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config.md

~/.bashrc

```
setxkbmap -option caps:escape
```

~/.vimrc

```
set nu rnu sw=4 ts=4 ai ls=2
```

template.cpp

```
#include <bits/stdc++.h>
using namespace std;
#define fastio ios base::sync with stdio(0);cin.tie(0)
#define pb push_back
#define mp make_pair
#define sz(x) int(x.size())
#define trace(x) cerr << #x << ": " << x << endl;
typedef long long 11;
const ll N = 1e6;
const ll INF = 1LL << 61;</pre>
const 11 \text{ MOD} = 1e9 + 7;
void solve() {
signed main() {
    fastio;
    solve();
    return 0;
}
```

Makefile

reference.md

bash

 Make CapsLock work as Esc setxkbmap -option caps:escape

```
• Remove gdb warranty message
```

alias gdb="gdb --silent"

• Copy file to clipboard

xclip -sel c < file

• Change keyboard language setxkbmap us|br

vim

commands

• Compile the program. If there are any errors, the cursor is moved to the first one

:make file

• Go to the next error

:cnext

• Show full error message

:cc

• Show the list of errors

:clist

Split horizontaly

:splitCtrl-w s

• Split verticaly

:vsplitCtrl-w v

options

```
" show line numbers (nu)
set number
" display relative line numbers instead of absolute (rnu)
set relativenumber
" the width of indent for operations like << and >> (sw)
set shiftwidth=4
" display size of the tab character (ts)
set tabstop=4
" use the indentation of the current line when creating a new one (ai)
set autoindent
" constantly show the file name (ls)
set laststatus=2
"" not used
" show folds in a column (fdc)
set foldcolumn=4
" fold by indent (fdm)
set foldmethod=indent
" set vim colorscheme (colo)
colorscheme delek
" autoindent when typing { and }
set smartindent
" number of spaces that a tab counts when typing \TAB> (sts)
set softtabstop=4
" inserted tabs turn into spaces (et)
set expandtab
" show information about the current position (activated by default, but just
set ruler
" show tabs and trailing spaces
set list
```

Data structures

```
segment_tree_beats_1.cpp, iterative_segment_tree.cpp, persistent_segment_tree.cpp,
iterative_lazy_segment_tree.cpp, iterative_segment_tree_2d.cpp, heavy_light_decomposition.cpp,
ordered_set.cpp, median.cpp, search_buckets.cpp, sparse_table.cpp, fenwick_tree.cpp,
implicit_lazy_treap.cpp, mos.cpp, monotonic_convex_hull_trick.cpp
```

segment_tree_beats_1.cpp

```
/* Level 1 of Segment tree beats. Can do range updates that assign to
 * every element in the range the minimum between its current value
 ^{\star} and a given value x
* A - (input) initial array
 * Complexity: O(logn) per query
                amortized O(logn) per update
struct node{
    11 maxi_count, maxi, second, sum;
    bool leaf, lazy;
} t[N*4];
int A[N];
void build(int node) {
    t[node].maxi = max(t[node<<1].maxi, t[node<<1|1].maxi);
    t[node].sum = t[node << 1].sum + t[node << 1|1].sum;
    if(t[node<<1].maxi == t[node<<1|1].maxi) {</pre>
        t[node].maxi_count = t[node<<1].maxi_count + t[node<<1|1].maxi_count;
        t[node].second = max(t[node<<1].second, t[node<<1|1].second);</pre>
    } else if(t[node<<1].maxi > t[node<<1|1].maxi) {</pre>
        t[node].maxi count = t[node<<1].maxi count;
        t[node].second = max(t[node<<1|1].maxi, t[node<<1].second);
    } else {
        t[node].maxi count = t[node<<1|1].maxi count;
        t[node].second = max(t[node<<1].maxi, t[node<<1|1].second);
}
void init(int 1, int r, int node=1) {
    t[node].lazy = 0;
    t[node].leaf = 0;
    if(1 == r) {
        t[node].maxi_count = 1;
        t[node].sum = t[node].maxi = A[1];
        t[node].second = -1;
        t[node].leaf = 1;
        return;
    int m = (1 + r) >> 1;
    init(1, m, node << 1);
init(m + 1, r, node << 1 | 1);</pre>
    build(node);
void putTag(int node, ll x){
    t[node].lazy = 1;
t[node].sum -= (t[node].maxi - x) * t[node].maxi_count;
    t[node].maxi = x;
void propagate(int node) {
    if(!t[node].leaf){
        if(t[node << 1].maxi > t[node].maxi)
            putTag(node << 1, t[node].maxi);</pre>
        if(t[node << 1 | 1].maxi > t[node].maxi)
            putTag(node <<1 | 1, t[node].maxi);</pre>
    t[node].lazy = 0;
// Queries maximum in closed range [ll, rr] \,
int queMax(int ll, int rr, int l=0, int r=n-1, int node=1) {
    if(ll <= l && r <= rr) return t[node].maxi;</pre>
    if(rr < 1 || r < 11) return -1;
    if(t[node].lazy) propagate(node);
```

```
int m = (1 + r) >> 1;
    return max(queMax(11, rr, 1, m, node << 1), queMax(11, rr, m+1, r, node << 1 | 1));
// Queries sum of closed range [11, rr] ll queSum(int ll, int rr, int l=0, int r=n-1, int node=1) {
    if(ll <= 1 && r <= rr) return t[node].sum;</pre>
    if(rr < 1 || r < 11) return 0;
    if(t[node].lazy) propagate(node);
    int m = (1 + r) >> 1;
    return queSum(ll, rr, l, m, node << 1) + queSum(ll, rr, m + 1, r, node << 1 | 1);
// Updates to the minimum between x and current value in the closed range [11, rr]
void upd(int ll, int rr, long long int x, int l=0, int r=n-1, int node=1) {
    if(rr < 1 || r < 11 || t[node].maxi <= x)
         return;
    if(ll \leq 1 && r \leq rr && t[node].second < x) {
         putTag(node, x);
         return;
    if(t[node].lazy) propagate(node);
int m = (1 + r) >> 1;
    upd(11, rr, x, 1, m, node << 1);
upd(11, rr, x, m+1, r, node << 1 | 1);
    build(node);
```

