Nearest neighbor for handwritten digit recognition In this notebook we will build a classifier that takes an image of a handwritten digit and outputs a label 0-9. We will look at a particularly simple strategy for this problem known as the **nearest neighbor classifier**. To run this notebook you should have the following Python packages installed: numpy matplotlib sklearn 1. The MNIST dataset MNIST is a classic dataset in machine learning, consisting of 28x28 gray-scale images handwritten digits. The original training set contains 60,000 examples and the test set contains 10,000 examples. In this notebook we will be working with a subset of this data: a training set of 7,500 examples and a test set of 1,000 examples. %matplotlib inline import numpy as np import matplotlib.pyplot as plt import time ## Load the training set train data = np.load('MNIST/train data.npy') train labels = np.load('MNIST/train labels.npy') ## Load the testing set test data = np.load('MNIST/test data.npy') test labels = np.load('MNIST/test labels.npy') ## Print out their dimensions print("Training dataset dimensions: ", np.shape(train\_data)) print("Number of training labels: ", len(train labels)) print("Testing dataset dimensions: ", np.shape(test\_data)) print("Number of testing labels: ", len(test\_labels)) Training dataset dimensions: (7500, 784) Number of training labels: 7500 Testing dataset dimensions: (1000, 784) Number of testing labels: 1000 ## Compute the number of examples of each digit train\_digits, train\_counts = np.unique(train\_labels, return\_counts=True) print("Training set distribution:") print(dict(zip(train\_digits, train\_counts))) test digits, test\_counts = np.unique(test\_labels, return\_counts=True) print("Test set distribution:") print(dict(zip(test\_digits, test\_counts))) Training set distribution: {0: 750, 1: 750, 2: 750, 3: 750, 4: 750, 5: 750, 6: 750, 7: 750, 8: 750, 9: 750} Test set distribution: {0: 100, 1: 100, 2: 100, 3: 100, 4: 100, 5: 100, 6: 100, 7: 100, 8: 100, 9: 100} 2. Visualizing the data Each data point is stored as 784-dimensional vector. To visualize a data point, we first reshape it to a 28x28 image. In [4]: ## Define a function that displays a digit given its vector representation def show\_digit(x): plt.axis('off') plt.imshow(x.reshape((28,28)), cmap=plt.cm.gray) plt.show() return ## Define a function that takes an index into a particular data set ("train" or "test") and displays that im def vis\_image(index, dataset="train"): if (dataset=="train"): show digit(train data[index,]) label = train\_labels[index] else: show\_digit(test\_data[index,]) label = test\_labels[index] print("Label " + str(label)) return ## View the first data point in the training set vis\_image(0, "train") ## Now view the first data point in the test set vis\_image(0, "test") Label 9 Label 0 3. Squared Euclidean distance To compute nearest neighbors in our data set, we need to first be able to compute distances between data points. A natural distance function is *Euclidean distance*: for two vectors  $x,y\in\mathbb{R}^d$  , their Euclidean distance is defined as  $\|x-y\|=\sqrt{\sum_{i=1}^d(x_i-y_i)^2}.$ Often we omit the square root, and simply compute squared Euclidean distance:  $\|x-y\|^2 = \sum_{i=1}^d (x_i-y_i)^2.$ For the purposes of nearest neighbor computations, the two are equivalent: for three vectors  $x,y,z\in\mathbb{R}^d$  , we have  $||x - y|| \le ||x - z||$  if and only if  $||x - y||^2 \le ||x - z||^2$ . Now we just need to be able to compute squared Euclidean distance. The following function does so. In [5]: ## Computes squared Euclidean distance between two vectors. **def** squared dist(x,y): return np.sum(np.square(x-y)) ## Compute distance between a seven and a one in our training set. print("Distance from 7 to 1: ", squared dist(train data[4,],train data[5,])) ## Compute distance between a seven and a two in our training set. print("Distance from 7 to 2: ", squared dist(train data[4,],train data[1,])) ## Compute distance