```
In [1]: # 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 unzip
        # assignment folder, e.g. 'cs6353/assignments/assignment1/'
        FOLDERNAME = 'MS/CS6353/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/$FOLDERNAME/cs6353/datasets/
        !bash get datasets.sh
        %cd /content/drive/My\ Drive/$FOLDERNAME
        # Install requirements from colab requirements.txt
        # TODO: Please change your path below to the colab requirements.txt file
        #! python -m pip install -r /content/drive/My\ Drive/$FOLDERNAME/colab re
        Drive already mounted at /content/drive; to attempt to forcibly remount,
        call drive.mount("/content/drive", force remount=True).
        /content/drive/My Drive/MS/CS6353/assignment1/cs6353/datasets
        --2024-09-08 05:09:59-- http://www.cs.toronto.edu/~kriz/cifar-10-python.
        tar.qz
        Resolving www.cs.toronto.edu (www.cs.toronto.edu)... 128.100.3.30
        Connecting to www.cs.toronto.edu (www.cs.toronto.edu) | 128.100.3.30 | :80...
        connected.
        HTTP request sent, awaiting response... 200 OK
        Length: 170498071 (163M) [application/x-gzip]
        Saving to: 'cifar-10-python.tar.gz'
        cifar-10-python.tar 100%[==========] 162.60M 31.2MB/s
                                                                           in 5.
        1 s
        2024-09-08 05:10:04 (31.7 MB/s) - 'cifar-10-python.tar.gz' saved [1704980
        71/1704980711
        cifar-10-batches-py/
        cifar-10-batches-py/data batch 4
        cifar-10-batches-py/readme.html
        cifar-10-batches-py/test batch
        cifar-10-batches-py/data batch 3
        cifar-10-batches-py/batches.meta
        cifar-10-batches-py/data batch 2
        cifar-10-batches-py/data batch 5
        cifar-10-batches-py/data batch 1
        /content/drive/My Drive/MS/CS6353/assignment1
```

# 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 transferring the labels of the k most similar training examples
- · The value of k is cross-validated

In this exercise you will implement these steps and understand the basic Image Classification pipeline, cross-validation, and gain proficiency in writing efficient, vectorized code.

```
In [2]: # Run some setup code for this notebook.
        from future import print function
        import random
        import numpy as np
        from cs6353.data utils import load CIFAR10
        import matplotlib.pyplot as plt
        import copy
        # This is a bit of magic to make matplotlib figures appear inline in the
        # 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 module
        # see http://stackoverflow.com/questions/1907993/autoreload-of-modules-in
        %load ext autoreload
        %autoreload 2
In [3]: # Load the raw CIFAR-10 data.
        cifar10 dir = 'cs6353/datasets/cifar-10-batches-py'
        # Cleaning up variables to prevent loading data multiple times (which may
        try:
           del X_train, y_train
           del X test, y test
           print('Clear previously loaded data.')
        except:
           pass
        X train, y train, X test, y test = load CIFAR10(cifar10 dir)
        # As a sanity check, we print out the size of the training and test data.
        print('Training data shape: ', X_train.shape)
        print('Training labels shape: ', y train.shape)
        print('Test data shape: ', X test.shape)
        print('Test labels shape: ', y test.shape)
        Training data shape: (50000, 32, 32, 3)
        Training labels shape: (50000,)
        Test data shape: (10000, 32, 32, 3)
        Test labels shape: (10000,)
```

```
In [4]: # 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()
                        bird
         plane
                                       deer
                                              dog
                                                                    ship
                 car
                                cat
                                                     frog
                                                            horse
                                                                           truck
         🚾 🚰 👺 🤔 🍇 🌌 🥦
In [5]: # 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]
In [6]: # Reshape the image data into rows
        X train = np.reshape(X train, (X train.shape[0], -1))
        X test = np.reshape(X test, (X test.shape[0], -1))
        print(X train.shape, X test.shape)
```

```
(5000, 3072) (500, 3072)
In [7]: from cs6353.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 i-th train example.

First, open cs6353/classifiers/k\_nearest\_neighbor.py and implement the function compute\_distances\_two\_loops that uses a (very inefficient) double loop over all pairs of (test, train) examples and computes the distance matrix one element at a time.

