```
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 [28]: # Run some setup code for this notebook.
          from __future__ import print function
          import random
          import numpy as np
          from cs682.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
          The autoreload extension is already loaded. To reload it, use:
            %reload_ext autoreload
In [29]: # Load the raw CIFAR-10 data.
          cifar10 dir = 'cs682/datasets/cifar-10-batches-py'
          # Cleaning up variables to prevent loading data multiple times (which may cause memory issue)
          try:
             del X_train, y_train
             del X test, y test
             print('Clear previously loaded data.')
          except:
            pass
          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)
          Clear previously loaded data.
          Training data shape: (50000, 32, 32, 3)
          Training labels shape: (50000,)
          Test data shape: (10000, 32, 32, 3)
          Test labels shape: (10000,)
In [30]: # 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()
                                                      frog
                                                            horse
                                 cat
          In [31]: # 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_{\text{test}} = X_{\text{test}}[\text{mask}]
          y_test = y_test[mask]
In [32]: # 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 [33]: from cs682.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 cs682/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 [34]: # Open cs682/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 [11]: # 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()
           250
                          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:

    distinctly bright rows are caused by a particular test image being similar to most training images.

    distinctly bright columns are caused by a particular training image being similar to many test images.

In [35]: # 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 [36]: 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.
          Inline Question 2 We can also other distance metrics such as L1 distance. The performance of a Nearest Neighbor classifier that uses L1 distance will not
          change if (Select all that apply.):
          1. The data is preprocessed by subtracting the mean.
           2. The data is preprocessed by subtracting the mean and dividing by the standard deviation.
           3. The coordinate axes for the data are rotated.
           4. None of the above. (Mean and standard deviation in (1) and (2) are vectors and can be different across dimensions)
          Your Answer: A KNN Classifier that uses L1 norm's performance will not change if (3) the axes are rotated.
          Your explanation: This is because for L1 norm the absolute value of the values for each axis are just added, so it shouldn't matter which axis is which.
In [22]: # 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:</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 [46]: # 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:</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 [48]: # 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 21.526645 seconds
          One loop version took 33.403456 seconds
          No loop version took 0.247977 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 [47]: num folds = 5
          k 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))
          y_train_folds = np.array(np.array_split(y_train, num_folds))
          END OF YOUR CODE
          # A dictionary holding the accuracies for different values of k that we find
          # when running cross-validation. After running cross-validation,
          # k to accuracies[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 = {}
          # TODO:
          # 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 fold in range(num folds):
              x_te = X_train_folds[fold]
             y_te = y_train_folds[fold]
             x_tr = np.concatenate(np.array([x for i, x in enumerate(X_train_folds) if i != fold]), axis=0)
              y_tr = np.concatenate(np.array([y for i, y in enumerate(y_train_folds) if i != fold]), axis=0)
              classifier = KNearestNeighbor()
              classifier.train(x_tr, y_tr)
              for k_choice in k_choices:
                  y_test_pred = classifier.predict(x_te, k=k_choice)
                  num_correct = np.sum(y_test_pred == y_te)
                  accuracy = float(num_correct) / y_te.shape[0]
                  if k_choice in k_to_accuracies.keys():
                      k_to_accuracies[k_choice].append(accuracy)
                  else:
                      k_to_accuracies[k_choice] = [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))
          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]: # 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
            0.29
            0.28
          Cross-validation accuracy
            0.27
            0.26
            0.25
            0.24
                              20
                                                                                100
                                                       60
                                                                    80
In [27]: \# 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
          # 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
          Inline Question 3 Which of the following statements about k-Nearest Neighbor (k-NN) are true in a classification setting, and for all k? Select all that apply.
          1. The training error of a 1-NN will always be better than or equal to that of 5-NN.
          True for training error on 1-NN it will always find the distance to itself == 0. With a higher k, other different class neighbors around it could increase error.
          1. The test error of a 1-NN will always be better than that of a 5-NN.
          False depends on layout of training and test data
```

1. The decision boundary of the k-NN classifier is linear.

False arbitrary boundaries depending on closest neighbors at all points. Not necessarily linear

True As training data increases, you need to compute distances and compare for more points.

1. The time needed to classify a test example with the k-NN classifier grows with the size of the training set.

k-Nearest Neighbor (kNN) exercise

more details see the assignments page on the course website.