iterative segment_tree.cpp

```
/* Simple iterative segment tree, with point update and range queries.
\mbox{\ensuremath{^{\star}}} The operator needs to be commutative for this implementation.
* Time complexity: O(n) for building and O(log n) for updates and queries.
* Space complexity: O(n)
struct segTree {
    vector<ll> st;
   const ll NEUT = 0; // TODO define neutral element
    // combine two elements, doesn't need to be commutative
    inline ll combine(ll a, ll b) {
        return a + b; // TODO define merge operator
    // build the tree with vector v
    void build(vector<ll> &v) {
        for (int i = 0; i < n; i++) {
            st[n + i] = v[i];
        for (int i = n-1; i >= 1; i--) {
            st[i] = combine(st[i << 1], st[i << 1 | 1]);
    public:
    segTree() {}
    // initialize with neutral elements
    segTree(int n) {
        resize(n);
    // initialize with vector
    segTree(vector<ll> &v) : segTree(v.size()) {
        build(v);
    void resize(int s) {
        n = s;
        st.assign(2*s, NEUT);
    // add x to position i
    void update(int i, ll x) {
```

```
st[i += n] += x; // TODO change update operation
while (i > 1) {
    i >>= 1;
    st[i] = combine(st[i << 1], st[i << 1 | 1]);
}

// query from 1 to r, inclusive

ll query(int 1, int r) {
    ll resl = NEUT, resr = NEUT;
    for (1 += n, r += n+1; 1 < r; 1 >>= 1, r >>= 1) {
        if (1 & 1) resl = combine(resl, st[l++]);
        if (r & 1) resr = combine(st[--r], resr);
    }

    return combine(resl, resr);
}
```

persistent_segment_tree.cpp

```
/* Persistent segment tree. This example is for queries of sum in range
 * and updates of sum in position, but any query or update can be achieved
 * changing the NEUT value, and functions updNode and merge.
^{\star} The version 0 of the persistent segTree has implicitly an array of
* length n full of NEUT values.
* It's recommend to set n as the actual length of the array.
* int n; cin >> n;
* segTree::n = n;
 * Complexity: O(logn) memory and time per query/update
const int NEUT = 0;
struct segTree {
    vector<int> t = vector<int>(1, NEUT);
    vector<int> left = vector<int>(1, 0), right = vector<int>(1, 0);
    static int n;
    int newNode(int v, int l=0, int r=0) {
        t.pb(v), left.pb(l), right.pb(r);
        return sz(t) - 1;
    int merge(int a, int b) {
        return a + b;
    // Initializes a segTree with the values of the array A of length n
    int init(int* A, int L=0, int R=n) {
        if(L + 1 == R) return newNode(A[L]);
        int M = (L + R) >> 1;
        int l = init(A, L, M), r = init(A, M, R);
        return newNode(merge(t[l], t[r]), l , r);
    int updNode(int cur value, int upd value) {
        return cur_value + upd_value;
    // updates the position pos of version k with the value v
    int upd(int k, int pos, int v, int L=0, int R=n) {
   int nxt = newNode(t[k], left[k], right[k]);
        if(L + 1 == R) t[nxt] = updNode(t[nxt], v);
        else {
            int M = (L + R) >> 1;
            int temp;
            if(pos < M) temp = upd(left[nxt], pos, v, L, M), left[nxt] = temp;</pre>
            else temp = upd(right[nxt], pos, v, M, R), right[nxt] = temp;
            t[nxt] = merge(t[left[nxt]], t[right[nxt]]);
        return nxt;
    // query in the range [l, r) of version \boldsymbol{k}
    int que(int k, int l, int r, int L=0, int R=n) {
        if (r \le L \mid \mid R \le 1) return NEUT;
        if(l <= L && R <= r) return t[k];
        int M = (L + R) >> 1;
        return merge(que(left[k], 1, r, L, M), que(right[k], 1, r, M, R));
int segTree::n = N;
```

iterative_lazy_segment_tree.cpp

```
/* Iterative segment tree with lazy propagation. Supports range updates and queries.
* This example is for querying the maximum value in a range and updating with sum in
 * a range.
* Changes must be done in the struct node operators nad NEUT value. For more complicated
* lazy values, change the `apply` method in segTree too.
* Time complexity: O(n) for building and O(\log n) for updates and queries.
* Space complexity: O(n)
const ll NEUT = -INF;
struct node {
    node() : val(0) {} // initial value
    node(ll val) : val(val) {}
    // combine two nodes
    node operator+(const node& other) {
        return node(max(val, other.val));
    // update a node by the lazy value
    void operator+=(ll x) {
        val += x;
};
struct segTree {
    int n, h;
    vector<ll> d;
    vector<node> t;
    segTree(int n) : n(n), t(n << 1) {
    d.resize(n, 0);</pre>
        h = sizeof(int) * 8 - __builtin clz(n);
    void apply(int p, ll x) {
        t[p] += x;
        if(p < n) d[p] += x;
    void push(int p) {
        for(int s = h; s > 0; s--) {
   int i = p >> s;
            if(d[i]) {
                apply(i << 1, d[i]);
                apply(i << 1 | 1, d[i]);
                d[i] = 0;
            }
        }
    void build(int p) {
        for(; p >>= 1;) {
            t[p] = t[p << 1] + t[p << 1 | 1];
            t[p] += d[p];
    void update(int 1, int r, ll x){
        1 += n, r += n;
        int 10 = 1, r0 = r;
        push(1);
        push(r - 1);
for(; 1 < r; 1 >>= 1, r >>= 1) {
            if(l \& 1) apply(l++, x);
            if(r & 1) apply(--r, x);
        build(10);
        build(r0 - 1);
    node query(int 1, int r) {
       1 += n, r += n;
        push(1);
        push(r - 1);
        node ans (NEUT);
        for(; 1 < r; 1 >>= 1, r >>= 1){
            if(1 \& 1) ans = ans + t[1++];
            if(r & 1) ans = ans + t[--r];
        return ans;
    }
} ;
```

