knn

September 29, 2021

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.

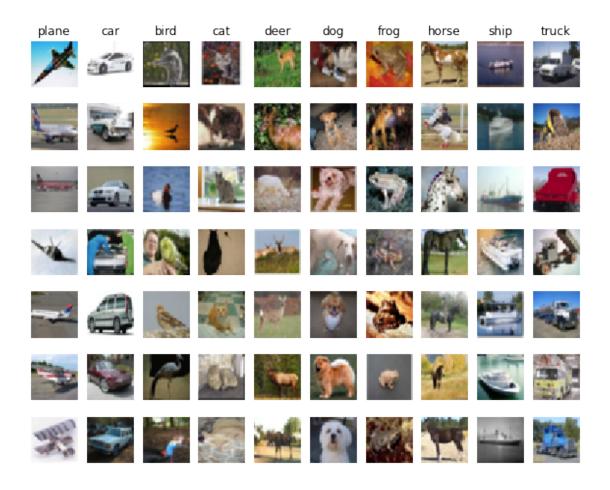
The kNN classifier consists of two stages:

- 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.

```
[1]: # 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
      \rightarrownotebook
     # 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/
     \rightarrow autoreload-of-modules-in-ipython
     %load ext autoreload
     %autoreload 2
```

```
[2]: # 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,)
[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', _
     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()
```



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

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

```
[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)

```
[6]: 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 \mathbf{Ntr} training examples and \mathbf{Nte} test examples, this stage should result in a $\mathbf{Nte} \times \mathbf{Ntr}$ matrix where each element (i,j) is the distance between the i-th test and j-th train example.

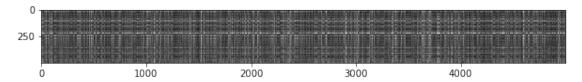
First, open cs682/classifiers/k_nearest_neighbor.py and implement the function compute_distances_two_loops that uses a (very inefficient) double loop over all pairs of (test, train) examples and computes the distance matrix one element at a time.

```
[7]: # 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)

```
[8]: # 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()
```



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.

- What in the data is the cause behind the distinctly bright rows? The test datas that are very different to all train data.
- What causes the columns? The train datas that are very different to all test data.

```
[9]: # 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:

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

Got 139 / 500 correct => accuracy: 0.278000

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

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: 1, 2, 3

Your explanation:

Because with k-nn we are only finding the k nearest datas, and with all three conditions, even though the value will change, the relativity would stay the same, we will still get the same k nearest datas.

```
[11]: # 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,
\( \to reshape \)
# 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')</pre>
```

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

```
[12]: # 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')
else:
    print('Uh-oh! The distance matrices are different')</pre>
```

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

```
[13]: # Let's compare how fast the implementations are
      def time_function(f, *args):
          11 11 11
          Call a function f with args and return the time (in seconds) that it took_{\sqcup}
       \rightarrow 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 \underline{\mbox{\ }} \rightarrow implementation
```

```
Two loop version took 33.363664 seconds One loop version took 25.649612 seconds No loop version took 0.124133 seconds
```