```
In [8]: # Open cs6353/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)
       # We can visualize the distance matrix: each row is a single test example
        # its distances to training examples
        plt.imshow(dists, interpolation='none')
        plt.show()
          0
        250
            0
                        1000
                                       2000
                                                     3000
                                                                    4000
```

**Inline Question #1:** Notice the structured patterns in the distance matrix, where some rows or columns are visible brighter. (Note that with the default color scheme black indicates low distances while white indicates high distances.)

- What in the data is the cause behind the distinctly bright rows?
- · What causes the columns?

#### Your Answer:

- For brighter rows, it indicates that the test image distance is relatively high from all the training images. It suggests that there is low similarity between all the classes in the training labels to the test image.
- For brighter columns, it indicates that that particular training image has a relatively high distance to all test images. It suggests that there is low similarity between all the test images in the testing set to that training image.

```
In [10]: # 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, acc
Got 137 / 500 correct => accuracy: 0.274000
```

You should expect to see approximately 27% accuracy. Now lets try out a larger k, say k = 5:

```
In [11]: 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, acc
    Got 139 / 500 correct => accuracy: 0.278000
```

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

**Inline Question 2** We can also other distance metrics such as L1 distance. The performance of a Nearest Neighbor classifier that uses L1 distance will not change if (Select all that apply.):

- 1. The data is preprocessed by subtracting the mean.
- 2. The data is preprocessed by subtracting the mean and dividing by the standard deviation.
- 3. The coordinate axes for the data are rotated.
- 4. None of the above. (Mean and standard deviation in (1) and (2) are vectors and can be different across dimensions)

Your Answer.

## 1 and 2

Your explanation:

 The mean will cancel out when computing distance as shown in the equation below because it is the same constant. By subtracting the mean to each data points, we are just shifting the data to center around the origin. Hence, no change in the distance.

```
|(p_i - mean) - (q_i - mean)| = |p_i - mean - q_i + mean| = |p_i - q_i|.
```

The standard deviation (sd) will only scale the distance by some constant. Each data point is scaled by the same sd constant, hence there is no difference in the distance comparison.

```
|[(p_i - mean)]/sd - [(q_i - mean)]/sd| = |p_i - q_i|/sd.
```

3. L1 distance computation considers the axis orientation. Hence, if the data matrix is rotated, then the data matrix is transformed and it will more likely to produce different value for the distance.

```
In [12]: # Now lets speed up distance matrix computation by using partial vectoriz
         # with one loop. Implement the function compute distances one loop and ru
         # code below:
         dists one = classifier.compute distances one loop(X test)
         # To ensure that our vectorized implementation is correct, we make sure t
         # agrees with the naive implementation. There are many ways to decide whe
         # two matrices are similar; one of the simplest is the Frobenius norm. In
         \# you haven't seen it before, the Frobenius norm of two matrices is the s
         # root of the squared sum of differences of all elements; in other words,
         # the matrices into vectors and compute the Euclidean distance between th
         difference = np.linalg.norm(dists - dists one, ord='fro')
         print('Difference was: %f' % (difference, ))
         if difference < 0.001:</pre>
            print('Good! The distance matrices are the same')
         else:
            print('Uh-oh! The distance matrices are different')
         Difference was: 0.000000
         Good! The distance matrices are the same
In [13]: # Now implement the fully vectorized version inside compute distances no
         # and run the code
         dists two = classifier.compute distances no loops(X test)
         # check that the distance matrix agrees with the one we computed before:
         difference = np.linalg.norm(dists - dists two, ord='fro')
         print('Difference was: %f' % (difference, ))
         if difference < 0.001:</pre>
            print('Good! The distance matrices are the same')
         else:
            print('Uh-oh! The distance matrices are different')
         Difference was: 0.000000
         Good! The distance matrices are the same
In [14]: # Let's compare how fast the implementations are
         def time function(f, *args):
             0.00
             Call a function f with args and return the time (in seconds) that it
             import time
             tic = time.time()
             f(*args)
             toc = time.time()
             return toc - tic
         two loop time = time function(classifier.compute distances two loops, X t
```