iterative segment tree 2d.cpp

```
/* 2d iterative segment tree, with point update and range queries.
 ^{\star} The operator needs to be commutative for this implementation.
* From: https://github.com/mhunicken/icpc-team-notebook-el-
vasito/blob/master/data structures/segment tree 2d.cpp
 * Time complexity: O(n*m) for building and O(\log n * \log m) for updates and queries.
 * Space complexity: O(n*m)
struct segTree {
    int n, m;
    vector<vector<ll>>> st;
    const ll NEUT = 0; // TODO define neutral element
    // combine two elements, needs to be commutative
inline ll combine(ll a, ll b) {
   return a + b; // TODO define merge operator
    // build the tree with matriz mat
    void build(vector<vector<ll>> &mat) {
        for (int i = 0; i < n; i++) for (int j = 0; j < m; j++)
             st[n + i][m + j] = mat[i][j];
        for (int i = 0; i < n; i++) for (int j = m-1; j >= 1; j--)
             st[n + i][j] = combine(st[n + i][j << 1], st[n + i][j << 1 | 1]);
         for (int i = n-1; i >= 1; i--) for (int j = 0; j < 2*m; j++)
             st[i][j] = combine(st[i << 1][j], st[i << 1 | 1][j]);
    public:
    segTree() {}
    // initialize with neutral elements
    segTree(int n, int m) {
        resize(n, m);
    // initialize with matrix
    segTree(vector<vector<11>> &m) : segTree(m.size(), m.front().size()) {
        build (m);
    void resize(int new_n, int new_m) {
        n = new n;
        m = new m;
        st.assign(2*n, vector<ll>(2*m, NEUT));
    // set position (x, y) to k
    void update(int x, int y, ll k) {
        st[n + x][m + y] = k; // TODO change update operation
         for (int j = m + y; j > 1; j >>= 1)
             st[n + x][j >> 1] = combine(st[n + x][j], st[n + x][j ^ 1]);
        for (int i = n + x; i > 1; i >>= 1) for (int j = m + y; j >= 1; j >>= 1) st[i >> 1][j] = combine(st[i][j], st[i ^1][j]);
    }
    // query in the rectangle (is, js) (ie, je), INCLUSIVE !!!
    ll query(int is, int js, int ie, int je) {
         11 \text{ res} = \text{NEUT};
         for (int i0 = n + is, i1 = n + ie + 1; i0 < i1; i0 >>= 1, i1 >>= 1) {
             11 t[2], q = 0;
             if (i0 & 1) t[q++] = i0++;
             if (i1 & 1) t[q++] = --i1;
             for (int k = 0; k < q; k++)
                 for (int j0 = m + js, j1 = m + je + 1; j0 < j1; j0 >>= 1, j1 >>= 1) {
   if (j0 & 1) res = combine(res, st[t[k]][j0++]);
                      if (j1 & 1) res = combine(res, st[t[k]][--j1]);
         return res;
```

};

heavy_light_decomposition.cpp

```
/st Data structure to answer queries on paths of a tree. It divides the tree in chains
  and for each chain saves a data structure that answers the query as if it were done
 ^{\star} in an array. In this template, this data structure is assumed to be a segment tree
* but that's not mandatory.
* Complexity: O(T(n) * logn) time per query/update, where T(n) is the time of
                query/update of the inherent data structure.
struct hld {
    int gid, r;
    vector<int> tam, id, p, d, rt;
    vector<vector<int> > adj;
    segTree st;
    hld(int n, int root=0)
       : r(root), gid(0), tam(n), id(n), p(n), d(n), rt(n), adj(n), st(n) {}
    void addEdge(int u, int v) {
       adj[u].pb(v);
        adj[v].pb(u);
    int prec(int v, int par=-1, int depth=0) {
        tam[v] = 1;
        p[v] = par;
        d[v] = depth;
        for(int u: adj[v]) if(u != par)
            tam[v] += prec(u, v, depth + 1);
        return tam[v];
    void build(int v, int root) {
        id[v] = gid++;
        rt[v] = root;
        if(sz(adj[v]) > 1 && adj[v][0] == p[v])
           swap(adj[v][0], adj[v][1]);
        for (auto &u: adj[v]) if (u != p[v] && tam[u] > tam[adj[v][0]])
            swap(adj[v][0], u);
        for(auto u: adj[v]) if(u != p[v])
            build(u, u == adj[v][0] ? root : u);
    void init() {
        prec(r);
        build(r, r);
    void updateVertex(int u, int x) {
        st.update(id[u], x);
    void updateEdge(int u, int v, int x) {
        st.update(max(id[u], id[v]), x);
    // for queries on the edges of the path set for_edge to true
    // this code assumes that the segment tree queries are right-exclusive
    int query(int u, int v, bool for edge=false){
    auto oper = (ll a, ll b) { // you probably will only need to change this
            return max(a, b);
        ll ans = -INF;
        while (rt[u] != rt[v]) {
            if(d[rt[u]] > d[rt[v]]) swap(u, v);
            ans = oper(ans, st.query(id[rt[v]], id[v] + 1));
            v = p[rt[v]];
        int a = id[u], b = id[v];
        ans = oper(ans, st.query(min(a, b) + for edge, max(a, b) + 1));
        return ans:
};
```

ordered_set.cpp

```
/* The ordered set data structure works just like the std::set, but it has additional
* information for every node. Namely, it can retrieve the order for a given key and also
* the key of a node which is in a given position.
*
* Methods:
```

```
* - find_by_order(pos)
*    returns the iterator of the node in the `pos` position.
* - order_of_key(key)
*    returns the order in which a node of key `key` would be in the ordered set
* The other methods are the same of set's methods.
*
* Time complexity: O(logn) for insert, erase and queries
* Space complexity: O(n)
*/

#include <ext/pb_ds/assoc_container.hpp>
#include <ext/pb_ds/tree_policy.hpp>
using namespace __gnu_pbds;

template<typename T>
using ordered_set = tree<T, null_type, less<T>, rb_tree_tag, tree_order_statistics_node_update>;
```