1.0.1 Cross-validation

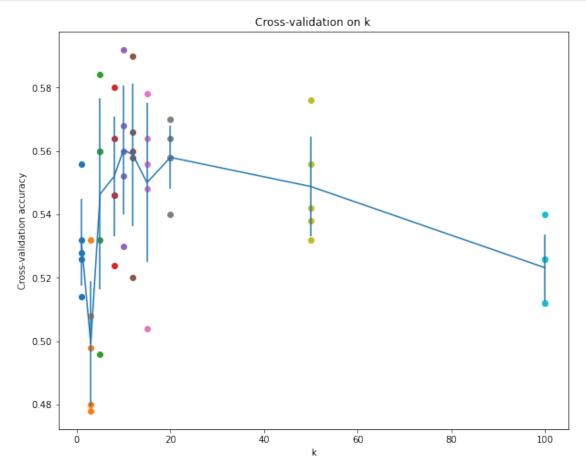
We have implemented the k-Nearest Neighbor classifier but we set the value k=5 arbitrarily. We will now determine the best value of this hyperparameter with cross-validation.

```
[14]: num folds = 5
    k_{choices} = [1, 3, 5, 8, 10, 12, 15, 20, 50, 100]
    X_train_folds = []
    y_train_folds = []
    # TODO:
    # 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.
    # Your code
    train_data = np.c_[X_train, y_train]
    train_data_folds = np.array_split(train_data, num_folds)
    for i in range(num folds):
       X_train_folds.append(train_data_folds[i][:, :-1])
       y_train_folds.append(train_data_folds[i][:, -1].astype(int))
    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.
# Your code
for k in k_choices:
    accuracies = []
    for i in range(num_folds):
       X_train_temp = X_train_folds[:]
       X_test_temp = X_train_temp.pop(i)
       X_train_temp = np.concatenate(X_train_temp)
       y_train_temp = y_train_folds[:]
       y_test_temp = y_train_temp.pop(i)
       y_train_temp = np.concatenate(y_train_temp)
       classifier = KNearestNeighbor()
       classifier.train(X_train_temp, y_train_temp)
       dists_temp = classifier.compute_distances_no_loops(X_test_temp)
       y_test_temp_pred = classifier.predict_labels(dists_temp, k)
       num_correct = np.sum(y_test_temp_pred == y_test_temp)
       accuracy = float(num_correct) / num_test
       accuracies.append(accuracy)
    k to accuracies[k] = accuracies
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 = 0.526000
k = 1, accuracy = 0.514000
k = 1, accuracy = 0.528000
k = 1, accuracy = 0.556000
k = 1, accuracy = 0.532000
k = 3, accuracy = 0.478000
k = 3, accuracy = 0.498000
k = 3, accuracy = 0.480000
k = 3, accuracy = 0.532000
k = 3, accuracy = 0.508000
k = 5, accuracy = 0.496000
k = 5, accuracy = 0.532000
k = 5, accuracy = 0.560000
k = 5, accuracy = 0.584000
```

```
k = 5, accuracy = 0.560000
     k = 8, accuracy = 0.524000
     k = 8, accuracy = 0.564000
     k = 8, accuracy = 0.546000
     k = 8, accuracy = 0.580000
     k = 8, accuracy = 0.546000
     k = 10, accuracy = 0.530000
     k = 10, accuracy = 0.592000
     k = 10, accuracy = 0.552000
     k = 10, accuracy = 0.568000
     k = 10, accuracy = 0.560000
     k = 12, accuracy = 0.520000
     k = 12, accuracy = 0.590000
     k = 12, accuracy = 0.558000
     k = 12, accuracy = 0.566000
     k = 12, accuracy = 0.560000
     k = 15, accuracy = 0.504000
     k = 15, accuracy = 0.578000
     k = 15, accuracy = 0.556000
     k = 15, accuracy = 0.564000
     k = 15, accuracy = 0.548000
     k = 20, accuracy = 0.540000
     k = 20, accuracy = 0.558000
     k = 20, accuracy = 0.558000
     k = 20, accuracy = 0.564000
     k = 20, accuracy = 0.570000
     k = 50, accuracy = 0.542000
     k = 50, accuracy = 0.576000
     k = 50, accuracy = 0.556000
     k = 50, accuracy = 0.538000
     k = 50, accuracy = 0.532000
     k = 100, accuracy = 0.512000
     k = 100, accuracy = 0.540000
     k = 100, accuracy = 0.526000
     k = 100, accuracy = 0.512000
     k = 100, accuracy = 0.526000
[15]: # 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()
```



```
[16]: # 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. 2. The test error of a 1-NN will always be better than that of a 5-NN. 3. The decision boundary of the k-NN classifier is linear. 4. The time needed to classify a test example with the k-NN classifier grows with the size of the training set. 5. None of the above.

Your Answer:

1, 4

Your explanation:

- 1. Since when training 1-NN, the nearest point would always be itself, the error would always be zero, while 5-NN would have an error that is greater or equal to 0.
- 2. This is because in order to find the k nearest neighbors, we would have to go through every single data in the training set.

svm

September 29, 2021

1 Multiclass Support Vector Machine 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.

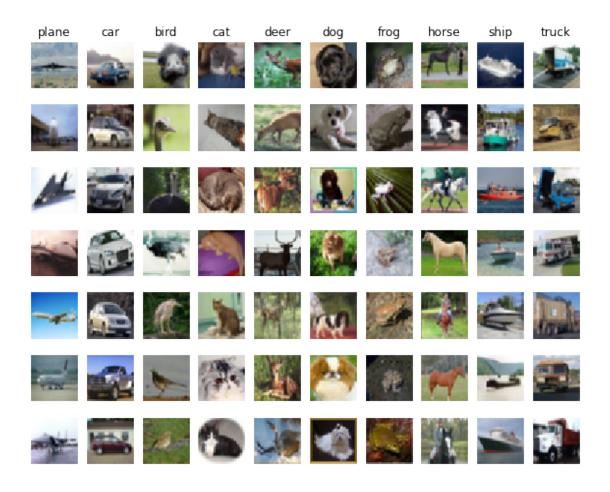
In this exercise you will:

- implement a fully-vectorized loss function for the SVM
- implement the fully-vectorized expression for its analytic gradient
- check your implementation using 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

```
[1]: # 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/
     \rightarrow autoreload-of-modules-in-ipython
     %load_ext autoreload
     %autoreload 2
```

1.1 CIFAR-10 Data Loading and Preprocessing

```
[2]: # 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,)
[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', u
     ⇔'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()
```



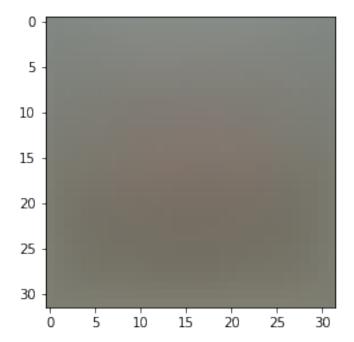
```
[4]: | # Split the data into train, val, and test sets. In addition we will
     # create a small development set as a subset of the training data;
     # we can use this for development so our code runs faster.
     num_training = 49000
     num_validation = 1000
     num_test = 1000
     num_dev = 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,)
[5]: # 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)
[6]: # 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]



```
[7]: # 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
```

```
[8]: # 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)

1.2 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 svm_loss_naive which uses for loops to evaluate the multiclass SVM loss function.

