

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 q and a_j , and replace the larger by $|q - a_j|$ while replacing the smaller by 0. The intuition is that if we decide to put q and a_j in different sets, then it is as though we have one element of size $|q - a_j|$ 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 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 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

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:

```

 $A = (10, 8, 7, 6, 5) \rightarrow A' = (10, 15, 0, 6, 5)$ 
 $(10, 15, 0, 6, 5) \rightarrow (0, 5, 0, 6, 5)$ 
 $\rightarrow (0, 0, 0, 1, 5)$ 
 $\rightarrow (0, 0, 0, 0, 4)$ 

```

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 $p_i \neq p_j$ and set p_i to p_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  $S$ 
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  $\exp(-(res(S')-res(S))/T(iter))$ 
return  $S$ 

```

Note that one could also have the code return the best solution seen thus far; in practice the two procedures are essentially the same if the process is run for enough iterations.

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.

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. Compare the results and discuss.

For the simulated annealing algorithm, you must choose a *cooling schedule*. That is, you must choose a function $T(iter)$. We suggest $T(iter) = 10^{10}(0.8)^{\lfloor iter/300 \rfloor}$ 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.)