# knn

June 8, 2024

```
[]: # This mounts your Google Drive to the Colab VM.
     from google.colab import drive
     drive.mount('/content/drive')
     # TODO: Enter the foldername in your Drive where you have saved the unzipped
     # assignment folder, e.g. 'cs231n/assignments/assignment1/'
     FOLDERNAME = None
     assert FOLDERNAME is not None, "[!] Enter the foldername."
     # Now that we've mounted your Drive, this ensures that
     # the Python interpreter of the Colab VM can load
     # python files from within it.
     import sys
     sys.path.append('/content/drive/My Drive/{}'.format(FOLDERNAME))
     # This downloads the CIFAR-10 dataset to your Drive
     # if it doesn't already exist.
     %cd /content/drive/My\ Drive/$FOLDERNAME/cs231n/datasets/
     !bash get datasets.sh
     %cd /content/drive/My\ Drive/$FOLDERNAME
```

# 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.

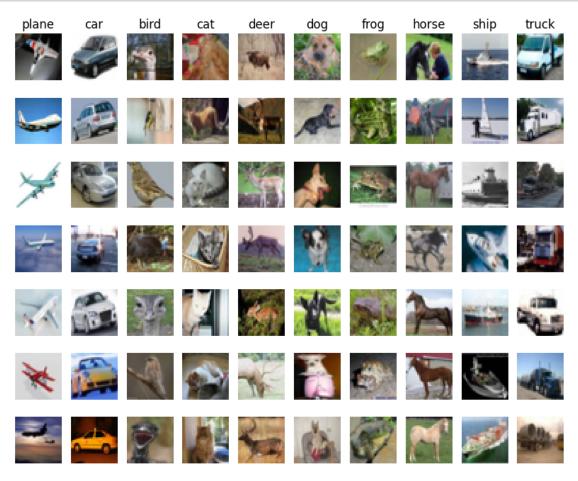
The kNN classifier consists of two stages:

- 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.

```
[3]: # Run some setup code for this notebook.
     import random
     import numpy as np
     from cs231n.data_utils import load_CIFAR10
     import matplotlib.pyplot as plt
     # This is a bit of magic to make matplotlib figures appear inline in the
      \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
[4]: # Load the raw CIFAR-10 data.
     cifar10_dir = 'cs231n/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,)
```

```
[5]: # 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()
```



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

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

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

(5000, 3072) (500, 3072)

```
[7]: from cs231n.classifiers import KNearestNeighbor

# Create a kNN classifier instance.

# Remember that training a kNN classifier is a noop:

# the Classifier simply remembers the data and does no further processing classifier = KNearestNeighbor() classifier.train(X_train, y_train)
```

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

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

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

Note: For the three distance computations that we require you to implement in this notebook, you may not use the np.linalg.norm() function that numpy provides.

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

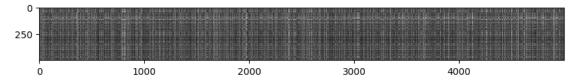
```
[11]: # Open cs231n/classifiers/k_nearest_neighbor.py and implement
    # compute_distances_two_loops.

# Test your implementation:
dists = classifier.compute_distances_two_loops(X_test)
```

```
print(dists.shape)
```

(500, 5000)

```
[12]: # 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 visibly 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?

### YourAnswer:

We know that each row is a single test example and each column is a single train example:

- If a test image is very different from all the training images, the row will be bright.
- If all the test images are very different from a single training image, the column will be bright.

```
[14]: # 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:

```
[16]: 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 use other distance metrics such as L1 distance. For pixel values  $p_{ij}^{(k)}$  at location (i,j)of some image  $I_k$ ,

the mean  $\mu$  across all pixels over all images is

$$\mu = \frac{1}{nhw} \sum_{k=1}^{n} \sum_{i=1}^{h} \sum_{j=1}^{w} p_{ij}^{(k)}$$

And the pixel-wise mean  $\mu_{ij}$  across all images is

$$\mu_{ij} = \frac{1}{n} \sum_{k=1}^{n} p_{ij}^{(k)}.$$

The general standard deviation  $\sigma$  and pixel-wise standard deviation  $\sigma_{ij}$  is defined similarly.

Which of the following preprocessing steps will not change the performance of a Nearest Neighbor classifier that uses L1 distance? Select all that apply. To clarify, both training and test examples are preprocessed in the same way.

- 1. Subtracting the mean  $\mu$   $(\tilde{p}_{ij}^{(k)} = p_{ij}^{(k)} \mu.)$
- 2. Subtracting the per pixel mean  $\mu_{ij}$   $(\tilde{p}_{ij}^{(k)} = p_{ij}^{(k)} \mu_{ij})$ . Subtracting the mean  $\mu$  and dividing by the standard deviation  $\sigma$ .
- 4. Subtracting the pixel-wise mean  $\mu_{ij}$  and dividing by the pixel-wise standard deviation  $\sigma_{ij}$ .
- 5. Rotating the coordinate axes of the data, which means rotating all the images by the same angle. Empty regions in the image caused by rotation are padded with a same pixel value and no interpolation is performed.

#### Your Answer:

# 1, 2, and 3

# Your Explanation:

- 1. Mean dist
- 2. SDpixel-wise SD dist dist

performance

3. L1 dist L1

Detailed Explanation from mantatsu

First, we need to make some assumptions. Consider  $\mathbf{p}^{(k)}$  is our pixel vector which is just a flattened matrix of pixel values of some image  $I_k$ . There are a total of  $S = h \times w$  values where each entry  $p_m^{(k)}$  in  $\mathbf{p}^{(k)}$  corresponds to each entry  $p_{ij}^{(k)}$  in  $I_k$ .

Let's remember how L1 is calculated between 2 pixel vectors belonging to different images:

$$d_1\left(\mathbf{p}^{(k)},\mathbf{p}^{(k')}\right) = \left|\left|\mathbf{p}^{(k)} - \mathbf{p}^{(k')}\right|\right|_1$$

where L1 norm is expressed as  $||-||_1$  and calculated as follows:

$$\left|\left|\mathbf{p}^{(k)} - \mathbf{p}^{(k')}\right|\right|_1 = \sum_{m=1}^{S} \left|p_m^{(k)} - p_m^{(k')}\right|$$

Since distances are computed between pixels element-wise, it can be derived that only the last preprocessing step could change the L1 distance, thus the classifier performance: 1. Mean subtraction \* Formula:  $\sum_{m=1}^{S} \left| \left( p_m^{(k)} - \mu \right) - \left( p_m^{(k')} - \mu \right) \right| * Explanation$ : it can be seen that  $\mu$  will cancel out during distance computation because both pixel values are modified by the same constant.

2. Per-pixel mean subtraction \* Formula:  $\sum_{m=1}^{S} \left| \left( p_m^{(k)} - \mu_m \right) - \left( p_m^{(k')} - \mu_m \right) \right| * Explanation$ : same as 1 3. Standartization \* Formula:  $\sum_{m=1}^{S} \left| \frac{p_m^{(k)} - \mu}{\sigma} - \frac{p_m^{(k')} - \mu}{\sigma} \right| * Explanation:$  it can be seen that  $\sigma^{-1}$  is a constant by which each absolute difference is weighted. The constant for all the differences is the same, therefore, each distance is scaled by the same amount. This means that the performance is unaffected because the distance comparisons will remain the same (just on a different scale) 4. **Per-pixel standartization** \* Formula:  $\sum_{m=1}^{S} \left| \frac{p_m^{(k)} - \mu_m}{\sigma_m} - \frac{p_m^{(k')} - \mu_m}{\sigma_m} \right|$  \* Explanation: it can be seen that  $\sigma_m^{-1}$  is a constant by which each absolute difference is weighted. The constant is different for every difference, therefore, some differences can become more prominent than they originally were while computing some L1, whereas the same constants may have not much affect on the differences when computing some other L1. \* Example: assume  $\mathbf{p}^{(k)} = \begin{bmatrix} 0 \\ 0 \end{bmatrix}$ ,  $\mathbf{p}^{(k')} = \begin{bmatrix} 0 \\ 0.9 \end{bmatrix}$ ,

$$\mathbf{p}^{(k'')} = \begin{bmatrix} 1 \\ 0.1 \end{bmatrix}$$
, and  $\sigma = \begin{bmatrix} 1 \\ 0.1 \end{bmatrix}$ .

The following distances are acquired before scaling:

- \*  $d_1\left(\mathbf{p}^{(k)}, \mathbf{p}^{(k')}\right) = 0 + 0.9 = 0.9$
- \*  $d_1\left(\mathbf{p}^{(k)}, \mathbf{p}^{(k'')}\right) = 1 + 0.1 = 1.1$

And the following distances are acquired after scaling:

- \*  $d_1\left(\mathbf{p}^{(k)}, \mathbf{p}^{(k')}\right) = 1 \cdot 0 + 10 \cdot 0.9 = 10$
- \*  $d_1\left(\mathbf{p}^{(k)}, \mathbf{p}^{(k'')}\right) = 1 \cdot 1 + 10 \cdot 0.1 = 2$

And we can see that before scaling  $d_1\left(\frac{p}^{(k)}, \mathcal{p}^{(k')}\right) < 0$ 

#### 5. Rotation

- Formula:  $||R\mathbf{p}^{(k)} R\mathbf{p}^{(k')}||_1$
- Explanation: in Manhattan space, each absolute difference takes into account axis orientation which changes as the space is rotated, resulting in relative difference changes (not scaled by the same constant). Therefore, the L1 distance between 2 points of 2 different vectors in the same dimension after the transformation will differ because the (unnormalized) basis for the 2 vector spaces are (usually) different. Note that, in Euclidean space, the rotated basis would also be different but that will not change the distances (L2 in this case) between points since per-axis squared differences lose the information
- Example: assume  $\mathbf{p}^{(k)} = \begin{bmatrix} 0 \\ 0 \end{bmatrix}$ ,  $\mathbf{p}^{(k')} = \begin{bmatrix} 1 \\ 0 \end{bmatrix}$ ,  $\mathbf{p}^{(k'')} = \begin{bmatrix} 1 \\ 1 \end{bmatrix}$ , and  $R = \begin{bmatrix} \cos(\frac{\pi}{4}) & -\sin(\frac{\pi}{4}) \\ \sin(\frac{\pi}{4}) & \cos(\frac{\pi}{4}) \end{bmatrix} = \mathbf{p}^{(k')}$

```
-d_1(\mathbf{p}^{(k)},\mathbf{p}^{(k'')}) = 1+1=2
                 And the following distances are acquired after rotating:
                  -d_1\left(\mathbf{p}^{(k)},\mathbf{p}^{(k')}\right) = \frac{\sqrt{2}}{2} + \frac{\sqrt{2}}{2} = \sqrt{2}
                  -d_1(\mathbf{p}^{(k)},\mathbf{p}^{(k'')}) = 0 + \sqrt{2} = \sqrt{2}
                 And we can see that before rotating d_1\left(\mathbf{p}^{(k)},\mathbf{p}^{(k')}\right) < d_1\left(\mathbf{p}^{(k)},\mathbf{p}^{(k'')}\right) but after rotating
                 d_1\left(\mathbf{p}^{(k)},\mathbf{p}^{(k')}\right) = d_1\left(\mathbf{p}^{(k)},\mathbf{p}^{(k'')}\right).
[27]: # 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, \Box
        \hookrightarrow reshape
       # the matrices into vectors and compute the Euclidean distance between them.
       difference = np.linalg.norm(dists - dists_one, ord='fro')
       print('One loop 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')
      One loop difference was: 0.000000
      Good! The distance matrices are the same
[22]: # 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('No loop 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')
      No loop difference was: 0.000000
      Good! The distance matrices are the same
```

The following distances are acquired before rotating:

 $-d_1(\mathbf{p}^{(k)}, \mathbf{p}^{(k')}) = 1 + 0 = 1$ 