```
[9]: # 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) * 0.0001

loss, grad = svm_loss_naive(W, X_dev, y_dev, 0.000005)
    print('loss: %f' % (loss, ))
```

loss: 9.234847

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 svm_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:

```
[10]: # 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 |
      \rightarrow 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)
      print("now w/ regularization")
      # 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: 0.910133 analytic: 0.910133, relative error: 6.432312e-10 numerical: -24.208321 analytic: -24.208321, relative error: 1.568762e-11 numerical: 7.626115 analytic: 7.626115, relative error: 8.662189e-11

```
numerical: 11.575777 analytic: 11.575777, relative error: 1.079016e-11
numerical: 22.141280 analytic: 22.141280, relative error: 1.893267e-11
numerical: 0.590871 analytic: 0.590871, relative error: 5.162115e-10
numerical: 36.153483 analytic: 36.153483, relative error: 4.742293e-12
numerical: 3.947682 analytic: 3.947682, relative error: 6.284088e-11
numerical: -3.631061 analytic: -3.631061, relative error: 6.378823e-11
numerical: 13.608739 analytic: 13.608739, relative error: 1.422914e-11
now w/ regularization
numerical: -8.377741 analytic: -8.377741, relative error: 3.035389e-11
numerical: 16.888644 analytic: 16.888644, relative error: 7.118199e-12
numerical: 26.754491 analytic: 26.754491, relative error: 1.798586e-11
numerical: -24.216629 analytic: -24.216629, relative error: 1.694647e-11
numerical: 13.463277 analytic: 13.463277, relative error: 6.773779e-12
numerical: 33.023519 analytic: 33.023519, relative error: 3.667877e-12
numerical: -44.411010 analytic: -44.411010, relative error: 1.062990e-12
numerical: 4.296020 analytic: 4.296020, relative error: 5.511394e-11
numerical: -9.370645 analytic: -9.370645, relative error: 5.694974e-11
numerical: 21.512107 analytic: 21.512107, relative error: 2.287321e-13
```

1.2.1 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: fill this in.

Yes it is possible. When x would let f(x) be exactly, which is when $s_j - s_j - s_j = -1$. Because at that point the analytic gradient would be 0, while the numerical would be a little greater than zero, due to the fact that f(x+h) would be a bit greater than zero, while f(x-h) be zero.

Naive loss: 9.234847e+00 computed in 0.134170s Vectorized loss: 9.234847e+00 computed in 0.005889s difference: 0.000000

```
[12]: # Complete the implementation of sum 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)
      toc = time.time()
      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 0.125564s Vectorized loss and gradient: computed in 0.001873s difference: 0.000000

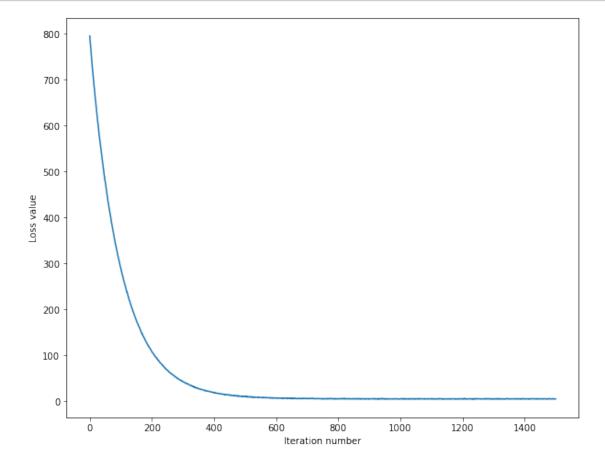
1.2.2 Stochastic Gradient Descent

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.

iteration 0 / 1500: loss 795.237026
iteration 100 / 1500: loss 289.900307
iteration 200 / 1500: loss 107.701745
iteration 300 / 1500: loss 42.840852
iteration 400 / 1500: loss 18.888930

```
iteration 500 / 1500: loss 10.398202 iteration 600 / 1500: loss 7.375045 iteration 700 / 1500: loss 5.986475 iteration 800 / 1500: loss 5.628352 iteration 900 / 1500: loss 5.183372 iteration 1000 / 1500: loss 5.421224 iteration 1100 / 1500: loss 5.267346 iteration 1200 / 1500: loss 5.807447 iteration 1300 / 1500: loss 5.337107 iteration 1400 / 1500: loss 5.207017 That took 4.512303s
```

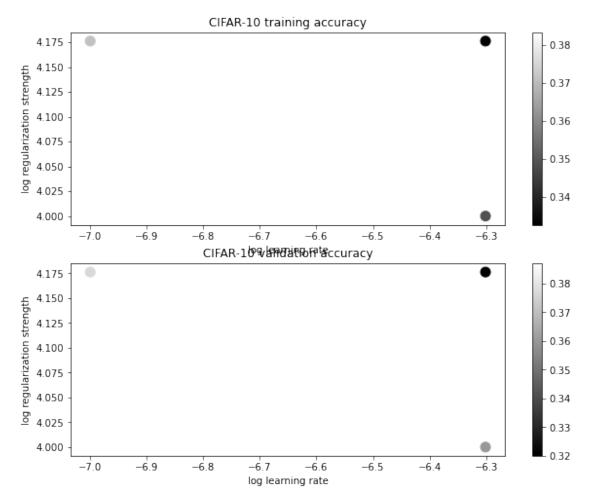
```
[17]: # A useful debugging strategy is to plot the loss as a function of
    # iteration number:
    plt.plot(loss_hist)
    plt.xlabel('Iteration number')
    plt.ylabel('Loss value')
    plt.show()
```



```
[21]: # 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: 0.365939
     validation accuracy: 0.377000
[63]: # 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 = [8e-8, 9e-8, 1e-7]
     regularization_strengths = [9e3, 1e4, 1.5e4]
     # 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
      \rightarrow 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.
     # Your code
     for lr in learning_rates:
         for reg in regularization_strengths:
             svm = LinearSVM()
             svm.train(X_train, y_train, lr, reg, num_iters=3000)
             y_train_pred = svm.predict(X_train)
             train_accuracy = np.mean(y_train == y_train_pred)
```

