knn

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1 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 assignments page 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
        # This is a bit of magic to make matplotlib figures appear inline in the notebook
        # 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-ipython
        %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.
        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 [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()
         plane
                 car
                         bird
                                 cat
                                        deer
                                                 dog
                                                        frog
                                                                horse
                                                                        ship
                                                                                truck
```

```
In [4]: # Subsample the data for more efficient code execution in this exercise
    num_training = 5000
    mask = range(num_training)
    X_train = X_train[mask]
    y_train = y_train[mask]

    num_test = 500
    mask = 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)
```

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 [36]: 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)
In [37]: # 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 [38]: # We can visualize the distance matrix: each row is a single test example and
         # its distances to training examples
         plt.imshow(dists, interpolation='none')
         plt.show()
     100
     200
     300
     400
        0
                      1000
                                     2000
                                                     3000
                                                                    4000
```

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 [39]: # Now implement the function predict_labels and run the code below:
         # We use k = 1 (which is Nearest Neighbor).
         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
         y_test_pred = classifier.predict_labels(dists, k=1)
         print y_test_pred.shape
         print y_test.shape
         # 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)
Training data shape: (5000, 3072)
Training labels shape: (5000,)
Test data shape: (500, 3072)
Test labels shape: (500,)
(500,)
(500.)
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 [40]: 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 [41]: # Now lets speed up distance matrix computation by using partial vectorization
         # 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 that it
         # agrees with the naive implementation. There are many ways to decide whether
         # two matrices are similar; one of the simplest is the Frobenius norm. In case
         # you haven't seen it before, the Frobenius norm of two matrices is the square
         # root of the squared sum of differences of all elements; in other words, reshape
```

```
# 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:
           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 [42]: # Now implement the fully vectorized version inside compute_distances_no_loops
         # 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: 44278155389.041199
Uh-oh! The distance matrices are different
In [43]: # 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 took 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_test)
         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 23.161831 seconds
```

One loop version took 24.999564 seconds No loop version took 0.153357 seconds

1.0.1 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 [57]: num_folds = 5
       k_{\text{choices}} = [1, 3, 5, 8, 10, 12, 15, 20, 50, 100]
       X_train_folds = []
       y_train_folds = []
       # 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(np.array_split(X_train, num_folds))
       print(X_train_folds.shape)
       y_train_folds = np.array(np.array_split(y_train, num_folds))
       END OF YOUR CODE 1
       # A dictionary holding the accuracies for different values of k that we find
       # when running cross-validation. After running cross-validation,
       \# k_{to} = ccuracies[k] should be a list of length num_folds giving the different
       \# accuracy values that we found when using that value of k.
       k to accuracies = {}
       \# Perform k-fold cross validation t50 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 in k_choices:
          for n in xrange(num_folds):
             combinat = [x for x in xrange(num_folds) if x != n] # n'th fold is used for testing
             # print(combinat)
             x_training_dat = np.concatenate(X_train_folds[combinat]) # (4000, 3072)
             y_training_dat = np.concatenate(y_train_folds[combinat]) # (4000, )
             # print(x_training_dat.shape)
             # print(y_training_dat.shape)
             classifier_k = KNearestNeighbor()
             classifier_k.train(x_training_dat, y_training_dat)
             y_cross_validation_pred = classifier_k.predict(X_train_folds[n], k) # pass n'th fold
             num_of_accuracy_test_per_iteration = y_cross_validation_pred.shape[0]
             # print(num_of_accuracy_test_per_iteration)
             num_correct = np.sum(y_cross_validation_pred == y_train_folds[n]) #
             # print(num_correct)
             # print(num_test)
             accuracy = float(num_correct) / num_of_accuracy_test_per_iteration
             k_to_accuracies.setdefault(k, []).append(accuracy)
```

```
END OF YOUR CODE
         # Print out the computed accuracies
        for k in sorted(k_to_accuracies):
            for accuracy in k_to_accuracies[k]:
                print 'k = \%d, accuracy = \%f' \% (k, accuracy)
(5, 1000, 3072)
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 [59]: # plot the raw observations
         for k in k_choices:
           accuracies = k_to_accuracies[k]
           plt.scatter([k] * len(accuracies), accuracies)
         # plot the trend line with error bars that correspond to standard deviation
         accuracies_mean = np.array([np.mean(v) for k,v in sorted(k_to_accuracies.items())])
         accuracies_std = np.array([np.std(v) for k,v in sorted(k_to_accuracies.items())])
         plt.errorbar(k_choices, accuracies_mean, yerr=accuracies_std)
         plt.title('Cross-validation on k')
         plt.xlabel('k')
         plt.ylabel('Cross-validation accuracy')
         plt.show()
                                         Cross-validation on k
        0.30
        0.29
        0.28
     Cross-validation accuracy
        0.27
        0.26
        0.25
        0.24
       0.23 L
-20
```

In [61]: # 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 test

0

20

60

100

120

80

```
# 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 []:
```