median.cpp

```
struct Median {
    multiset<11, greater<11>> sm;
multiset<11> gt;
    int size = 0;
    void balance() {
        if (sm.size() < (size + 1) / 2) {
             11 x = *(gt.begin());
             gt.erase(gt.begin());
             sm.insert(x);
         } else if (sm.size() > (size + 1) / 2) {
             11 x = *(sm.begin());
             sm.erase(sm.begin());
             gt.insert(x);
    void insert(ll v) {
        if (size == 0 || v <= median first()) sm.insert(v);</pre>
        else gt.insert(v);
        size++:
        balance();
    bool remove(ll v) {
   if (size == 0) return false;
         if (v <= median first()) {</pre>
             auto it = \overline{sm.find(v)};
             if (it == sm.end()) return false;
             sm.erase(it);
         } else {
             auto it = gt.find(v);
             if (it == gt.end()) return false;
             gt.erase(it);
         size--;
        balance();
        return true;
    11 median first() {
        if (\overline{\text{size}} \% 2 == 1) \text{ return } *(\text{sm.begin}());
        return *(sm.begin());
    double median() {
        if (size % 2 == 1) return *(sm.begin());
        return (*(sm.begin()) + *(gt.begin())) / 2.0;
    string median string() {
        if (size \frac{1}{8} 2 == 1) return to string(*(sm.begin()));
        11 a = *(sm.begin()) + *(gt.\overline{begin()});
        string s;
        if (a < 0) s += '-';
         s += to_string(abs(a) / 2);
        if (abs(a) % 2 == 1) s += ".5";
        return s;
};
```

search buckets.cpp

```
#include <algorithm>
#include <cassert>
#include <cmath>
#include <iostream>
#include <vector>
using namespace std;
// search_buckets provides two operations on an array:
// 1) set array[i] = x
// 2) count how many i in [start, end) satisfy array[i] < value</pre>
// Both operations take sqrt(N log N) time. Amazingly, because of the cache efficiency this is
faster than the
// (log N)^2 algorithm until N = 2-5 million.
template<typename T>
struct search buckets {
    // values are just the values in order. buckets are sorted in segments of BUCKET SIZE (last
segment may be smaller)
    int N, BUCKET_SIZE;
    vector<T> values, buckets;
    search buckets(const vector<T> &initial = {}) {
        init(initial);
    int get bucket end(int bucket start) const {
        return min(bucket_start + BUCKET_SIZE, N);
    void init(const vector<T> &initial) {
        values = buckets = initial;
        N = values.size();
        BUCKET SIZE = 3 * sqrt(N * log(N + 1)) + 1;
        cerr << "Bucket size: " << BUCKET SIZE << endl;
        for (int start = 0; start < N; start += BUCKET SIZE)</pre>
            sort(buckets.begin() + start, buckets.begin() + get_bucket_end(start));
    }
    int bucket_less_than(int bucket_start, T value) const {
        auto begin = buckets.begin() + bucket_start;
        auto end = buckets.begin() + get_bucket_end(bucket_start);
        return lower bound(begin, end, value) - begin;
    int less_than(int start, int end, T value) const {
        int \overline{\text{count}} = 0;
        int bucket start = start - start % BUCKET SIZE;
        int bucket end = min(get bucket end(bucket start), end);
        if (start - bucket_start < bucket_end - start) {</pre>
            while (start > bucket_start)
                count -= values[--start] < value;</pre>
            while (start < bucket end)
                count += values[start++] < value;</pre>
        if (start == end)
            return count;
        bucket start = end - end % BUCKET SIZE;
        bucket end = get bucket end(bucket start);
        if (end - bucket_start < bucket_end - end) {
    while (end > bucket_start)
                count += values[--end] < value;</pre>
        } else {
            while (end < bucket end)
                count -= values[end++] < value;</pre>
        while (start < end && get bucket end(start) <= end) {
            count += bucket less than(start, value);
            start = get_bucket_end(start);
        assert(start == end);
        return count;
    }
```

```
int prefix_less_than(int n, T value) const {
        return less_than(0, n, value);
    void modify(int index, T value) {
        int bucket_start = index - index % BUCKET_SIZE;
        int old_pos = bucket_start + bucket_less_than(bucket_start, values[index]);
        int new pos = bucket start + bucket less than(bucket start, value);
        if (old_pos < new_pos) {</pre>
            copy(buckets.begin() + old_pos + 1, buckets.begin() + new_pos, buckets.begin() +
old_pos);
            // memmove(&buckets[old pos], &buckets[old pos + 1], (new pos - old pos) * sizeof(T));
        } else {
            copy backward(buckets.begin() + new pos, buckets.begin() + old pos, buckets.begin() +
old pos + 1);
            // memmove(&buckets[new pos + 1], &buckets[new pos], (old pos - new pos) * sizeof(T));
        buckets[new_pos] = value;
        values[index] = value;
};
int main() {
   int N, M;
    scanf("%d %d", &N, &M);
    vector<int> A(N), B(N);
    vector<int> location(N + 1);
    for (int i = 0; i < N; i++) {
        scanf("%d", &A[i]);
        location[A[i]] = i;
    for (int &b : B) {
        scanf("%d", &b);
        b = location[b];
    search buckets<int> buckets(B);
    for (int i = 0; i < M; i++) {
        int type;
        scanf("%d", &type);
        if (type == 1) {
            int LA, RA, LB, RB;
            scanf("%d %d %d %d", &LA, &RA, &LB, &RB);
            LA--; LB--;
            printf("%d\n", buckets.less\_than(LB, RB, RA) - buckets.less\_than(LB, RB, LA));
        } else if (type == 2) {
            int x, y;
scanf("%d %d", &x, &y);
            x--; y--;
            buckets.modify(x, B[y]);
            buckets.modify(y, B[x]);
            swap(B[x], B[y]);
        } else {
            assert(false);
   }
}
```

sparse_table.cpp

```
/* Sparse table. Useful for queries of idempotent functions in a range.
  * Change the oper method accordingly.
  *
  * Time complexity:
  * - O(n logn) for building the structure
  * - O(1) for queries of idempotent functions.
  * Space complexity: O(n logn)
  */

struct sparseTable {
  int n, logn;
  vector<vector<ll>> t;
  int log_floor(int tam) {
    return 31 - __builtin_clz(tam);
}
```

```
11 oper(11 a, 11 b) { // example with minimum in range
        return min(a, b);
    sparseTable() {}
    sparseTable(vector<ll>& v) {
        n = sz(v);
        logn = log floor(n) + 1;
        t.resize(logn);
        t[0].resize(n);
        for (int i = 0; i < n; i++)
            t[0][i] = v[i];
        for (int k = 1; k < logn; k++) {
            t[k].resize(n);
            for(int i = 0; i + (1 << k) <= n; i++)
                t[k][i] = oper(t[k-1][i], t[k-1][i+(1 << (k-1))]);
    Il que(int 1, int r) { // queries in the semi-open interval [1, r) int k = log_floor(r - 1);
        return oper(t[k][1], t[k][r - (1 << k)]);
};
```