```
y_val_pred = svm.predict(X_val)
            val_accuracy = np.mean(y_val == y_val_pred)
            results[(lr, reg)] = (train_accuracy, val_accuracy)
             if val_accuracy > best_val:
                best_svm = svm
                best_val = val_accuracy
     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)
    lr 8.000000e-08 reg 9.000000e+03 train accuracy: 0.386143 val accuracy: 0.385000
    lr 8.000000e-08 reg 1.000000e+04 train accuracy: 0.384878 val accuracy: 0.388000
    lr 8.000000e-08 reg 1.500000e+04 train accuracy: 0.383388 val accuracy: 0.396000
    lr 9.000000e-08 reg 9.000000e+03 train accuracy: 0.383143 val accuracy: 0.384000
    lr 9.000000e-08 reg 1.000000e+04 train accuracy: 0.384633 val accuracy: 0.389000
    lr 9.000000e-08 reg 1.500000e+04 train accuracy: 0.377061 val accuracy: 0.392000
    lr 1.000000e-07 reg 9.000000e+03 train accuracy: 0.390327 val accuracy: 0.391000
    lr 1.000000e-07 reg 1.000000e+04 train accuracy: 0.381612 val accuracy: 0.399000
    lr 1.000000e-07 reg 1.500000e+04 train accuracy: 0.380327 val accuracy: 0.388000
    best validation accuracy achieved during cross-validation: 0.399000
[58]: # 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()
```



```
[59]: # Evaluate the best sum 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: 0.380000





1.2.3 Inline question 2:

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

Your answer: fill this in

Since we define our scores being a dot product between the data and weights, in order to get the highest scores for the correct class, the weight vector would be some form of parallel vector to the train data. As a result each weight for each class would look like the class data it wants to predict. Which could be observed in the visualizations above, we could see how both the plane and ship have large amounts of blue, since plane and ship usually have the sky and sea as background. We could also see how the car and truck do look alike while the car has a more rounded feature while the truck has a more square ish appearance.

softmax

September 29, 2021

1 Softmax 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.

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

```
[2]: def get_CIFAR10_data(num_training=49000, num_validation=1000, num_test=1000, u → 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.
```

```
HHHH
    # 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_dev = np.reshape(X_dev, (X_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,)

1.1 Softmax Classifier

Your code for this section will all be written inside cs682/classifiers/softmax.py.

```
[3]: # 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: 2.373042

sanity check: 2.302585

1.2 Inline Question 1:

Why do we expect our loss to be close to $-\log(0.1)$? Explain briefly.**

Your answer: Fill this in

This is because we initiate our weights randomly, and we have ten different classes. Since the weights are random, it is most likely that every class would have the same score before training. This means the probabilities would be 1/num_classes before train, which in our case is 1/10=0.1.

```
[4]: # 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: 0.371524 analytic: 0.371524, relative error: 1.641621e-07
numerical: 0.664583 analytic: 0.664583, relative error: 3.825252e-08
numerical: -0.670274 analytic: -0.670274, relative error: 8.670363e-09
numerical: 5.662179 analytic: 5.662178, relative error: 2.007403e-08
numerical: -1.291537 analytic: -1.291537, relative error: 2.540536e-08
numerical: -1.068029 analytic: -1.068029, relative error: 9.433747e-09
numerical: 0.889227 analytic: 0.889226, relative error: 9.961711e-08
numerical: -0.699912 analytic: -0.699913, relative error: 5.477711e-08
numerical: 0.102961 analytic: 0.102961, relative error: 9.655173e-08
numerical: 0.414729 analytic: 0.414729, relative error: 7.764006e-08
numerical: -0.363680 analytic: -0.363680, relative error: 3.198507e-08
numerical: 0.465865 analytic: 0.465865, relative error: 8.059718e-08
numerical: -1.091263 analytic: -1.091263, relative error: 9.133470e-08
numerical: -1.925673 analytic: -1.925674, relative error: 1.588992e-08
numerical: -3.747741 analytic: -3.747741, relative error: 1.872167e-08
numerical: 0.269944 analytic: 0.269944, relative error: 1.539578e-07
numerical: 0.448560 analytic: 0.448560, relative error: 2.354892e-08
numerical: 1.426694 analytic: 1.426694, relative error: 4.527782e-08
numerical: -3.059476 analytic: -3.059476, relative error: 2.512482e-09
```

numerical: 0.287101 analytic: 0.287101, relative error: 2.864537e-07

```
[5]:  # 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
\rightarrowshould 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: 2.373042e+00 computed in 0.093381s vectorized loss: 2.373042e+00 computed in 0.004147s

Loss difference: 0.000000 Gradient difference: 0.000000

```
[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 = [8e-8, 9e-8, 1e-7]
    regularization_strengths = [9e3, 1e4, 1.5e4]
    # 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.
    # Your code
    for lr in learning_rates:
       for reg in regularization_strengths:
          softmax = Softmax()
```

```
softmax.train(X_train, y_train, lr, reg, num_iters=3000)
           y_train_pred = softmax.predict(X_train)
           train_accuracy = np.mean(y_train == y_train_pred)
           y_val_pred = softmax.predict(X_val)
           val_accuracy = np.mean(y_val == y_val_pred)
           results[(lr, reg)] = (train_accuracy, val_accuracy)
           if val_accuracy > best_val:
               best softmax = softmax
               best_val = val_accuracy
    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)
   lr 8.000000e-08 reg 9.000000e+03 train accuracy: 0.359837 val accuracy: 0.370000
   lr 8.000000e-08 reg 1.000000e+04 train accuracy: 0.357898 val accuracy: 0.368000
   lr 8.000000e-08 reg 1.500000e+04 train accuracy: 0.346408 val accuracy: 0.363000
   lr 9.000000e-08 reg 9.000000e+03 train accuracy: 0.357245 val accuracy: 0.377000
   lr 9.000000e-08 reg 1.000000e+04 train accuracy: 0.358612 val accuracy: 0.372000
   lr 9.000000e-08 reg 1.500000e+04 train accuracy: 0.337245 val accuracy: 0.351000
   lr 1.000000e-07 reg 9.000000e+03 train accuracy: 0.361286 val accuracy: 0.372000
   lr 1.000000e-07 reg 1.000000e+04 train accuracy: 0.358367 val accuracy: 0.371000
   lr 1.000000e-07 reg 1.500000e+04 train accuracy: 0.344653 val accuracy: 0.355000
   best validation accuracy achieved during cross-validation: 0.377000
[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: 0.370000