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 **CIFAR-10 Data Loading and Preprocessing** In [6]: # Load the raw CIFAR-10 data. cifar10 dir = 'cs682/datasets/cifar-10-batches-py' # Cleaning up variables to prevent loading data multiple times (which may cause memory issue) try: del X train, y_train del X test, y test print('Clear previously loaded data.') except: pass 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 [7]: # 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 [8]: # Split the data into train, val, and test sets. In addition we will # create a small development set as a subset of the training data; # we can use this for development so our code runs faster. num training = 49000 num validation = 1000 num test = 1000num dev = 500# Our validation set will be num validation points from the original # training set. mask = range(num training, num training + num validation) X_val = X_train[mask] y_val = y_train[mask] # Our training set will be the first num train points from the original # training set. mask = range(num training) X_train = X_train[mask] y train = y train[mask] # We will also make a development set, which is a small subset of # the training set. mask = np.random.choice(num_training, num_dev, replace=False) X dev = X train[mask] y dev = y train[mask] # We use the first num test points of the original test set as our # test set. mask = range(num_test) X_test = X_test[mask] y test = y test[mask] print('Train data shape: ', X_train.shape) print('Train labels shape: ', y_train.shape) print('Validation data shape: ', X_val.shape) print('Validation labels shape: ', y_val.shape) print('Test data shape: ', X test.shape) print('Test labels shape: ', y_test.shape) Train data shape: (49000, 32, 32, 3) Train labels shape: (49000,) Validation data shape: (1000, 32, 32, 3) Validation labels shape: (1000,) Test data shape: (1000, 32, 32, 3) Test labels shape: (1000,) In [9]: # Preprocessing: reshape the image data into rows X train = np.reshape(X_train, (X_train.shape[0], -1)) X_val = np.reshape(X_val, (X_val.shape[0], -1)) X_test = np.reshape(X_test, (X_test.shape[0], -1)) X_dev = np.reshape(X_dev, (X_dev.shape[0], -1)) # As a sanity check, print out the shapes of the data print('Training data shape: ', X_train.shape) print('Validation data shape: ', X val.shape) print('Test data shape: ', X_test.shape) print('dev data shape: ', X_dev.shape) Training data shape: (49000, 3072) Validation data shape: (1000, 3072) Test data shape: (1000, 3072) dev data shape: (500, 3072) In [10]: # Preprocessing: subtract the mean image # first: compute the image mean based on the training data mean_image = np.mean(X_train, axis=0) print(mean image[:10]) # print a few of the elements plt.figure(figsize=(4,4)) plt.imshow(mean_image.reshape((32,32,3)).astype('uint8')) # visualize the mean image plt.show() [130.64189796 135.98173469 132.47391837 130.05569388 135.34804082 131.75402041 130.96055102 136.14328571 132.47636735 131.48467347] 0 5 10 15 20 25 30 5 10 15 20 25 30 In [11]: # second: subtract the mean image from train and test data X train -= mean image X val -= mean image X test -= mean image X dev -= mean image In [12]: # third: append the bias dimension of ones (i.e. bias trick) so that our SVM # only has to worry about optimizing a single weight matrix W. X train = np.hstack([X train, np.ones((X train.shape[0], 1))]) X_val = np.hstack([X_val, np.ones((X_val.shape[0], 1))]) X test = np.hstack([X test, np.ones((X test.shape[0], 1))]) X dev = np.hstack([X dev, np.ones((X dev.shape[0], 1))]) print(X train.shape, X val.shape, X test.shape, X dev.shape) (49000, 3073) (1000, 3073) (1000, 3073) (500, 3073) **SVM Classifier** Your code for this section will all be written inside cs682/classifiers/linear_svm.py. As you can see, we have prefilled the function sym loss naive which uses for loops to evaluate the multiclass SVM loss function. In [13]: # Evaluate the naive implementation of the loss we provided for you: from cs682.classifiers.linear_svm import svm loss naive import time # generate a random SVM weight matrix of small numbers W = np.random.randn(3073, 10) * 0.0001loss, grad = svm loss naive(W, X dev, y dev, 0.000005) print('loss: %f' % (loss,)) loss: 8.606381 The grad returned from the function above is right now all zero. Derive and implement the gradient for the SVM cost function and implement it inline inside the function sym loss naive. You will find it helpful to interleave your new code inside the existing function. To check that you have correctly implemented the gradient correctly, you can numerically estimate the gradient of the loss function and compare the numeric estimate to the gradient that you computed. We have provided code that does this for you: In [14]: # Once you've implemented the gradient, recompute it with the code below # and gradient check it with the function we provided for you # Compute the loss and its gradient at W. loss, grad = svm loss naive(W, X dev, y dev, 0.0) # Numerically compute the gradient along several randomly chosen dimensions, and # compare them with your analytically computed gradient. The numbers should match # almost exactly along all dimensions. from cs682.gradient check import grad check sparse f = lambda w: svm_loss_naive(w, X_dev, y_dev, 0.0)[0] grad numerical = grad check sparse(f, W, grad) # do the gradient check once again with regularization turned on # you didn't forget the regularization gradient did you? loss, grad = svm_loss_naive(W, X_dev, y_dev, 5e1) f = lambda w: svm_loss_naive(w, X_dev, y_dev, 5e1)[0] grad numerical = grad check sparse(f, W, grad) numerical: 32.662110 analytic: 32.662110, relative error: 1.090752e-11 numerical: 1.663175 analytic: 1.663175, relative error: 3.914669e-11 numerical: -39.167971 analytic: -39.167971, relative error: 9.851151e-12 numerical: 26.852921 analytic: 26.852921, relative error: 3.227127e-12 numerical: -14.226808 analytic: -14.226808, relative error: 1.909513e-11 numerical: 20.146966 analytic: 20.146966, relative error: 3.368709e-12 numerical: -7.677977 analytic: -7.677977, relative error: 5.282179e-11 numerical: -7.721987 analytic: -7.721987, relative error: 3.408445e-11 numerical: 13.315709 analytic: 13.315709, relative error: 3.016378e-12 numerical: -7.827410 analytic: -7.827410, relative error: 4.291792e-11 numerical: 7.752431 analytic: 7.752431, relative error: 1.451516e-11 numerical: 0.760045 analytic: 0.760045, relative error: 4.056273e-10 numerical: -5.437939 analytic: -5.437939, relative error: 5.611144e-11 numerical: 30.272562 analytic: 30.272562, relative error: 8.595795e-12 numerical: 12.690939 analytic: 12.690939, relative error: 2.237610e-11 numerical: -8.214033 analytic: -8.214033, relative error: 5.088807e-11 numerical: -53.618353 analytic: -53.618353, relative error: 6.217904e-12 numerical: -46.852087 analytic: -46.852087, relative error: 4.444220e-12 numerical: -24.371575 analytic: -24.371575, relative error: 1.975149e-12 numerical: 13.691466 analytic: 13.691466, relative error: 1.827882e-11 **Inline Question 1:** It is possible that once in a while a dimension in the gradcheck will not match exactly. What could such a discrepancy be caused by? Is it a reason for concern? What is a simple example in one dimension where a gradient check could fail? How would change the margin affect of the frequency of this happening? Hint: the SVM loss function is not strictly speaking differentiable Your Answer: Yes, it is possible that a dimension of grad_check may be different. Since grad_check computes the gradient by simply adding a small value to X and checking the function values instead of taking the actual derivative it could be a different value. Also if there is a point where the loss isn't differentiable (like where the hinge loss goes from 0 to above 0) the numerical derivative will return a correct value but the analytical derivative may not. In [15]: # Next implement the function svm loss vectorized; for now only compute the loss; # we will implement the gradient in a moment. tic = time.time() loss naive, grad naive = svm loss naive(W, X dev, y dev, 0.000005) toc = time.time() print('Naive loss: %e computed in %fs' % (loss_naive, toc - tic)) from cs682.classifiers.linear svm import svm loss vectorized tic = time.time() loss_vectorized, _ = svm_loss_vectorized(W, X_dev, y_dev, 0.000005) toc = time.time() print('Vectorized loss: %e computed in %fs' % (loss_vectorized, toc - tic)) # The losses should match but your vectorized implementation should be much faster. print('difference: %f' % (loss_naive - loss_vectorized)) Naive loss: 8.606381e+00 computed in 0.140188s Vectorized loss: 8.606381e+00 computed in 0.005475s difference: -0.000000 In [16]: # Complete the implementation of svm loss vectorized, and compute the gradient # of the loss function in a vectorized way. # The naive implementation and the vectorized implementation should match, but # the vectorized version should still be much faster. tic = time.time() _, grad_naive = svm_loss_naive(W, X_dev, y_dev, 0.000005) toc = time.time() print('Naive loss and gradient: computed in %fs' % (toc - tic)) tic = time.time() _, grad_vectorized = svm_loss_vectorized(W, X_dev, y_dev, 0.000005) print('Vectorized loss and gradient: computed in %fs' % (toc - tic)) # The loss is a single number, so it is easy to compare the values computed # by the two implementations. The gradient on the other hand is a matrix, so # we use the Frobenius norm to compare them. difference = np.linalg.norm(grad_naive - grad_vectorized, ord='fro') print('difference: %f' % difference) Naive loss and gradient: computed in 0.104291s Vectorized loss and gradient: computed in 0.004068s difference: 0.000000 **Stochastic Gradient Descent** We now have vectorized and efficient expressions for the loss, the gradient and our gradient matches the numerical gradient. We are therefore ready to do SGD to minimize the loss. In [17]: # In the file linear classifier.py, implement SGD in the function # LinearClassifier.train() and then run it with the code below. from cs682.classifiers import LinearSVM svm = LinearSVM() tic = time.time() loss_hist = svm.train(X_train, y_train, learning_rate=1e-7, reg=2.5e4, num iters=1500, verbose=True) toc = time.time() print('That took %fs' % (toc - tic)) iteration 0 / 1500: loss 796.140715 iteration 100 / 1500: loss 290.013147 iteration 200 / 1500: loss 109.450232 iteration 300 / 1500: loss 42.935666 iteration 400 / 1500: loss 19.140520 iteration 500 / 1500: loss 10.010751 iteration 600 / 1500: loss 7.287787 iteration 700 / 1500: loss 5.990759 iteration 800 / 1500: loss 5.262746 iteration 900 / 1500: loss 5.252881 iteration 1000 / 1500: loss 4.651852 iteration 1100 / 1500: loss 5.187224 iteration 1200 / 1500: loss 4.735986 iteration 1300 / 1500: loss 6.000257 iteration 1400 / 1500: loss 5.278133 That took 4.907961s In [18]: # A useful debugging strategy is to plot the loss as a function of # iteration number: plt.plot(loss hist) plt.xlabel('Iteration number') plt.ylabel('Loss value') plt.show() 800 700 600 500 oss value 300 200 100 0 200 1400 400 800 1200 1000 Iteration number In [19]: # Write the LinearSVM.predict function and evaluate the performance on both the # training and validation set y train pred = svm.predict(X train) print('training accuracy: %f' % (np.mean(y train == y train pred),)) y val pred = svm.predict(X val) print('validation accuracy: %f' % (np.mean(y_val == y_val_pred),)) training accuracy: 0.363776 validation accuracy: 0.377000 In [21]: # Use the validation set to tune hyperparameters (regularization strength and # learning rate). You should experiment with different ranges for the learning # rates and regularization strengths; if you are careful you should be able to # get a classification accuracy of about 0.4 on the validation set. learning rates = [2e-8, 1e-7]regularization_strengths = [2.5e4, 1e4] # results is dictionary mapping tuples of the form # (learning rate, regularization strength) to tuples of the form # (training accuracy, validation accuracy). The accuracy is simply the fraction # of data points that are correctly classified. results = {} best val = -1 # The highest validation accuracy that we have seen so far. best_svm = None # The LinearSVM object that achieved the highest validation rate. # TODO: # Write code that chooses the best hyperparameters by tuning on the validation # # set. For each combination of hyperparameters, train a linear SVM on the # training set, compute its accuracy on the training and validation sets, and # # store these numbers in the results dictionary. In addition, store the best # # validation accuracy in best val and the LinearSVM object that achieves this # # accuracy in best svm. # Hint: You should use a small value for num iters as you develop your # validation code so that the SVMs don't take much time to train; once you are # # confident that your validation code works, you should rerun the validation # # code with a larger value for num iters. for lr in learning rates: for reg in regularization strengths: print("Testing LR=", lr, " reg=", reg) svm = LinearSVM() svm.train(X train, y train, learning rate=lr, reg=reg, num_iters=10000, verbose=False) y train pred = svm.predict(X train) tr acc = np.mean(y train == y train pred) print('training accuracy: %f' % tr acc) y val pred = svm.predict(X val) val_acc = np.mean(y_val == y_val_pred) print('validation accuracy: %f' % val acc) results[(lr, reg)] = tr acc, val acc, svm END OF YOUR CODE # Print out results. for lr, reg in sorted(results): train accuracy, val accuracy, svm = results[(lr, reg)] if val accuracy > best val: best val = val accuracy best svm = svm print('lr %e reg %e train accuracy: %f val accuracy: %f' % (lr, reg, train_accuracy, val_accuracy)) print('best validation accuracy achieved during cross-validation: %f' % best_val) Testing LR= 2e-08 reg= 25000.0 training accuracy: 0.374898 validation accuracy: 0.394000 Testing LR= 2e-08 reg= 10000.0 training accuracy: 0.389592 validation accuracy: 0.403000 Testing LR= 1e-07 reg= 25000.0 training accuracy: 0.367000 validation accuracy: 0.373000 Testing LR= 1e-07 reg= 10000.0 training accuracy: 0.384204 validation accuracy: 0.382000 lr 2.000000e-08 reg 1.000000e+04 train accuracy: 0.389592 val accuracy: 0.403000 lr 2.000000e-08 reg 2.500000e+04 train accuracy: 0.374898 val accuracy: 0.394000 lr 1.000000e-07 reg 1.000000e+04 train accuracy: 0.384204 val accuracy: 0.382000 lr 1.000000e-07 reg 2.500000e+04 train accuracy: 0.367000 val accuracy: 0.373000 best validation accuracy achieved during cross-validation: 0.403000 In [22]: # Visualize the cross-validation results import math $x_scatter = [math.log10(x[0]) for x in results]$ y scatter = [math.log10(x[1]) for x in results] # plot training accuracy marker size = 100 colors = [results[x][0] for x in results] plt.subplot(2, 1, 1)plt.scatter(x scatter, y scatter, marker size, c=colors) plt.colorbar() plt.xlabel('log learning rate') plt.ylabel('log regularization strength') plt.title('CIFAR-10 training accuracy') # plot validation accuracy colors = [results[x][1] for x in results] # default size of markers is 20 plt.subplot(2, 1, 2)plt.scatter(x_scatter, y_scatter, marker_size, c=colors) plt.colorbar() plt.xlabel('log learning rate') plt.ylabel('log regularization strength') plt.title('CIFAR-10 validation accuracy') plt.show() CIFAR-10 training accuracy 4.40 4.35 4.30 4.25 4.20 4.15 4.10 4.05 - 0.385 0.380 0.375 0.370 4.00 -7.0 -7.6-7.5-7.4-7.3-7.2-7.1-7.7CIFAR-10% MONTH TO FIRE CCUracy 4.40 0.400 4.35 - 0.395 4.30 4.25 0.3904.20 0.385 4.15 4.10 0.380 4.05 0.375 4.00 -7.7-7.6-7.5-7.4-7.3-7.2-7.1log learning rate In [24]: # Evaluate the best svm on test set y_test_pred = best_svm.predict(X_test) test_accuracy = np.mean(y_test == y_test_pred) print('linear SVM on raw pixels final test set accuracy: %f' % test_accuracy) linear SVM on raw pixels final test set accuracy: 0.382000 In [25]: # Visualize the learned weights for each class. # Depending on your choice of learning rate and regularization strength, these may # or may not be nice to look at. w = best_svm.W[:-1,:] # strip out the bias w = w.reshape(32, 32, 3, 10) w_{\min} , $w_{\max} = np.min(w)$, np.max(w)classes = ['plane', 'car', 'bird', 'cat', 'deer', 'dog', 'frog', 'horse', 'ship', 'truck'] for i in range (10): plt.subplot(2, 5, i + 1)# Rescale the weights to be between 0 and 255 wimg = $255.0 * (w[:, :, i].squeeze() - w_min) / (w_max - w_min)$ plt.imshow(wimg.astype('uint8')) plt.axis('off') plt.title(classes[i]) plane bird cat deer car