fenwick_tree.cpp

```
/\star Fenwick tree, with point update and range queries.
* The operator needs to be commutative for this implementation.
* Time complexity: O(n \log(n)) for building and O(\log n) for updates and queries.
* Space complexity: O(n)
struct Fenwick {
   vector<ll> bit;
   int n;
    void build(vector<ll> &v) {
        for (size_t i = 0; i < v.size(); i++) {
            update(i, v[i]);
    Fenwick() {}
    Fenwick(int n) : n(n) {
        resize(n);
    Fenwick(vector<ll> &v) : Fenwick(v.size()) {
       build(v);
    void resize(int n) {
        bit.assign(n, 0);
    11 query(int r) {
        11 \text{ ret} = 0;
        for (; r \ge 0; r = (r \& (r + 1)) - 1) {
           ret += bit[r];
        return ret;
    }
    11 query(int 1, int r) {
        return query(r) - query(1-1);
    void update(int i, ll v) {
       for (; i < n; i = i | (i + 1)) {
            bit[i] += v;
};
```

implicit_lazy_treap.cpp

```
/* Implicit treap with lazy propagation. It is easy to adapt to versions without
 * lazy propagation or that are not implicit.
 * For non-lazy versions, just delete the functions apply and push.
 * For non-implicit versions, use an actual key in the split instead of the implicit key.
 ^{\star} This example uses lazy propagation for reversing a range in implicit treaps.
 * Time complexity: O(log n) for merge and split.
 * Space complexity: O(n)
mt19937 rng(chrono::steady_clock::now().time_since_epoch().count());
typedef struct item* pitem;
struct item {
    int cnt, pr, val;
    bool lazy;
    pitem 1, r;
    item(int val)
        : pr(rng()), val(val), cnt(1), lazy(0), l(nullptr), r(nullptr) {}
};
int cnt(pitem t) {
    return t != nullptr ? t->cnt : 0;
void upd(pitem t) {
    if(t != nullptr)
        t->cnt = cnt(t->1) + 1 + cnt(t->r);
void apply(pitem t, bool d) {
    if(t != nullptr) {
        if (d.first) swap(t->1, t->r);
        t\rightarrow lazy ^= d;
}
void push(pitem t) {
    if(t != nullptr) {
        apply(t->1, t->lazy);
apply(t->r, t->lazy);
t->lazy = {0, 0};
    }
}
void merge(pitem& t, pitem l, pitem r) {
   if(l == nullptr || r == nullptr)
        t = 1 != nullptr ? 1 : r;
    else if(l->pr > r->pr) {
        push(1);
        merge(1->r, 1->r, r), t = 1;
    } else {
        push(r);
        merge(r->1, 1, r->1), t = r;
    upd(t);
}
void split(pitem t, pitem& l, pitem& r, int key, int add=0) {
    if(t == nullptr) {
        1 = r = nullptr; return;
    push(t);
    int cur_key = add + cnt(t->1);
if(key <= cur key)</pre>
        split(t->1, 1, t->1, key, add), r = t;
    else
         split(t->r, t->r, r, key, add + 1 + cnt(t->1)), 1 = t;
    upd(t);
void insert(pitem& t, int pos) {
    pitem it = new item(pos);
    pitem t1, t2;
    split(t, t1, t2, pos);
    merge(t1, t1, it);
merge(t, t1, t2);
}
```

mos.cpp

```
/\!\!\!\!\!\!^\star Optimization for problems with queries that can be answered offline
 * Complexity: O(q * (BLOCK SIZE) + n * (n / BLOCK SIZE))
               O((q + n) * sqrt(n)) if BLOCK SIZE ~ sqrt(n)
const int BLOCK SIZE = 700;
struct Query {
    int 1, r, idx;
    bool operator<(Query other) const
        return make pair(1 / BLOCK SIZE, r) <
               make_pair(other.l / BLOCK SIZE, other.r);
};
struct Mos {
    // variables specific for the problem
    11 sum;
    vector<11> v;
    // stores the answers to the queries in the original order
    void exec(vector<Query> &queries, vector<ll> &answers) {
        answers.resize(queries.size());
        sort(queries.begin(), queries.end());
        int cur l = 0;
        int cur_r = -1;
        for (Query q : queries) {
    while (cur_l > q.l) {
                cur 1--;
                add(cur_l);
            while (cur r < q.r) {
                cur r++;
                add(cur r);
            while (cur_l < q.l) {
                remove(cur_1);
                 cur 1++;
            while (cur_r > q.r) {
                remove(cur_r);
                 cur_r--;
            answers[q.idx] = get answer(cur 1, cur r);
        }
    // functions below are specific for the problem
    Mos(vector<11> &v) : sum(0), v(v) {}
    void add(int i) {
        sum += v[i];
    void remove(int i) {
        sum -= v[i];
    11 get answer(int 1, int r) {
       return sum;
};
```

monotonic_convex_hull_trick.cpp

```
/* Convex hull trick version for monotonic slopes and queries, this is,
  * the inserted lines have only increasing/decreasing slopes and the
  * queries are done only in increasing/decreasing x-coordinates.
  *
  * This template assumes non-decreasing slopes in inserted lines
  * and queries of maximization in non-decreasing x-coordinates.
  *
  * Complexity: amortized O(1) for each insertion and query
  */
typedef long double ld;
```