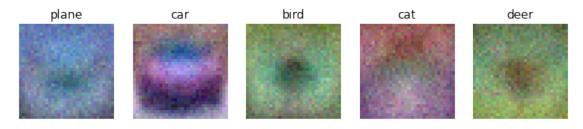
Inline Question - True or False

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: True

Your explanation:

We could find data that when in SVM it's score for all other classes would lead to $S_j - S_yi < -1$, which then the loss would only be 0. However, when in softmax, since e^x could never be zero, $e^(s_j)$ could never be zero which means $e^{(s_yi)/\text{sum}_j(e}(s_j))$ would always be < 1, therefore $-\log(e^{(s_yi)/\text{sum}_j(e}(s_j)))$ would never be 0.





two_layer_net

September 29, 2021

1 Implementing a Neural Network

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.

```
[1]: # A bit of setup
     from __future__ import print_function
     import numpy as np
     import matplotlib.pyplot as plt
     from cs682.classifiers.neural net import TwoLayerNet
     %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 external modules
     # see http://stackoverflow.com/questions/1907993/
     \rightarrow autoreload-of-modules-in-ipython
     %load ext autoreload
     %autoreload 2
     def rel_error(x, y):
         """ returns relative error """
         return np.max(np.abs(x - y) / (np.maximum(1e-8, np.abs(x) + np.abs(y))))
```

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 model that we will use to develop your implementation.

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

input_size = 4
```

```
hidden_size = 10
num_classes = 3
num_inputs = 5

def init_toy_model():
    np.random.seed(0)
    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, y = init_toy_data()
```

2 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.

```
[3]: 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
```

3 Forward pass: compute loss

In the same function, implement the second part that computes the data and regularizaion loss.

```
[4]: loss, _ = net.loss(X, y, reg=0.05)
    correct_loss = 1.30378789133

# should be very small, we get < 1e-12
    print('Difference between your loss and correct loss:')
    print(np.sum(np.abs(loss - correct_loss)))</pre>
```

Difference between your loss and correct loss: 1.794120407794253e-13

4 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:

```
[5]: 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], u
→verbose=False)

print('%s max relative error: %e' % (param_name, rel_error(param_grad_num, u
→grads[param_name])))

W1 max relative error: 3.561318e-09
b1 max relative error: 2.738423e-09
W2 max relative error: 3.440708e-09
```

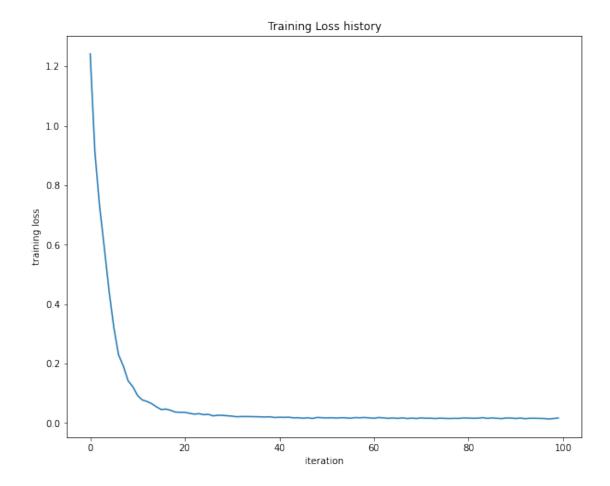
5 Train the network

b2 max relative error: 3.865039e-11

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.

Final training loss: 0.01714960793873202



6 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.

```
[7]: 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 data snape: (49000, 3072)
Train labels shape: (49000,)
Validation data shape: (1000, 3072)
Validation labels shape: (1000,)

```
Test data shape: (1000, 3072)
Test labels shape: (1000,)
```

7 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.

```
iteration 0 / 1000: loss 2.302954
iteration 100 / 1000: loss 2.302550
iteration 200 / 1000: loss 2.297648
iteration 300 / 1000: loss 2.259602
iteration 400 / 1000: loss 2.204170
iteration 500 / 1000: loss 2.118565
iteration 600 / 1000: loss 2.051535
iteration 700 / 1000: loss 1.988466
iteration 800 / 1000: loss 2.006591
iteration 900 / 1000: loss 1.951473
Validation accuracy: 0.287
```

8 Debug the training

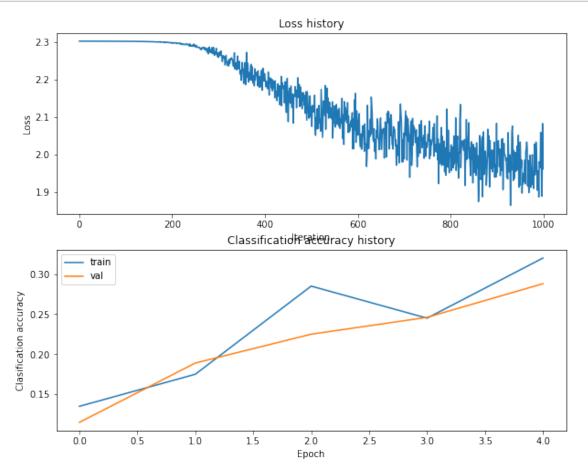
With the default parameters we provided above, you should get a validation accuracy of about 0.29 on the validation set. This isn't very good.