```
[29]: # 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 _{\! \sqcup}
       \hookrightarrow to execute.
           11 11 11
          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
       \rightarrow implementation!
      # NOTE: depending on what machine you're using,
      # you might not see a speedup when you go from two loops to one loop,
      # and might even see a slow-down.
```

Two loop version took 83.989085 seconds One loop version took 83.330165 seconds No loop version took 0.576901 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.

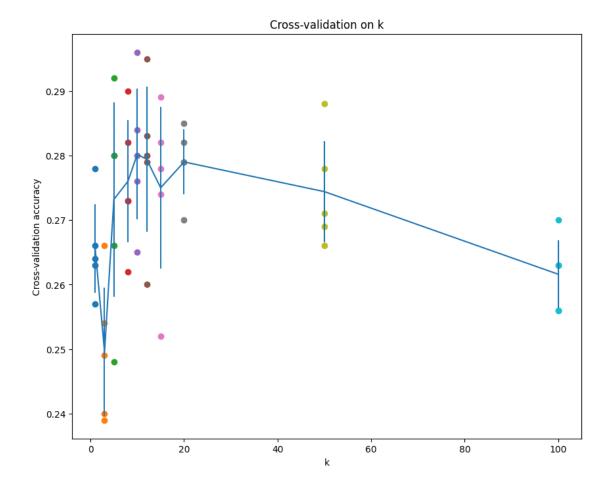
```
# *****START OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)****
X_train_folds = np.array_split(X_train, num_folds)
y_train_folds = np.array_split(y_train, num_folds)
# *****END OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)****
# A dictionary holding the accuracies for different values of k that we 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.
# *****START OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)****
for k in k choices:
   k_to_accuracies[k] = []
   for i in range(num_folds):
       # Create num folds-1 of the data and labels as the training samples
      X_train_temp = np.concatenate(np.compress(
          np.arange(num_folds) != i, X_train_folds, axis=0))
       y_train_temp = np.concatenate(np.compress(
          np.arange(num_folds) != i, y_train_folds, axis=0))
       # Train the classifier based on the training data
       classifier.train(X_train_temp, y_train_temp)
       # Predict using the remaining fold representing validation data
      y pred temp = classifier.predict(X train folds[i], k=k, num loops=0)
       # Compute the accuracy of the predicted label
      num_correct = np.sum(y_pred_temp == y_train_folds[i])
      k_to_accuracies[k].append(num_correct / len(y_pred_temp))
# *****END OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)****
```

```
# 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
[31]: # 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()
```



```
[33]: # 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 = k_choices[accuracies_mean.argmax()]

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 decision boundary of the k-NN classifier is

linear. 2. The training error of a 1-NN will always be lower than or equal to that of 5-NN. 3. The test error of a 1-NN will always be lower than that of a 5-NN. 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:

# Your Explanation:

1. **FALSE**: n hyperplane

2. TRUE: 1-NN training error 5-NN 1-NN 5-NN 5

3. FALSE: 1-NN 5-NN 5-NN 5-NN 1-NN

4. **TRUE**:

#### svm

June 8, 2024

```
[2]: # This mounts your Google Drive to the Colab VM.
     from google.colab import drive
     drive.mount('/content/drive')
     # TODO: Enter the foldername in your Drive where you have saved the unzipped
     # assignment folder, e.g. 'cs231n/assignments/assignment1/'
     FOLDERNAME = 'CS231n Assignments/assignment1/'
     assert FOLDERNAME is not None, "[!] Enter the foldername."
     # Now that we've mounted your Drive, this ensures that
     # the Python interpreter of the Colab VM can load
     # python files from within it.
     import sys
     sys.path.append('/content/drive/My Drive/{}'.format(FOLDERNAME))
     # This downloads the CIFAR-10 dataset to your Drive
     # if it doesn't already exist.
     %cd /content/drive/My\ Drive/$FOLDERNAME/cs231n/datasets/
     !bash get datasets.sh
     %cd /content/drive/My\ Drive/$FOLDERNAME
```

Mounted at /content/drive /content/drive/My Drive/CS231n Assignments/assignment1/cs231n/datasets /content/drive/My Drive/CS231n Assignments/assignment1

# 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

```
[20]: # Run some setup code for this notebook.
      import random
      import numpy as np
      from cs231n.data_utils import load_CIFAR10
      import matplotlib.pyplot as plt
      # This is a bit of magic to make matplotlib figures appear inline in the
      # notebook rather than in a new window.
      %matplotlib inline
      plt.rcParams['figure.figsize'] = (10.0, 8.0) # set default size of plots
      plt.rcParams['image.interpolation'] = 'nearest'
      plt.rcParams['image.cmap'] = 'gray'
      # Some more magic so that the notebook will reload external python modules;
      # see http://stackoverflow.com/questions/1907993/
       \rightarrow autoreload-of-modules-in-ipython
      %load ext autoreload
      %autoreload 2
```

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

# 1.1 CIFAR-10 Data Loading and Preprocessing

```
[21]: # Load the raw CIFAR-10 data.
      cifar10_dir = 'cs231n/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)
```

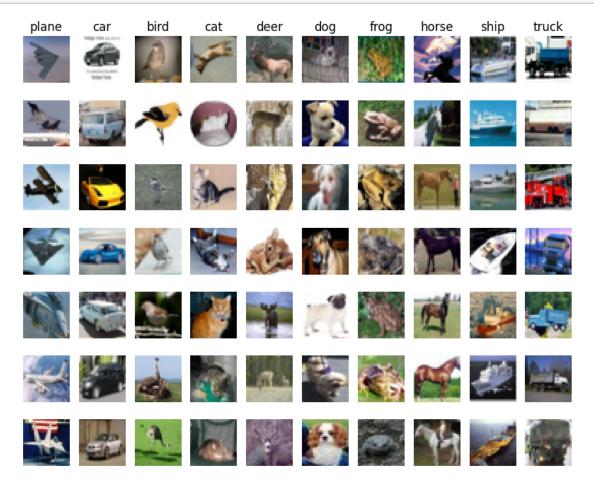
Clear previously loaded data.

Training data shape: (50000, 32, 32, 3)

Training labels shape: (50000,)

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

```
[22]: # 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()
```



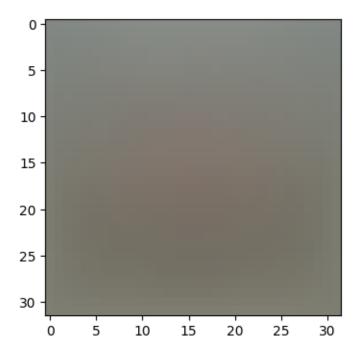
```
[23]: # Split the data into train, val, and test sets. In addition we will
      # create a small development set as a subset of the training data;
      # we can use this for development so our code runs faster.
      num_training = 49000
      num_validation = 1000
      num_test = 1000
      num_dev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

```
[24]: # 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)
[25]: # 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_i
       \hookrightarrow image
      plt.show()
      # 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
      # 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))])
```

[130.64189796 135.98173469 132.47391837 130.05569388 135.34804082 131.75402041 130.96055102 136.14328571 132.47636735 131.48467347]

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 cs231n/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.

```
[26]: # Evaluate the naive implementation of the loss we provided for you:
    from cs231n.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: 9.380900

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, 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:

```
[27]: # 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 cs231n.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: 18.884505 analytic: 18.884505, relative error: 4.024633e-12
numerical: 0.022778 analytic: 0.022778, relative error: 8.107635e-09
numerical: -6.771037 analytic: -6.771037, relative error: 2.097558e-11
numerical: 2.463959 analytic: 2.463959, relative error: 2.051104e-10
numerical: 20.563848 analytic: 20.563848, relative error: 9.726066e-12
numerical: 0.970949 analytic: 0.970949, relative error: 4.125456e-10
numerical: 9.104800 analytic: 9.104800, relative error: 1.076627e-11
numerical: -1.258153 analytic: -1.258825, relative error: 2.672292e-04
numerical: -17.047273 analytic: -17.047273, relative error: 9.208210e-12
numerical: 11.682778 analytic: 11.682778, relative error: 2.628170e-12
numerical: -12.317642 analytic: -12.319517, relative error: 7.613245e-05
numerical: 5.429396 analytic: 5.429033, relative error: 3.348866e-05
numerical: -44.567230 analytic: -44.679777, relative error: 1.261079e-03
numerical: -21.654647 analytic: -21.654647, relative error: 1.291212e-11
numerical: 7.161745 analytic: 7.161745, relative error: 1.053909e-11
numerical: 13.260400 analytic: 13.260400, relative error: 1.631660e-11
numerical: 0.097807 analytic: 0.097807, relative error: 3.478209e-09
numerical: 16.463840 analytic: 16.438967, relative error: 7.559699e-04
numerical: 20.938328 analytic: 20.937656, relative error: 1.606197e-05
numerical: 14.443883 analytic: 14.443883, relative error: 3.204939e-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

#### YourAnswer:

#### Good answer from Mantasu

First, we need to make some assumptions. To compute our **SVM loss**, we use **Hinge loss** which takes the form  $\max(0, -)$ . For 1D case, we can define it as follows  $(\hat{y} - \text{score}, i - \text{any class}, c - \text{correct class}, \Delta - \text{margin})$ :

$$f(x) = \max(0, x), \text{ where } x = \hat{y}_i - \hat{y}_c + \Delta$$

Let's now see how our max function fits the definition of computing the gradient. It is the formula we use for computing the gradient numerically when, instead of implementing the limit approaching to 0, we choose some arbitrary small h:

$$\frac{df(x)}{dx} = \lim_{h \to 0} \frac{\max(0, x+h) - \max(0, x)}{h}$$

Now we can talk about the possible mismatches between numeric and analytic gradient computation: 1. Cause of mismatch \* Relative error - the discrepancy is caused due to arbitrary choice of small values of h because by definition it should approach 0. Analytic computation produces an exact result (as precise as computation precision allows) while numeric solution only approximates the result. \* Kinks - max only has a subgradient because when both values in max are equal, its gradient is undefined, therefore, not smooth. Such parts, referred to as kinks, may cause numeric gradient to produce different results from analytic computation due to (again) arbitrary choice of h. 2. Concerns \* When comparing analytic and numeric methods, kinks are more dangerous than small inaccuracies where the gradient is smooth. Small derivative inaccuracies still change the weight by approximately the same amount but kinks may cause unintentional updates as seen in an example below. If the unintentional values would have a noticable affect on parameter updates, it is a reason for concern. 3. 1D example of numeric gradient fail \* Assume  $x = -10^{-9}$ . Then the analytic computation of the derivative of  $\max(0, x)$  would yield 0. However, if we choose our  $h = 10^{-8}$ , then the numeric computation would yield 0.9. 4. Relation between margin and mismatch \* Assuming all other parameters remain unchanged, increasing  $\Delta$  will lower the frequency of kinks. This is because higher  $\Delta$  will cause more x to be positive, thus reducing the probability of kinks. In reality though, it would not have a big effect - if we increase the margin  $\Delta$ , the **SVM** will only learn to increase the (negative) gap between  $\hat{y}_i - \hat{y}_c$  and 0 (when  $i \neq c$ ). But that still means, if we add  $\Delta$ , there is the same chance for x to result on the edge.