Dynamic programming

longest_increasing_subsequence.cpp, knuth.cpp, divide_and_conquer.cpp

longest_increasing_subsequence.cpp

```
/* Computes the longest (weakly) increasing subsequence in a vector.
 * Is easy to change the code below for the strictly increasing version of the problem.
 * v - (input) vector for which the longest increasing subsequence will be computed.
 * aux - one of the longest increasing subsequences of v
 *
 * Complexity: O(n logn) time
 * O(n) memory
 */

vector<int> lis(vector<int>& v) {
    vector<int> aux;
    for(auto x: v) {
        // Change to lower_bound for strictly increasing version
        auto it = upper_bound(aux.begin(), aux.end(), x);
        if(it == aux.end()) aux.pb(x);
        else *it = x;
    }
    return aux;
}
```

knuth.cpp

```
/\,^\star A dp of the form
      dp[l][r] = min_{1 < m < r}(dp[l][m] + dp[m][r]) + cost(l, r)
* can be solved in O(n^2) with Knuth opmitization if we have that
       opt[1][r - 1] \le opt[1][r] \le opt[1 + 1][r]
  where
       dp[1][r] = dp[1][opt[1][r]] + dp[opt[1][r]][r] + cost(1, r).
* Other sufficient condition (that implies the previous one) is
     given a \leq b \leq c \leq d, we have:
         - quadrangle inequality: cost(a, c) + cost(b, d) \le cost(a, d) + cost(b, c)
          - monotonicity: cost(b, c) <= cost(a, d)
* Complexity: O(n^2) time
               O(n^2) memory
   vector<vector<ll>>> dp = vector<vector<ll>>>(n, vector<ll>(n));
   vector<vector<int>> opt = vector<vector<int>> (n, vector<int>(n));
    for (int k = 0; k \le n; k++) {
        for(int l = 0; l + k \le n; l++) {
            int r = 1 + k;
            if(k < 2) {
                dp[1][r] = 0; // base case
                opt[1][r] = 1;
            dp[1][r] = INF;
            for (int m = opt[1][r - 1]; m \le opt[1 + 1][r]; m++) {
                ll cur = dp[1][m] + dp[m][r] + cost(l, r); // must define O(1) cost function
                if(cur < dp[l][r]) {
                    dp[l][r] = cur;
                    opt[l][r] = m;
                }
            }
        }
    return dp[0][n];
}
```

divide_and_conquer.cpp

```
/* A dp of the form

* dp[i][j] = min_{k < j}(dp[i - 1][k] + cost(k, j))
```

```
\ast can be solved in O(m n logn) with divide and conquer optimization if we have that
       opt[i][j] <= opt[i][j + 1]
       dp[i][j] = dp[i - 1][opt[i][j]] + cost(opt[i][j], j).
 * Complexity: O(m \ n \ logn) time (for a partition in m subarrays of an array of size n)
                O(n) memory
ll dp[N][2];
void go(int k, int l, int r, int optl, int optr) {
    if(1 > r) return;
    int opt, m = (1 + r) >> 1;
dp[m][k&1] = INF;
    for(int i = optl; i <= min(m, optr + 1); i++) {
    ll cur = dp[i][~k&1] + cost(i, m); // must define O(1) cost function</pre>
         if(cur < dp[m][k&1]) {
             dp[m][k&1] = cur;
             opt = i;
        }
    go(k, l, m - 1, optl, opt);
    go(k, m + 1, r, opt, optr);
}
11 dc(int n, int m) {
   dp[0][0] = dp[0][1] = 0;
    for(int i = 0; i < n; i++) dp[i][0] = INF;
for(int i = 1; i <= m; i++) go(i, i, n, i - 1, n - 1);
    return dp[n - 1][m & 1];
}
```

Flows and matchings

bipartite_matching.cpp, dinic.cpp, hungarian.cpp, stable_matching.cpp, flow_with_demands.cpp,
min cost max flow dijkstra.cpp, min cost max flow spfa.cpp

bipartite_matching.cpp

```
/* Solves the bipartite matching problem for color classes X and Y.
 ^{\star} n - (input) size of color class {\rm X}
 * m - (input) size of color class Y
 * g - (input) g[x] contains the neighbours of x \in X
 * mat - mat[y] is the vertex matched with y \in Y or -1 if there's no such vertex * arcs - set of arcs of the bipartite matching. Each arc is a pair such that the first
           element is in X and the second in Y.
 * Complexity: O(V * (V + E)) where V is the number of vertex and E the number of edges
                                  in the bipartite graph
 * /
vector<int> q[N];
int mat[N];
bool vis[N];
int n, m;
int match(int x) {
    if(vis[x]) return 0;
    vis[x] = true;
    for (int y: g[x]) if (mat[y] < 0 || match (mat[y])) {
        mat[y] = x;
    return 0;
vector<pair<int, int> > maxMatching() {
    vector<pair<int, int> > arcs;
    for (int i = 0; i < m; i++)
        mat[i] = -1;
    for(int i = 0; i < n; i++) {
        for(int j = 0; j < n; j++)
            vis[j] = 0;
         match(i);
    for (int i = 0; i < m; i++)
        if(mat[i] >= 0) arcs.pb({mat[i], i});
    return arcs;
}
```

dinic.cpp

```
/* Fast max flow algorithm.
* Constructor:
 * dinic(n, s, t)
* n - number of nodes in the flow network.
* s - source of the flow network.
^{\star} t - sink of the flow network.
* Methods:
  - addEdge(u, v, cap)
    adds a directed edge from `u` to `v` with capacity `cap`.
* - getFlow()
     returns the maximum flow of the network.
* Complexity: In general, the time complexity of getFlow is O(E\ V^2), but there are better
                upper bounds for bipartite graphs (O(E sqrt(V))) and networks with unit
                capacities (O(E sqrt(E))).
const int INF = 1 << 29;
struct dinic {
    ll n, s, t;
    vector<ll> dist, q, work;
```

```
struct edge {
        11 to, rev, f, cap;
    vector<vector<edge> > g;
    \label{eq:dinic(int n, int s, int t) : n(n), s(s), t(t), g(n), dist(n), q(n), work(n) {} \{\}
    void addEdge(int u, int v, ll cap) {
        g[u].pb((edge){v, sz(g[v]), 0, cap});
        g[v].pb((edge)\{u, sz(g[u]) - 1, 0, 0\});
    bool bfs() {
        for(int i = 0; i < n; i++) dist[i] = -1;
        dist[s] = 0;
        int qt = 0;
        q[qt++] = s;
        for(int qh = 0; qh < qt; qh++) {
            int u = q[qh];
            for(int i = 0; i < sz(g[u]); i++) {
                edge &e = g[u][i];
                int v = g[u][i].to;
                if(dist[v] < 0 && e.f < e.cap) {
                     dist[v] = dist[u] + 1;
                     q[qt++]=v;
            }
        return dist[t] >= 0;
    ll dfs(int u, ll f) {
        if(u == t) return f;
        for(ll &i = work[u]; i < sz(g[u]); i++) {
            edge &e = g[u][i];
            if(e.cap <= e.f) continue;
            int v = e.to;
            if(dist[v] == dist[u] + 1) {
                11 df = dfs(v, min(f, e.cap - e.f));
                 if(df > 0){
                     e.f += df;
                     q[v][e.rev].f -= df;
                     return df;
                 }
            }
        return 0;
    11 getFlow() {
        11 \text{ res} = 0;
        while(bfs()) {
            for (int i = 0; i < n; i++) work [i] = 0;
            while(ll delta = dfs(s, INF))
                res += delta;
        return res;
};
```