```
print('Two loop version took %f seconds' % two_loop_time)

one_loop_time = time_function(classifier.compute_distances_one_loop, X_te
print('One loop version took %f seconds' % one_loop_time)

no_loop_time = time_function(classifier.compute_distances_no_loops, X_tes
print('No loop version took %f seconds' % no_loop_time)

# you should see significantly faster performance with the fully vectoriz
```

Two loop version took 36.838056 seconds One loop version took 37.831194 seconds No loop version took 0.562591 seconds

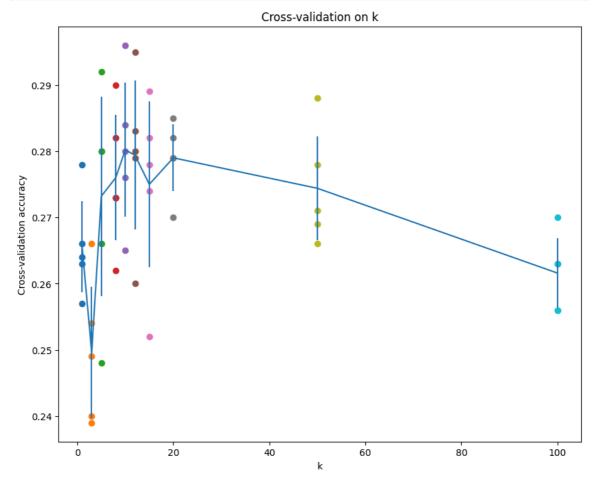
## Cross-validation

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

```
In [15]: num folds = 5
       k \text{ choices} = [1, 3, 5, 8, 10, 12, 15, 20, 50, 100]
       X train folds = []
       y train folds = []
       # TODO:
       # Split up the training data into folds. After splitting, X train folds a
       # y train folds should each be lists of length num folds, where
       # y train folds[i] is the label vector for the points in X train folds[i]
       # Hint: Look up the numpy array split function.
       **
       # Your code
       X train folds = np.array split(X train, num folds)
       y train folds = np.array split(y train, num folds)
       END OF YOUR CODE
       \# A dictionary holding the accuracies for different values of k that we f
       # when running cross-validation. After running cross-validation,
       # k to accuracies[k] should be a list of length num folds giving the diff
       # 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 tim
       # where in each case you use all but one of the folds as training data an
       \# last fold as a validation set. Store the accuracies for all fold and al
       # values of k in the k to accuracies dictionary.
       # Your code
       for k in k choices:
        accuracies = []
        for i in range(num folds):
           classifier = KNearestNeighbor()
           X train without valid set = np.copy(X train folds)
```