dog

Inline question 2:

Your answer:

frog

horse

ship

Describe what your visualized SVM weights look like, and offer a brief explanation for why they look they way that they do.

because their is a large variety of different images per class and the sym learns the features most common to each class.

truck

The SVM weights look like fairly random mixed colors. In some you can make out what looks like images like the truck, car, horse and frog. They look so mixed

Multiclass Support Vector Machine exercise

more details see the assignments page on the course website.

• implement a fully-vectorized loss function for the SVM

• check your implementation using numerical gradient

• optimize the loss function with SGD

• **visualize** the final learned weights

In [5]: # Run some setup code for this notebook.

from future import print function

from cs682.data utils import load CIFAR10

• implement the fully-vectorized expression for its **analytic gradient**

• use a validation set to tune the learning rate and regularization strength

In this exercise you will:

import random

import numpy as np

```
more details see the assignments page on the course website.
         This exercise is analogous to the SVM exercise. You will:
           • implement a fully-vectorized loss function for the Softmax classifier
           • implement the fully-vectorized expression for its analytic gradient
           • check your implementation with numerical gradient
           • use a validation set to tune the learning rate and regularization strength

    optimize the loss function with SGD

           • visualize the final learned weights
 In [3]: from __future__ import print_function
         import random
         import numpy as np
         from cs682.data_utils import load_CIFAR10
         import matplotlib.pyplot as plt
         %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'
         # for auto-reloading extenrnal modules
         # see http://stackoverflow.com/questions/1907993/autoreload-of-modules-in-ipython
         %load_ext autoreload
         %autoreload 2
 In [4]: def get_CIFAR10_data(num_training=49000, num_validation=1000, num_test=1000, num_dev=500):
             Load the CIFAR-10 dataset from disk and perform preprocessing to prepare
             it for the linear classifier. These are the same steps as we used for the
             SVM, but condensed to a single function.
             # Load the raw CIFAR-10 data
             cifar10_dir = 'cs682/datasets/cifar-10-batches-py'
             X train, y train, X test, y test = load CIFAR10(cifar10 dir)
             # subsample the data
             mask = list(range(num_training, num_training + num validation))
             X_val = X_train[mask]
             y_val = y_train[mask]
             mask = list(range(num training))
             X train = X train[mask]
             y_train = y_train[mask]
             mask = list(range(num_test))
             X test = X test[mask]
             y_test = y_test[mask]
             mask = np.random.choice(num_training, num_dev, replace=False)
             X_dev = X_train[mask]
             y_dev = y_train[mask]
             # Preprocessing: reshape the image data into rows
             X train = np.reshape(X train, (X train.shape[0], -1))
             X_val = np.reshape(X_val, (X_val.shape[0], -1))
             X_test = np.reshape(X_test, (X_test.shape[0], -1))
             X \text{ dev} = \text{np.reshape}(X \text{ dev}, (X \text{ dev.shape}[0], -1))
             # Normalize the data: subtract the mean image
             mean_image = np.mean(X_train, axis = 0)
             X_train -= mean_image
             X_val -= mean_image
             X_test -= mean_image
             X dev -= mean image
             # add bias dimension and transform into columns
             X_train = np.hstack([X_train, np.ones((X_train.shape[0], 1))])
             X_val = np.hstack([X_val, np.ones((X_val.shape[0], 1))])
             X_test = np.hstack([X_test, np.ones((X_test.shape[0], 1))])
             X_dev = np.hstack([X_dev, np.ones((X_dev.shape[0], 1))])
             return X_train, y_train, X_val, y_val, X_test, y_test, X_dev, y_dev
         # Cleaning up variables to prevent loading data multiple times (which may cause memory issue)
         try:
            del X_train, y_train
            del X_test, y_test
            print('Clear previously loaded data.')
         except:
            pass
         # Invoke the above function to get our data.
         X_train, y_train, X_val, y_val, X_test, y_test, X_dev, y_dev = get_CIFAR10_data()
         print('Train data shape: ', X_train.shape)
         print('Train labels shape: ', y train.shape)
         print('Validation data shape: ', X_val.shape)
         print('Validation labels shape: ', y_val.shape)
         print('Test data shape: ', X_test.shape)
         print('Test labels shape: ', y_test.shape)
         print('dev data shape: ', X dev.shape)
         print('dev labels shape: ', y_dev.shape)
         Train data shape: (49000, 3073)
         Train labels shape: (49000,)
         Validation data shape: (1000, 3073)
         Validation labels shape: (1000,)
         Test data shape: (1000, 3073)
         Test labels shape: (1000,)
         dev data shape: (500, 3073)
         dev labels shape: (500,)
         Softmax Classifier
         Your code for this section will all be written inside cs682/classifiers/softmax.py.
 In [5]: # First implement the naive softmax loss function with nested loops.
         # Open the file cs682/classifiers/softmax.py and implement the
         # softmax loss naive function.
         from cs682.classifiers.softmax import softmax loss naive
         import time
         # Generate a random softmax weight matrix and use it to compute the loss.
         W = np.random.randn(3073, 10) * 0.0001
         loss, grad = softmax_loss_naive(W, X_dev, y_dev, 0.0)
         # As a rough sanity check, our loss should be something close to -\log(0.1).
         print('loss: %f' % loss)
         print('sanity check: %f' % (-np.log(0.1)))
         loss: 2.329227
         sanity check: 2.302585
         Inline Question 1:
         Why do we expect our loss to be close to -log(0.1)? Explain briefly.**
         Your answer:
         The loss should be around -log(0.1 because we have 10 classes and assuming the scores are all approximately the same, the loss should be around -
         \log(e^score / (10 * e^score)) = -\log(1/10).
 In [5]: # Complete the implementation of softmax loss naive and implement a (naive)
         # version of the gradient that uses nested loops.
         loss, grad = softmax loss naive(W, X dev, y dev, 0.0)
         # As we did for the SVM, use numeric gradient checking as a debugging tool.
         # The numeric gradient should be close to the analytic gradient.
         from cs682.gradient_check import grad check sparse
         f = lambda w: softmax_loss_naive(w, X_dev, y_dev, 0.0)[0]
         grad numerical = grad check sparse(f, W, grad, 10)
          # similar to SVM case, do another gradient check with regularization
         loss, grad = softmax loss naive(W, X dev, y dev, 5e1)
         f = lambda w: softmax loss naive(w, X_dev, y_dev, 5e1)[0]
         grad numerical = grad check sparse(f, W, grad, 10)
         numerical: -1.544645 analytic: -1.544645, relative error: 5.320434e-08
         numerical: -2.178610 analytic: -2.178610, relative error: 5.554907e-09
         numerical: 0.361697 analytic: 0.361697, relative error: 1.125100e-07
         numerical: 0.076178 analytic: 0.076178, relative error: 5.267031e-07
         numerical: -0.248585 analytic: -0.248585, relative error: 2.649104e-08
         numerical: 0.309712 analytic: 0.309712, relative error: 5.420240e-08
         numerical: 1.630478 analytic: 1.630478, relative error: 1.205564e-08
         numerical: -0.452599 analytic: -0.452599, relative error: 7.546752e-08
         numerical: -0.276927 analytic: -0.276928, relative error: 2.554157e-07
         numerical: -3.078943 analytic: -3.078943, relative error: 1.107484e-08
         numerical: -1.695526 analytic: -1.695526, relative error: 2.552773e-08
         numerical: -6.392450 analytic: -6.392450, relative error: 9.196914e-09