```
# The losses should match but your vectorized implementation should be muchusfaster.

print('difference: %f' % (loss_naive - loss_vectorized))
```

Naive loss: 9.380900e+00 computed in 0.071088s Vectorized loss: 9.380900e+00 computed in 0.015362s difference: 0.000000

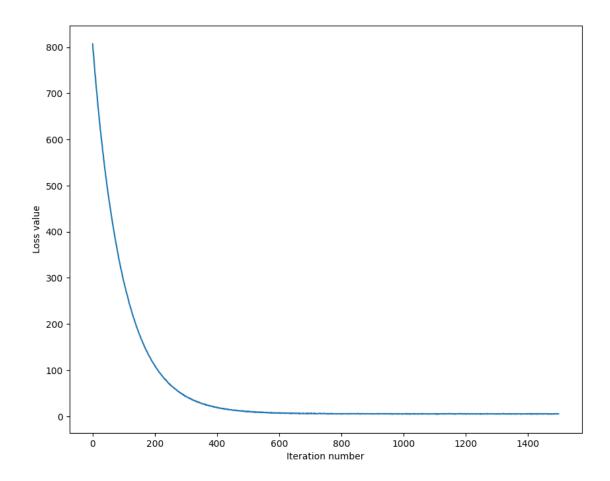
```
[29]: # Complete the implementation of sum_loss_vectorized, and compute the gradient
      # of the loss function in a vectorized way.
      # The naive implementation and the vectorized implementation should match, but
      # the vectorized version should still be much faster.
      tic = time.time()
      _, 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 0.245092s Vectorized loss and gradient: computed in 0.004506s difference: 0.000000

### 1.2.1 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. Your code for this part will be written inside cs231n/classifiers/linear\_classifier.py.

```
print('That took %fs' % (toc - tic))
     iteration 0 / 1500: loss 806.897213
     iteration 100 / 1500: loss 291.059551
     iteration 200 / 1500: loss 109.538098
     iteration 300 / 1500: loss 43.138137
     iteration 400 / 1500: loss 19.037643
     iteration 500 / 1500: loss 10.793517
     iteration 600 / 1500: loss 7.042191
     iteration 700 / 1500: loss 5.822904
     iteration 800 / 1500: loss 4.821756
     iteration 900 / 1500: loss 4.885208
     iteration 1000 / 1500: loss 5.388546
     iteration 1100 / 1500: loss 5.218789
     iteration 1200 / 1500: loss 4.548064
     iteration 1300 / 1500: loss 5.004626
     iteration 1400 / 1500: loss 5.720977
     That took 9.160720s
[31]: # 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()
```



training accuracy: 0.368449 validation accuracy: 0.371000

```
[33]: # 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 0.39 (> 0.385) on the validation set.

# Note: you may see runtime/overflow warnings during hyper-parameter search.
# This may be caused by extreme values, and is not a bug.

# results is dictionary mapping tuples of the form
```

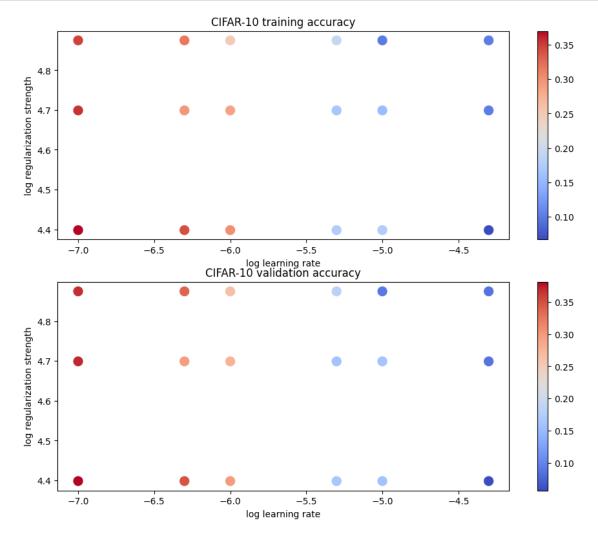
```
# (learning rate, regularization strength) to tuples of the form
# (training accuracy, validation accuracy). The accuracy is simply the fraction
# of data points that are correctly classified.
results = {}
best_val = -1  # The highest validation accuracy that we have seen so far.
best_svm = None # The LinearSVM object that achieved the highest validation
 -rate.
# 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.
# Provided as a reference. You may or may not want to change these
\rightarrowhyperparameters
learning_rates = [1e-7, 1e-6, 5e-7, 5e-6, 1e-5, 5e-5]
regularization strengths = [2.5e4, 5e4, 7.5e4]
# *****START OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)****
for lr in learning_rates:
   for reg in regularization_strengths:
       svm = LinearSVM()
       svm.train(X_train, y_train, lr, reg, num_iters=1000)
       y_train_pred, y_val_pred = svm.predict(X_train), svm.predict(X_val)
       results[(lr, reg)] = np.mean(
           y_train == y_train_pred), np.mean(y_val == y_val_pred)
       if results[(lr, reg)][1] > best_val:
           best_val = results[(lr, reg)][1]
           best_svm = svm
# *****END OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)*****
# 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-07 reg 2.500000e+04 train accuracy: 0.370020 val accuracy: 0.381000
     lr 1.000000e-07 reg 5.000000e+04 train accuracy: 0.355714 val accuracy: 0.370000
     lr 1.000000e-07 reg 7.500000e+04 train accuracy: 0.347878 val accuracy: 0.368000
     lr 5.000000e-07 reg 2.500000e+04 train accuracy: 0.340449 val accuracy: 0.348000
     lr 5.000000e-07 reg 5.000000e+04 train accuracy: 0.295714 val accuracy: 0.297000
     lr 5.000000e-07 reg 7.500000e+04 train accuracy: 0.317367 val accuracy: 0.339000
     lr 1.000000e-06 reg 2.500000e+04 train accuracy: 0.300449 val accuracy: 0.298000
     lr 1.000000e-06 reg 5.000000e+04 train accuracy: 0.286980 val accuracy: 0.278000
     lr 1.000000e-06 reg 7.500000e+04 train accuracy: 0.248551 val accuracy: 0.264000
     lr 5.000000e-06 reg 2.500000e+04 train accuracy: 0.174673 val accuracy: 0.164000
     lr 5.000000e-06 reg 5.000000e+04 train accuracy: 0.167388 val accuracy: 0.159000
     lr 5.000000e-06 reg 7.500000e+04 train accuracy: 0.193163 val accuracy: 0.182000
     lr 1.000000e-05 reg 2.500000e+04 train accuracy: 0.174510 val accuracy: 0.159000
     lr 1.000000e-05 reg 5.000000e+04 train accuracy: 0.154837 val accuracy: 0.160000
     lr 1.000000e-05 reg 7.500000e+04 train accuracy: 0.099061 val accuracy: 0.091000
     lr 5.000000e-05 reg 2.500000e+04 train accuracy: 0.066673 val accuracy: 0.057000
     lr 5.000000e-05 reg 5.000000e+04 train accuracy: 0.100265 val accuracy: 0.087000
     lr 5.000000e-05 reg 7.500000e+04 train accuracy: 0.100265 val accuracy: 0.087000
     best validation accuracy achieved during cross-validation: 0.381000
[34]: # Visualize the cross-validation results
      import math
      import pdb
      # pdb.set trace()
      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.tight_layout(pad=3)
      plt.scatter(x_scatter, y_scatter, marker_size, c=colors, cmap=plt.cm.coolwarm)
      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, cmap=plt.cm.coolwarm)
plt.colorbar()
plt.xlabel('log learning rate')
plt.ylabel('log regularization strength')
plt.title('CIFAR-10 validation accuracy')
plt.show()
```



```
[35]: # 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: 0.371000





## Inline question 2

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

### Your Answer:

Course Video

# softmax

June 8, 2024

```
[]: # This mounts your Google Drive to the Colab VM.
     from google.colab import drive
     drive.mount('/content/drive')
     # TODO: Enter the foldername in your Drive where you have saved the unzipped
     # assignment folder, e.g. 'cs231n/assignments/assignment1/'
     FOLDERNAME = None
     assert FOLDERNAME is not None, "[!] Enter the foldername."
     # Now that we've mounted your Drive, this ensures that
     # the Python interpreter of the Colab VM can load
     # python files from within it.
     import sys
     sys.path.append('/content/drive/My Drive/{}'.format(FOLDERNAME))
     # This downloads the CIFAR-10 dataset to your Drive
     # if it doesn't already exist.
     %cd /content/drive/My\ Drive/$FOLDERNAME/cs231n/datasets/
     !bash get datasets.sh
     %cd /content/drive/My\ Drive/$FOLDERNAME
```

### 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

```
[1]: import random import numpy as np
```

```
[2]: def get_CIFAR10_data(num_training=49000, num_validation=1000, num_test=1000,
      \rightarrownum dev=500):
         11 11 11
         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.
         11 11 11
         # Load the raw CIFAR-10 data
         cifar10_dir = 'cs231n/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)
         # 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
# 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,)
```

```
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 cs231n/classifiers/softmax.py.

```
[18]: # First implement the naive softmax loss function with nested loops.
# Open the file cs231n/classifiers/softmax.py and implement the
# softmax_loss_naive function.

from cs231n.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: 2.396957

sanity check: 2.302585

# Inline Question 1

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

### Your Answer:

10 0.1

0.1

```
[19]: # 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 cs231n.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: 0.994886 analytic: 0.994886, relative error: 4.040906e-08 numerical: -0.110419 analytic: -0.110419, relative error: 5.890618e-08 numerical: 2.673210 analytic: 2.673210, relative error: 3.454302e-08 numerical: -0.554525 analytic: -0.554525, relative error: 9.031858e-08 numerical: -0.064739 analytic: -0.064739, relative error: 4.892423e-07 numerical: 0.437539 analytic: 0.437539, relative error: 1.084708e-08 numerical: 2.071003 analytic: 2.071003, relative error: 1.726489e-08
```

```
numerical: -0.661887 analytic: -0.661887, relative error: 8.552310e-08
     numerical: 2.238773 analytic: 2.238772, relative error: 4.553863e-08
     numerical: -1.071853 analytic: -1.071853, relative error: 1.000644e-09
     numerical: -0.529727 analytic: -0.529727, relative error: 1.408935e-07
     numerical: -1.015710 analytic: -1.015710, relative error: 3.644312e-08
     numerical: 3.409544 analytic: 3.409544, relative error: 1.114026e-08
     numerical: 1.811444 analytic: 1.811444, relative error: 1.141884e-10
     numerical: -0.211312 analytic: -0.211312, relative error: 1.126727e-07
     numerical: -0.107410 analytic: -0.107410, relative error: 6.769405e-07
     numerical: -0.524265 analytic: -0.524265, relative error: 6.052658e-08
     numerical: -0.466537 analytic: -0.466538, relative error: 1.119220e-07
     numerical: 3.848656 analytic: 3.848656, relative error: 9.182922e-09
     numerical: -1.667093 analytic: -1.667093, relative error: 3.443010e-08
[28]: # 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 \Box
       ⇔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 cs231n.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: 2.396957e+00 computed in 0.058194s vectorized loss: 2.396957e+00 computed in 0.003887s