hungarian.cpp

```
/* Hungarian method algorithm that solves the bipartite perfect matching of minimum cost.
 \ensuremath{^{\star}} Can solve the problem of maximum cost with minor changes.
* Constructor:
* hungarian(n, m)
* n - (input) size of color class X
* m - (input) size of color class Y
* Methods:
 * - set(x, y, c)
    sets the cost c for the edge between x \in X to y \in Y
 * - assign()
     returns the cost of an optimal matching and fills the vectors matchl and matchr
     with the assignment of such matching
 * Complexity: O(V^3) where V is the number of vertex of the graph
typedef long double ld;
const ld INF = 1e100; // for maximization set INF to 0 and negate costs
bool zero(ld x) {
    return fabs(x) < 1e-9; // change to x == 0 for integer types
struct hungarian {
```

```
int n;
    vector<vector<ld> > cs;
    vector<int> matchl, matchr;
    hungarian(int _n, int _m) : n(max(_n, _m)), cs(n, vector<ld>(n)), match1(n), matchr(n) {
    for(int x = 0; x < _n; x++)
        for(int y = 0; y < _m; y++)
                cs[x][y] = INF;
    void set(int x, int y, ld c) {
        cs[x][y] = c;
    ld assign() {
       int mat = 0;
        vector < ld > ds(n), y(n), z(n);
        vector < int > dad(n), vis(n);
        for(int i = 0; i < n; i++) {
            matchl[i] = matchr[i] = -1;
            y[i] = *min element(cs[i].begin(), cs[i].end());
        for(int j = 0; j < n; j++) {
            z[j] = cs[0][j] - y[0];
             for (int i = 1; i < n; i++)
                z[j] = min(z[j], cs[i][j] - y[i]);
        for(int i = 0; i < n; i++)
            matchl[i] = j;
                     matchr[j] = i;
                     mat++;
                     break;
        for(;mat < n; mat++) {</pre>
            int s = 0, j, i;
            while (matchl[s] != -1) s++;
             for (int i = 0; i < n; i++) {
                dad[i] = -1;
                 vis[i] = 0;
            for (int k = 0; k < n; k++)
                ds[k] = cs[s][k] - y[s] - z[k];
            while(1) {
                 for (int k = 0; k < n; k++)
                     if(!vis[k] \&\& (j == -1 || ds[k] < ds[j]))
                        j = k;
                 vis[j] = 1;
                 i = matchr[j];
                 if(i == -1)
                     break;
                 for (int k = 0; k < n; k++) if (!vis[k]) {
                     auto new_ds = ds[j] + cs[i][k] - y[i] - z[k];
if(ds[k] > new_ds) {
   ds[k] = new_ds;
                         dad[k] = j;
                     }
                 }
            for (int k = 0; k < n; k++) if (k != j \&\& vis[k]) {
                auto w = ds[k] - ds[j];
                 z[k] += w;
                y[matchr[k]] -= w;
            y[s] += ds[j];
            while (dad[j] != -1) {
                matchl[matchr[j] = matchr[dad[j]]] = j;
                j = dad[j];
            matchr[j] = s;
            matchl[s] = j;
        1d value = 0;
        for(int i = 0; i < n; i++)
            value += cs[i][matchl[i]];
        return value;
};
```

stable_matching.cpp

/* Solves the rural hospital version of the stable matching problem.

```
* This is, you have a set of doctors and hospitals and you want to
 * assign at most cap[i] doctors to the i-th hospital and each doctor
 * to at most one hospital. Also, each doctor has a list of candidate
 * hospitals, sorted in decreasing order of preference, and each hospital
 * has a list of candidate doctors, sorted in decreasing order of
 * preference too. Find an assigment of doctors to hospitals such that
  there is no doctor D assigned to hospital H in which *both* D and H
 * would rather be matched with other candidates (consider the 'empty'
  assignment as the least prefered one).
 * Remark: any stable matching has the same set of doctors assigned to
 * any hospital, and the same capactity used in every hospital.
^{\star} For the less constrained version of the problem, just set all the
* capacities to 1.
\star n - (input) number of doctors.
* m - (input) number of hospitals.
 * cap - (input) array of hospitals' capacities. In the end of the algorithm,
        it will have the remaining capacity of each hospital.
 * doc - (input) list of hospital candidates of each doctor sorted by preference.
 ^{\star} hos - (input) list of doctor candidates of each hospital sorted by preference.
* assign - assign[i] is the hospital assigned to doctor i or -1 if it hasn't any
           hospital assigned.
 * Complexity: O(L), where L is the total size of the lists of candidates.
int n, m;
vector<int> doc[N], hos[N];
int pos[N], cap[N], last[N];
unordered map<int, int> prior[N];
vector<int> galeShapley() {
    for (int i = 0; i < m; i++) {
        last[i] = sz(hos[i]) - 1;
        for(int j = 0; j < sz(hos[i]); j++)
            prior[i][hos[i][j]] = j;
    vector<int> assign(n, -1);
    queue<int> Q;
    auto tryPush = [&] (int i) {
        if(pos[i] < sz(doc[i]))
            Q.push(i);
    for (int i = 0; i < n; i++)
        tryPush(i);
    while(!Q.empty()) {
        int u = Q.front();
        Q.pop();
        int cur = doc[u][pos[u]++];
        if(cap[cur]) {
            cap[cur]--;
            assign[u] = cur;
        } else {
            if(prior[cur][u] <= last[cur]) {</pre>
                assign[u] = cur;
                while(assign[hos[cur][last[cur]]] != cur)
                    last[cur]--;
                assign[hos[cur][last[cur]]] = -1;
                tryPush(hos[cur][last[cur]]);
            } else tryPush(u);
        }
    return assign;
}
```