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.

```
[9]: # 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()
```

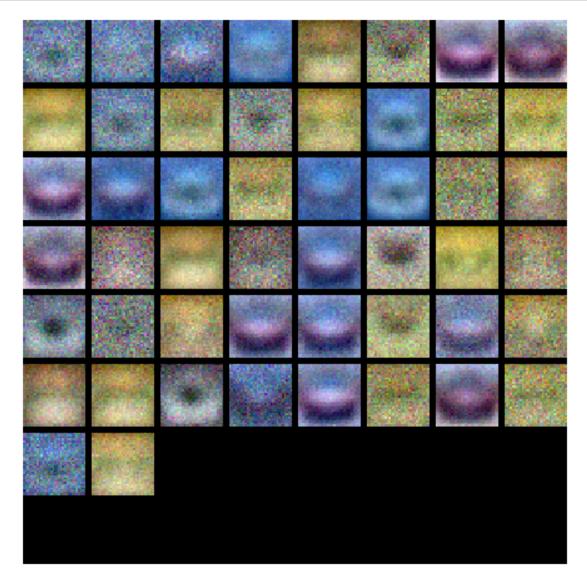


```
[10]: from cs682.vis_utils import visualize_grid

# 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)
```



9 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.

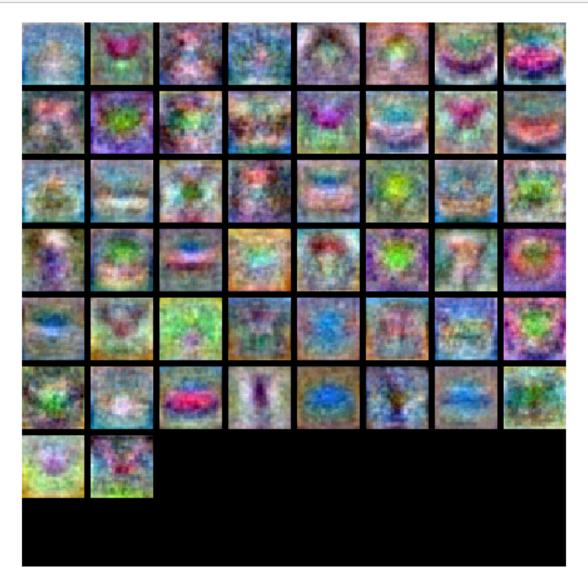
Experiment: You goal in this exercise is to get as good of a result on CIFAR-10 as you can, with a fully-connected Neural Network. Feel free implement your own techniques (e.g. PCA to reduce dimensionality, or adding dropout, or adding features to the solver, etc.).

```
[11]: best_net = None # store the best model into this
     # TODO: Tune hyperparameters using the validation set. Store your best trained \Box
     →#
     # model in best net.
     →#
     →#
     # To help debug your network, it may help to use visualizations similar to the \Box
     # ones we used above; these visualizations will have significant qualitative
     →#
     # differences from the ones we saw above for the poorly tuned network.
     →#
     #
                                                                     ш
     →#
     # Tweaking hyperparameters by hand can be fun, but you might find it useful to \Box
     →#
     # write code to sweep through possible combinations of hyperparameters
     →#
     # automatically like we did on the previous exercises.
                                                                     Ш
     # Your code
```

```
input_size = 32 * 32 * 3
hidden_size = 50
num_classes = 10
results = {}
best_val = -1
best net = None
learning_rates = [1e-4, 1e-3]
regularization strengths = [0.15, 0.25]
learning_rate_decay = 0.95
for lr in learning_rates:
    for reg in regularization_strengths:
       net = TwoLayerNet(input_size, hidden_size, num_classes)
       # Train the network
       stats = net.train(X_train, y_train, X_val, y_val, lr,__
 →learning_rate_decay, reg,
                       num_iters=1000, batch_size=200, verbose=False)
       y train pred = net.predict(X train)
       train_accuracy = np.mean(y_train == y_train_pred)
       y val pred = net.predict(X val)
       val_accuracy = np.mean(y_val == y_val_pred)
       results[(lr, reg)] = (train_accuracy, val_accuracy)
       if val_accuracy > best_val:
           best_net = net
           best_val = val_accuracy
# 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' %u
 →best val)
#
                             END OF YOUR CODE
 →#
lr 1.000000e-04 reg 1.500000e-01 train accuracy: 0.281143 val accuracy: 0.285000
lr 1.000000e-04 reg 2.500000e-01 train accuracy: 0.282592 val accuracy: 0.280000
```

```
lr 1.000000e-04 reg 1.500000e-01 train accuracy: 0.281143 val accuracy: 0.285000
lr 1.000000e-04 reg 2.500000e-01 train accuracy: 0.282592 val accuracy: 0.280000
lr 1.000000e-03 reg 1.500000e-01 train accuracy: 0.490816 val accuracy: 0.451000
lr 1.000000e-03 reg 2.500000e-01 train accuracy: 0.483551 val accuracy: 0.482000
best validation accuracy achieved during cross-validation: 0.482000
```

[12]: # visualize the weights of the best network show_net_weights(best_net)



10 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%.

```
[13]: test_acc = (best_net.predict(X_test) == y_test).mean()
print('Test accuracy: ', test_acc)
```

Test accuracy: 0.456

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: 1, 3
Your explanation:

When our testing accuracy is much lower than the training accuracy, it is likely that our classifier is overfitting to our training data.