Loss difference: 0.000000 Gradient difference: 0.000000

[38]: # 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 cs231n.classifiers import Softmax
results = {}
best_val = -1
best_softmax = 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 softmax classifer in best softmax.
# Provided as a reference. You may or may not want to change these
 ⇔hyperparameters
learning_rates = [2e-7, 0.75e-7, 1.5e-7, 1.25e-7, 0.75e-7]
regularization_strengths = [3e4, 3.25e4, 3.5e4,
                         3.75e4, 4e4, 4.25e4, 4.5e4, 4.75e4, 5e4]
# *****START OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)****
for lr in learning rates:
   for reg in regularization_strengths:
       softmax = Softmax()
       softmax.train(X_train, y_train, lr, reg, num_iters=1000)
       y_train_pred, y_val_pred = softmax.predict(
          X_train), softmax.predict(X_val)
       results[(lr, reg)] = np.mean(
          y_train == y_train_pred), np.mean(y_val == y_val_pred)
       if results[(lr, reg)][1] > best_val:
          best_val = results[(lr, reg)][1]
          best_softmax = softmax
# *****END OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)****
# 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 7.500000e-08 reg 3.000000e+04 train accuracy: 0.329673 val accuracy: 0.342000
lr 7.500000e-08 reg 3.250000e+04 train accuracy: 0.321510 val accuracy: 0.337000

```
1r 7.500000e-08 reg 3.500000e+04 train accuracy: 0.317959 val accuracy: 0.332000
lr 7.500000e-08 reg 3.750000e+04 train accuracy: 0.318286 val accuracy: 0.336000
lr 7.500000e-08 reg 4.000000e+04 train accuracy: 0.307612 val accuracy: 0.322000
lr 7.500000e-08 reg 4.250000e+04 train accuracy: 0.314265 val accuracy: 0.331000
1r 7.500000e-08 reg 4.500000e+04 train accuracy: 0.309429 val accuracy: 0.329000
1r 7.500000e-08 reg 4.750000e+04 train accuracy: 0.303592 val accuracy: 0.322000
1r 7.500000e-08 reg 5.000000e+04 train accuracy: 0.308327 val accuracy: 0.324000
lr 1.250000e-07 reg 3.000000e+04 train accuracy: 0.315878 val accuracy: 0.332000
lr 1.250000e-07 reg 3.250000e+04 train accuracy: 0.329592 val accuracy: 0.346000
lr 1.250000e-07 reg 3.500000e+04 train accuracy: 0.316469 val accuracy: 0.331000
lr 1.250000e-07 reg 3.750000e+04 train accuracy: 0.313388 val accuracy: 0.324000
lr 1.250000e-07 reg 4.000000e+04 train accuracy: 0.320082 val accuracy: 0.335000
lr 1.250000e-07 reg 4.250000e+04 train accuracy: 0.312694 val accuracy: 0.329000
lr 1.250000e-07 reg 4.500000e+04 train accuracy: 0.309224 val accuracy: 0.323000
lr 1.250000e-07 reg 4.750000e+04 train accuracy: 0.306224 val accuracy: 0.323000
lr 1.250000e-07 reg 5.000000e+04 train accuracy: 0.305714 val accuracy: 0.325000
lr 1.500000e-07 reg 3.000000e+04 train accuracy: 0.324347 val accuracy: 0.339000
lr 1.500000e-07 reg 3.250000e+04 train accuracy: 0.315755 val accuracy: 0.335000
lr 1.500000e-07 reg 3.500000e+04 train accuracy: 0.314959 val accuracy: 0.330000
lr 1.500000e-07 reg 3.750000e+04 train accuracy: 0.316796 val accuracy: 0.328000
lr 1.500000e-07 reg 4.000000e+04 train accuracy: 0.316408 val accuracy: 0.328000
lr 1.500000e-07 reg 4.250000e+04 train accuracy: 0.313347 val accuracy: 0.338000
lr 1.500000e-07 reg 4.500000e+04 train accuracy: 0.301306 val accuracy: 0.317000
lr 1.500000e-07 reg 4.750000e+04 train accuracy: 0.310082 val accuracy: 0.331000
lr 1.500000e-07 reg 5.000000e+04 train accuracy: 0.312041 val accuracy: 0.322000
lr 2.000000e-07 reg 3.000000e+04 train accuracy: 0.325612 val accuracy: 0.341000
lr 2.000000e-07 reg 3.250000e+04 train accuracy: 0.313347 val accuracy: 0.328000
lr 2.000000e-07 reg 3.500000e+04 train accuracy: 0.311633 val accuracy: 0.326000
lr 2.000000e-07 reg 3.750000e+04 train accuracy: 0.313388 val accuracy: 0.325000
lr 2.000000e-07 reg 4.000000e+04 train accuracy: 0.307449 val accuracy: 0.331000
lr 2.000000e-07 reg 4.250000e+04 train accuracy: 0.312286 val accuracy: 0.335000
lr 2.000000e-07 reg 4.500000e+04 train accuracy: 0.318714 val accuracy: 0.328000
lr 2.000000e-07 reg 4.750000e+04 train accuracy: 0.306367 val accuracy: 0.321000
lr 2.000000e-07 reg 5.000000e+04 train accuracy: 0.314122 val accuracy: 0.327000
best validation accuracy achieved during cross-validation: 0.346000
```

```
[41]: # 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: 0.339000

### Inline Question 2 - True or False

Suppose the overall training loss is defined as the sum of the per-datapoint loss over all training examples. It is 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:

SVM

SVM

Detailed Explanation from Mantasu

First, let's revise how individual loss functions look ( $\hat{y}$  - score, i - any class label, c - true class label):

$$\text{SVM loss:} \ \ L_n = \sum_{i \neq c}^I \max(0, \hat{y}_i - \hat{y}_c + \Delta)$$

Softmax loss: 
$$L_n = -\log \left( \frac{e^{\hat{y}_c}}{\sum_{i=1}^{I} e^{\hat{y}_i}} \right)$$

Explanation: \* The nature of **SVM loss** is to teach the model to assign a high enough score for the correct label. The total loss is unaffected if the scores for the incorrect labels are already lower by some margin than the score for a correct label because nothing is added to the overall loss. \* The nature of **Softmax loss** is to make the model raise the probability for the correct label to 1 while the probabilities for other labels decrease to 0. The loss is always affected because no matter what score is produced for an incorrect label, it still induces some probability which does not allow the true label to reach the probability of 1. \* Aside: theoretically, if all the scores for incorrect classes were  $-\infty$ , the probability of the correct class would be 1 so the individual loss would be  $-\log 1 = 0$  and that would not affect the overall loss \* Assuming the overall loss is just the sum of every individual loss, the data point which produces safe score differences between the incorrect labels and a correct one is a data point that does not affect the **SVM loss**, however, it is guaranteed that it will affect the **Softmax loss** by some (however tiny) amount.

Example: \* Suppose we get a score vector  $\mathbf{\hat{y}} = \begin{bmatrix} 1 \\ 0 \end{bmatrix}$  and that the first entry represents the score for a correct class, i.e.,  $\hat{y}_1 = 1$ . \* Assuming our margin for **SVM loss** is 1:  $\Delta = 1$ , the individual loss for  $\mathbf{\hat{y}}$  is 0 because it satisfies the margin. However, the individual **Softmax loss** will produce a value of  $\log(1 + e) - 1$  which will affect the overall loss.

Hence it is possible to add a datapoint which would not affect the (described) overall **SVM loss** but would affect the overall **Softmax loss**.

```
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])
```





[]:

# two\_layer\_net

June 8, 2024

```
[]: # This mounts your Google Drive to the Colab VM.
     from google.colab import drive
     drive.mount('/content/drive')
     # TODO: Enter the foldername in your Drive where you have saved the unzipped
     # assignment folder, e.g. 'cs231n/assignments/assignment1/'
     FOLDERNAME = None
     assert FOLDERNAME is not None, "[!] Enter the foldername."
     # Now that we've mounted your Drive, this ensures that
     # the Python interpreter of the Colab VM can load
     # python files from within it.
     import sys
     sys.path.append('/content/drive/My Drive/{}'.format(FOLDERNAME))
     # This downloads the CIFAR-10 dataset to your Drive
     # if it doesn't already exist.
     %cd /content/drive/My\ Drive/$FOLDERNAME/cs231n/datasets/
     !bash get datasets.sh
     %cd /content/drive/My\ Drive/$FOLDERNAME
```

# 1 Fully-Connected Neural Nets

In this exercise we will implement fully-connected networks using a modular approach. For each layer we will implement a forward and a backward function. The forward function will receive inputs, weights, and other parameters and will return both an output and a cache object storing data needed for the backward pass, like this:

```
def layer_forward(x, w):
    """ Receive inputs x and weights w """
    # Do some computations ...
    z = # ... some intermediate value
    # Do some more computations ...
    out = # the output

cache = (x, w, z, out) # Values we need to compute gradients
    return out, cache
```

The backward pass will receive upstream derivatives and the cache object, and will return gradients with respect to the inputs and weights, like this:

```
def layer_backward(dout, cache):
    """
    Receive dout (derivative of loss with respect to outputs) and cache,
    and compute derivative with respect to inputs.
    """
    # Unpack cache values
    x, w, z, out = cache

# Use values in cache to compute derivatives
    dx = # Derivative of loss with respect to x
    dw = # Derivative of loss with respect to w

return dx, dw
```

After implementing a bunch of layers this way, we will be able to easily combine them to build classifiers with different architectures.

```
[2]: # As usual, a bit of setup
     from __future__ import print_function
     import time
     import numpy as np
     import matplotlib.pyplot as plt
     from cs231n.classifiers.fc net import *
     from cs231n.data_utils import get_CIFAR10_data
     from cs231n.gradient_check import eval_numerical_gradient,__
      ⇔eval_numerical_gradient_array
     from cs231n.solver import Solver
     %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))))
```

```
[3]: # Load the (preprocessed) CIFAR10 data.
```

```
data = get_CIFAR10_data()
for k, v in list(data.items()):
    print(('%s: ' % k, v.shape))

('X_train: ', (49000, 3, 32, 32))
('y_train: ', (49000,))
('X_val: ', (1000, 3, 32, 32))
```

# 2 Affine layer: forward

('X\_test: ', (1000, 3, 32, 32))

('y\_val: ', (1000,))

('y\_test: ', (1000,))

Open the file cs231n/layers.py and implement the affine\_forward function.

Once you are done you can test your implementaion by running the following:

```
[4]: # Test the affine_forward function
     num inputs = 2
     input\_shape = (4, 5, 6)
     output dim = 3
     input_size = num_inputs * np.prod(input_shape)
     weight_size = output_dim * np.prod(input_shape)
     x = np.linspace(-0.1, 0.5, num=input_size).reshape(num_inputs, *input_shape)
     w = np.linspace(-0.2, 0.3, num=weight_size).reshape(np.prod(input_shape),__
      →output_dim)
     b = np.linspace(-0.3, 0.1, num=output dim)
     out, _ = affine_forward(x, w, b)
     correct_out = np.array([[ 1.49834967, 1.70660132, 1.91485297],
                             [ 3.25553199, 3.5141327, 3.77273342]])
     # Compare your output with ours. The error should be around e-9 or less.
     print('Testing affine_forward function:')
     print('difference: ', rel_error(out, correct_out))
```

Testing affine\_forward function: difference: 9.769849468192957e-10

# 3 Affine layer: backward

Now implement the affine\_backward function and test your implementation using numeric gradient checking.

```
[5]: # Test the affine backward function
    np.random.seed(231)
     x = np.random.randn(10, 2, 3)
     w = np.random.randn(6, 5)
     b = np.random.randn(5)
     dout = np.random.randn(10, 5)
     dx_num = eval_numerical_gradient_array(lambda x: affine_forward(x, w, b)[0], x,_
     dw_num = eval_numerical_gradient_array(lambda w: affine_forward(x, w, b)[0], w,__
      ⊸dout)
     db_num = eval_numerical_gradient_array(lambda b: affine_forward(x, w, b)[0], b,_

dout)
     _, cache = affine_forward(x, w, b)
     dx, dw, db = affine_backward(dout, cache)
     # The error should be around e-10 or less
     print('Testing affine_backward function:')
     print('dx error: ', rel_error(dx_num, dx))
     print('dw error: ', rel_error(dw_num, dw))
     print('db error: ', rel_error(db_num, db))
```

Testing affine\_backward function: dx error: 5.399100368651805e-11 dw error: 9.904211865398145e-11 db error: 2.4122867568119087e-11

## 4 ReLU activation: forward

Implement the forward pass for the ReLU activation function in the relu\_forward function and test your implementation using the following:

```
Testing relu_forward function: difference: 4.999999798022158e-08
```

## 5 ReLU activation: backward

Now implement the backward pass for the ReLU activation function in the relu\_backward function and test your implementation using numeric gradient checking:

```
[7]: np.random.seed(231)
    x = np.random.randn(10, 10)
    dout = np.random.randn(*x.shape)

    dx_num = eval_numerical_gradient_array(lambda x: relu_forward(x)[0], x, dout)

_, cache = relu_forward(x)
    dx = relu_backward(dout, cache)

# The error should be on the order of e-12
    print('Testing relu_backward function:')
    print('dx error: ', rel_error(dx_num, dx))
```

Testing relu\_backward function: dx error: 3.2756349136310288e-12

## 5.1 Inline Question 1:

We've only asked you to implement ReLU, but there are a number of different activation functions that one could use in neural networks, each with its pros and cons. In particular, an issue commonly seen with activation functions is getting zero (or close to zero) gradient flow during backpropagation. Which of the following activation functions have this problem? If you consider these functions in the one dimensional case, what types of input would lead to this behaviour? 1. Sigmoid 2. ReLU 3. Leaky ReLU

#### 5.2 Answer:

```
Answer: 1, 2.

