sheet2

October 27, 2017

1 Exercise Sheet 2: Timing, Numpy, Plotting

The previous exercise sheet introduced several methods for classification: decision trees, nearest neighbors, and nearest means. Of those, the one that could learn from the data, and that also offered enough complexity to produce an accurate decision function was k-nearest neighbors. However, nearest neighbors can be slow when implemented in pure Python (i.e. with loops). This is especially the case when the number of data points or input dimensions is large.

In this exercise sheet, we will speed up nearest neighbors by utilizing numpy and scipy packages. Your task will be to replace list-based operations by vector-based operations between numpy arrays. The speed and correctness of the implementations will then be tested. In particular, performance graphs will be drawn using the library matplotlib.

1.1 Python Nearest Neighbor

The most basic element of computation of nearest neighbors is its distance function relating two arbitrary data points x1 and x2. We assume that these points are iterable (i.e. we can use a loop over their dimensions). One way among others to compute the square Euclidean distance between two points is by computing the sum of the component-wise distances

where the prefix "py-" of the function indicates that the latter makes use of Python instead of numpy. Once the distance matrix has been implemented, the nearest neighbor for a given unlabeled point u that we would like to classify is obtained by iterating over all points in the training set (X,Y), selecting the point with smallest distance to u, and returning its corresponding label. Here X denotes the list of inputs in the training set and Y denotes the list of labels.

```
In [136]: def pynearest(u,X,Y,distance=pydistance):
    xbest = None
    ybest = None
    dbest = float('inf')

for x,y in zip(X,Y):
    d = distance(u,x)
    if d < dbest:
        ybest = y</pre>
```

```
xbest = x
dbest = d
return ybest
```

Note that this function either uses function pydistance (given as default if the argument distance is not specified). Or one could specify as argument a more optimized function for distance computation, for example, one that uses numpy. Finally, one might not be interested in classifying a single point, but many of them. The method below receives a collection of such unlabeled test points stored in the variable U. The function returns a list of predictions associated to each test point.

Again, such function uses by default the Python nearest neighbor search (with a specified distance function). However, we can also specified a more optimized nearest neighbor function, for example, based on numpy. Finally, one could consider an alternative function to pybatch that would use numpy from the beginning to the end. The implementation of such more optimized functions, and the testing of their correct behavior and higher performance will be the object of this exercise sheet.

1.2 Testing and correctness

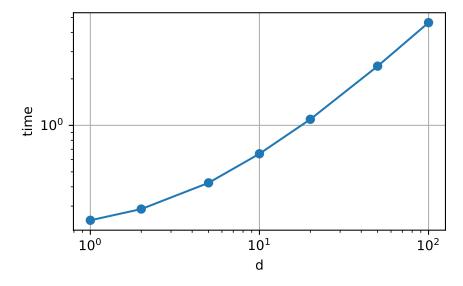
As a starting point, the code below tests the output of the nearest neighbor algorithm for some toy dataset with fixed parameters. In particular, the function $\mathtt{data.toy}(\texttt{M}, \texttt{N}, \texttt{d})$ generates a problem with M unlabeled test points stored in a matrix U of size (M x d), then N labeled training points stored in a matrix X of size (N x d) and the output label is stored in a vector Y of size N composed of zeros and ones encoding the two possible classes. The variable d denotes the number of dimensions of each point. The toy dataset is pseudo-random, that is, for fixed parameters, it produce a random-looking dataset, but every time the method is called with the same parameters, the dataset is the same. The pseudo-randomness property will be useful to verify that each nearest neighbor implementation performs the same overall computation.

In particular, the output of this function will help us to verify that the more optimized numpy-based versions of nearest neighbor are still valid.

1.3 Plotting and performance

We now describe how to build a plot that relates a certain parameter of the dataset (e.g. the number of input dimensions d to the time required for the computation. We first initialize the basic plotting environment.

The command "%matplotlib inline" tells IPython notebook that the plots should be rendered inside the notebook. The following code plots the computation time of predicting 100 points from the test set using a training set of size 100, and where we vary the number of input dimensions.