```
k = 1, accuracy = 0.263000
         k = 1, accuracy = 0.257000
         k = 1, accuracy = 0.264000
         k = 1, accuracy = 0.278000
         k = 1, accuracy = 0.266000
         k = 3, accuracy = 0.239000
         k = 3, accuracy = 0.249000
         k = 3, accuracy = 0.240000
         k = 3, accuracy = 0.266000
         k = 3, accuracy = 0.254000
         k = 5, accuracy = 0.248000
         k = 5, accuracy = 0.266000
         k = 5, accuracy = 0.280000
         k = 5, accuracy = 0.292000
         k = 5, accuracy = 0.280000
         k = 8, accuracy = 0.262000
         k = 8, accuracy = 0.282000
         k = 8, accuracy = 0.273000
         k = 8, accuracy = 0.290000
         k = 8, accuracy = 0.273000
         k = 10, accuracy = 0.265000
         k = 10, accuracy = 0.296000
         k = 10, accuracy = 0.276000
         k = 10, accuracy = 0.284000
         k = 10, accuracy = 0.280000
         k = 12, accuracy = 0.260000
         k = 12, accuracy = 0.295000
         k = 12, accuracy = 0.279000
         k = 12, accuracy = 0.283000
         k = 12, accuracy = 0.280000
         k = 15, accuracy = 0.252000
         k = 15, accuracy = 0.289000
         k = 15, accuracy = 0.278000
         k = 15, accuracy = 0.282000
         k = 15, accuracy = 0.274000
         k = 20, accuracy = 0.270000
         k = 20, accuracy = 0.279000
         k = 20, accuracy = 0.279000
         k = 20, accuracy = 0.282000
         k = 20, accuracy = 0.285000
         k = 50, accuracy = 0.271000
         k = 50, accuracy = 0.288000
         k = 50, accuracy = 0.278000
         k = 50, accuracy = 0.269000
         k = 50, accuracy = 0.266000
         k = 100, accuracy = 0.256000
         k = 100, accuracy = 0.270000
         k = 100, accuracy = 0.263000
         k = 100, accuracy = 0.256000
         k = 100, accuracy = 0.263000
In [16]: # 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 deviati
         accuracies_mean = np.array([np.mean(v) for k, v in sorted(k_to_accuracies
         accuracies_std = np.array([np.std(v) for k,v in sorted(k_to_accuracies.it
         plt.errorbar(k choices, accuracies mean, yerr=accuracies std)
         plt.title('Cross-validation on k')
         plt.xlabel('k')
```





```
In [17]: # Based on the cross-validation results above, choose the best value for
    # retrain the classifier using all the training data, and test it on the
    # 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, acc
Got 141 / 500 correct => accuracy: 0.282000
```

**Inline Question 3** Which of the following statements about \$k\$-Nearest Neighbor (\$k\$-NN) are true in a classification setting, and for all \$k\$? Select all that apply.

- 1. The training accuracy of a 1-NN (nearest neighbor) will always be better than or equal to that of 5-NN.
- 2. The test accuracy of a 1-NN will always be better than that of a 5-NN.
- 3. The decision boundary of the k-NN classifier is linear.
- 4. The time needed to classify a test example with the k-NN classifier grows with the size of the training set.
- 5. It remembers all of the training data.
- 6. remembers all of the test data.

- 7. KNN cannot be applied for distance metrics other than L1 and L2.
- 8. KNN does more computation during test time than training time.
- 9. KNN does more computation during training time than test time.
- 10. KNN does equal computation during training time and test time.
- 11. KNN does equal computation during training time and test time.

### Your Answer.

**True**: 1, 4, 5, 8

## Your explanation:

- 1. *True*: Because in 1-NN all data points in training set are classified to the single closest neighbour which is itself. Hence, there is no error.
- 2. False: Because there is a chance of misclassification in the testing phase. 1-NN has a higher possibility of misclassification than 5-NN. Because in 5-NN it has more options to consider so it might get the wrong label for the closest image but correct class from the other closest images.
- 3. *False*: Because the data is not always linearly seperable. Since it is n-dimensional space, it may have a complex space and can not be represented by a hyperplane.
- 4. *True*: Because when kNN model is classifying for a test example, it went through all the training examples and do the computation to classify. Hence, more data in training set, more time for computation.
- 5. *True*: Because in the training phase, it stores all the data but no computation.
- 6. *False*: Because the test data is not used in the training phase and in testing phase, it just do the computation to classify the images but not storing them.
- 7. False: Because L1 and L2 are just two techniques for measuring distance. Other techniques for measuring distance can still be applied to kNN algorithm concept.
- 8. True: Because all computation is done in testing phase.
- 9. False: Because there is no computation in training phase.
- 10. *False*: Because all computation is done in testing phase and there is no computation in training phase.

In [24]: !jupyter nbconvert --to html /content/drive/MyDrive/MS/CS6353/assignment1

 $[{\tt NbConvertApp}] \ {\tt Converting \ notebook / content/drive/MyDrive/MS/CS6353/assignment1/knn.ipynb \ to \ html}$ 

[NbConvertApp] Writing 1165688 bytes to /content/drive/MyDrive/MS/CS6353/assignment1/knn.html