- Train on a larger dataset: introduce more new data to make our classifier more general.
- Increase the regularization strength: adding more penalty to our weights to again make our classifier more general

features

September 29, 2021

1 Image features 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.

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.

```
[1]: 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
```

1.1 Load data

Similar to previous exercises, we will load CIFAR-10 data from disk.

```
[2]: 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)
try:
  del X train, y train
  del X test, y test
  print('Clear previously loaded data.')
except:
  pass
X_train, y_train, X_val, y_val, X_test, y_test = get_CIFAR10_data()
```

1.2 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.

```
[3]: 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
Done extracting features for 13000 / 49000 images
Done extracting features for 14000 / 49000 images
Done extracting features for 15000 / 49000 images
Done extracting features for 16000 / 49000 images
Done extracting features for 17000 / 49000 images
Done extracting features for 18000 / 49000 images
Done extracting features for 19000 / 49000 images
Done extracting features for 20000 / 49000 images
Done extracting features for 21000 / 49000 images
Done extracting features for 22000 / 49000 images
Done extracting features for 23000 / 49000 images
Done extracting features for 24000 / 49000 images
```

```
Done extracting features for 25000 / 49000 images
Done extracting features for 26000 / 49000 images
Done extracting features for 27000 / 49000 images
Done extracting features for 28000 / 49000 images
Done extracting features for 29000 / 49000 images
Done extracting features for 30000 / 49000 images
Done extracting features for 31000 / 49000 images
Done extracting features for 32000 / 49000 images
Done extracting features for 33000 / 49000 images
Done extracting features for 34000 / 49000 images
Done extracting features for 35000 / 49000 images
Done extracting features for 36000 / 49000 images
Done extracting features for 37000 / 49000 images
Done extracting features for 38000 / 49000 images
Done extracting features for 39000 / 49000 images
Done extracting features for 40000 / 49000 images
Done extracting features for 41000 / 49000 images
Done extracting features for 42000 / 49000 images
Done extracting features for 43000 / 49000 images
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
```

1.3 Train SVM on features

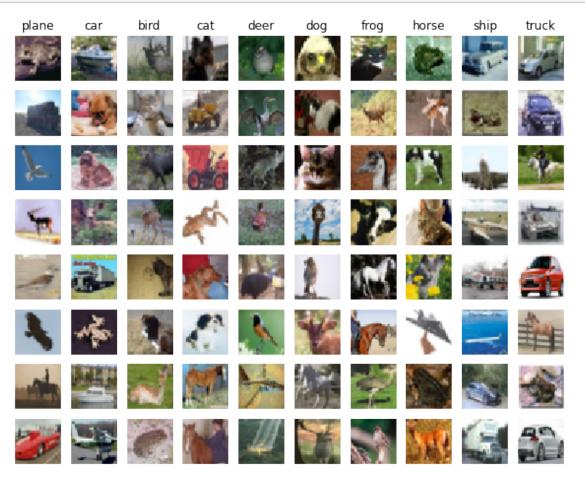
Using the multiclass SVM code developed earlier in the assignment, train SVMs on top of the features extracted above; this should achieve better results than training SVMs directly on top of raw pixels.

```
[4]: # Use the validation set to tune the learning rate and regularization strength
    from cs682.classifiers.linear_classifier import LinearSVM
    learning_rates = [1e-9, 1e-8, 1e-7]
    regularization_strengths = [5e4, 5e5, 5e6]
    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 sum. 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.
    # Your code
    for lr in learning_rates:
       for reg in regularization_strengths:
           svm = LinearSVM()
           svm.train(X_train_feats, y_train, lr, reg, num_iters=1500)
           y train pred = svm.predict(X train feats)
           train_accuracy = np.mean(y_train == y_train_pred)
           y val pred = svm.predict(X val feats)
           val_accuracy = np.mean(y_val == y_val_pred)
           results[(lr, reg)] = (train_accuracy, val_accuracy)
           if val_accuracy > best_val:
               best_svm = svm
               best_val = val_accuracy
    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)
   lr 1.000000e-09 reg 5.000000e+04 train accuracy: 0.092776 val accuracy: 0.089000
   lr 1.000000e-09 reg 5.000000e+05 train accuracy: 0.120327 val accuracy: 0.119000
   lr 1.000000e-09 reg 5.000000e+06 train accuracy: 0.413959 val accuracy: 0.418000
   lr 1.000000e-08 reg 5.000000e+04 train accuracy: 0.114265 val accuracy: 0.126000
   lr 1.000000e-08 reg 5.000000e+05 train accuracy: 0.411918 val accuracy: 0.413000
   lr 1.000000e-08 reg 5.000000e+06 train accuracy: 0.402061 val accuracy: 0.384000
   lr 1.000000e-07 reg 5.000000e+04 train accuracy: 0.414061 val accuracy: 0.410000
   lr 1.000000e-07 reg 5.000000e+05 train accuracy: 0.409837 val accuracy: 0.417000
   lr 1.000000e-07 reg 5.000000e+06 train accuracy: 0.323714 val accuracy: 0.312000
   best validation accuracy achieved during cross-validation: 0.418000
[5]: # 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)
```

0.418

```
[6]: # 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', _
     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()
```



1.3.1 Inline question 1:

Describe the misclassification results that you see. Do they make sense?

Just by looking at the pictures, I think the easiest to observe is that most images that were miss classified as planes or ships were images with a large blue background. Which makes sense because that was what we observed when we were looking at the visualization of the weights.