The time on the vertical axis is in seconds. Note that the exact computation time depends on the speed of your computer. As expected, the computation time increases with the number of input dimensions. Unfortunately, for the small dataset considered here (100 training and test points of 100 dimensions each), the algorithm already takes more than one second to execute. Thus, it is necessary for practical applications (e.g. the digit recognition task that we will consider at the end of this exercise sheet) to accelerate this nearest neighbor algorithm.

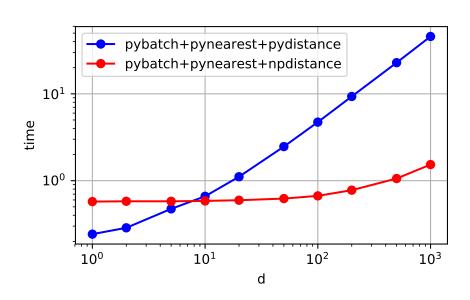
1.4 Accelerating the distance computation (25 P)

In this first exercise, we would like to accelerate the function that compute pairwise distances.

- Create a new function npdistance(x1,x2) with the same output as pydistance(x1,x2), but that computes the squared Euclidean distance using numpy operations.
- Print its output for the same toy example with parameters M=20, N=100, d=50 considered before (i.e. data.toy(20,100,50)). Verify that in both cases (i.e. using either npdistance or pydistance in the function pybatch) the output remains the same.
- Create a plot similar to the one above, but where the computation time required by both methods are shown in a superposed manner. Here, we fix M=100, N=100, and we let d vary from 1 to 1000, taking the list of values [1, 2, 5, 10, 20, 50, 100, 200, 500, 1000].

```
In [190]: ### Replace by your own code
          #import solutions
          #solutions.exercise1()
          import numpy
          def npdistance(x1,x2):
              if (len(x1.shape) == 1):
                  return numpy.sum(numpy.power(x1-x2,2))
              else:
                  return numpy.sum(numpy.power(x1-x2,2),1)
          def plot1():
              # Values for the number of dimensions d to test
              dlist = [1,2,5,10,20,50,100,200,500,1000]
              # Measure the computation time for each choice of number of dimensions d
              tlistPy = []
              tlistNp = []
              for d in dlist:
                  U, X, Y = data.toy(100, 100, d)
                  #Python
                  a = time.clock()
                  pybatch(U,X,Y)
                  b = time.clock()
                  tlistPy += [b-a]
                  #Numpy
```

```
a = time.clock()
                 pybatch(U,X,Y,distance=npdistance)
                  b = time.clock()
                  tlistNp += [b-a]
              # Plot the results in a graph
             plt.figure(figsize=(5,3))
             plt.plot(dlist,tlistPy, "-o", color='blue', label="pybatch+pynearest+pydistance")
             plt.plot(dlist,tlistNp, "-o", color='red', label="pybatch+pynearest+npdistance")
             plt.xscale('log');plt.yscale('log'); plt.xlabel('d'); plt.ylabel('time'); plt.grid
             plt.legend()
             plt.show()
          #Testing
         U, X, Y = data.toy(20, 100, 50)
          print('comparison to baseline implementation')
         print(('pybatch+pynearest+pydistance', pybatch(U,X,Y)))
          print(('pybatch+pynearest+npdistance',pybatch(U,X,Y,distance=npdistance)))
         plot1()
comparison to baseline implementation
('pybatch+pynearest+pydistance', [1, 1, 1, 0, 0, 1, 0, 1, 1, 0, 1, 1, 0, 1, 1, 1, 0])
('pybatch+pynearest+npdistance', [1, 1, 1, 0, 0, 1, 0, 1, 1, 0, 1, 1, 0, 1, 1, 0])
```



• Based on your results, explain what kind of speedup numpy provides, and in what regime do you expect the speedup to be the most important.

As one can see in the plot below from about dimension 10 numpy provides an speed up in the performance of the programm. For lower dimensions it would be more efficient to use only Python, but the higher the dimension gets, the better it becomes to use numpy. As we are concered about reading heaps of data in real time with huge dimensions, numpy gives us a great performance for many real life applications

1.5 Accelerating the nearest neighbor search (25 P)

Motivated by the success of the numpy optimized distance computation, we would like further accelerate the code by performing nearest neighbor search directly in numpy.

