# **ECE180 HW2**

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# k-Nearest Neighbor (kNN) exercise

\*Complete and hand in this completed worksheet (including its outputs) with your assignment submission.

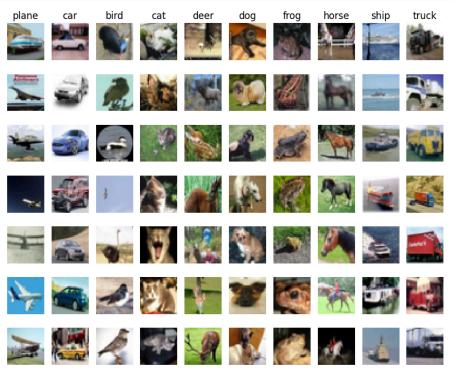
The kNN classifier consists of two stages:

- · 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.

```
# Run some setup code for this notebook.
from builtins import range
from builtins import object
import random
import numpy as np
import matplotlib.pyplot as plt
from __future__ import print_function
# 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'
from keras.datasets import cifar10
# load dataset
(X_train, y_train), (X_test, y_test) = cifar10.load_data()
y_train = y_train.squeeze()
y_test = y_test.squeeze()
# 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,)
```

```
# 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()
```



```
# Subsample the data for more efficient code execution in this exercise
# Take 5,000 samples from training set and 500 samples from test set.
num\_training = 5000
num\_test = 500
#
# Randomly select 5000 and 500 samples from train and test set.
                                                          #
\# Overwrite them to X_train, y_train, X_test, y_test.
                                                          #
# Use the above visualization to double-check images and labels.
# *****START OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE) ****
indices_train = np.random.choice(X_train.shape[0], num_training, replace=False)
X_train = X_train[indices_train]
y_train = y_train[indices_train]
indices_test = np.random.choice(X_test.shape[0], num_test, replace=False)
X_test = X_test[indices_test]
y_test = y_test[indices_test]
# *****END OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)*****
```

```
# Reshape the image data into rows
# Output should be (5000, 3072) (500, 3072)
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)

```
class KNearestNeighbor(object):
   """ a kNN classifier """
   def __init__(self):
       pass
   def train(self, X, y):
       Train the classifier. For k-nearest neighbors this is just
       memorizing the training data.
       Inputs:
       - X: A numpy array of shape (num_train, D) containing the training data
        consisting of num_train samples each of dimension D.
       - y: A numpy array of shape (N,) containing the training labels, where
            y[i] is the label for X[i].
       self.X_train = X
       self.y_train = y
   def predict(self, X, k=1):
       Predict labels for test data using this classifier.
       Inputs:
       - X: A numpy array of shape (num_test, D) containing test data consisting
            of num_test samples each of dimension D.
       - k: The number of nearest neighbors that vote for the predicted labels.
       Returns:
       - y: A numpy array of shape (num_test,) containing predicted labels for the
         test data, where y[i] is the predicted label for the test point X[i].
       dists = self.compute_distances(X)
       return self.predict_labels(dists, k=k)
   def compute_distances(self, X):
       Compute the distance between each test point in X and each training point
       in self.X_train
       X: A numpy array of shape (num_test, D) containing test data.
       Returns:
       - dists: A numpy array of shape (num_test, num_train) where dists[i, j]
         is the distance between the ith test point and the jth training
         point.
       num_test = X.shape[0]
       num_train = self.X_train.shape[0]
       dists = np.zeros((num_test, num_train))
       # TOD0:
       # Compute the 12 distance between the ith test point and the jth
       # training point, and store the result in dists[i, j]. You should
       # not use a loop over dimension, nor use np.linalg.norm().
       # Try to use numpy function to compute dists by a single function
       # call, comparing a matrix and a matrix. (deduction 0)
       # You can iterate for test samples to compute distances between a
       # vector (a test sample) and a matrix (train samples). (deduction 2)#
       # Or you can have two iterations (two for) to compute distances
       # between a vector (a test sample) and a vector (a train sample),
       # but that will be slow and your cross-validation will take a long #
       # time.
       # *****START OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)****
       # for i in range(num_test):
       # dists[i] = np.sqrt(np.sum(np.square(X[i] - self.X_train), axis = 1))
       test_square = np.sum(X**2, axis=1, keepdims=True)
       train_square = np.sum(self.X_train**2, axis=1, keepdims=True).T
       dists = np.sqrt(test_square + train_square - 2 * np.dot(X, self.X_train.T))
       # *****END OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)****
```