         numerical: 0.545001 analytic: 0.545001, relative error: 1.411039e-07
         numerical: 2.092425 analytic: 2.092425, relative error: 2.684974e-08
         numerical: 0.603232 analytic: 0.603232, relative error: 8.809838e-08
         numerical: -2.125793 analytic: -2.125793, relative error: 1.717946e-09
         numerical: 0.652094 analytic: 0.652094, relative error: 9.618587e-08
         numerical: -6.011416 analytic: -6.011416, relative error: 7.344141e-11
         numerical: -3.414886 analytic: -3.414886, relative error: 2.819845e-09
         numerical: 1.816323 analytic: 1.816322, relative error: 5.299007e-08
 In [6]: # Now that we have a naive implementation of the softmax loss function and its gradient,
         # implement a vectorized version in softmax loss vectorized.
         # The two versions should compute the same results, but the vectorized version should be
         # much faster.
         tic = time.time()
         loss naive, grad naive = softmax loss naive(W, X dev, y dev, 0.000005)
         toc = time.time()
         print('naive loss: %e computed in %fs' % (loss naive, toc - tic))
         from cs682.classifiers.softmax import softmax loss vectorized
         tic = time.time()
         loss vectorized, grad vectorized = softmax loss vectorized(W, X dev, y dev, 0.000005)
         toc = time.time()
         print('vectorized loss: %e computed in %fs' % (loss vectorized, toc - tic))
         # As we did for the SVM, we use the Frobenius norm to compare the two versions
         # of the gradient.
         grad difference = np.linalg.norm(grad naive - grad vectorized, ord='fro')
         print('Loss difference: %f' % np.abs(loss naive - loss vectorized))
         print('Gradient difference: %f' % grad_difference)
         naive loss: 2.418885e+00 computed in 0.148879s
         vectorized loss: 2.418885e+00 computed in 0.005713s
         Loss difference: 0.000000
         Gradient difference: 0.000000
 In [6]: # Use the validation set to tune hyperparameters (regularization strength and
         # learning rate). You should experiment with different ranges for the learning
         # rates and regularization strengths; if you are careful you should be able to
         # get a classification accuracy of over 0.35 on the validation set.
         from cs682.classifiers import Softmax
         results = {}
         best_val = -1
         best softmax = None
         learning rates = [5e-8, 1e-7, 5e-7]
         regularization strengths = [5e3, 1.5e4, 2.5e4, 5e4]
         # TODO:
         # Use the validation set to set the learning rate and regularization strength. #
         # This should be identical to the validation that you did for the SVM; save #
         # the best trained softmax classifer in best softmax.
         for lr in learning rates:
             for reg in regularization_strengths:
                 print("Testing LR=", lr, " reg=", reg)
                 softmax = Softmax()
                 softmax.train(X train, y train, learning rate=lr, reg=reg,
                                num iters=6000, verbose=False)
                 y_train_pred = softmax.predict(X_train)
                 tr acc = np.mean(y train == y train pred)
                 print('training accuracy: %f' % tr acc)
                 y_val_pred = softmax.predict(X_val)
                 val acc = np.mean(y val == y val pred)
                 print('validation accuracy: %f' % val_acc)
                 results[(lr, reg)] = tr acc, val acc, softmax
          END OF YOUR CODE
         # Print out results.
         for lr, reg in sorted(results):
             train_accuracy, val_accuracy, softmax = results[(lr, reg)]
             if val accuracy > best val:
                 best val = val accuracy
                 best softmax = softmax
             print('lr %e reg %e train accuracy: %f val accuracy: %f' % (
                         lr, reg, train accuracy, val accuracy))
         print('best validation accuracy achieved during cross-validation: %f' % best val)
         Testing LR= 5e-08 reg= 5000.0
         training accuracy: 0.374224
         validation accuracy: 0.379000
         Testing LR= 5e-08 reg= 15000.0
         training accuracy: 0.346388
         validation accuracy: 0.365000
         Testing LR= 5e-08 reg= 25000.0
         training accuracy: 0.332714
         validation accuracy: 0.347000
         Testing LR= 5e-08 reg= 50000.0
         training accuracy: 0.314714
         validation accuracy: 0.332000
         Testing LR= 1e-07 reg= 5000.0
         training accuracy: 0.377020
         validation accuracy: 0.386000
         Testing LR= 1e-07 reg= 15000.0
         training accuracy: 0.342898
         validation accuracy: 0.360000
         Testing LR= 1e-07 reg= 25000.0
         training accuracy: 0.330020
         validation accuracy: 0.346000
         Testing LR= 1e-07 reg= 50000.0
         training accuracy: 0.306755
         validation accuracy: 0.318000
         Testing LR= 5e-07 reg= 5000.0
         training accuracy: 0.371041
         validation accuracy: 0.382000
         Testing LR= 5e-07 reg= 15000.0
         training accuracy: 0.338633
         validation accuracy: 0.350000
         Testing LR= 5e-07 reg= 25000.0
         training accuracy: 0.337837
         validation accuracy: 0.345000
         Testing LR= 5e-07 reg= 50000.0
         training accuracy: 0.306224
         validation accuracy: 0.321000
         1r 5.000000e-08 reg 5.000000e+03 train accuracy: 0.374224 val accuracy: 0.379000
         lr 5.000000e-08 reg 1.500000e+04 train accuracy: 0.346388 val accuracy: 0.365000
         lr 5.000000e-08 reg 2.500000e+04 train accuracy: 0.332714 val accuracy: 0.347000
         lr 5.000000e-08 reg 5.000000e+04 train accuracy: 0.314714 val accuracy: 0.332000
         lr 1.000000e-07 reg 5.000000e+03 train accuracy: 0.377020 val accuracy: 0.386000
         lr 1.000000e-07 reg 1.500000e+04 train accuracy: 0.342898 val accuracy: 0.360000
         lr 1.000000e-07 reg 2.500000e+04 train accuracy: 0.330020 val accuracy: 0.346000
         lr 1.000000e-07 reg 5.000000e+04 train accuracy: 0.306755 val accuracy: 0.318000
         lr 5.000000e-07 reg 5.000000e+03 train accuracy: 0.371041 val accuracy: 0.382000
         lr 5.000000e-07 reg 1.500000e+04 train accuracy: 0.338633 val accuracy: 0.350000
         lr 5.000000e-07 reg 2.500000e+04 train accuracy: 0.337837 val accuracy: 0.345000
         lr 5.000000e-07 reg 5.000000e+04 train accuracy: 0.306224 val accuracy: 0.321000
         best validation accuracy achieved during cross-validation: 0.386000
 In [7]: # evaluate on test set
         # Evaluate the best softmax on test set
         y_test_pred = best_softmax.predict(X_test)
         test_accuracy = np.mean(y_test == y_test_pred)
         print('softmax on raw pixels final test set accuracy: %f' % (test accuracy, ))
         softmax on raw pixels final test set accuracy: 0.383000
         Inline Question - True or False
         It's possible to add a new datapoint to a training set that would leave the SVM loss unchanged, but this is not the case with the Softmax classifier loss.
         Your answer:
         This is because sym loss uses a hinge loss type function and margins so if the new added datapoint's score is outside the margin for it's class the loss
         wouldn't change however softmax loss takes all data points into account no matter what their score is so the loss would change.
In [44]: # Visualize the learned weights for each class
         w = best softmax.W[:-1,:] # strip out the bias
         w = w.reshape(32, 32, 3, 10)
         w \min, w \max = np.\min(w), np.\max(w)
         classes = ['plane', 'car', 'bird', 'cat', 'deer', 'dog', 'frog', 'horse', 'ship', 'truck']
         for i in range(10):
             plt.subplot(2, 5, i + 1)
             # Rescale the weights to be between 0 and 255
             wimg = 255.0 * (w[:, :, i].squeeze() - w_min) / (w_max - w_min)
             plt.imshow(wimg.astype('uint8'))
             plt.axis('off')
             plt.title(classes[i])
                                          bird
                                                                      deer
              plane
                                                         cat
```