between two seven's in our training set. print("Distance from 7 to 7: ", squared dist(train data[4,],train data[7,])) Distance from 7 to 1: 5357193.0 Distance from 7 to 2: 12451684.0 Distance from 7 to 7: 5223403.0 4. Computing nearest neighbors Now that we have a distance function defined, we can now turn to nearest neighbor classification. ## Takes a vector x and returns the index of its nearest neighbor in train data **def** find NN(x): # Compute distances from x to every row in train data distances = [squared dist(x,train data[i,]) for i in range(len(train labels))] # Get the index of the smallest distance return np.argmin(distances) ## Takes a vector x and returns the class of its nearest neighbor in train data def NN classifier(x): # Get the index of the the nearest neighbor index = find NN(x)# Return its class return train labels[index] ## A success case: print("A success case:") print("NN classification: ", NN classifier(test data[0,])) print("True label: ", test\_labels[0]) print("The test image:") vis image(0, "test") print("The corresponding nearest neighbor image:") vis image(find NN(test data[0,]), "train") A success case: NN classification: 0 True label: 0 The test image: Label 0 The corresponding nearest neighbor image: Label 0 test\_data.shape Out[8]: (1000, 784) ## A failure case: print("A failure case:") print("NN classification: ", NN\_classifier(test\_data[39,])) print("True label: ", test\_labels[39]) print("The test image:") vis\_image(39, "test") print("The corresponding nearest neighbor image:") vis\_image(find\_NN(test\_data[39,]), "train") A failure case: NN classification: 2 True label: 3 The test image: The corresponding nearest neighbor image: Label 2 5. Processing the full test set Now let's apply our nearest neighbor classifier over the full data set. Note that to classify each test point, our code takes a full pass over each of the 7500 training examples. Thus we should not expect testing to be very fast. The following code takes about 100-150 seconds on 2.6 GHz Intel Core i5. ## Predict on each test data point (and time it!) t before = time.time()  $\texttt{test\_predictions} = \texttt{[NN\_classifier(test\_data[i,])} \ \ \textbf{for} \ \ i \ \ \textbf{in} \ \ \texttt{range(len(test\_labels))]}$ t\_after = time.time() ## Compute the error err\_positions = np.not\_equal(test\_predictions, test\_labels) error = float(np.sum(err\_positions))/len(test\_labels) print("Error of nearest neighbor classifier: ", error) print("Classification time (seconds): ", t after - t before) Error of nearest neighbor classifier: 0.046 Classification time (seconds): 49.61181592941284 6. Faster nearest neighbor methods Performing nearest neighbor classification in the way we have presented requires a full pass through the training set in order to classify a single point. If there are N training points in  $\mathbb{R}^d$ , this takes O(Nd) time. Fortunately, there are faster methods to perform nearest neighbor look up if we are willing to spend some time preprocessing the training set. scikit-learn has fast implementations of two useful nearest neighbor data structures: the ball tree and the k-d tree. from sklearn.neighbors import BallTree ## Build nearest neighbor structure on training data t before = time.time() ball tree = BallTree(train data) t after = time.time() ## Compute training time t training = t after - t before print("Time to build data structure (seconds): ", t training) ## Get nearest neighbor predictions on testing data t before = time.time() test neighbors = np.squeeze(ball tree.query(test data, k=1, return distance=False)) ball tree predictions = train labels[test neighbors] t after = time.time() ## Compute testing time t testing = t after - t before print("Time to classify test set (seconds): ", t testing) ## Verify that the predictions are the same print("Ball tree produces same predictions as above? ", np.array equal(test predictions, ball tree prediction Time to build data structure (seconds): 0.40031003952026367 