1_norm, L2_norm, Linf_norm]
        # YOUR CODE HERE:
          Calculate the cross-validation error for each norm in norms, testing
          the trained model on each of the 5 folds. Average these errors
           together and make a plot of the norm used vs the cross-validation error
           Use the best cross-validation k from the previous part.
           Feel free to use the compute_distances function. We're testing just
           three norms, but be advised that this could still take some time.
           You're welcome to write a vectorized form of the L1- and Linf- norms
           to speed this up, but it is not necessary.
        # ------ #
        norm_error_list = []
        for norm in norms:
           norm\_error = 0
           for i in np.arange(num_folds):
              X fold train = np.concatenate(X train folds[:i] + X train folds[i + 1:], axis=0)
              y_fold_train = np.concatenate(y_train_folds[:i] + y_train_folds[i + 1:], axis=0)
              X_fold_val = X_train_folds[i]
              y_fold_val = y_train_folds[i]
               knn.train(X_fold_train, y_fold_train)
              cur_pred = knn.predict_labels(knn.compute_distances(X_fold_val, norm), best_k)
              norm_error += (1 - np.mean(np.equal(y_fold_val, cur_pred)))
           norm_error_list.append(norm_error / num_folds)
        # ============ #
        # END YOUR CODE HERE
        # ------ #
        print('Computation time: %.2f'%(time.time()-time start))
```

Computation time: 362.95

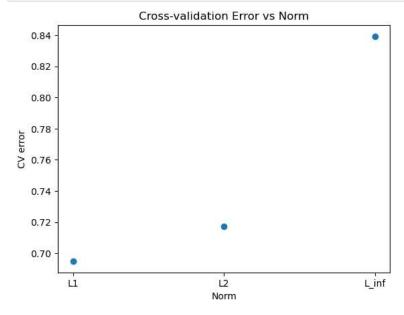
```
In [84]: ns = ["L1", "L2", "L_inf"]

plt.scatter(["L1", "L2", "L_inf"], norm_error_list)
plt.title('Cross-validation Error vs Norm')
plt.ylabel('CV error')
plt.xlabel('Norm')
plt.show()

for i in range(len(norm_error_list)):
    print(ns[i] + ' error: ' + str(norm_error_list[i]))

print()

print()
best_error = min(norm_error_list)
norm_idx = norm_error_list.index(best_error)
print("Best_norm, error: " + ns[norm_idx] + ', ' + str(best_error))
```



L1 error: 0.695 L2 error: 0.7172 L_inf error: 0.8392 Best norm, error: L1, 0.695

Questions:

- (1) What norm has the best cross-validation error?
- (2) What is the cross-validation error for your given norm and k?

Answers:

- (1) L1 norm
- (2) 0.695

Evaluating the model on the testing dataset.

Now, given the optimal \boldsymbol{k} and norm you found in earlier parts, evaluate the testing error of the k-nearest neighbors model.

Error rate achieved: 0.718

Question:

How much did your error improve by cross-validation over naively choosing k=1 and using the L2-norm?

Answer:

Error with L2-norm and k=1: 0.726

Error with L1-norm and k=15: 0.718

Improvement: 0.008