## k-Nearest Neighbor (kNN) exercise

Complete and hand in this completed worksheet (including its outputs and any supporting code outside of the worksheet) with your assignment submission. For more details see the <u>assignments page (http://vision.stanford.edu/teaching/cs231n/assignments.html)</u> on the course website.

The kNN classifier consists of two stages:

- During training, the classifier takes the training data and simply remembers it
- During testing, kNN classifies every test image by comparing to all training images and transfering the labels of the k
  most similar training examples
- The value of k is cross-validated

In this exercise you will implement these steps and understand the basic Image Classification pipeline, cross-validation, and gain proficiency in writing efficient, vectorized code.

```
In [1]: # Run some setup code for this notebook.
        import random
        import numpy as np
        from cs231n.data utils import load CIFAR10
        import matplotlib.pyplot as plt
        from __future__ import print_function
        # This is a bit of magic to make matplotlib figures appear inline in the no
        tebook
        # rather than in a new window.
        %matplotlib inline
        plt.rcParams['figure.figsize'] = (10.0, 8.0) # set default size of plots
        plt.rcParams['image.interpolation'] = 'nearest'
        plt.rcParams['image.cmap'] = 'gray'
        # Some more magic so that the notebook will reload external python modules;
        # see http://stackoverflow.com/questions/1907993/autoreload-of-modules-in-i
        python
        %load_ext autoreload
        %autoreload 2
In [2]: # Load the raw CIFAR-10 data.
        cifar10 dir = 'cs231n/datasets/cifar-10-batches-py'
        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.
```

```
cifar10_dir = 'cs231n/datasets/cifar-10-batches-py'
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,)

Test labels shape: (10000,)
```

```
In [3]: # Visualize some examples from the dataset.
        # We show a few examples of training images from each class.
        classes = ['plane', 'car', 'bird', 'cat', 'deer', 'dog', 'frog', 'horse', '
        ship', 'truck']
        num_classes = len(classes)
        samples_per_class = 7
        for y, cls in enumerate(classes):
            idxs = np.flatnonzero(y_train == y)
            idxs = np.random.choice(idxs, samples_per_class, replace=False)
            for i, idx in enumerate(idxs):
                plt idx = i * num classes + y + 1
                plt.subplot(samples_per_class, num_classes, plt_idx)
                plt.imshow(X_train[idx].astype('uint8'))
                plt.axis('off')
                if i == 0:
                    plt.title(cls)
        plt.show()
```



```
In [4]: # 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]
```

```
In [5]: # 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)

In [6]: from cs231n.classifiers import KNearestNeighbor

# Create a kNN classifier instance.
# Remember that training a kNN classifier is a noop:
# the Classifier simply remembers the data and does no further processing classifier = KNearestNeighbor()
    classifier.train(X_train, y_train)
```

We would now like to classify the test data with the kNN classifier. Recall that we can break down this process into two steps:

- 1. First we must compute the distances between all test examples and all train examples.
- 2. Given these distances, for each test example we find the k nearest examples and have them vote for the label

Lets begin with computing the distance matrix between all training and test examples. For example, if there are **Ntr** training examples and **Nte** test examples, this stage should result in a **Nte** x **Ntr** matrix where each element (i,j) is the distance between the i-th test and j-th train example.

First, open cs231n/classifiers/k\_nearest\_neighbor.py and implement the function compute\_distances\_two\_loops that uses a (very inefficient) double loop over all pairs of (test, train) examples and computes the distance matrix one element at a time.

```
In [7]: # Open cs231n/classifiers/k nearest neighbor.py and implement
        # compute_distances_two_loops.
        # Test your implementation:
        dists = classifier.compute_distances_two_loops(X_test)
        print(dists.shape)
        (500, 5000)
In [8]: # We can visualize the distance matrix: each row is a single test example a
        # its distances to training examples
        plt.imshow(dists, interpolation='none')
        plt.show()
           0
         250
         500
                        1000
                                      2000
                                                   3000
                                                                4000
                                                                              5000
```

**Inline Question #1:** Notice the structured patterns in the distance matrix, where some rows or columns are visible brighter. (Note that with the default color scheme black indicates low distances while white indicates high distances.)

- What in the data is the cause behind the distinctly bright rows?
- · What causes the columns?