Time to classify test set (seconds): 7.583951950073242 Ball tree produces same predictions as above? True from sklearn.neighbors import KDTree ## Build nearest neighbor structure on training data t before = time.time() kd tree = KDTree(train data) t after = time.time() ## Compute training time t training = t after - t before print("Time to build data structure (seconds): ", t training) ## Get nearest neighbor predictions on testing data t before = time.time() test neighbors = np.squeeze(kd tree.query(test data, k=1, return distance=False)) kd tree predictions = train labels[test neighbors] t after = time.time() ## Compute testing time t testing = t after - t before print("Time to classify test set (seconds): ", t testing) ## Verify that the predictions are the same print("KD tree produces same predictions as above? ", np.array\_equal(test\_predictions, kd\_tree\_predictions)) Time to build data structure (seconds): 0.3857090473175049 Time to classify test set (seconds): 8.6266348361969 KD tree produces same predictions as above? True Worksheet 1 Question 7 (a) For test point 100, print its image as well as the image of its nearest neighbor in the training set. Put these images in your writeup. Is this test point classified correctly? ## Now view the 100th data point in the test set vis image(99, "test") print("The corresponding nearest neighbor image:") vis image(find NN(test data[99,]), "train") The corresponding nearest neighbor image: Conclusion: Yes, this test point classified correctly. (b) The confusion matrix for the classifier is a 10  $\times$  10 matrix Nij with 0  $\leq$  i, j  $\leq$  9, where Nij is the number of test points whose true label is i but which are classified as j. Thus, if all test points are correctly classified, the off-diagonal entries of the matrix will be zero. Compute the matrix N for the 1-NN classifier and print it out. Which digit is misclassified most often? Least often? In [14]: from sklearn.metrics import confusion matrix cm = confusion\_matrix(test\_predictions, test\_labels) print('Confusion matrix of the classifier:\n', cm) Confusion matrix of the classifier:  $[[ 99 \quad 0 \quad 0 \quad 0 \quad 0 \quad 1 \quad 0 \quad 0 \quad 2 \quad 1]$  $[ \quad 0 \quad 100 \quad 1 \quad 0 \quad 0 \quad 0 \quad 0 \quad 4 \quad 0 \quad 1]$  $[ \quad 0 \quad 0 \quad 94 \quad 2 \quad 0 \quad 0 \quad 0 \quad 0 \quad 1 \quad 1]$ [ 0 0 1 91 0 0 0 1 1][ 0 0 0 2 97 0 0 1 1 [ 1 0 0 4 0 98 1 0 0 [ 0 0 0 0 0 0 0 99 0 1 [ 0 0 3 0 0 0 0 94 1 1] 3]  $[ 0 \quad 0 \quad 0 \quad 0 \quad 3 \quad 1 \quad 0 \quad 1 \quad 1 \quad 90]]$ x = cm[0][0]**for** i **in** range(0, 10): **for** j **in** range (0, 10): **if** i == j: **if** cm[i][j] <= x: x = cm[i][j]most\_often = np.where(cm == x) most often print('Digits that misclassified most often:\n', most often[0]) least often = np.where(cm == cm.max()) least often[0] print('\nDigits that misclassified least often:\n', least often[0]) Digits that misclassified most often: [9] Digits that misclassified least often: (c) For each digit  $0 \le i \le 9$ : look at all training instances of image i, and compute their mean. This average is a 784-dimensional vector. Use the show digit routine to print out these 10 average-digits. zero = np.mean(train\_data[train\_labels == 0,], axis = 0) one = np.mean(train\_data[train\_labels == 1,], axis = 0) two = np.mean(train\_data[train\_labels == 2,], axis = 0) three = np.mean(train\_data[train\_labels == 3,], axis = 0) four = np.mean(train\_data[train\_labels == 4,], axis = 0) five = np.mean(train\_data[train\_labels == 5,], axis = 0) six = np.mean(train\_data[train\_labels == 6,], axis = 0) seven = np.mean(train\_data[train\_labels == 7,], axis = 0) eight = np.mean(train\_data[train\_labels == 8,], axis = 0) nine = np.mean(train\_data[train\_labels == 9,], axis = 0) show\_digit(zero) show\_digit(one) show\_digit(two) show\_digit(three) show\_digit(four) show\_digit(five) show\_digit(six) show\_digit(seven) show\_digit(eight) show\_digit(nine)