horse

dog

frog

ship

truck

Softmax exercise ¶

return TwoLayerNet(input size, hidden size, num classes, std=1e-1) def init toy data(): np.random.seed(1) X = 10 * np.random.randn(num inputs, input size) y = np.array([0, 1, 2, 2, 1])return X, y net = init toy model() X_{i} y = init toy data() Forward pass: compute scores Open the file cs682/classifiers/neural net.py and look at the method TwoLayerNet.loss. This function is very similar to the loss functions you have written for the SVM and Softmax exercises: It takes the data and weights and computes the class scores, the loss, and the gradients on the parameters. Implement the first part of the forward pass which uses the weights and biases to compute the scores for all inputs. In [4]: scores = net.loss(X) print('Your scores:') print(scores) print() print('correct scores:') correct scores = np.asarray([[-0.81233741, -1.27654624, -0.70335995],[-0.17129677, -1.18803311, -0.47310444],[-0.51590475, -1.01354314, -0.8504215],[-0.15419291, -0.48629638, -0.52901952],[-0.00618733, -0.12435261, -0.15226949]])print(correct scores) print() # The difference should be very small. We get < 1e-7 print('Difference between your scores and correct scores:') print(np.sum(np.abs(scores - correct_scores))) Your scores: [[-0.81233741 -1.27654624 -0.70335995][-0.17129677 -1.18803311 -0.47310444][-0.51590475 -1.01354314 -0.8504215][-0.15419291 -0.48629638 -0.52901952][-0.00618733 -0.12435261 -0.15226949]]correct scores: [[-0.81233741 -1.27654624 -0.70335995][-0.17129677 -1.18803311 -0.47310444][-0.51590475 -1.01354314 -0.8504215][-0.15419291 -0.48629638 -0.52901952][-0.00618733 -0.12435261 -0.15226949]]Difference between your scores and correct scores: 3.6802720496109664e-08 Forward pass: compute loss In the same function, implement the second part that computes the data and regularization loss. In [5]: loss, _ = net.loss(X, y, reg=0.05) correct loss = 1.30378789133# should be very small, we get < 1e-12</pre> print('Difference between your loss and correct loss:') print(np.sum(np.abs(loss - correct loss))) Difference between your loss and correct loss: 1.794120407794253e-13 **Backward pass** Implement the rest of the function. This will compute the gradient of the loss with respect to the variables W1, b1, W2, and b2. Now that you (hopefully!) have a correctly implemented forward pass, you can debug your backward pass using a numeric gradient check: In [6]: from cs682.gradient check import eval numerical gradient # Use numeric gradient checking to check your implementation of the backward pass. # If your implementation is correct, the difference between the numeric and # analytic gradients should be less than 1e-8 for each of W1, W2, b1, and b2. loss, grads = net.loss(X, y, reg=0.05) # these should all be less than 1e-8 or so for param name in grads: f = lambda W: net.loss(X, y, reg=0.05)[0]param grad num = eval numerical gradient(f, net.params[param name], verbose=False) print('%s max relative error: %e' % (param name, rel error(param grad num, grads[param name]))) W2 max relative error: 3.440708e-09 b2 max relative error: 3.865091e-11 W1 max relative error: 3.561318e-09 b1 max relative error: 2.738421e-09 Train the network To train the network we will use stochastic gradient descent (SGD), similar to the SVM and Softmax classifiers. Look at the function TwoLayerNet.train and fill in the missing sections to implement the training procedure. This should be very similar to the training procedure you used for the SVM and Softmax classifiers. You will also have to implement TwoLayerNet.predict, as the training process periodically performs prediction to keep track of accuracy over time while the network trains. Once you have implemented the method, run the code below to train a two-layer network on toy data. You should achieve a training loss less than 0.2. In [7]: net = init toy model() stats = net.train(X, y, X, y, learning rate=1e-1, reg=5e-6, num iters=100, verbose=False, batch size=10) print('Final training loss: ', stats['loss_history'][-1]) # plot the loss history plt.plot(stats['loss history']) plt.xlabel('iteration') plt.ylabel('training loss') plt.title('Training Loss history') plt.show() Final training loss: 0.01925119514188341 Training Loss history 1.2 1.0 0.8 training loss 0.4 0.2 0.0 20 60 80 100 iteration Load the data Now that you have implemented a two-layer network that passes gradient checks and works on toy data, it's time to load up our favorite CIFAR-10 data so we can use it to train a classifier on a real dataset. In [8]: from cs682.data utils import load CIFAR10 def get CIFAR10 data(num training=49000, num validation=1000, num test=1000): Load the CIFAR-10 dataset from disk and perform preprocessing to prepare it for the two-layer neural net classifier. These are the same steps as we used for the SVM, but condensed to a single function. # Load the raw CIFAR-10 data cifar10 dir = 'cs682/datasets/cifar-10-batches-py' X_train, y_train, X_test, y_test = load_CIFAR10(cifar10_dir) # Subsample the data mask = list(range(num training, num training + num validation)) X_val = X_train[mask] y val = y train[mask] mask = list(range(num training)) X train = X train[mask] y train = y train[mask] mask = list(range(num test)) X_test = X_test[mask] y test = y test[mask] # Normalize the data: subtract the mean image mean_image = np.mean(X_train, axis=0) X train -= mean image X_val -= mean_image X_test -= mean_image # Reshape data to rows X train = X train.reshape(num training, -1) X val = X val.reshape(num validation, -1) X test = X test.reshape(num test, -1)return X train, y train, X val, y val, X test, y test # Cleaning up variables to prevent loading data multiple times (which may cause memory issue) try: del X_train, y_train del X_test, y_test print('Clear previously loaded data.') except: pass # Invoke the above function to get our data. X train, y train, X val, y val, X test, y test = get CIFAR10 data() print('Train data shape: ', X train.shape) print('Train labels shape: ', y_train.shape) print('Validation data shape: ', X_val.shape) print('Validation labels shape: ', y_val.shape) print('Test data shape: ', X test.shape) print('Test labels shape: ', y test.shape) Train data shape: (49000, 3072) Train labels shape: (49000,) Validation data shape: (1000, 3072) Validation labels shape: (1000,) Test data shape: (1000, 3072) Test labels shape: (1000,) Train a network To train our network we will use SGD. In addition, we will adjust the learning rate with an exponential learning rate schedule as optimization proceeds; after each epoch, we will reduce the learning rate by multiplying it by a decay rate. In [9]: input size = 32 * 32 * 3 hidden size = 50 num classes = 10 net = TwoLayerNet(input size, hidden size, num classes) # Train the network stats = net.train(X train, y train, X val, y val, num_iters=1000, batch_size=200, learning_rate=1e-4, learning_rate_decay=0.95, reg=0.25, verbose=**True**) # Predict on the validation set val acc = (net.predict(X val) == y val).mean() print('Validation accuracy: ', val acc) iteration 0 / 1000: loss 2.302963 iteration 100 / 1000: loss 2.302517 iteration 200 / 1000: loss 2.297871 iteration 300 / 1000: loss 2.264625 iteration 400 / 1000: loss 2.199637 iteration 500 / 1000: loss 2.103959 iteration 600 / 1000: loss 2.027170 iteration 700 / 1000: loss 2.017522 iteration 800 / 1000: loss 1.972401 iteration 900 / 1000: loss 1.924231 Validation accuracy: 0.288 **Debug the training** With the default parameters we provided above, you should get a validation accuracy of about 0.29 on the validation set. This isn't very good. One strategy