" " course notes 0

• Sigmoid. Sigmoid 0 1

• Relu. Relu
```

# 6 "Sandwich" layers

There are some common patterns of layers that are frequently used in neural nets. For example, affine layers are frequently followed by a ReLU nonlinearity. To make these common patterns easy, we define several convenience layers in the file cs231n/layer utils.py.

For now take a look at the affine\_relu\_forward and affine\_relu\_backward functions, and run the following to numerically gradient check the backward pass:

```
[8]: from cs231n.layer_utils import affine_relu_forward, affine_relu_backward
     np.random.seed(231)
     x = np.random.randn(2, 3, 4)
     w = np.random.randn(12, 10)
     b = np.random.randn(10)
     dout = np.random.randn(2, 10)
     out, cache = affine_relu_forward(x, w, b)
     dx, dw, db = affine relu backward(dout, cache)
     dx num = eval numerical gradient array(lambda x: affine relu forward(x, w, )
      \rightarrowb)[0], x, dout)
     dw num = eval_numerical_gradient_array(lambda w: affine relu_forward(x, w,__
      \hookrightarrowb)[0], w, dout)
     db num = eval numerical gradient array(lambda b: affine relu forward(x, w, ...
      →b)[0], b, dout)
     # Relative error should be around e-10 or less
     print('Testing affine relu forward and affine relu backward:')
     print('dx error: ', rel_error(dx_num, dx))
     print('dw error: ', rel_error(dw_num, dw))
     print('db error: ', rel_error(db_num, db))
```

Testing affine\_relu\_forward and affine\_relu\_backward:

dx error: 2.299579177309368e-11
dw error: 8.162011105764925e-11
db error: 7.826724021458994e-12

# 7 Loss layers: Softmax and SVM

Now implement the loss and gradient for softmax and SVM in the softmax\_loss and svm\_loss function in cs231n/layers.py. These should be similar to what you implemented in cs231n/classifiers/softmax.py and cs231n/classifiers/linear\_svm.py.

You can make sure that the implementations are correct by running the following:

```
[9]: np.random.seed(231)
num_classes, num_inputs = 10, 50
x = 0.001 * np.random.randn(num_inputs, num_classes)
y = np.random.randint(num_classes, size=num_inputs)

dx_num = eval_numerical_gradient(lambda x: svm_loss(x, y)[0], x, verbose=False)
loss, dx = svm_loss(x, y)

# Test svm_loss function. Loss should be around 9 and dx error should be around_u
the order of e-9
print('Testing svm_loss:')
```

Testing svm\_loss:

loss: 8.999602749096233

dx error: 1.4021566006651672e-09

Testing softmax\_loss:

loss: 2.3025458445007376

dx error: 8.234144091578429e-09

## 8 Two-layer network

Open the file cs231n/classifiers/fc\_net.py and complete the implementation of the TwoLayerNet class. Read through it to make sure you understand the API. You can run the cell below to test your implementation.

```
[10]: np.random.seed(231)
      N, D, H, C = 3, 5, 50, 7
      X = np.random.randn(N, D)
      y = np.random.randint(C, size=N)
      std = 1e-3
      model = TwoLayerNet(input_dim=D, hidden_dim=H, num_classes=C, weight_scale=std)
      print('Testing initialization ... ')
      W1_std = abs(model.params['W1'].std() - std)
      b1 = model.params['b1']
      W2_std = abs(model.params['W2'].std() - std)
      b2 = model.params['b2']
      assert W1_std < std / 10, 'First layer weights do not seem right'
      assert np.all(b1 == 0), 'First layer biases do not seem right'
      assert W2_std < std / 10, 'Second layer weights do not seem right'
      assert np.all(b2 == 0), 'Second layer biases do not seem right'
      print('Testing test-time forward pass ... ')
      model.params['W1'] = np.linspace(-0.7, 0.3, num=D*H).reshape(D, H)
```

```
model.params['b1'] = np.linspace(-0.1, 0.9, num=H)
model.params['W2'] = np.linspace(-0.3, 0.4, num=H*C).reshape(H, C)
model.params['b2'] = np.linspace(-0.9, 0.1, num=C)
X = np.linspace(-5.5, 4.5, num=N*D).reshape(D, N).T
scores = model.loss(X)
correct_scores = np.asarray(
  [[11.53165108, 12.2917344, 13.05181771, 13.81190102, 14.57198434, 15.
 →33206765, 16.09215096],
   [12.05769098, 12.74614105, 13.43459113, 14.1230412, 14.81149128, 15.
 →49994135, 16.18839143],
   [12.58373087, 13.20054771, 13.81736455, 14.43418138, 15.05099822, 15.
 →66781506, 16.2846319 ]])
scores_diff = np.abs(scores - correct_scores).sum()
assert scores_diff < 1e-6, 'Problem with test-time forward pass'
print('Testing training loss (no regularization)')
y = np.asarray([0, 5, 1])
loss, grads = model.loss(X, y)
correct_loss = 3.4702243556
assert abs(loss - correct_loss) < 1e-10, 'Problem with training-time loss'</pre>
model.reg = 1.0
loss, grads = model.loss(X, y)
correct_loss = 26.5948426952
assert abs(loss - correct_loss) < 1e-10, 'Problem with regularization loss'
# Errors should be around e-7 or less
for reg in [0.0, 0.7]:
  print('Running numeric gradient check with reg = ', reg)
  model.reg = reg
  loss, grads = model.loss(X, y)
  for name in sorted(grads):
    f = lambda _: model.loss(X, y)[0]
    grad_num = eval_numerical_gradient(f, model.params[name], verbose=False)
    print('%s relative error: %.2e' % (name, rel_error(grad_num, grads[name])))
Testing initialization ...
Testing test-time forward pass ...
Testing training loss (no regularization)
Running numeric gradient check with reg = 0.0
W1 relative error: 1.83e-08
W2 relative error: 3.20e-10
b1 relative error: 9.83e-09
b2 relative error: 4.33e-10
Running numeric gradient check with reg = 0.7
W1 relative error: 2.53e-07
```

W2 relative error: 2.85e-08 b1 relative error: 1.56e-08 b2 relative error: 9.09e-10

## 9 Solver

Open the file cs231n/solver.py and read through it to familiarize yourself with the API. After doing so, use a Solver instance to train a TwoLayerNet that achieves about 36% accuracy on the validation set.

```
[11]: input size = 32 * 32 * 3
   hidden_size = 50
   num_classes = 10
   model = TwoLayerNet(input_size, hidden_size, num_classes)
   solver = None
   # TODO: Use a Solver instance to train a TwoLayerNet that achieves about 36% #
   # accuracy on the validation set.
   # *****START OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE) *****
   solver = Solver(model, data, optim_config={'learning_rate': 1e-3})
   solver.train()
   # *****END OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)****
   END OF YOUR CODE
```

```
(Iteration 1 / 4900) loss: 2.300089
(Epoch 0 / 10) train acc: 0.171000; val acc: 0.170000
(Iteration 11 / 4900) loss: 2.258416
(Iteration 21 / 4900) loss: 2.160948
(Iteration 31 / 4900) loss: 2.016287
(Iteration 41 / 4900) loss: 2.138642
(Iteration 51 / 4900) loss: 2.059372
(Iteration 61 / 4900) loss: 1.897230
(Iteration 71 / 4900) loss: 1.801362
(Iteration 81 / 4900) loss: 1.816051
(Iteration 91 / 4900) loss: 1.767155
(Iteration 101 / 4900) loss: 1.782419
(Iteration 111 / 4900) loss: 1.861558
(Iteration 121 / 4900) loss: 1.650334
(Iteration 131 / 4900) loss: 1.698891
(Iteration 141 / 4900) loss: 1.833884
(Iteration 151 / 4900) loss: 1.763588
(Iteration 161 / 4900) loss: 1.635848
```

```
(Iteration 171 / 4900) loss: 1.962973
(Iteration 181 / 4900) loss: 1.716530
(Iteration 191 / 4900) loss: 1.652857
(Iteration 201 / 4900) loss: 1.803466
(Iteration 211 / 4900) loss: 1.646484
(Iteration 221 / 4900) loss: 1.753302
(Iteration 231 / 4900) loss: 1.665357
(Iteration 241 / 4900) loss: 1.558509
(Iteration 251 / 4900) loss: 1.553072
(Iteration 261 / 4900) loss: 1.875181
(Iteration 271 / 4900) loss: 1.571899
(Iteration 281 / 4900) loss: 1.671443
(Iteration 291 / 4900) loss: 1.563806
(Iteration 301 / 4900) loss: 1.712676
(Iteration 311 / 4900) loss: 1.687622
(Iteration 321 / 4900) loss: 1.695280
(Iteration 331 / 4900) loss: 1.704362
(Iteration 341 / 4900) loss: 1.763766
(Iteration 351 / 4900) loss: 1.485922
(Iteration 361 / 4900) loss: 1.552275
(Iteration 371 / 4900) loss: 1.833757
(Iteration 381 / 4900) loss: 1.760161
(Iteration 391 / 4900) loss: 1.494989
(Iteration 401 / 4900) loss: 1.693946
(Iteration 411 / 4900) loss: 1.547319
(Iteration 421 / 4900) loss: 1.333266
(Iteration 431 / 4900) loss: 1.526631
(Iteration 441 / 4900) loss: 1.568662
(Iteration 451 / 4900) loss: 1.543955
(Iteration 461 / 4900) loss: 1.413793
(Iteration 471 / 4900) loss: 1.519715
(Iteration 481 / 4900) loss: 1.548025
(Epoch 1 / 10) train acc: 0.399000; val_acc: 0.428000
(Iteration 491 / 4900) loss: 1.662400
(Iteration 501 / 4900) loss: 1.712717
(Iteration 511 / 4900) loss: 1.563596
(Iteration 521 / 4900) loss: 1.573986
(Iteration 531 / 4900) loss: 1.608347
(Iteration 541 / 4900) loss: 1.693400
(Iteration 551 / 4900) loss: 1.612438
(Iteration 561 / 4900) loss: 1.741232
(Iteration 571 / 4900) loss: 1.707330
(Iteration 581 / 4900) loss: 1.492098
(Iteration 591 / 4900) loss: 1.617551
(Iteration 601 / 4900) loss: 1.460141
(Iteration 611 / 4900) loss: 1.392740
(Iteration 621 / 4900) loss: 1.578618
(Iteration 631 / 4900) loss: 1.568944
```

```
(Iteration 641 / 4900) loss: 1.571395
(Iteration 651 / 4900) loss: 1.506858
(Iteration 661 / 4900) loss: 1.584271
(Iteration 671 / 4900) loss: 1.525855
(Iteration 681 / 4900) loss: 1.505960
(Iteration 691 / 4900) loss: 1.517432
(Iteration 701 / 4900) loss: 1.563845
(Iteration 711 / 4900) loss: 1.443370
(Iteration 721 / 4900) loss: 1.691380
(Iteration 731 / 4900) loss: 1.569986
(Iteration 741 / 4900) loss: 1.611344
(Iteration 751 / 4900) loss: 1.426493
(Iteration 761 / 4900) loss: 1.686239
(Iteration 771 / 4900) loss: 1.582190
(Iteration 781 / 4900) loss: 1.605619
(Iteration 791 / 4900) loss: 1.489545
(Iteration 801 / 4900) loss: 1.535489
(Iteration 811 / 4900) loss: 1.618398
(Iteration 821 / 4900) loss: 1.546814
(Iteration 831 / 4900) loss: 1.524163
(Iteration 841 / 4900) loss: 1.560121
(Iteration 851 / 4900) loss: 1.728512
(Iteration 861 / 4900) loss: 1.451588
(Iteration 871 / 4900) loss: 1.512195
(Iteration 881 / 4900) loss: 1.714012
(Iteration 891 / 4900) loss: 1.733377
(Iteration 901 / 4900) loss: 1.359039
(Iteration 911 / 4900) loss: 1.581966
(Iteration 921 / 4900) loss: 1.428448
(Iteration 931 / 4900) loss: 1.561284
(Iteration 941 / 4900) loss: 1.568419
(Iteration 951 / 4900) loss: 1.518509
(Iteration 961 / 4900) loss: 1.494198
(Iteration 971 / 4900) loss: 1.328252
(Epoch 2 / 10) train acc: 0.468000; val acc: 0.444000
(Iteration 981 / 4900) loss: 1.446261
(Iteration 991 / 4900) loss: 1.393918
(Iteration 1001 / 4900) loss: 1.339096
(Iteration 1011 / 4900) loss: 1.344386
(Iteration 1021 / 4900) loss: 1.256831
(Iteration 1031 / 4900) loss: 1.422381
(Iteration 1041 / 4900) loss: 1.738975
(Iteration 1051 / 4900) loss: 1.553349
(Iteration 1061 / 4900) loss: 1.603838
(Iteration 1071 / 4900) loss: 1.486361
(Iteration 1081 / 4900) loss: 1.417496
(Iteration 1091 / 4900) loss: 1.429808
(Iteration 1101 / 4900) loss: 1.397059
```