```
def predict_labels(self, dists, k=1):
       Given a matrix of distances between test points and training points,
       predict a label for each test point.
       - dists: A numpy array of shape (num_test, num_train) where dists[i, j]
        gives the distance betwen the ith test point and the jth training point.
       - y: A numpy array of shape (num_test,) containing predicted labels for the
        test data, where y[i] is the predicted label for the test point X[i].
       num_test = dists.shape[0]
       y_pred = np.zeros(num_test, dtype=int) # Ensure dtype is int for indexing
       for i in range(num_test):
          # A list of length k storing the labels of the k nearest neighbors to
          # the ith test point.
          # TOD0:
          # Use the distance matrix to find the k nearest neighbors of the ith
          # testing point, and use self.y_train to find the labels of these
                                                                           #
          # neighbors. Store these labels in closest_y.
          # Hint: Look up the function numpy.argsort.
          # *****START OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)****
          closest_y_indices = np.argsort(dists[i])[:k]
          closest_y = self.y_train[closest_y_indices]
          # *****END OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)*****
          # TODO:
          # Now that you have found the labels of the k nearest neighbors, you
                                                                           #
          # need to find the most common label in the list closest_y of labels.
          # Store this label in y_pred[i]. Break ties by choosing the smaller
          # label.
          # *****START OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)****
          # Voting mechanism: Find the most frequent label
          unique, counts = np.unique(closest_y, return_counts=True)
          y_pred[i] = unique[np.argmax(counts)]
          # *****END OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)****
       return y_pred
# 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)
# Now implement the function predict_labels and run the code below:
# We use k = 1 (which is Nearest Neighbor).
dists = classifier.compute_distances(X_test)
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 52 / 500 correct => accuracy: 0.104000
```

You should expect to see approximately 20% accuracy. Now lets try out a larger k, say k = 5:

```
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 52 / 500 correct => accuracy: 0.104000

#### 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.

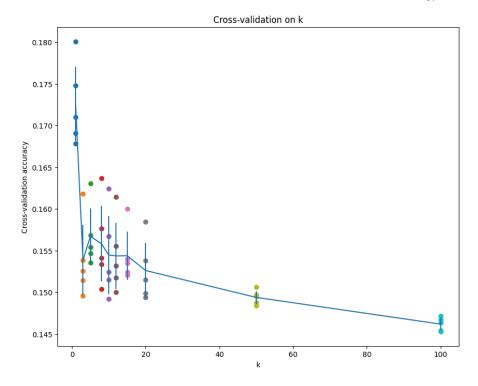
```
num_folds = 5
k_{c} hoices = [1, 3, 5, 8, 10, 12, 15, 20, 50, 100] #You can reduce the k's if the running time is too long.
\# k_choices = [1, 3, 5, 8, 10, 12]
X_train_folds = []
y_train_folds = []
# Split up the training data into folds. After splitting, X_train_folds and
                                                                      #
# y_train_folds should each be lists of length num_folds, where
# y_train_folds[i] is the label vector for the points in X_train_folds[i].
                                                                      #
# Hint: Look up the numpy array_split function.
# *****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 curr_k in k_choices:
   k_to_accuracies[curr_k] = []
   for curr_fold in range(num_folds):
       classifier = KNearestNeighbor()
       classifier.train(np.concatenate(X_train_folds[:curr_fold] + X_train_folds[curr_fold + 1:]), np.concatenate(y_train_fold
       y_pred_fold = classifier.predict(X_train_folds[curr_fold], k = curr_k)
       k_to_accuracies[curr_k].append(np.mean(y_pred_fold == y_train_folds[curr_fold]))
# ****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 = 0.167857
    k = 1, accuracy = 0.169082
    k = 1, accuracy = 0.174796
    k = 1, accuracy = 0.180102
    k = 1, accuracy = 0.171020
    k = 3, accuracy = 0.151429
    k = 3, accuracy = 0.149592
    k = 3, accuracy = 0.153878
    k = 3, accuracy = 0.161837
    k = 3, accuracy = 0.152551
    k = 5, accuracy = 0.154694
    k = 5, accuracy = 0.153571
```