1.4 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.

```
[7]: # 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)

```
best val = -1
learning_rates = [1e-1, 5e-1, 9e-1]
regularization_strengths = [0.025, 0.05, 0.15]
learning_rate_decay = 0.95
for lr in learning_rates:
    for reg in regularization_strengths:
       net = TwoLayerNet(input_dim, hidden_dim, num_classes)
        # Train the network
        stats = net.train(X_train_feats, y_train, X_val_feats, y_val, lr,__
 →learning_rate_decay, reg,
                        num_iters=1500, batch_size=200, verbose=False)
       y_train_pred = net.predict(X_train_feats)
       train_accuracy = np.mean(y_train == y_train_pred)
       y_val_pred = net.predict(X_val_feats)
        val_accuracy = np.mean(y_val == y_val_pred)
       results[(lr, reg)] = (train_accuracy, val_accuracy)
        if val_accuracy > best_val:
           best_net = net
           best_val = val_accuracy
# 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' %u
 →best_val)
END OF YOUR CODE
lr 1.000000e-01 reg 2.500000e-02 train accuracy: 0.484449 val accuracy: 0.485000
lr 1.000000e-01 reg 5.000000e-02 train accuracy: 0.437061 val accuracy: 0.430000
lr 1.000000e-01 reg 1.500000e-01 train accuracy: 0.250592 val accuracy: 0.259000
lr 5.000000e-01 reg 2.500000e-02 train accuracy: 0.451816 val accuracy: 0.435000
lr 5.000000e-01 reg 5.000000e-02 train accuracy: 0.397878 val accuracy: 0.403000
lr 5.000000e-01 reg 1.500000e-01 train accuracy: 0.257816 val accuracy: 0.236000
lr 9.000000e-01 reg 2.500000e-02 train accuracy: 0.432776 val accuracy: 0.443000
lr 9.000000e-01 reg 5.000000e-02 train accuracy: 0.402939 val accuracy: 0.389000
lr 9.000000e-01 reg 1.500000e-01 train accuracy: 0.242796 val accuracy: 0.242000
best validation accuracy achieved during cross-validation: 0.485000
```

```
[9]: # Run your best neural net classifier on the test set. You should be able
# to get more than 55% accuracy.

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

0.472

cool bonus

September 29, 2021

1 Cool Bonus: Dynamic Mini-batch

This is mainly testing a question I had during class. Which would having different batch sizes during the training help improve the process? Since in class the batch size was always the same during the same training, I wanted to test out whether having three different batch-size(small, medium, large) had any effect.

The specific question I wish to answer in the following experient are: - Would a dynamic batch be faster than a large or slower than a small batch-size? If so, by how much? - Would a dynamic batch have a better than a small or worse than a large batch-size? If so, by how much?

```
[1]: # 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/
     \rightarrow autoreload-of-modules-in-ipython
     %load_ext autoreload
     %autoreload 2
```

1.1 CIFAR-10 Data Loading and Preprocessing

```
[2]: 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_dev = np.reshape(X_dev, (X_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 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,)

```
[3]: from cs682.classifiers import LinearSVM
import time

results = {}
best_val = -1
best_svm = None
batch_sizes = [100, [100, 300, 500], 500]

for batch_size in batch_sizes:

   tic = time.time()

   svm = LinearSVM()
   svm.train(X_train, y_train, batch_size=batch_size, learning_rate=1e-7, usine content of the co
```

```
toc = time.time()
         cost_time = toc-tic
         results[str(batch_size)] = (train_accuracy, val_accuracy, cost_time)
         if val_accuracy > best_val:
             best_svm = svm
             best_val = val_accuracy
     # Print out results.
     for batch_size in sorted(results):
         train_accuracy, val_accuracy, cost_time = results[str(batch_size)]
         print('batch size %s train accuracy: %f val accuracy: %f took %f seconds' %⊔
     → (
                     batch_size, train_accuracy, val_accuracy, cost_time))
     print('best validation accuracy achieved during cross-validation: %f' %⊔
     →best_val)
    batch size 100 train accuracy: 0.371224 val accuracy: 0.384000 took 1.818013
    seconds
    batch size 500 train accuracy: 0.385857 val accuracy: 0.399000 took 8.026039
    batch size [100, 300, 500] train accuracy: 0.387633 val accuracy: 0.394000 took
    4.874423 seconds
    best validation accuracy achieved during cross-validation: 0.399000
[4]: from cs682.classifiers import Softmax
     results = {}
     best_val = -1
     best_softmax = None
     batch_sizes = [100, [100, 300, 500], 500]
     for batch_size in batch_sizes:
         tic = time.time()
         softmax = Softmax()
         softmax.train(X_train, y_train, learning_rate=8e-8, reg=9e3,_
     →batch_size=batch_size, num_iters=1500)
         y_train_pred = softmax.predict(X_train)
         train_accuracy = np.mean(y_train == y_train_pred)
         y_val_pred = softmax.predict(X_val)
         val_accuracy = np.mean(y_val == y_val_pred)
         toc = time.time()
```

```
batch size 100 train accuracy: 0.337184 val accuracy: 0.353000 took 1.148682 seconds
batch size 500 train accuracy: 0.343367 val accuracy: 0.372000 took 8.106762 seconds
batch size [100, 300, 500] train accuracy: 0.339429 val accuracy: 0.352000 took 4.760601 seconds
best validation accuracy achieved during cross-validation: 0.372000
```

The results I got for both SVM and Softmax were interesting.

Dynamic mini-batch was faster in both cases! However the val accuracy for Softmax seems the same as the small batch-size, while for SVM the dynamic mini-batch was much closer to large batch-size and only cost 4.87s when the large batch-size cost 8.02.

I think it would definitely need more experiments to make sure of the benefits for dynamic minibatch. But the result I got on SVM looks promising!