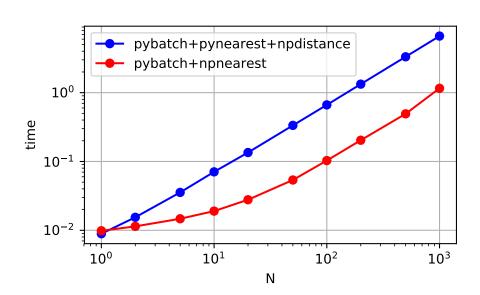
- Create a new function npnearest(u, X, Y) as an alternative to the function pynearest(u, X, Y, distance=npdistance) that we have used in the previous exercise.
- Print its output for the same toy example as before (i.e. data.toy(20,100,50)). Verify that the output remains the same compared to the implementation of the previous exercise.
- Create a plot similar to the one above, where the new method is compared to the previous one. Here, we fix M=100, d=100, and we let N take different values [1, 2, 5, 10, 20, 50, 100, 200, 500, 1000].

```
In [182]: ### Replace by your own code
          #import solutions
          #solutions.exercise2()
          def npnearest(u, X, Y, distance=npdistance):
              return Y[numpy.argmin(distance(X,u))]
          def plot2():
              # Values for the number of dimensions d to test
              Nlist = [1,2,5,10,20,50,100,200,500,1000]
              # Measure the computation time for each choice of number of dimensions d
              tlistPy = []
              tlistNp = []
              for N in Nlist:
                  U, X, Y = data.toy(100, N, 100)
                  #Python
                  a = time.clock()
                  pybatch(U,X,Y,nearest=pynearest,distance=npdistance)
                  b = time.clock()
                  tlistPy += [b-a]
                  #Numpy
                  a = time.clock()
                  pybatch(U,X,Y,nearest=npnearest, distance=npdistance)
                  b = time.clock()
                  tlistNp += [b-a]
              # Plot the results in a graph
```

```
plt.figure(figsize=(5,3))
    plt.plot(Nlist,tlistPy, "-o", color='blue', label="pybatch+pynearest+npdistance")
    plt.plot(Nlist,tlistNp, "-o", color='red', label="pybatch+npnearest")
    plt.xscale('log');plt.yscale('log'); plt.xlabel('N'); plt.ylabel('time'); plt.grid
    plt.legend()
    plt.show()

U,X,Y = data.toy(20,100,50)
    print('comparison to baseline implementation')
    print(('pybatch+pynearest+npdistance', pybatch(U,X,Y,nearest=pynearest, distance=npdist print(('pybatch+npnearest', pybatch(U,X,Y,nearest=npnearest, distance=npdistance)))
    plot2()

comparison to baseline implementation
('pybatch+pynearest+npdistance', [1, 1, 1, 0, 0, 1, 0, 1, 1, 0, 1, 1, 1, 0, 1, 1, 1, 0])
('pybatch+npnearest', [1, 1, 1, 0, 0, 1, 0, 1, 1, 1, 0, 1, 0, 1, 1, 1, 0])
```



• Based on your results, explain what kind of speedup this further optimization provides, and in what regime the speedup is the most significant.

The only difference between the blue and the red curve is the use of npnearest and pynearest. Since the python implementation uses a for loop and the numpy variant internally works with c, it is clear that npnearest is faster. In particular by increasing N we also enlarge X and Y (i.e. the for loop lasts longer) in every step, which slows down pynearest. Considering the graphic the most significant speedup is in the first and the second regime.

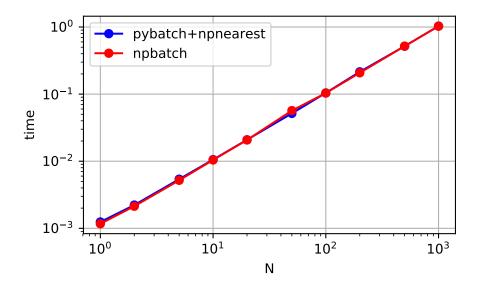
1.6 Accelerating the processing of multiple test points (25 P)

Not yet fully happy with the performance of the algorithm, we would like to further optimize it by avoiding performing a loop on the test points, and instead, classify them all at once.

- Create a new function npbatch(U,X,Y) as a replacement of the implementation pybatch(U,X,Y,nearest=npnearest) that we have built in the previous exercise.
- Print its output for the same dataset data.toy(20,100,50) and verify that the output remains the same as for the previous implementation.
- Create a plot comparing the computation time of the new implementation compared to the previous one. Here, we fix N=100, d=100, and we let M vary from 1 to 1000 with values [1, 2, 5, 10, 20, 50, 100, 200, 500, 1000].

```
In [186]: ### Replace by your own code
          def npbatch(U,X,Y,distance=npdistance):
              return [npnearest(u, X, Y, distance=distance) for u in U]
          def plot3():
              # Values for the number of dimensions d to test
              Mlist = [1, 2, 5, 10, 20, 50, 100, 200, 500, 1000]
              # Measure the computation time for each choice of number of dimensions d
              tlistPy = []
              tlistNp = []
              for M in Mlist:
                  U,X,Y = data.toy(M,100,100)
                  #Python
                  a = time.clock()
                  pybatch(U,X,Y,distance=npdistance, nearest=npnearest)
                  b = time.clock()
                  tlistPy += [b-a]
                  #Numpy
                  a = time.clock()
                  npbatch(U,X,Y)
                  b = time.clock()
                  tlistNp += [b-a]
              # Plot the results in a graph
              plt.figure(figsize=(5,3))
              plt.plot(Mlist,tlistPy, "-o", color='blue', label="pybatch+npnearest")
              plt.plot(Mlist,tlistNp, "-o", color='red', label="npbatch")
              plt.xscale('log');plt.yscale('log'); plt.xlabel('N'); plt.ylabel('time'); plt.grid
              plt.legend()
              plt.show()
          U, X, Y = data.toy(20, 100, 50)
          print('comparison to baseline implementation')
          print(('pybatch+npnearest', pybatch(U,X,Y,distance=npdistance, nearest=npnearest)))
          print(('npbatch', npbatch(U,X,Y)))
          plot3()
comparison to baseline implementation
```

('pybatch+npnearest', [1, 1, 1, 0, 0, 1, 0, 1, 1, 0, 1, 1, 0, 1, 0, 1, 1, 1, 0])

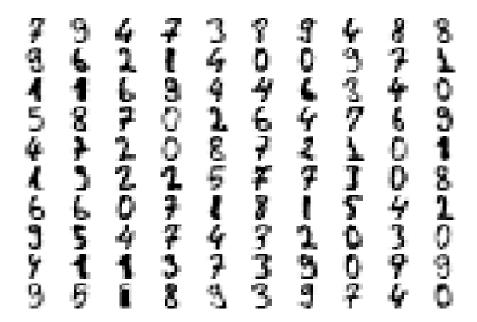


1.7 Application to real data (25 P)

Having now implemented an efficient K-nearest neighbor classifier, we can test it on real problems with many data points and dimensions. We consider a small handwritten digits recognition dataset, that can be directly obtained from the library scikit-learn. This dataset consists of handwritten digits of size 8 x 8 flattened into arrays of size 64, with class between 0 and 9. We use a function data.digits() to load the data and arrange data points in some predefined order.

```
In [188]: X,Y = data.digits()
```

Using the function imshow of matplotlib to visualize the first 100 digits of the dataset.