for getting insight into what's wrong is to plot the loss function and the accuracies on the training and validation sets during optimization. Another strategy is to visualize the weights that were learned in the first layer of the network. In most neural networks trained on visual data, the first layer weights typically show some visible structure when visualized. In [10]: # Plot the loss function and train / validation accuracies plt.subplot(2, 1, 1)plt.plot(stats['loss_history']) plt.title('Loss history') plt.xlabel('Iteration') plt.ylabel('Loss') plt.subplot(2, 1, 2) plt.plot(stats['train acc history'], label='train') plt.plot(stats['val acc history'], label='val') plt.title('Classification accuracy history') plt.xlabel('Epoch') plt.ylabel('Clasification accuracy') plt.legend() plt.show() Loss history 2.3 2.2 2.1 2.0 1.9 200 400 600 800 1000 Classification action of the Classification 0.300 train 0.275 0.250 0.225 0.200 0.175 0.150 0.125 0.100 1.5 2.5 4.0 0.0 0.5 1.0 2.0 3.0 3.5 Epoch In [11]: from cs682.vis_utils import visualize_grid # Visualize the weights of the network def show net weights(net): W1 = net.params['W1'] W1 = W1.reshape(32, 32, 3, -1).transpose(3, 0, 1, 2)plt.imshow(visualize_grid(W1, padding=3).astype('uint8')) plt.gca().axis('off') plt.show() show_net_weights(net) **Tune your hyperparameters** What's wrong?. Looking at the visualizations above, we see that the loss is decreasing more or less linearly, which seems to suggest that the learning rate may be too low. Moreover, there is no gap between the training and validation accuracy, suggesting that the model we used has low capacity, and that we should increase its size. On the other hand, with a very large model we would expect to see more overfitting, which would manifest itself as a very large gap between the training and validation accuracy. **Tuning**. Tuning the hyperparameters and developing intuition for how they affect the final performance is a large part of using Neural Networks, so we want you to get a lot of practice. Below, you should experiment with different values of the various hyperparameters, including hidden layer size, learning rate, numer of training epochs, and regularization strength. You might also consider tuning the learning rate decay, but you should be able to get good performance using the default value. Approximate results. You should be aim to achieve a classification accuracy of greater than 48% on the validation set. Our best network gets over 52% on the validation set. **Experiment**: You goal in this exercise is to get as good of a result on CIFAR-10 as you can, with a fully-connected Neural Network. Feel free implement your own techniques (e.g. PCA to reduce dimensionality, or adding dropout, or adding features to the solver, etc.). In [12]: best net = None # store the best model into this # TODO: Tune hyperparameters using the validation set. Store your best trained # # model in best net. # To help debug your network, it may help to use visualizations similar to the # # ones we used above; these visualizations will have significant qualitative # differences from the ones we saw above for the poorly tuned network. # Tweaking hyperparameters by hand can be fun, but you might find it useful to # # write code to sweep through possible combinations of hyperparameters # automatically like we did on the previous exercises. learning rates = [5e-4, 1e-3, 2.5e-3]regularization_strengths = [0.25, 0.5] batch_sizes = [100, 200] hidden_sizes = [50] best val = -1.0for lr in learning_rates: for reg in regularization strengths: for batch size in batch sizes: for hidden_size in hidden_sizes: print("Testing lr =", lr, " reg =", reg, "batch size =", batch size, "hidden size =", hidden size) net = TwoLayerNet(input_size, hidden_size, num_classes) # Train the network stats = net.train(X_train, y_train, X_val, y_val, num_iters=2000, batch_size=batch_size, learning rate=lr, learning rate decay=0.95, reg=reg, verbose=False) # Predict on the validation set val acc = (net.predict(X val) == y val).mean() y train pred = net.predict(X train) tr acc = np.mean(y train == y train pred) print('training accuracy: %f' % tr_acc) y val pred = net.predict(X val) val_acc = np.mean(y_val == y_val_pred) print('validation accuracy: %f' % val acc) if val acc > best val: print("---> Found new best net with params lr =", lr, " reg =", reg, "batch_size =", batch_size, "hidden size =", hidden size) best val = val acc best net = net END OF YOUR CODE Testing lr = 0.0005 reg = 0.25 batch_size = 100 hidden_size = 50 training accuracy: 0.491347 validation accuracy: 0.471000 ---> Found new best net with params lr = 0.0005 reg = 0.25 batch size = 100 hidden size = 50 Testing lr = 0.0005 reg = 0.25 batch size = 200 hidden size = 50 training accuracy: 0.496837 validation accuracy: 0.470000 Testing lr = 0.0005 reg = 0.5 batch_size = 100 hidden_size = 50 training accuracy: 0.475980 validation accuracy: 0.456000 Testing lr = 0.0005 reg = 0.5 batch_size = 200 hidden_size = 50 training accuracy: 0.488694 validation accuracy: 0.468000 Testing lr = 0.001 reg = 0.25 batch size = 100 hidden size = 50 training accuracy: 0.507204 validation accuracy: 0.464000 Testing lr = 0.001 reg = 0.25 batch size = 200 hidden size = 50 training accuracy: 0.519061 validation accuracy: 0.479000 ---> Found new best net with params lr = 0.001 reg = 0.25 batch_size = 200 hidden_size = 50 Testing lr = 0.001 reg = 0.5 batch size = 100 hidden size = 50 training accuracy: 0.488429 validation accuracy: 0.471000 Testing lr = 0.001 reg = 0.5 batch size = 200 hidden size = 50 training accuracy: 0.512898 validation accuracy: 0.488000 ---> Found new best net with params lr = 0.001 reg = 0.5 batch size = 200 hidden size = 50 Testing lr = 0.0025 reg = 0.25 batch_size = 100 hidden_size = 50 training accuracy: 0.427918 validation accuracy: 0.416000 Testing lr = 0.0025 reg = 0.25 batch_size = 200 hidden_size = 50 training accuracy: 0.492878 validation accuracy: 0.483000 Testing lr = 0.0025 reg = 0.5 batch_size = 100 hidden_size = 50 training accuracy: 0.441408 validation accuracy: 0.439000 Testing lr = 0.0025 reg = 0.5 batch_size = 200 hidden_size = 50 training accuracy: 0.498367 validation accuracy: 0.485000 In [125]: # visualize the weights of the best network show_net_weights(best_net) Run on the test set When you are done experimenting, you should evaluate your final trained network on the test set; you should get above 48%. In [124]: test_acc = (best_net.predict(X_test) == y_test).mean() print('Test accuracy: ', test_acc) Test accuracy: 0.498 **Inline Question** Now that you have trained a Neural Network classifier, you may find that your testing accuracy is much lower than the training accuracy. In what ways can we decrease this gap? Select all that apply. 1. Train on a larger dataset. 2. Add more hidden units. 3. Increase the regularization strength. 4. None of the above. Your answer: While my testing accuracy was similar to my training accuracy, these strategies all have the ability to increase testing accuracy by reducing bias. 1. Train on a larger dataset would make the nets parameters less dependant on each individual training example which would make the net more applicable to a larger variety of test data. 2. Adding more hidden units is less likely than the other strategies to make a difference on the test data, however it could possibly help testing accuracy by allowing the net to learn more parameters. 3. Increasing the regularization strength could help increase training accuracy because regularization is used to reduce bias by making it less dependant on training data.