k = 5, accuracy = 0.155408

```
k = 5, accuracy = 0.163061
k = 5, accuracy = 0.156837
k = 8, accuracy = 0.150408
k = 8, accuracy = 0.154082
k = 8, accuracy = 0.157653
k = 8, accuracy = 0.163673
k = 8, accuracy = 0.153367
k = 10, accuracy = 0.151531
k = 10, accuracy = 0.152449
k = 10, accuracy = 0.156735
k = 10, accuracy = 0.162449
k = 10, accuracy = 0.149184
k = 12, accuracy = 0.153163
k = 12, accuracy = 0.151735
k = 12, accuracy = 0.155510
k = 12, accuracy = 0.161429
k = 12, accuracy = 0.150000
k = 15, accuracy = 0.152041
k = 15, accuracy = 0.153469
k = 15, accuracy = 0.153980
k = 15, accuracy = 0.160000
k = 15, accuracy = 0.152449
k = 20, accuracy = 0.149898
k = 20, accuracy = 0.153776
k = 20, accuracy = 0.151531
k = 20, accuracy = 0.158469
k = 20, accuracy = 0.149388
k = 50, accuracy = 0.149490
k = 50, accuracy = 0.149694
k = 50, accuracy = 0.148776
k = 50, accuracy = 0.150612
k = 50, accuracy = 0.148367
k = 100, accuracy = 0.145306
k = 100, accuracy = 0.146735
k = 100, accuracy = 0.146327
k = 100, accuracy = 0.147143
k = 100, accuracy = 0.145408
```

```
# 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()
```



```
# 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 184 / 500 correct => accuracy: 0.368000

# Softmax exercise

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

```
def get_CIFAR10_data(num_training=49000, num_validation=1000, num_test=1000, num_dev=500):
    Load the CIFAR-10 dataset from disk and perform preprocessing to prepare
    it for the linear classifier. These are the same steps as we used for the
    SVM, but condensed to a single function.
    # Load the raw CIFAR-10 data
    (X_train, y_train), (X_test, y_test) = cifar10.load_data()
    y_train = y_train.squeeze()
    y_test = y_test.squeeze()
    # 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_{\text{test}} = y_{\text{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\_original = np.mean(X\_train, axis = 0)
    mean_image = mean_image_original.astype(np.uint8)
    X_train -= mean_image
    X_val -= mean_image
    X_test -= mean_image
    X_dev -= mean_image
    # add bias dimension and transform into columns
    X_train = np.hstack([X_train, np.ones((X_train.shape[0], 1))])
    X_val = np.hstack([X_val, np.ones((X_val.shape[0], 1))])
    X_test = np.hstack([X_test, np.ones((X_test.shape[0], 1))])
    X_dev = np.hstack([X_dev, np.ones((X_dev.shape[0], 1))])
    return X_train, y_train, X_val, y_val, X_test, y_test, X_dev, y_dev
# Cleaning up variables to prevent loading data multiple times (which may cause memory issue)
try:
   del X_train, y_train
   del X_test, y_test
   print('Clear previously loaded data.')
except:
   pass
# Invoke the above function to get our data.
X_train, y_train, X_val, y_val, X_test, y_test, X_dev, y_dev = get_CIFAR10_data()
print('Train data shape: ', X_train.shape)
print('Train labels shape: ', y_train.shape)
print('Validation data shape: ', X_val.shape)
print('Validation labels shape: ', y_val.shape)
print('Test data shape: ', X_test.shape)
print('Test labels shape: ', y_test.shape)
print('dev data shape: ', X_dev.shape)
print('dev labels shape: ', y_dev.shape)
     Clear previously loaded data.
```

```
Clear previously loaded data.
Train data shape: (49000, 3073)
Train labels shape: (49000,)
Validation data shape: (1000, 3073)
Validation labels shape: (1000,)
Test data shape: (1000, 3073)
Test labels shape: (1000,)
dev data shape: (500, 3073)
dev labels shape: (500,)
```

## Softmax Classifier

```
def softmax_loss(W, X, y, reg):
   Softmax loss function, vectorized version.
   Inputs and outputs are the same as softmax_loss_naive.
   # Initialize the loss and gradient to zero.
   loss = 0.0
   dW = np.zeros_like(W)
   # TODO: Compute the softmax loss and its gradient using no explicit loops. #
   # Store the loss in loss and the gradient in dW. If you are not careful
   # here, it is easy to run into numeric instability. Don't forget the
                                                                      #
   # regularization!
   # *****START OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE) ****
   num_classes = W.shape[1]
   num_train = X.shape[0]
   scores = X.dot(W)
   scores -= scores.max(axis = 1, keepdims = True)
   probs = np.exp(scores)/np.sum(np.exp(scores), axis = 1, keepdims = True)
   loss = -np.log(probs[np.arange(num_train), y])
   loss = np.sum(loss)
   dscores = probs.reshape(num_train, -1)
   dscores[np.arange(num\_train), y] = 1
   dW = np.dot(X.T.reshape(X.shape[1], num_train), dscores)
   loss /= num_train
   dW /= num_train
   loss += reg * np.sum(W * W)
   dW += 2 * reg * W
   # *****END OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)*****
   return loss, dW
```

