

## Edit Peptide

Palindromic peptides prevail! At least according to Anna, but what else can you expect from a chemist with a palindromic name. Peptides are chemical compounds consisting of  $N \geq 2$  amino acids linked in a chain. We number the amino acids from 0 to  $N - 1$  according to their position in the chain.

The reason why Anna is so enthusiastic about palindromic peptides is that their special spatial structure allows chemists to induce a reaction which swaps a pair of amino acids in the chain: acid 0 can be swapped with acid  $N - 1$ , acid 2 can be swapped with acid  $N - 2$ , and so on. However, editing the molecule structure too much may make the peptide unstable, thus it is not allowed to perform more than  $K$  swaps.

Each amino acid has some potential energy  $P_i$ , represented by a nonnegative integer. Anna wants to perform  $Q$  experiments. In each experiment, she wants to maximize the total potential energy of the amino acids over some contiguous segment of the chain. That is, for a given interval  $[l, r]$  ( $0 \leq l \leq r < N$ ), she wants to maximize the sum  $P_l + P_{l+1} + \dots + P_r$  by performing at most  $K$  palindromic swaps.

Anna starts each experiment with a new peptide molecule, and she uses the same type of molecule for each experiment. Given the description of the peptide and  $Q$  intervals, your task is to compute the maximum potential energy that Anna can achieve in each case. You have to compute the answers *online*, that is, you have to solve the problem for an interval first before learning the next interval (see the Input section for details).

### Input

The first line of the input contains three integers  $N$ ,  $K$  and  $Q$  ( $2 \leq N, Q \leq 200\,000$  and  $0 \leq K \leq N$ ), the size of the peptide chain, the maximum number of swaps and the number of experiments.

The second line contains  $N$  integers  $P_i$  ( $0 \leq P_i \leq 10^9$ ), the potential energies of the acids.

Each of the next  $Q$  lines contains two integers.

- The first line contains the boundaries  $l$  and  $r$  for the first experiment ( $0 \leq l \leq r < N$ ).
- In the  $j$ -th line ( $2 \leq j \leq Q$ ) there are two integers  $s_j$  and  $t_j$  ( $0 \leq s_j, t_j \leq 10^{18}$ ). Let  $a_{j-1}$  denote the correct answer to the  $(j - 1)$ -th question. Then, the value of  $l$  and  $r$  for the  $j$ -th experiment are  $s_j \oplus a_{j-1}$  and  $t_j \oplus a_{j-1}$ , respectively.

Here,  $\oplus$  denotes the binary XOR operator.

### Output

Print  $Q$  lines, where the  $j$ -th line ( $1 \leq j \leq Q$ ) contains the integer  $a_j$ , the maximum sum of potentials in the  $j$ -th experiment.

## Examples

input	output
7 1 5 4 4 0 6 1 1 0 0 2 8 15 18 22 12 10 20 20	9 16 12 16 1

## Explanation

The first interval is  $[l, r] = [0, 2]$ . The sum of potentials without any swaps is  $4 + 4 + 0 = 8$  over this segment, which can be improved to  $a_1 = 4 + 4 + 1 = 9$  by swapping acids 2 and 4.

The second interval is  $[l, r] = [8 \oplus 9, 15 \oplus 9] = [1, 6]$ . The sum of potentials without swaps is 12 which can be improved to  $a_2 = 16$  by swapping acids 0 and 6.

The third interval is  $[l, r] = [18 \oplus 16, 22 \oplus 16] = [2, 6]$ . The sum of potentials without swaps is 8. There are two ways to improve this, either by swapping acids 0 and 6 or by swapping 1 and 5. Since  $K = 1$ , we are only allowed to make one swap, and swapping 0 and 6 is better, so  $a_3 = 12$ .

The fourth interval is  $[l, r] = [12 \oplus 12, 10 \oplus 12] = [0, 6]$ , the whole peptide chain. The sum of potentials is 16 which cannot be improved by swapping a pair of acids, so  $a_4 = 16$ .

The fifth interval is  $[l, r] = [20 \oplus 16, 20 \oplus 16] = [4, 4]$ . The potential of acid 4 is 1, which cannot be improved by swapping acids 4 and 2. So we have  $a_5 = 1$ .