Implementing a Neural Network

from cs682.classifiers.neural net import TwoLayerNet

plt.rcParams['image.interpolation'] = 'nearest'

plt.rcParams['figure.figsize'] = (10.0, 8.0) # set default size of plots

see http://stackoverflow.com/questions/1907993/autoreload-of-modules-in-ipython

return np.max(np.abs(x - y) / (np.maximum(1e-8, np.abs(x) + np.abs(y))))

from __future__ import print function

import matplotlib.pyplot as plt

plt.rcParams['image.cmap'] = 'gray'

for auto-reloading external modules

""" returns relative error """

model that we will use to develop your implementation.

In [3]: # Create a small net and some toy data to check your implementations. # Note that we set the random seed for repeatable experiments.

In [2]: # A bit of setup

import numpy as np

%matplotlib inline

%load ext autoreload

def rel error(x, y):

%autoreload 2

input_size = 4
hidden_size = 10
num_classes = 3
num_inputs = 5

def init_toy_model():

np.random.seed(0)

In this exercise we will develop a neural network with fully-connected layers to perform classification, and test it out on the CIFAR-10 dataset.

We will use the class TwoLayerNet in the file cs682/classifiers/neural net.py to represent instances of our network. The network parameters are

stored in the instance variable self.params where keys are string parameter names and values are numpy arrays. Below, we initialize toy data and a toy

```
We have seen that we can achieve reasonable performance on an image classification task by training a linear classifier on the pixels of the input image. In this
          exercise we will show that we can improve our classification performance by training linear classifiers not on raw pixels but on features that are computed from
          the raw pixels.
          All of your work for this exercise will be done in this notebook.
In [105]: from __future__ import print_function
          import random
          import numpy as np
          from cs682.data_utils import load_CIFAR10
          import matplotlib.pyplot as plt
          %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'
          # for auto-reloading extenrnal modules
          # 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
          Load data
          Similar to previous exercises, we will load CIFAR-10 data from disk.
In [106]: from cs682.features import color_histogram_hsv, hog_feature
          def get_CIFAR10_data(num_training=49000, num_validation=1000, num_test=1000):
              # Load the raw CIFAR-10 data
              cifar10_dir = 'cs682/datasets/cifar-10-batches-py'
              X_train, y_train, X_test, y_test = load_CIFAR10(cifar10_dir)
              # Subsample the data
              mask = list(range(num_training, num_training + num_validation))
              X_val = X_train[mask]
              y_val = y_train[mask]
              mask = list(range(num_training))
              X_train = X_train[mask]
              y_train = y_train[mask]
              mask = list(range(num_test))
              X_test = X_test[mask]
              y test = y test[mask]
              return X_train, y_train, X_val, y_val, X_test, y_test
          # Cleaning up variables to prevent loading data multiple times (which may cause memory issue)
             del X train, y train
             del X_test, y_test
             print('Clear previously loaded data.')
             pass
          X_train, y_train, X_val, y_val, X_test, y_test = get_CIFAR10_data()
          Clear previously loaded data.
          Extract Features
          For each image we will compute a Histogram of Oriented Gradients (HOG) as well as a color histogram using the hue channel in HSV color space. We form our
          final feature vector for each image by concatenating the HOG and color histogram feature vectors.
          Roughly speaking, HOG should capture the texture of the image while ignoring color information, and the color histogram represents the color of the input
          image while ignoring texture. As a result, we expect that using both together ought to work better than using either alone. Verifying this assumption would be a
          good thing to try for your interests.
          The hog_feature and color_histogram_hsv functions both operate on a single image and return a feature vector for that image. The extract_features
          function takes a set of images and a list of feature functions and evaluates each feature function on each image, storing the results in a matrix where each
          column is the concatenation of all feature vectors for a single image.
In [107]: from cs682.features import *
          num color bins = 10 # Number of bins in the color histogram
          feature_fns = [hog_feature, lambda img: color_histogram_hsv(img, nbin=num_color_bins)]
          X_train_feats = extract_features(X_train, feature_fns, verbose=True)
          X val feats = extract features(X val, feature fns)
          X test feats = extract features(X test, feature fns)
          # Preprocessing: Subtract the mean feature
          mean_feat = np.mean(X_train_feats, axis=0, keepdims=True)
          X train feats -= mean feat
          X val feats -= mean feat
          X_test_feats -= mean_feat
          # Preprocessing: Divide by standard deviation. This ensures that each feature
          # has roughly the same scale.
          std feat = np.std(X train feats, axis=0, keepdims=True)
          X train feats /= std feat
          X val feats /= std feat
          X test feats /= std feat
          # Preprocessing: Add a bias dimension
          X train feats = np.hstack([X train feats, np.ones((X_train_feats.shape[0], 1))])
          X val feats = np.hstack([X val feats, np.ones((X val feats.shape[0], 1))])
          X test feats = np.hstack([X test feats, np.ones((X test feats.shape[0], 1))])
          Done extracting features for 1000 / 49000 images
          Done extracting features for 2000 / 49000 images
          Done extracting features for 3000 / 49000 images
          Done extracting features for 4000 / 49000 images
          Done extracting features for 5000 / 49000 images
          Done extracting features for 6000 / 49000 images
          Done extracting features for 7000 / 49000 images
          Done extracting features for 8000 / 49000 images
          Done extracting features for 9000 / 49000 images
          Done extracting features for 10000 / 49000 images
          Done extracting features for 11000 / 49000 images
          Done extracting features for 12000 / 49000 images
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          Done extracting features for 44000 / 49000 images
          Done extracting features for 45000 / 49000 images
          Done extracting features for 46000 / 49000 images
          Done extracting features for 47000 / 49000 images
          Done extracting features for 48000 / 49000 images
          Train SVM on features
          Using the multiclass SVM code developed earlier in the assignment, train SVMs on top of the features extracted above; this should achieve better results than
          training SVMs directly on top of raw pixels.
In [130]: # Use the validation set to tune the learning rate and regularization strength
          from cs682.classifiers.linear_classifier import LinearSVM
          learning_rates = [5e-7, 7.5e-7]
          regularization strengths = [1.5e4, 2.5e4]
          results = {}
          best val = -1
          best_svm = None
          # TODO:
          # Use the validation set to set the learning rate and regularization strength. #
          # This should be identical to the validation that you did for the SVM; save
          # the best trained classifer in best_svm. You might also want to play
          # with different numbers of bins in the color histogram. If you are careful
          # you should be able to get accuracy of near 0.44 on the validation set.
          for lr in learning_rates:
              for reg in regularization strengths:
                  print("Testing LR=", lr, " reg=", reg)
                  svm = LinearSVM()
                  svm.train(X train_feats, y_train, learning_rate=lr, reg=reg,
                                num_iters=2000, verbose=False)
                  y train pred = svm.predict(X train feats)
                  tr_acc = np.mean(y_train == y_train_pred)
                  print('training accuracy: %f' % tr_acc)
                  y_val_pred = svm.predict(X_val_feats)
                  val_acc = np.mean(y_val == y_val_pred)
                  print('validation accuracy: %f' % val acc)
                  if val_acc > best_val:
                      best val = val acc
                      best svm = svm
                  results[(lr, reg)] = tr_acc, val_acc
          END OF YOUR CODE
          # Print out results.
          for lr, reg in sorted(results):
              train accuracy, val accuracy = results[(lr, reg)]
              print('lr %e reg %e train accuracy: %f val accuracy: %f' % (
                          lr, reg, train_accuracy, val_accuracy))
          print('best validation accuracy achieved during cross-validation: %f' % best val)
          Testing LR= 5e-07 reg= 15000.0
          training accuracy: 0.409592
          validation accuracy: 0.403000
          Testing LR= 5e-07 reg= 25000.0
          training accuracy: 0.413633
          validation accuracy: 0.424000
          Testing LR= 7.5e-07 reg= 15000.0
          training accuracy: 0.415755
          validation accuracy: 0.411000
          Testing LR= 7.5e-07 reg= 25000.0
          training accuracy: 0.412224
          validation accuracy: 0.413000
          lr 5.000000e-07 reg 1.500000e+04 train accuracy: 0.409592 val accuracy: 0.403000
          lr 5.000000e-07 reg 2.500000e+04 train accuracy: 0.413633 val accuracy: 0.424000
          lr 7.500000e-07 reg 1.500000e+04 train accuracy: 0.415755 val accuracy: 0.411000
          lr 7.500000e-07 reg 2.500000e+04 train accuracy: 0.412224 val accuracy: 0.413000
          best validation accuracy achieved during cross-validation: 0.424000
In [131]: # Evaluate your trained SVM on the test set
          y_test_pred = best_svm.predict(X_test_feats)
          test_accuracy = np.mean(y_test == y_test_pred)
          print(test accuracy)
          0.425
 In [ ]:
In [132]: # An important way to gain intuition about how an algorithm works is to
          # visualize the mistakes that it makes. In this visualization, we show examples
          # of images that are misclassified by our current system. The first column
          # shows images that our system labeled as "plane" but whose true label is
          # something other than "plane".
          examples per class = 8
          classes = ['plane', 'car', 'bird', 'cat', 'deer', 'dog', 'frog', 'horse', 'ship', 'truck']
          for cls, cls name in enumerate(classes):
              idxs = np.where((y_test != cls) & (y_test_pred == cls))[0]
              idxs = np.random.choice(idxs, examples per class, replace=False)
              for i, idx in enumerate(idxs):
                  plt.subplot(examples_per_class, len(classes), i * len(classes) + cls + 1)
                  plt.imshow(X_test[idx].astype('uint8'))
                  plt.axis('off')
                  if i == 0:
                      plt.title(cls_name)
          plt.show()
                                                                         truck
            plane
                                       deer
                                              dog
                                                     frog
                                                           horse
                                                                   ship
                 Inline question 1:
          Describe the misclassification results that you see. Do they make sense?
          While some of the misclassifications are more difficult to explain, many of them have features similar to what the class looks like which could cause them to be
          misclassified. For example some birds are misclassified as planes because they have similar texture features (such as outspread wings) that could be confused
          for a plane. Similar features like this can cause the scores for a specific class to be higher than the correct class.
          Neural Network on image features
          Earlier in this assignment we saw that training a two-layer neural network on raw pixels achieved better classification performance than linear classifiers on raw
          pixels. In this notebook we have seen that linear classifiers on image features outperform linear classifiers on raw pixels.
          For completeness, we should also try training a neural network on image features. This approach should outperform all previous approaches: you should easily
          be able to achieve over 55% classification accuracy on the test set; our best model achieves about 60% classification accuracy.
In [30]: # Preprocessing: Remove the bias dimension
          # Make sure to run this cell only ONCE
          print(X train feats.shape)
          X_train_feats = X_train_feats[:, :-1]
          X val feats = X val feats[:, :-1]
          X test feats = X test feats[:, :-1]
          print(X_train_feats.shape)
          (49000, 155)
          (49000, 154)
In [76]: from cs682.classifiers.neural net import TwoLayerNet
          input dim = X train feats.shape[1]
          hidden dim = 50
          num_classes = 10
          net = TwoLayerNet(input_dim, hidden_dim, num_classes)
          best net = None
          # TODO: Train a two-layer neural network on image features. You may want to
          # cross-validate various parameters as in previous sections. Store your best #
          # model in the best net variable.
          best_net = TwoLayerNet(input_dim, hidden_dim, num_classes)
          # Train the network
          stats = best net.train(X train feats, y train, X val feats, y val,
                  num_iters=4000, batch_size=400,
                  learning_rate=0.15, learning_rate_decay=0.98,
                  reg=1e-5, verbose=True)
          y_train_pred = best_net.predict(X_train_feats)
          tr acc = np.mean(y train == y train pred)
          print('training accuracy: %f' % tr_acc)
          y val pred = best net.predict(X val feats)
          val_acc = np.mean(y_val == y_val_pred)
          print('validation accuracy: %f' % val acc)
          END OF YOUR CODE
          iteration 0 / 4000: loss 2.302585
          iteration 100 / 4000: loss 2.300256
          iteration 200 / 4000: loss 1.867220
          iteration 300 / 4000: loss 1.617228
          iteration 400 / 4000: loss 1.407140
          iteration 500 / 4000: loss 1.415947
          iteration 600 / 4000: loss 1.359692
          iteration 700 / 4000: loss 1.439941
          iteration 800 / 4000: loss 1.307052
          iteration 900 / 4000: loss 1.289611
          iteration 1000 / 4000: loss 1.359644
          iteration 1100 / 4000: loss 1.289393
          iteration 1200 / 4000: loss 1.265420
          iteration 1300 / 4000: loss 1.224796
          iteration 1400 / 4000: loss 1.202857
          iteration 1500 / 4000: loss 1.225035
          iteration 1600 / 4000: loss 1.196401
          iteration 1700 / 4000: loss 1.274015
          iteration 1800 / 4000: loss 1.180586
          iteration 1900 / 4000: loss 1.183681
          iteration 2000 / 4000: loss 1.237354
          iteration 2100 / 4000: loss 1.217919
          iteration 2200 / 4000: loss 1.099059
          iteration 2300 / 4000: loss 1.178842
          iteration 2400 / 4000: loss 1.159751
          iteration 2500 / 4000: loss 1.107960
          iteration 2600 / 4000: loss 1.069555
```

iteration 2700 / 4000: loss 1.085661
iteration 2800 / 4000: loss 1.165164
iteration 2900 / 4000: loss 1.111434
iteration 3000 / 4000: loss 1.127369
iteration 3100 / 4000: loss 1.124869
iteration 3200 / 4000: loss 1.170346
iteration 3300 / 4000: loss 1.140856
iteration 3400 / 4000: loss 1.081912
iteration 3500 / 4000: loss 1.113889
iteration 3600 / 4000: loss 1.121366
iteration 3700 / 4000: loss 1.121366
iteration 3800 / 4000: loss 1.065460
iteration 3900 / 4000: loss 1.071608

training accuracy: 0.620898 validation accuracy: 0.574000

print(test_acc)

0.565

to get more than 55% accuracy.

In [75]: # Run your best neural net classifier on the test set. You should be able

test_acc = (best_net.predict(X_test_feats) == y_test).mean()

Image features exercise

more details see the assignments page on the course website.