Your Answer: fill this in.

```
In [16]: # Now implement the function predict_labels and run the code below:
    # We use k = 1 (which is Nearest Neighbor).
    y_test_pred = classifier.predict_labels(dists, k=1)

# Compute and print the fraction of correctly predicted examples
    num_correct = np.sum(y_test_pred == y_test)
    accuracy = float(num_correct) / num_test
    print('Got %d / %d correct => accuracy: %f' % (num_correct, num_test, accuracy))

Got 137 / 500 correct => accuracy: 0.274000
```

You should expect to see approximately 27% accuracy. Now lets try out a larger k, say k = 5:

```
In [10]: y_test_pred = classifier.predict_labels(dists, k=5)
num_correct = np.sum(y_test_pred == y_test)
accuracy = float(num_correct) / num_test
print('Got %d / %d correct => accuracy: %f' % (num_correct, num_test, accuracy))
Got 139 / 500 correct => accuracy: 0.278000
```

You should expect to see a slightly better performance than with k = 1.

```
In [11]: # Now lets speed up distance matrix computation by using partial vectorizat
         # with one loop. Implement the function compute_distances_one_loop and run
         the
         # code below:
         dists one = classifier.compute distances one loop(X test)
         # To ensure that our vectorized implementation is correct, we make sure tha
         t it
         # agrees with the naive implementation. There are many ways to decide wheth
         er
         # two matrices are similar; one of the simplest is the Frobenius norm. In c
         ase
         # you haven't seen it before, the Frobenius norm of two matrices is the squ
         are
         # root of the squared sum of differences of all elements; in other words, r
         # the matrices into vectors and compute the Euclidean distance between them
         difference = np.linalg.norm(dists - dists one, ord='fro')
         print('Difference was: %f' % (difference, ))
         if difference < 0.001:</pre>
             print('Good! The distance matrices are the same')
         else:
             print('Uh-oh! The distance matrices are different')
```

Difference was: 0.000000 Good! The distance matrices are the same

```
In [12]: # Now implement the fully vectorized version inside compute_distances_no_lo
         # and run the code
         dists_two = classifier.compute_distances_no_loops(X_test)
         # check that the distance matrix agrees with the one we computed before:
         difference = np.linalg.norm(dists - dists_two, ord='fro')
         print('Difference was: %f' % (difference, ))
         if difference < 0.001:
             print('Good! The distance matrices are the same')
             print('Uh-oh! The distance matrices are different')
         Difference was: 0.000000
         Good! The distance matrices are the same
In [13]: # Let's compare how fast the implementations are
         def time_function(f, *args):
             Call a function f with args and return the time (in seconds) that it to
         ok to execute.
             import time
             tic = time.time()
             f(*args)
             toc = time.time()
             return toc - tic
         two_loop_time = time_function(classifier.compute_distances_two_loops, X_tes
         print('Two loop version took %f seconds' % two_loop_time)
         one_loop_time = time_function(classifier.compute_distances_one_loop, X_test
         print('One loop version took %f seconds' % one_loop_time)
         no_loop_time = time_function(classifier.compute_distances_no_loops, X_test)
         print('No loop version took %f seconds' % no_loop_time)
         # you should see significantly faster performance with the fully vectorized
         implementation
         Two loop version took 16.554962 seconds
```

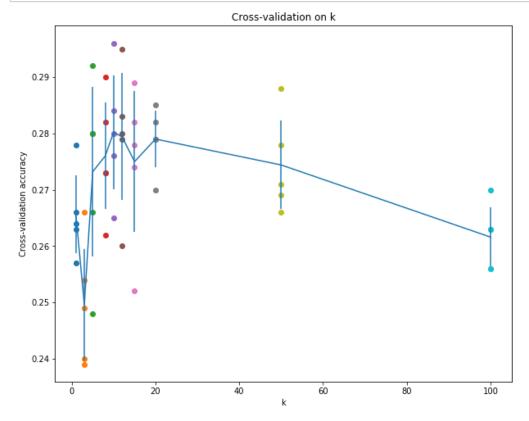
One loop version took 27.078910 seconds No loop version took 0.198320 seconds