```
class Softmax(object):
    def __init__(self):
        self.W = None
    def train(
        self,
        Χ,
        у,
        learning_rate=1e-3,
        reg=1e-5,
        num_iters=100,
        batch_size=200,
        verbose=False,
    ):
        Train this linear classifier using stochastic gradient descent.
        Inputs:
        - X: A numpy array of shape (N, D) containing training data; there are N
         training samples each of dimension D.
        - y: A numpy array of shape (N,) containing training labels; y[i] = c
          means that X[i] has label 0 \ll c \ll C for C classes.
        - learning_rate: (float) learning rate for optimization.
        - reg: (float) regularization strength.
        - num_iters: (integer) number of steps to take when optimizing
        - batch_size: (integer) number of training examples to use at each step.
        - verbose: (boolean) If true, print progress during optimization.
        A list containing the value of the loss function at each training iteration.
        num_train, dim = X.shape
        num_classes = (
           np.max(y) + 1
        ) # assume y takes values 0...K-1 where K is number of classes
        if self.W is None:
            # lazily initialize W
            self.W = 0.001 * np.random.randn(dim, num_classes)
        # Run stochastic gradient descent to optimize W
        loss_history = []
        for it in range(num_iters):
            X_batch = None
            y_batch = None
            indices = np.random.choice(num_train, batch_size)
            X_batch = X[indices]
            y_batch = y[indices]
            # evaluate loss and gradient
            loss, grad = self.loss(X_batch, y_batch, reg)
            loss_history.append(loss)
            self.W -= learning_rate * grad
            if verbose and it % 100 == 0:
                print("iteration %d / %d: loss %f" % (it, num_iters, loss))
        return loss_history
    def predict(self, X):
        Use the trained weights of this linear classifier to predict labels for
        data points.
        Inputs:
        - X: A numpy array of shape (N, D) containing training data; there are N
          training samples each of dimension \ensuremath{\text{D.}}
        Returns:
        - y_pred: Predicted labels for the data in X. y_pred is a 1-dimensional
          array of length N, and each element is an integer giving the predicted
          class.
        # *****START OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)****
        y_pred = np.argmax(X.dot(self.W), axis = 1)
```

```
# *****END OF YOUR CODE (DO NOT DELETE/MODIFY THIS LINE)*****
return y_pred

def loss(self, X_batch, y_batch, reg):
    """
    Compute the loss function and its derivative.
    Subclasses will override this.

Inputs:
    - X_batch: A numpy array of shape (N, D) containing a minibatch of N
    data points; each point has dimension D.
    - y_batch: A numpy array of shape (N,) containing labels for the minibatch.
    - reg: (float) regularization strength.

Returns: A tuple containing:
    - loss as a single float
    - gradient with respect to self.W; an array of the same shape as W
    """
    return softmax_loss(self.W, X_batch, y_batch, reg)
```

## Stochastic Gradient Descent

```
We now have vectorized and efficient expressions for the loss, the gradient and our gradient matches the numerical gradient. We are therefore
ready to do SGD to minimize the loss.
# In the file linear_classifier.py, implement SGD in the function
softmax = Softmax()
loss_hist = softmax.train(X_train, y_train, learning_rate=1e-7, reg=2.5e4,
                       num_iters=1500, verbose=True)
     iteration 0 / 1500: loss 786.942138
     iteration 100 / 1500: loss 286.234105 iteration 200 / 1500: loss 105.929184
     iteration 300 / 1500: loss 39.916208
     iteration 400 / 1500: loss 15.924812
     iteration 500 / 1500: loss 7.193934
     iteration 600 / 1500: loss 4.006732
     iteration 700 / 1500: loss 2.870526
     iteration 800 / 1500: loss 2.436226
     iteration 900 / 1500: loss 2.241114
     iteration 1000 / 1500: loss 2.256278
     iteration 1100 / 1500: loss 2.194643
     iteration 1200 / 1500: loss 2.153940 iteration 1300 / 1500: loss 2.191554
     iteration 1400 / 1500: loss 2.224378
# Write the Softmax.predict function and evaluate the performance on both the
# training and validation set
y_train_pred = softmax.predict(X_train)
print('training accuracy: %f' % (np.mean(y_train == y_train_pred), ))
y_val_pred = softmax.predict(X_val)
print('validation accuracy: %f' % (np.mean(y_val == y_val_pred), ))
     training accuracy: 0.256612
     validation accuracy: 0.266000
# 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()
```