```
(Iteration 1111 / 4900) loss: 1.653973
(Iteration 1121 / 4900) loss: 1.450812
(Iteration 1131 / 4900) loss: 1.437969
(Iteration 1141 / 4900) loss: 1.448549
(Iteration 1151 / 4900) loss: 1.273936
(Iteration 1161 / 4900) loss: 1.430609
(Iteration 1171 / 4900) loss: 1.559522
(Iteration 1181 / 4900) loss: 1.324973
(Iteration 1191 / 4900) loss: 1.679667
(Iteration 1201 / 4900) loss: 1.499237
(Iteration 1211 / 4900) loss: 1.509699
(Iteration 1221 / 4900) loss: 1.402940
(Iteration 1231 / 4900) loss: 1.277391
(Iteration 1241 / 4900) loss: 1.450792
(Iteration 1251 / 4900) loss: 1.351350
(Iteration 1261 / 4900) loss: 1.600031
(Iteration 1271 / 4900) loss: 1.534798
(Iteration 1281 / 4900) loss: 1.448485
(Iteration 1291 / 4900) loss: 1.116751
(Iteration 1301 / 4900) loss: 1.349809
(Iteration 1311 / 4900) loss: 1.348709
(Iteration 1321 / 4900) loss: 1.377703
(Iteration 1331 / 4900) loss: 1.702516
(Iteration 1341 / 4900) loss: 1.418050
(Iteration 1351 / 4900) loss: 1.537872
(Iteration 1361 / 4900) loss: 1.391911
(Iteration 1371 / 4900) loss: 1.354068
(Iteration 1381 / 4900) loss: 1.705837
(Iteration 1391 / 4900) loss: 1.600954
(Iteration 1401 / 4900) loss: 1.449001
(Iteration 1411 / 4900) loss: 1.453994
(Iteration 1421 / 4900) loss: 1.489336
(Iteration 1431 / 4900) loss: 1.495201
(Iteration 1441 / 4900) loss: 1.226016
(Iteration 1451 / 4900) loss: 1.263878
(Iteration 1461 / 4900) loss: 1.477084
(Epoch 3 / 10) train acc: 0.466000; val acc: 0.429000
(Iteration 1471 / 4900) loss: 1.400271
(Iteration 1481 / 4900) loss: 1.309975
(Iteration 1491 / 4900) loss: 1.410661
(Iteration 1501 / 4900) loss: 1.319916
(Iteration 1511 / 4900) loss: 1.393971
(Iteration 1521 / 4900) loss: 1.445102
(Iteration 1531 / 4900) loss: 1.489275
(Iteration 1541 / 4900) loss: 1.500323
(Iteration 1551 / 4900) loss: 1.613116
(Iteration 1561 / 4900) loss: 1.215678
(Iteration 1571 / 4900) loss: 1.375205
```

```
(Iteration 1581 / 4900) loss: 1.250286
(Iteration 1591 / 4900) loss: 1.442918
(Iteration 1601 / 4900) loss: 1.441210
(Iteration 1611 / 4900) loss: 1.584907
(Iteration 1621 / 4900) loss: 1.555882
(Iteration 1631 / 4900) loss: 1.538311
(Iteration 1641 / 4900) loss: 1.350846
(Iteration 1651 / 4900) loss: 1.586443
(Iteration 1661 / 4900) loss: 1.244578
(Iteration 1671 / 4900) loss: 1.529737
(Iteration 1681 / 4900) loss: 1.490461
(Iteration 1691 / 4900) loss: 1.429346
(Iteration 1701 / 4900) loss: 1.351760
(Iteration 1711 / 4900) loss: 1.439255
(Iteration 1721 / 4900) loss: 1.404750
(Iteration 1731 / 4900) loss: 1.478821
(Iteration 1741 / 4900) loss: 1.645068
(Iteration 1751 / 4900) loss: 1.452901
(Iteration 1761 / 4900) loss: 1.591717
(Iteration 1771 / 4900) loss: 1.594355
(Iteration 1781 / 4900) loss: 1.531285
(Iteration 1791 / 4900) loss: 1.314133
(Iteration 1801 / 4900) loss: 1.430905
(Iteration 1811 / 4900) loss: 1.280109
(Iteration 1821 / 4900) loss: 1.330116
(Iteration 1831 / 4900) loss: 1.325624
(Iteration 1841 / 4900) loss: 1.394343
(Iteration 1851 / 4900) loss: 1.371149
(Iteration 1861 / 4900) loss: 1.345813
(Iteration 1871 / 4900) loss: 1.425325
(Iteration 1881 / 4900) loss: 1.410831
(Iteration 1891 / 4900) loss: 1.484850
(Iteration 1901 / 4900) loss: 1.347602
(Iteration 1911 / 4900) loss: 1.382680
(Iteration 1921 / 4900) loss: 1.733093
(Iteration 1931 / 4900) loss: 1.418113
(Iteration 1941 / 4900) loss: 1.313570
(Iteration 1951 / 4900) loss: 1.435940
(Epoch 4 / 10) train acc: 0.504000; val_acc: 0.451000
(Iteration 1961 / 4900) loss: 1.232119
(Iteration 1971 / 4900) loss: 1.611118
(Iteration 1981 / 4900) loss: 1.624994
(Iteration 1991 / 4900) loss: 1.439298
(Iteration 2001 / 4900) loss: 1.452116
(Iteration 2011 / 4900) loss: 1.377280
(Iteration 2021 / 4900) loss: 1.345182
(Iteration 2031 / 4900) loss: 1.723098
(Iteration 2041 / 4900) loss: 1.337191
```

```
(Iteration 2051 / 4900) loss: 1.367573
(Iteration 2061 / 4900) loss: 1.363604
(Iteration 2071 / 4900) loss: 1.373193
(Iteration 2081 / 4900) loss: 1.527777
(Iteration 2091 / 4900) loss: 1.294007
(Iteration 2101 / 4900) loss: 1.369872
(Iteration 2111 / 4900) loss: 1.252515
(Iteration 2121 / 4900) loss: 1.632426
(Iteration 2131 / 4900) loss: 1.221843
(Iteration 2141 / 4900) loss: 1.588294
(Iteration 2151 / 4900) loss: 1.382462
(Iteration 2161 / 4900) loss: 1.451325
(Iteration 2171 / 4900) loss: 1.232785
(Iteration 2181 / 4900) loss: 1.428744
(Iteration 2191 / 4900) loss: 1.528158
(Iteration 2201 / 4900) loss: 1.561029
(Iteration 2211 / 4900) loss: 1.545789
(Iteration 2221 / 4900) loss: 1.216724
(Iteration 2231 / 4900) loss: 1.368028
(Iteration 2241 / 4900) loss: 1.419180
(Iteration 2251 / 4900) loss: 1.338375
(Iteration 2261 / 4900) loss: 1.297813
(Iteration 2271 / 4900) loss: 1.293287
(Iteration 2281 / 4900) loss: 1.397591
(Iteration 2291 / 4900) loss: 1.448465
(Iteration 2301 / 4900) loss: 1.242521
(Iteration 2311 / 4900) loss: 1.520294
(Iteration 2321 / 4900) loss: 1.354138
(Iteration 2331 / 4900) loss: 1.309486
(Iteration 2341 / 4900) loss: 1.309089
(Iteration 2351 / 4900) loss: 1.420478
(Iteration 2361 / 4900) loss: 1.299870
(Iteration 2371 / 4900) loss: 1.374915
(Iteration 2381 / 4900) loss: 1.128089
(Iteration 2391 / 4900) loss: 1.195701
(Iteration 2401 / 4900) loss: 1.433959
(Iteration 2411 / 4900) loss: 1.271562
(Iteration 2421 / 4900) loss: 1.342458
(Iteration 2431 / 4900) loss: 1.389719
(Iteration 2441 / 4900) loss: 1.508970
(Epoch 5 / 10) train acc: 0.493000; val_acc: 0.465000
(Iteration 2451 / 4900) loss: 1.403120
(Iteration 2461 / 4900) loss: 1.362264
(Iteration 2471 / 4900) loss: 1.612289
(Iteration 2481 / 4900) loss: 1.489801
(Iteration 2491 / 4900) loss: 1.429105
(Iteration 2501 / 4900) loss: 1.229392
(Iteration 2511 / 4900) loss: 1.320781
```

```
(Iteration 2521 / 4900) loss: 1.348747
(Iteration 2531 / 4900) loss: 1.265773
(Iteration 2541 / 4900) loss: 1.437888
(Iteration 2551 / 4900) loss: 1.245283
(Iteration 2561 / 4900) loss: 1.283544
(Iteration 2571 / 4900) loss: 1.332414
(Iteration 2581 / 4900) loss: 1.376941
(Iteration 2591 / 4900) loss: 1.242511
(Iteration 2601 / 4900) loss: 1.527411
(Iteration 2611 / 4900) loss: 1.485586
(Iteration 2621 / 4900) loss: 1.393928
(Iteration 2631 / 4900) loss: 1.332372
(Iteration 2641 / 4900) loss: 1.302352
(Iteration 2651 / 4900) loss: 1.707333
(Iteration 2661 / 4900) loss: 1.350690
(Iteration 2671 / 4900) loss: 1.335264
(Iteration 2681 / 4900) loss: 1.427223
(Iteration 2691 / 4900) loss: 1.449200
(Iteration 2701 / 4900) loss: 1.262844
(Iteration 2711 / 4900) loss: 1.164357
(Iteration 2721 / 4900) loss: 1.364723
(Iteration 2731 / 4900) loss: 1.279289
(Iteration 2741 / 4900) loss: 1.248773
(Iteration 2751 / 4900) loss: 1.449679
(Iteration 2761 / 4900) loss: 1.137617
(Iteration 2771 / 4900) loss: 1.293644
(Iteration 2781 / 4900) loss: 1.296194
(Iteration 2791 / 4900) loss: 1.434576
(Iteration 2801 / 4900) loss: 1.540281
(Iteration 2811 / 4900) loss: 1.651625
(Iteration 2821 / 4900) loss: 1.333709
(Iteration 2831 / 4900) loss: 1.406024
(Iteration 2841 / 4900) loss: 1.375746
(Iteration 2851 / 4900) loss: 1.317436
(Iteration 2861 / 4900) loss: 1.410073
(Iteration 2871 / 4900) loss: 1.326769
(Iteration 2881 / 4900) loss: 1.210212
(Iteration 2891 / 4900) loss: 1.361641
(Iteration 2901 / 4900) loss: 1.061335
(Iteration 2911 / 4900) loss: 1.490168
(Iteration 2921 / 4900) loss: 1.531250
(Iteration 2931 / 4900) loss: 1.734393
(Epoch 6 / 10) train acc: 0.537000; val_acc: 0.488000
(Iteration 2941 / 4900) loss: 1.343652
(Iteration 2951 / 4900) loss: 1.309466
(Iteration 2961 / 4900) loss: 1.137099
(Iteration 2971 / 4900) loss: 1.404897
(Iteration 2981 / 4900) loss: 1.499112
```

```
(Iteration 2991 / 4900) loss: 1.182964
(Iteration 3001 / 4900) loss: 1.375969
(Iteration 3011 / 4900) loss: 1.475385
(Iteration 3021 / 4900) loss: 1.297671
(Iteration 3031 / 4900) loss: 1.455192
(Iteration 3041 / 4900) loss: 1.288831
(Iteration 3051 / 4900) loss: 1.252507
(Iteration 3061 / 4900) loss: 1.273133
(Iteration 3071 / 4900) loss: 1.245781
(Iteration 3081 / 4900) loss: 1.526020
(Iteration 3091 / 4900) loss: 1.244705
(Iteration 3101 / 4900) loss: 1.168029
(Iteration 3111 / 4900) loss: 1.199980
(Iteration 3121 / 4900) loss: 1.616235
(Iteration 3131 / 4900) loss: 1.228325
(Iteration 3141 / 4900) loss: 1.372394
(Iteration 3151 / 4900) loss: 1.261772
(Iteration 3161 / 4900) loss: 1.472383
(Iteration 3171 / 4900) loss: 1.129794
(Iteration 3181 / 4900) loss: 1.396840
(Iteration 3191 / 4900) loss: 1.405426
(Iteration 3201 / 4900) loss: 1.270886
(Iteration 3211 / 4900) loss: 1.421649
(Iteration 3221 / 4900) loss: 1.503245
(Iteration 3231 / 4900) loss: 1.568326
(Iteration 3241 / 4900) loss: 1.393653
(Iteration 3251 / 4900) loss: 1.318088
(Iteration 3261 / 4900) loss: 1.332783
(Iteration 3271 / 4900) loss: 1.606893
(Iteration 3281 / 4900) loss: 1.276238
(Iteration 3291 / 4900) loss: 1.216147
(Iteration 3301 / 4900) loss: 1.360948
(Iteration 3311 / 4900) loss: 1.328076
(Iteration 3321 / 4900) loss: 1.329054
(Iteration 3331 / 4900) loss: 1.150582
(Iteration 3341 / 4900) loss: 1.499849
(Iteration 3351 / 4900) loss: 1.232690
(Iteration 3361 / 4900) loss: 1.357872
(Iteration 3371 / 4900) loss: 1.577824
(Iteration 3381 / 4900) loss: 1.394309
(Iteration 3391 / 4900) loss: 1.458092
(Iteration 3401 / 4900) loss: 1.218180
(Iteration 3411 / 4900) loss: 1.264785
(Iteration 3421 / 4900) loss: 1.199487
(Epoch 7 / 10) train acc: 0.531000; val_acc: 0.468000
(Iteration 3431 / 4900) loss: 1.311152
(Iteration 3441 / 4900) loss: 1.364160
(Iteration 3451 / 4900) loss: 1.306556
```

```
(Iteration 3461 / 4900) loss: 1.399136
(Iteration 3471 / 4900) loss: 1.317526
(Iteration 3481 / 4900) loss: 1.274651
(Iteration 3491 / 4900) loss: 1.581665
(Iteration 3501 / 4900) loss: 1.492338
(Iteration 3511 / 4900) loss: 1.366844
(Iteration 3521 / 4900) loss: 1.406317
(Iteration 3531 / 4900) loss: 1.242057
(Iteration 3541 / 4900) loss: 1.591405
(Iteration 3551 / 4900) loss: 1.339095
(Iteration 3561 / 4900) loss: 1.088366
(Iteration 3571 / 4900) loss: 1.433389
(Iteration 3581 / 4900) loss: 1.272246
(Iteration 3591 / 4900) loss: 1.236100
(Iteration 3601 / 4900) loss: 1.193582
(Iteration 3611 / 4900) loss: 1.481540
(Iteration 3621 / 4900) loss: 1.297923
(Iteration 3631 / 4900) loss: 1.281803
(Iteration 3641 / 4900) loss: 1.311063
(Iteration 3651 / 4900) loss: 1.380828
(Iteration 3661 / 4900) loss: 1.321568
(Iteration 3671 / 4900) loss: 1.343057
(Iteration 3681 / 4900) loss: 1.389960
(Iteration 3691 / 4900) loss: 1.223307
(Iteration 3701 / 4900) loss: 1.478843
(Iteration 3711 / 4900) loss: 1.594951
(Iteration 3721 / 4900) loss: 1.250862
(Iteration 3731 / 4900) loss: 1.162776
(Iteration 3741 / 4900) loss: 1.340488
(Iteration 3751 / 4900) loss: 1.367186
(Iteration 3761 / 4900) loss: 1.366398
(Iteration 3771 / 4900) loss: 1.417237
(Iteration 3781 / 4900) loss: 1.344297
(Iteration 3791 / 4900) loss: 1.396264
(Iteration 3801 / 4900) loss: 1.238349
(Iteration 3811 / 4900) loss: 1.344179
(Iteration 3821 / 4900) loss: 1.219159
(Iteration 3831 / 4900) loss: 1.361244
(Iteration 3841 / 4900) loss: 1.443448
(Iteration 3851 / 4900) loss: 1.334387
(Iteration 3861 / 4900) loss: 1.360704
(Iteration 3871 / 4900) loss: 1.257395
(Iteration 3881 / 4900) loss: 1.345110
(Iteration 3891 / 4900) loss: 1.215053
(Iteration 3901 / 4900) loss: 1.245863
(Iteration 3911 / 4900) loss: 1.357334
(Epoch 8 / 10) train acc: 0.541000; val_acc: 0.463000
(Iteration 3921 / 4900) loss: 1.150966
```

```
(Iteration 3931 / 4900) loss: 1.319383
(Iteration 3941 / 4900) loss: 1.069093
(Iteration 3951 / 4900) loss: 1.462634
(Iteration 3961 / 4900) loss: 1.284100
(Iteration 3971 / 4900) loss: 1.159020
(Iteration 3981 / 4900) loss: 1.127782
(Iteration 3991 / 4900) loss: 1.110383
(Iteration 4001 / 4900) loss: 1.189466
(Iteration 4011 / 4900) loss: 1.271651
(Iteration 4021 / 4900) loss: 1.461067
(Iteration 4031 / 4900) loss: 1.365532
(Iteration 4041 / 4900) loss: 1.070091
(Iteration 4051 / 4900) loss: 1.431787
(Iteration 4061 / 4900) loss: 1.414741
(Iteration 4071 / 4900) loss: 1.361378
(Iteration 4081 / 4900) loss: 1.522182
(Iteration 4091 / 4900) loss: 1.287178
(Iteration 4101 / 4900) loss: 1.308431
(Iteration 4111 / 4900) loss: 1.589505
(Iteration 4121 / 4900) loss: 1.423470
(Iteration 4131 / 4900) loss: 1.724517
(Iteration 4141 / 4900) loss: 1.183873
(Iteration 4151 / 4900) loss: 1.029758
(Iteration 4161 / 4900) loss: 1.384384
(Iteration 4171 / 4900) loss: 1.469837
(Iteration 4181 / 4900) loss: 1.529209
(Iteration 4191 / 4900) loss: 1.190913
(Iteration 4201 / 4900) loss: 1.214756
(Iteration 4211 / 4900) loss: 1.133497
(Iteration 4221 / 4900) loss: 1.354617
(Iteration 4231 / 4900) loss: 1.273540
(Iteration 4241 / 4900) loss: 1.477614
(Iteration 4251 / 4900) loss: 1.584060
(Iteration 4261 / 4900) loss: 1.378007
(Iteration 4271 / 4900) loss: 1.200474
(Iteration 4281 / 4900) loss: 1.268054
(Iteration 4291 / 4900) loss: 1.154626
(Iteration 4301 / 4900) loss: 1.206561
(Iteration 4311 / 4900) loss: 1.067481
(Iteration 4321 / 4900) loss: 1.508419
(Iteration 4331 / 4900) loss: 1.348757
(Iteration 4341 / 4900) loss: 1.383121
(Iteration 4351 / 4900) loss: 1.300875
(Iteration 4361 / 4900) loss: 1.233478
(Iteration 4371 / 4900) loss: 1.221144
(Iteration 4381 / 4900) loss: 1.239890
(Iteration 4391 / 4900) loss: 1.323033
(Iteration 4401 / 4900) loss: 1.226766
```

```
(Epoch 9 / 10) train acc: 0.580000; val_acc: 0.469000
(Iteration 4411 / 4900) loss: 1.534061
(Iteration 4421 / 4900) loss: 1.022927
(Iteration 4431 / 4900) loss: 1.294994
(Iteration 4441 / 4900) loss: 1.360441
(Iteration 4451 / 4900) loss: 1.380112
(Iteration 4461 / 4900) loss: 1.143059
(Iteration 4471 / 4900) loss: 1.245423
(Iteration 4481 / 4900) loss: 1.408525
(Iteration 4491 / 4900) loss: 1.210972
(Iteration 4501 / 4900) loss: 1.394010
(Iteration 4511 / 4900) loss: 1.203786
(Iteration 4521 / 4900) loss: 1.060802
(Iteration 4531 / 4900) loss: 1.464206
(Iteration 4541 / 4900) loss: 1.260809
(Iteration 4551 / 4900) loss: 1.332313
(Iteration 4561 / 4900) loss: 1.226907
(Iteration 4571 / 4900) loss: 1.556329
(Iteration 4581 / 4900) loss: 1.326948
(Iteration 4591 / 4900) loss: 1.211093
(Iteration 4601 / 4900) loss: 1.254027
(Iteration 4611 / 4900) loss: 1.474277
(Iteration 4621 / 4900) loss: 1.288402
(Iteration 4631 / 4900) loss: 1.466134
(Iteration 4641 / 4900) loss: 1.088726
(Iteration 4651 / 4900) loss: 1.397102
(Iteration 4661 / 4900) loss: 1.441696
(Iteration 4671 / 4900) loss: 1.584400
(Iteration 4681 / 4900) loss: 1.289660
(Iteration 4691 / 4900) loss: 1.137633
(Iteration 4701 / 4900) loss: 1.267596
(Iteration 4711 / 4900) loss: 1.196905
(Iteration 4721 / 4900) loss: 1.246403
(Iteration 4731 / 4900) loss: 1.285876
(Iteration 4741 / 4900) loss: 1.326565
(Iteration 4751 / 4900) loss: 1.170252
(Iteration 4761 / 4900) loss: 1.246038
(Iteration 4771 / 4900) loss: 1.379887
(Iteration 4781 / 4900) loss: 1.176485
(Iteration 4791 / 4900) loss: 1.128133
(Iteration 4801 / 4900) loss: 1.317326
(Iteration 4811 / 4900) loss: 1.425641
(Iteration 4821 / 4900) loss: 1.359672
(Iteration 4831 / 4900) loss: 1.331216
(Iteration 4841 / 4900) loss: 1.210385
(Iteration 4851 / 4900) loss: 1.258650
(Iteration 4861 / 4900) loss: 1.214546
(Iteration 4871 / 4900) loss: 1.285254
```

```
(Iteration 4881 / 4900) loss: 1.383855
(Iteration 4891 / 4900) loss: 1.421827
(Epoch 10 / 10) train acc: 0.524000; val_acc: 0.456000
```

## 10 Debug the training

With the default parameters we provided above, you should get a validation accuracy of about 0.36 on the validation set. This isn't very good.

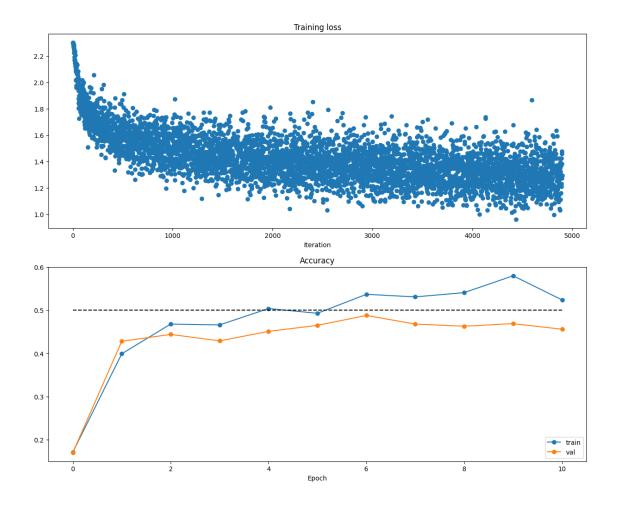
One strategy for getting insight into what's wrong is to plot the loss function and the accuracies on the training and validation sets during optimization.

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.

```
[12]: # Run this cell to visualize training loss and train / val accuracy

plt.subplot(2, 1, 1)
plt.title('Training loss')
plt.plot(solver.loss_history, 'o')
plt.xlabel('Iteration')

plt.subplot(2, 1, 2)
plt.title('Accuracy')
plt.plot(solver.train_acc_history, '-o', label='train')
plt.plot(solver.val_acc_history, '-o', label='val')
plt.plot([0.5] * len(solver.val_acc_history), 'k--')
plt.xlabel('Epoch')
plt.legend(loc='lower right')
plt.gcf().set_size_inches(15, 12)
plt.show()
```

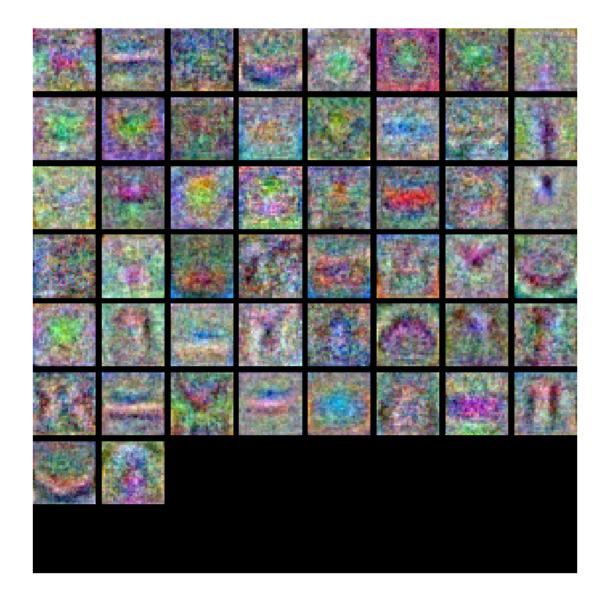


```
[13]: from cs231n.vis_utils import visualize_grid

# Visualize the weights of the network

def show_net_weights(net):
    W1 = net.params['W1']
    W1 = W1.reshape(3, 32, 32, -1).transpose(3, 1, 2, 0)
    plt.imshow(visualize_grid(W1, padding=3).astype('uint8'))
    plt.gca().axis('off')
    plt.show()

show_net_weights(model)
```



# 11 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 (52% could serve as a reference), 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.).

```
[15]: best_model = None
     # TODO: Tune hyperparameters using the validation set. Store your best trained \square
      →#
     # model in best_model.
                                                                          Ш
      →#
     #
                                                                          ш
      →#
     # 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 u
     # write code to sweep through possible combinations of hyperparameters
     # automatically like we did on thexs previous exercises.
                                                                          ш
     # *****START OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)****
     results = {}
     best_val = -1
     learning_rates = [2e-7, 0.75e-7, 1.5e-7, 1.25e-7, 0.75e-7]
     regularization_strengths = [3e4, 3.25e4, 3.5e4, 3.75e4, 4e4, 4.25e4, 4.5e4, 4.
      ⊶75e4, 5e4]
     learning_rates = np.geomspace(3e-4, 3e-2, 3)
     regularization_strengths = np.geomspace(1e-6, 1e-2, 5)
```

```
import itertools
for lr, reg in itertools.product(learning_rates, regularization_strengths):
    model = TwoLayerNet(hidden_dim=128, reg=reg)
    solver = Solver(model, data, optim_config={'learning_rate': lr},__
 →num_epochs=10, verbose=False)
    solver.train()
    results[(lr, reg)] = solver.best_val_acc
    if results[(lr, reg)] > best_val:
       best_val = results[(lr, reg)]
       best_model = model
for lr, reg in sorted(results):
    val_accuracy = results[(lr, reg)]
    print('lr %e reg %e val accuracy: %f' % (lr, reg, val_accuracy))
print('best validation accuracy achieved during cross-validation: %f' %11
 ⇒best_val)
# *****END OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)*****
END OF YOUR CODE
\CS231n Assignments\assignment1\cs231n\layers.py:821: RuntimeWarning:
g:\
divide by zero encountered in log
 loss = -np.log(P[range(N), y]).sum() / N # sum cross entropies as loss
    \CS231n Assignments\assignment1\cs231n\layers.py:817: RuntimeWarning:
overflow encountered in subtract
 P = np.exp(x - x.max(axis=1, keepdims=True))
    \CS231n Assignments\assignment1\cs231n\layers.py:32: RuntimeWarning:
overflow encountered in matmul
 out = x_reshaped @ w + b
    \CS231n Assignments\assignment1\cs231n\layers.py:817: RuntimeWarning:
invalid value encountered in subtract
 P = np.exp(x - x.max(axis=1, keepdims=True))
lr 3.000000e-04 reg 1.000000e-06 val accuracy: 0.519000
lr 3.000000e-04 reg 1.000000e-05 val accuracy: 0.511000
lr 3.000000e-04 reg 1.000000e-04 val accuracy: 0.528000
lr 3.000000e-04 reg 1.000000e-03 val accuracy: 0.521000
lr 3.000000e-04 reg 1.000000e-02 val accuracy: 0.533000
lr 3.000000e-03 reg 1.000000e-06 val accuracy: 0.217000
lr 3.000000e-03 reg 1.000000e-05 val accuracy: 0.375000
lr 3.000000e-03 reg 1.000000e-04 val accuracy: 0.189000
lr 3.000000e-03 reg 1.000000e-03 val accuracy: 0.231000
```

```
lr 3.000000e-03 reg 1.000000e-02 val accuracy: 0.384000
lr 3.000000e-02 reg 1.000000e-06 val accuracy: 0.132000
lr 3.000000e-02 reg 1.000000e-05 val accuracy: 0.172000
lr 3.000000e-02 reg 1.000000e-04 val accuracy: 0.113000
lr 3.000000e-02 reg 1.000000e-03 val accuracy: 0.145000
lr 3.000000e-02 reg 1.000000e-02 val accuracy: 0.118000
best validation accuracy achieved during cross-validation: 0.533000
```

## 12 Test your model!

Run your best model on the validation and test sets. You should achieve above 48% accuracy on the validation set and the test set.

```
[16]: y_val_pred = np.argmax(best_model.loss(data['X_val']), axis=1)
print('Validation set accuracy: ', (y_val_pred == data['y_val']).mean())
```

Validation set accuracy: 0.533

```
[17]: y_test_pred = np.argmax(best_model.loss(data['X_test']), axis=1)
print('Test set accuracy: ', (y_test_pred == data['y_test']).mean())
```

Test set accuracy: 0.52

## 12.1 Inline Question 2:

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.

#### YourAnswer:

1, 3.

### Your Explanation:

- 1. TRUE -
- 2. FALSE -
- 3. TRUE -

[]:

## features

June 8, 2024

```
[]: # This mounts your Google Drive to the Colab VM.
     from google.colab import drive
     drive.mount('/content/drive')
     # TODO: Enter the foldername in your Drive where you have saved the unzipped
     # assignment folder, e.g. 'cs231n/assignments/assignment1/'
     FOLDERNAME = None
     assert FOLDERNAME is not None, "[!] Enter the foldername."
     # Now that we've mounted your Drive, this ensures that
     # the Python interpreter of the Colab VM can load
     # python files from within it.
     import sys
     sys.path.append('/content/drive/My Drive/{}'.format(FOLDERNAME))
     # This downloads the CIFAR-10 dataset to your Drive
     # if it doesn't already exist.
     %cd /content/drive/My\ Drive/$FOLDERNAME/cs231n/datasets/
     !bash get datasets.sh
     %cd /content/drive/My\ Drive/$FOLDERNAME
```

# 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]: import random import numpy as np from cs231n.data_utils import load_CIFAR10 import matplotlib.pyplot as plt
```

#### 1.1 Load data

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

```
[2]: from cs231n.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 = 'cs231n/datasets/cifar-10-batches-py'
         # Cleaning up variables to prevent loading data multiple times (which may ...
      ⇒cause memory issue)
         trv:
            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)
         # 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
```

```
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 own interest.

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 cs231n.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
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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
Done extracting features for 49000 / 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
    import itertools
    from cs231n.classifiers.linear classifier import LinearSVM
    learning rates = np.geomspace(1e-4, 1e-2, 3)
    regularization_strengths = np.geomspace(0.01, 1, 3)
    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.
    # *****START OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)****
    for lr, reg in itertools.product(learning_rates, regularization_strengths):
        # Create SVM and train it
        svm = LinearSVM()
        svm.train(X_train_feats, y_train, lr, reg, num_iters=1500)
        # Compute training and validation sets accuracies and append to the
     \rightarrow dictionary
        y_train_pred, y_val_pred = svm.predict(
           X_train_feats), svm.predict(X_val_feats)
        results[(lr, reg)] = np.mean(
           y_train == y_train_pred), np.mean(y_val == y_val_pred)
        # Save if validation accuracy is the best
        if results[(lr, reg)][1] > best_val:
           best_val = results[(lr, reg)][1]
           best_svm = svm
    # *****END OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)****
```

```
lr 1.000000e-04 reg 1.000000e-02 train accuracy: 0.450429 val accuracy: 0.444000 lr 1.000000e-04 reg 1.000000e+00 train accuracy: 0.451020 val accuracy: 0.448000 lr 1.000000e-03 reg 1.000000e-02 train accuracy: 0.500694 val accuracy: 0.492000 lr 1.000000e-03 reg 1.000000e-01 train accuracy: 0.498408 val accuracy: 0.484000 lr 1.000000e-03 reg 1.000000e+00 train accuracy: 0.483796 val accuracy: 0.482000 lr 1.000000e-02 reg 1.000000e+00 train accuracy: 0.509122 val accuracy: 0.485000 lr 1.000000e-02 reg 1.000000e-01 train accuracy: 0.504531 val accuracy: 0.501000 lr 1.000000e-02 reg 1.000000e+00 train accuracy: 0.472959 val accuracy: 0.467000 best validation accuracy achieved: 0.501000
```

```
[5]: # Evaluate your trained SVM on the test set: you should be able to get at least
\( \times 0.40 \)

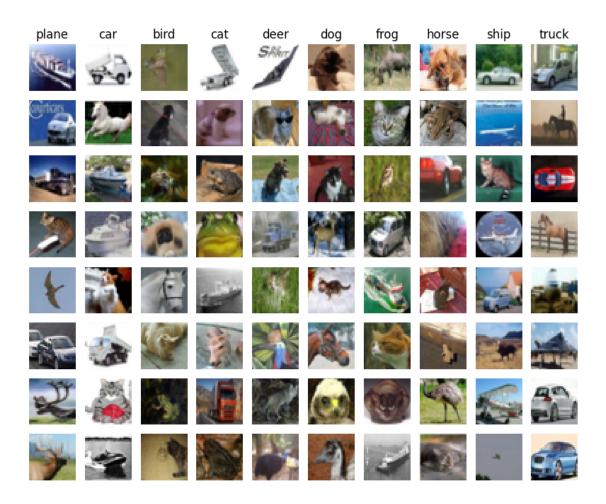
y_test_pred = best_svm.predict(X_test_feats)

test_accuracy = np.mean(y_test == y_test_pred)

print(test_accuracy)
```

#### 0.489

```
[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?

Your Answer:

## 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)
[8]: import itertools
    from cs231n.classifiers.fc_net import TwoLayerNet
    from cs231n.solver import Solver
    input_dim = X_train_feats.shape[1]
    hidden dim = 500
    num classes = 10
    data = {
       'X_train': X_train_feats,
       'y_train': y_train,
        'X_val': X_val_feats,
       'y_val': y_val,
       'X_test': X_test_feats,
       'y_test': y_test,
    }
    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.
    # *****START OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)****
    results = {}
    best_val = -1
    learning_rates = np.linspace(1e-2, 2.75e-2, 4)
    regularization_strengths = np.geomspace(1e-6, 1e-4, 3)
    data = {
       'X_train': X_train_feats,
       'X_val': X_val_feats,
        'X_test': X_test_feats,
```

```
'y_train': y_train,
         'y_val': y_val,
         'y_test': y_test
     }
     for lr, reg in itertools.product(learning_rates, regularization_strengths):
         # Create Two Layer Net and train it with Solver
         model = net
         solver = Solver(model, data, optim_config={
                         'learning rate': lr}, num epochs=15, verbose=False)
         solver.train()
         # Compute validation set accuracy and append to the dictionary
         results[(lr, reg)] = solver.best_val_acc
         # Save if validation accuracy is the best
         if results[(lr, reg)] > best_val:
             best_val = results[(lr, reg)]
             best_net = model
     # Print out results.
     for lr, reg in sorted(results):
         val accuracy = results[(lr, reg)]
         print('lr %e reg %e val accuracy: %f' % (lr, reg, val_accuracy))
     print('best validation accuracy achieved during cross-validation: %f' %u
      ⇒best_val)
     # *****END OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)****
    lr 1.000000e-02 reg 1.000000e-06 val accuracy: 0.522000
    lr 1.000000e-02 reg 1.000000e-05 val accuracy: 0.543000
    lr 1.000000e-02 reg 1.000000e-04 val accuracy: 0.574000
    lr 1.583333e-02 reg 1.000000e-06 val accuracy: 0.598000
    lr 1.583333e-02 reg 1.000000e-05 val accuracy: 0.610000
    lr 1.583333e-02 reg 1.000000e-04 val accuracy: 0.617000
    lr 2.166667e-02 reg 1.000000e-06 val accuracy: 0.614000
    lr 2.166667e-02 reg 1.000000e-05 val accuracy: 0.617000
    lr 2.166667e-02 reg 1.000000e-04 val accuracy: 0.615000
    lr 2.750000e-02 reg 1.000000e-06 val accuracy: 0.612000
    lr 2.750000e-02 reg 1.000000e-05 val accuracy: 0.613000
    lr 2.750000e-02 reg 1.000000e-04 val accuracy: 0.612000
    best validation accuracy achieved during cross-validation: 0.617000
[9]: # Run your best neural net classifier on the test set. You should be able
     # to get more than 55% accuracy.
```

```
y_test_pred = np.argmax(best_net.loss(data['X_test']), axis=1)
test_acc = (y_test_pred == data['y_test']).mean()
print(test_acc)
```

0.593