## **Cross-validation**

We have implemented the k-Nearest Neighbor classifier but we set the value k = 5 arbitrarily. We will now determine the best value of this hyperparameter with cross-validation.

```
In [27]:
       num folds = 5
       k_{choices} = [1, 3, 5, 8, 10, 12, 15, 20, 50, 100]
       X_{train_folds} = []
       y_train_folds = []
       #####
       # TOD0:
       # Split up the training data into folds. After splitting, X_train_folds and
       # y_train_folds should each be lists of length num_folds, where
       # y train folds[i] is the label vector for the points in X train folds[i].
       # Hint: Look up the numpy array split function.
       #####
       X_train_folds = np.array_split(X_train,num_folds)
       y_train_folds = np.array_split(y_train,num_folds)
       #####
       #
                                  END OF YOUR CODE
       #####
       # A dictionary holding the accuracies for different values of k that we fin
       # when running cross-validation. After running cross-validation,
       # k_to_accuracies[k] should be a list of length num_folds giving the differ
       ent
       # accuracy values that we found when using that value of k.
       k to accuracies = {}
       #####
       # TOD0:
       # Perform k-fold cross validation to find the best value of k. For each
       # possible value of k, run the k-nearest-neighbor algorithm num_folds times
       # where in each case you use all but one of the folds as training data and
       the #
       # last fold as a validation set. Store the accuracies for all fold and all
       # values of k in the k to accuracies dictionary.
       #####
       for k val in k choices:
         k to accuracies[k val]=[]
         for i in range(num folds):
          # print 'Cross-validation :'+ str(i)
          X train cycle = np.concatenate([f for j,f in enumerate(X train folds) i
       f j!=i ])
          y train cycle = np.concatenate([f for j,f in enumerate(y train folds) i
       f j!=i ])
          X_val_cycle = X_train_folds[i]
          v val cvcle = v train folds[i]
```

```
k = 1, accuracy = 0.263000
k = 1, accuracy = 0.257000
k = 1, accuracy = 0.264000
k = 1, accuracy = 0.278000
k = 1, accuracy = 0.266000
k = 3, accuracy = 0.239000
k = 3, accuracy = 0.249000
k = 3, accuracy = 0.240000
k = 3, accuracy = 0.266000
k = 3, accuracy = 0.254000
k = 5, accuracy = 0.248000
k = 5, accuracy = 0.266000
k = 5, accuracy = 0.280000
k = 5, accuracy = 0.292000
k = 5, accuracy = 0.280000
k = 8, accuracy = 0.262000
k = 8, accuracy = 0.282000
k = 8, accuracy = 0.273000
k = 8, accuracy = 0.290000
k = 8, accuracy = 0.273000
k = 10, accuracy = 0.265000
k = 10, accuracy = 0.296000
k = 10, accuracy = 0.276000
k = 10, accuracy = 0.284000
k = 10, accuracy = 0.280000
k = 12, accuracy = 0.260000
k = 12, accuracy = 0.295000
k = 12, accuracy = 0.279000
k = 12, accuracy = 0.283000
k = 12, accuracy = 0.280000
k = 15, accuracy = 0.252000
k = 15, accuracy = 0.289000
k = 15, accuracy = 0.278000
k = 15, accuracy = 0.282000
k = 15, accuracy = 0.274000
k = 20, accuracy = 0.270000
k = 20, accuracy = 0.279000
k = 20, accuracy = 0.279000
k = 20, accuracy = 0.282000
k = 20, accuracy = 0.285000
k = 50, accuracy = 0.271000
k = 50, accuracy = 0.288000
k = 50, accuracy = 0.278000
k = 50, accuracy = 0.269000
k = 50, accuracy = 0.266000
k = 100, accuracy = 0.256000
k = 100, accuracy = 0.270000
k = 100, accuracy = 0.263000
k = 100, accuracy = 0.256000
k = 100, accuracy = 0.263000
```



```
In [26]: # Based on the cross-validation results above, choose the best value for k,
# retrain the classifier using all the training data, and test it on the te
st
# data. You should be able to get above 28% accuracy on the test data.
best_k = 10

classifier = KNearestNeighbor()
classifier.train(X_train, y_train)
y_test_pred = classifier.predict(X_test, k=best_k)

# Compute and display the accuracy
num_correct = np.sum(y_test_pred == y_test)
accuracy = float(num_correct) / num_test
print('Got %d / %d correct => accuracy: %f' % (num_correct, num_test, accuracy))
```

Got 141 / 500 correct => accuracy: 0.282000

In [ ]: