

# CS 124 Programming Assignment 3: Spring 2024

**Your name(s) (up to two):**

**Collaborators:** (You shouldn't have any collaborators but the up-to-two of you, but tell us if you did.)

**No. of late days used on previous psets:**

**No. of late days used after including this pset:**

Homework is due Thursday 2024-04-18 at 11:59pm ET. You are allowed up to **twelve** late days through the semester, but the number of late days you take on each assignment must be a nonnegative integer at most **two**.

For this programming assignment, you will implement a number of heuristics for solving the NUMBER PARTITION problem, which is (of course) NP-complete. As input, the number partition problem takes a sequence  $A = (a_1, a_2, \dots, a_n)$  of non-negative integers. The output is a sequence  $S = (s_1, s_2, \dots, s_n)$  of signs  $s_i \in \{-1, +1\}$  such that the *residue*

$$u = \left| \sum_{i=1}^n s_i a_i \right|$$

is minimized. Another way to view the problem is the goal is to split the set (or multi-set) of numbers given by  $A$  into two subsets  $A_1$  and  $A_2$  with roughly equal sums. The absolute value of the difference of the sums is the residue.

As a warm-up exercise, you will first prove that even though Number Partition is NP-complete, it can be solved in pseudo-polynomial time. That is, suppose the sequence of terms in  $A$  sum up to some number  $b$ . Then each of the numbers in  $A$  has at most  $\log b$  bits, so a polynomial time algorithm would take time polynomial in  $n \log b$ . Instead you should find a dynamic programming algorithm that takes time polynomial in  $nb$ .

**Give a dynamic programming solution to the Number Partition problem.**

One deterministic heuristic for the Number Partition problem is the Karmarkar-Karp algorithm, or the KK algorithm. This approach uses *differencing*. The differencing idea is to take two elements from  $A$ , call them  $a_i$  and  $a_j$ , and replace the larger by  $|a_i - a_j|$  while replacing the smaller by 0. The intuition is that if we decide to put  $a_i$  and  $a_j$  in different sets, then it is as though we have one element of size  $|a_i - a_j|$  around. An algorithm based on differencing repeatedly takes two elements from  $A$  and performs a differencing until there is only one element left; this element equals an attainable residue. (A sequence of signs  $s_i$  that yields this residue can be determined from the differencing operations performed in linear time by two-coloring the graph  $(A, E)$  that arises, where  $E$  is the set of pairs  $(a_i, a_j)$  that are used in the differencing steps. You will not need to construct the  $s_i$  for this assignment.)

For the Karmarkar-Karp algorithm suggests repeatedly taking the largest two elements remaining in  $A$  at each step and differencing them. For example, if  $A$  is initially  $(10, 8, 7, 6, 5)$ , then the KK algorithm proceeds as follows:

$$\begin{aligned} (10, 8, 7, 6, 5) &\rightarrow (2, 0, 7, 6, 5) \\ &\rightarrow (2, 0, 1, 0, 5) \\ &\rightarrow (0, 0, 1, 0, 3) \\ &\rightarrow (0, 0, 0, 0, 2) \end{aligned}$$

Hence the KK algorithm returns a residue of 2. The best possible residue for the example is 0.

**Explain briefly how the Karmarkar-Karp algorithm can be implemented in  $O(n \log n)$  steps, assuming the values in  $A$  are small enough that arithmetic operations take one step.**

You will compare the Karmarkar-Karp algorithm and a variety of randomized heuristic algorithms on random input sets. Let us first discuss two ways to represent solutions to the problem and the state space based on these representations. Then we discuss heuristic search algorithms you will use.

The standard representation of a solution is simply as a sequence  $S$  of  $+1$  and  $-1$  values. A random solution can be obtained by generating a random sequence of  $n$  such values. Thinking of all possible solutions as a state space, a natural way to define neighbors of a solution  $S$  is as the set of all solutions that differ from  $S$  in either one or two places. This has a natural interpretation if we think of the  $+1$  and  $-1$  values as determining two subsets  $A_1$  and  $A_2$  of  $A$ . Moving from  $S$  to a neighbor is accomplished either by moving one or two elements from  $A_1$  to  $A_2$ , or moving one or two elements from  $A_2$  to  $A_1$ , or swapping a pair of elements where one is in  $A_1$  and one is in  $A_2$ .

A *random move* on this state space can be defined as follows. Choose two random indices  $i$  and  $j$  from  $[1, n]$  with  $i \neq j$ . Set  $s_i$  to  $-s_i$  and with probability  $1/2$ , set  $s_j$  to  $-s_j$ .

An alternative way to represent a solution called *prepartitioning* is as follows. We represent a solution by a sequence  $P = \{p_1, p_2, \dots, p_n\}$  where  $p_i \in \{1, \dots, n\}$ . The sequence  $P$  represents a prepartitioning of the elements of  $A$ , in the following way: if  $p_i = p_j$ , then we enforce the restriction that  $a_i$  and  $a_j$  have the same sign. Equivalently, if  $p_i = p_j$ , then  $a_i$  and  $a_j$  both lie in the same subset, either  $A_1$  or  $A_2$ .

We turn a solution of this form into a solution in the standard form using two steps:

- We derive a new sequence  $A'$  from  $A$  which enforces the prepartitioning from  $P$ . Essentially  $A'$  is derived by resetting  $a_i$  to be the sum of all values  $j$  with  $p_j = i$ , using for example the following pseudocode:

```

 $A' = (0, 0, \dots, 0)$ 
for  $j = 1$  to  $n$ 
     $a'_{p_j} = a'_{p_j} + a_j$ 

```

- We run the KK heuristic algorithm on the result  $A'$ .

For example, if  $A$  is initially  $(10, 8, 7, 6, 5)$ , the solution  $P = (1, 2, 2, 4, 5)$  corresponds to the following run of the KK algorithm:

$$\begin{aligned}
 A = (10, 8, 7, 6, 5) &\rightarrow A' = (10, 15, 0, 6, 5) \\
 (10, 15, 0, 6, 5) &\rightarrow (0, 5, 0, 6, 5) \\
 &\rightarrow (0, 0, 0, 1, 5) \\
 &\rightarrow (0, 0, 0, 0, 4)
 \end{aligned}$$

Hence in this case the solution  $P$  has a residue of 4.

Notice that all possible solution sequences  $S$  can be generated using this prepartition representation, as any split of  $A$  into sets  $A_1$  and  $A_2$  can be obtained by initially assigning  $p_i$  to 1 for all  $a_i \in A_1$  and similarly assigning  $p_i$  to 2 for all  $a_i \in A_2$ .

A random solution can be obtained by generating a sequence of  $n$  values in the range  $[1, n]$  and using this for  $P$ . Thinking of all possible solutions as a state space, a natural way to define neighbors of a solution  $P$  is as the set of all solutions that differ from  $P$  in just one place. The interpretation is that we change the prepartitioning by changing the partition of one element. A *random move* on this state space can be defined as follows. Choose two random indices  $i$  and  $j$  from  $[1, n]$  with  $p_i \neq j$  and set  $p_i$  to  $j$ .

You will try each of the following three algorithms for both representations.

- Repeated random: repeatedly generate random solutions to the problem, as determined by the representation.

```

Start with a random solution  $S$ 
for iter = 1 to max_iter
     $S' =$  a random solution
    if residue( $S'$ ) < residue( $S$ ) then  $S = S'$ 
return  $S$ 

```

- Hill climbing: generate a random solution to the problem, and then attempt to improve it through moves to better neighbors.

```

Start with a random solution  $S$ 
for iter = 1 to max_iter
     $S' =$  a random neighbor of  $S$ 
    if residue( $S'$ ) < residue( $S$ ) then  $S = S'$ 
return  $S$ 

```

- Simulated annealing: generate a random solution to the problem, and then attempt to improve it through moves to neighbors, that are not always better.

```

Start with a random solution  $S$ 
 $S'' = S$ 
for iter = 1 to max_iter
     $S' =$  a random neighbor of  $S$ 
    if residue( $S'$ ) < residue( $S$ ) then  $S = S'$ 
    else  $S = S'$  with probability  $\exp(-(res(S')-res(S))/T(iter))$ 
    if residue( $S$ ) < residue( $S''$ ) then  $S'' = S$ 
return  $S''$ 

```

Note that for simulated annealing we have the code return the best solution seen thus far.

You will run experiments on sets of 100 integers, with each integer being a random number chosen uniformly from the range  $[1, 10^{12}]$ . Note that these are big numbers. You should use 64 bit integers. Pay attention to things like whether your random number generator works on ranges this large!

Below is the main problem of the assignment.

First, write a routine that takes three arguments: a flag, an algorithm code (see Table 1), and an input file. We'll run typical commands to compile and execute your code, as in programming assignment 2; for example, for C/C++, the run command will look as follows:

```
$ ./partition flag algorithm inputfile
```

The flag is meant to provide you some flexibility; the autograder will only pass 0 as the flag but you may use other values for your own testing, debugging, or extensions. The algorithm argument is one of the values specified in Table 1. You can also assume the inputfile is a list of 100 (unsorted) integers, one per line. The desired output is the residue obtained by running the specified algorithm with these 100 numbers as input.

Code	Algorithm
0	Karmarkar-Karp
1	Repeated Random
2	Hill Climbing
3	Simulated Annealing
11	Prepartitioned Repeated Random
12	Prepartitioned Hill Climbing
13	Prepartitioned Simulated Annealing

Table 1: Algorithm command-line argument values

If you wish to use a programming language other than Python, C++, C, Java, and Go, please contact us first. As before, you should submit either 1) a single source file named one of partition.py, partition.c, partition.cpp, partition.java, Partition.java, or partition.go, or 2) possibly multiple source files named whatever you like, along with a Makefile (named makefile or Makefile).

Second, generate 50 random instances of the problem as described above. For each instance, find the result from using the Karmarkar-Karp algorithm. Also, for each instance, run a repeated random, a hill climbing, and a simulated annealing algorithm, using both representations, each for at least 25,000 iterations. Give tables and/or graphs clearly demonstrating the results. Compare the results and discuss.

For the simulated annealing algorithm, you must choose a *cooling schedule*. That is, you must choose a function  $T(\text{iter})$ . We suggest  $T(\text{iter}) = 10^{10}(0.8)^{\lfloor \text{iter}/300 \rfloor}$  for numbers in the range  $[1, 10^{12}]$ , but you can experiment with this as you please.

Note that, in our random experiments, we began with a random initial starting point.

Discuss briefly how you could use the solution from the Karmarkar-Karp algorithm as a starting point for the randomized algorithms, and suggest what effect that might have. (No experiments are necessary.)

Finally, the following is entirely optional; you'll get no credit for it. But if you want to try something else, it's interesting to do.

**Optional:** Can you design a BubbleSearch-based heuristic for this problem? The Karmarkar-Karp algorithm greedily takes the top two items at each step, takes their difference, and adds that difference back into the list of numbers. A BubbleSearch variant would not necessarily take the top two items in the list, but probabilistically take two items close to the top. (For instance, you might “flip coins” until the first heads; the number of flips (modulo the number of items) gives your first item. Then do the same, starting from where you left off, to obtain the second item. Once you’re down to a small number of numbers – five to ten – you might want to switch back to the standard Karmarkar-Karp algorithm.) Unlike the original Karmarkar-Karp algorithm, you can repeat this algorithm multiple times and get different answers, much like the Repeated Random algorithm you tried for the assignment. Test your BubbleSearch algorithm against the other algorithms you have tried. How does it compare?