- Partition the data into a "training" set and "test" set. The first one contains the 1000 first digits of X, and the second one contains the remaining ones.
- Assume that you don't know the labels for the test data and classify the test data using your efficient nearest neighbor implementation.
- Print the predicted labels for the test set.

```
8 2 6 1 1 3 8 2 3 5 5 8 0 5 4 8 0 7 3 6 4 0 8 9 4 8 9 9 7 4 4 6 8 4 5 2 9
9\; 4\; 0\; 5\; 8\; 5\; 2\; 2\; 7\; 6\; 4\; 8\; 3\; 0\; 7\; 6\; 5\; 6\; 1\; 0\; 9\; 3\; 5\; 6\; 3\; 6\; 3\; 3\; 0\; 0\; 1\; 4\; 1\; 1\; 9\; 3\; 8
8 8 8 2 0 7 6 5 6 8 2 0 6 8 6 0 0 0 6 9 3 7 0 1 8 9 9 9 1 7 0 5 5 5 6 4 1
4 8 6 6 8 3 1 0 5 2 2 6 8 4 2 1 0 4 6 9 9 6 1 7 2 3 4 0 5 5 7 4 8 1 1 7 8
7\ 1\ 7\ 5\ 1\ 2\ 1\ 3\ 2\ 2\ 9\ 8\ 7\ 8\ 2\ 7\ 2\ 7\ 1\ 0\ 9\ 2\ 8\ 4\ 2\ 1\ 0\ 4\ 2\ 7\ 2\ 6\ 9\ 2\ 1\ 2\ 5
4 7 1 6 3 4 4 7 0 0 9 9 9 9 9 1 1 9 5 7 3 8 4 8 6 6 3 1 8 6 8 4 3 6 2 3
2\ 1\ 1\ 8\ 1\ 9\ 4\ 4\ 9\ 0\ 1\ 7\ 9\ 8\ 3\ 6\ 2\ 2\ 5\ 4\ 1\ 2\ 6\ 1\ 1\ 2\ 3\ 6\ 7\ 8\ 3\ 4\ 4\ 4\ 6\ 5\ 9
4 6 6 1 3 2 6 5 7 9 4 7 6 8 5 6 1 9 1 9 7 5 7 8 8 5 1 0 7 2 6 9 1 0 7 3 2
2 7 3 0 7 0 9 3 8 8 3 6 2 4 2 8 5 9 6 2 7 6 2 9 5 5 3 8 2 6 2 7 5 3 8 3 0
3 0 4 0 2 8 5 1 0 7 7 3 2 7 6 8 8 3 7 4 9 5 0 4 1 1 2 1 3 6 5 0 5 1 2 6 1
5\; 4\; 7\; 4\; 8\; 4\; 1\; 8\; 8\; 8\; 8\; 8\; 7\; 1\; 8\; 9\; 3\; 1\; 2\; 1\; 2\; 3\; 0\; 1\; 6\; 1\; 3\; 5\; 3\; 7\; 1\; 3\; 6\; 9\; 3\; 6\; 6
2\; 3\; 2\; 0\; 0\; 1\; 9\; 6\; 5\; 9\; 5\; 9\; 4\; 4\; 2\; 5\; 4\; 9\; 0\; 0\; 6\; 2\; 5\; 7\; 2\; 1\; 7\; 0\; 3\; 0\; 6\; 2\; 1\; 6\; 9\; 9\; 5
8 3 3 6 6 1 6 5 7 2 8 2 3 1 0 9 3 9 7 9 3 6 9 8 7 4 6 9 7 6 8 3 5 0 3 0 4
4 3 1 4 2 6 5 6 1 7 5 6 6 3 2 9 2 3 1 1 1 2 3 3 2 0 9 4 9 8 3 2 0 0 9 5 1
3 5 5 3 6 0 3 2 0 1 9 4 0 9 6 7 8 5 0 6 5 9 3 3 9 5 5 5 6 4 3 5 6 1 4 6 8
7 1 8 6 5 2 4 6 1 1 8 2 5 7 0 6 4 3 0 9 7 0 6 5 7 7 7 5 5 4 3 5 0 3 2 5 9
3 8 1 0 6 7 3 5 2 8 9 0 8 5 2 0 7 6 5 1 6 9 9 9 4 2 6 6 4 4 4 2 3 2 0 9 2
5\; 3\; 9\; 3\; 3\; 6\; 0\; 0\; 2\; 3\; 9\; 5\; 5\; 6\; 6\; 4\; 7\; 0\; 0\; 9\; 4\; 7\; 9\; 7\; 5\; 4\; 1\; 0\; 8\; 1\; 1\; 6\; 9\; 2\; 9\; 7\; 8
5 1 0 5 6 6 9 9 6 3 4 5 1 5 6 9 1 3 4 2]
```

Finally, in order to determine the accuracy of the classifier, we would like to compare the predictions with the ground truth (i.e. the true labels from the test data).

• Compute the fraction of the time on the test set where the predictions of the nearest neighbor algorithm and labels disagree.