flow_with_demands.cpp

```
/* Finds *any* feasible solution (if exists) to a flow network with demands (also known as
* lower bounds). It's possible to find the minimum flow solution doing binary search with
* a minor change.
* It only does a graph transformation and applies a max flow algorithm on it, thus this
* template depends of another flow algorithm template. In this example, that algorithm
* is minimum-cost maximum-flow, but that's not always necessary. Adapting this template
* to other flow algorithms should be easy.

* * Constructor:
* demands(n, s, t)
* n - number of nodes in the original flow network.
* s - source of the original flow network.
* t - sink of the original flow network.
```

```
* Methods:
 * - addEdge(u, v, cap, dem, cost)
    adds an edge to the original flow network from u to v with capacity `cap`, flow
    demand `dem` and cost `cost`. For problems without cost, the last parameter should
    be erased.
 * - getFlow()
    finishes building the auxiliary graph and returns the result of applying the max flow algorithm on it. If and only if the maximum flow is exactly equal to the sum of all
     demands, then exists a feasible solution to this problem.
 \star Complexity: depends on the flow algorithm used, but take into account that this method
                adds O(V) edges to the network flow.
struct demands {
    vector<ll> din, dout;
    mcf g;
    int n, s, t;
    11 \text{ base} = 0;
    demands(int n, int _s, int _t) : n(n), s(n), t(n + 1), din(n, 0), dout(n, 0) { g = mcf(n + 2, s, t);
        g.addEdge(_t, _s, INF, 0); // for minimum solution, change INF according to the value
                                       // of the binary search
    void addEdge(int u, int v, ll cap, ll dem, ll cost) {
        din[v] += dem;
        dout[u] += dem;
base += dem * cost;
        g.addEdge(u, v, cap - dem, cost);
    pair<ll, ll> getFlow() {
        for(int i = 0; i < n; i++) {
             if(din[i] > 0) g.addEdge(s, i, din[i], 0);
             if(dout[i] > 0) g.addEdge(i, t, dout[i], 0);
        auto ans = g.getFlow();
        return {ans.first, base + ans.second};
};
```

min_cost_max_flow_dijkstra.cpp

```
/* Solves the minimum-cost maximum-flow problem using dijkstra for finding the incremental
* shortest paths.
* Constructor:
* mcf(n, s, t)
* n - number of nodes in the flow graph.
* s - source of the flow graph.
* t - sink of the flow graph.
* Methods:
  - addEdge(u, v, cap, cost)
    adds a directed edge from u to v with capacity `cap` and cost `cost`.
* - getFlow()
    returns a pair of integers in which the first value is the maximum flow and the
    second is the minimum cost to achieve this flow.
* Complexity: There are two upper bounds to the time complexity of getFlow
                - O(max_flow * (E log V))
- O(V * E * (E log V))
const ll INF = 1LL << 60;
struct mcf {
   int n, s, t;
   ll cost, fl;
   vector<int> first, prev;
   vector<ll> dist;
   struct edge {
        int to, next;
        edge(int to, ll cap, ll cost, int next)
            : to(_to), cap(_cap), cost(_cost), next(_next) {};
   } ;
   vector<edge> g;
   mcf() {}
   mcf(int n, int s, int t) : n(n), s(s), t(t), fl(0), cost(0) {
        first.resize(n, -1);
```

```
dist.resize(n);
        prev.resize(n);
         g.reserve(n*n);
    void addEdge(int u, int v, ll cap, ll cost) {
         g.pb(edge(v, cap, cost, first[u]));
         first[u] = sz(g) - 1;
         g.pb(edge(u, 0, -cost, first[v]));
         first[v] = sz(q) - 1;
    bool augment() {
        dist.assign(n, INF);
         dist[s] = 0;
        priority_queue<pair<11, int> > q;
q.push({0, s});
         while(!q.empty()) {
             if(dist[t] < INF) break;</pre>
             11 d, u;
             tie(d, u) = q.top();
d *= -1;
             q.pop();
             if(dist[u] < d) continue;</pre>
             for (int e = first[u]; e != -1; e = g[e].next) {
                 int v = g[e].to;
ll ndist = d + g[e].cost;
if(g[e].cap > 0 && ndist < dist[v]) {</pre>
                      dist[v] = ndist;
                      q.push({-ndist, v});
                      prev[v] = e;
                  }
             }
         return dist[t] < INF;</pre>
    pair<11, 11> getFlow() {
         while(augment()) {
             11 cur = t, curf = INF;
             while(cur != s) {
                 int e = prev[cur];
                  curf = min(curf, g[e].cap);
                 cur = g[e^1].to;
             fl += curf;
             cost += dist[t] * curf;
             cur = t;
             while(cur != s) {
                  int e = prev[cur];
                  g[e].cap -= curf;
                 g[e^1].cap += curf;
                  cur = g[e^1].to;
        return {fl, cost};
    }
};
```

min_cost_max_flow_spfa.cpp

```
/* Solves the minimum-cost maximum-flow problem using spfa for the finding the incremental
^{\star} shortest paths. Useful when the edges costs are negative.
* Constructor:
* mcf(n, s, t)
\mbox{\ensuremath{\star}} n - number of nodes in the flow graph.
* s - source of the flow graph.
* t - sink of the flow graph.
 * Methods:
* - addEdge(u, v, cap, cost)
    adds a directed edge from u to v with capacity `cap` and cost `cost`.
 * - getFlow()
     returns a pair of integers in which the first value is the maximum flow and the
    second is the minimum cost to achieve this flow.
* Complexity: There are two upper bounds to the time complexity of getFlow
                - O(max_flow * (E log V))
                 - O(V * E * (E log V))
 * /
const ll INF = 1LL << 60;
struct mcf {
```

```
int n, s, t;
    ll cost, fl;
    vector<int> first, prev;
    vector<ll> dist;
    vector<bool> queued;
    struct edge {
        int to, next;
        11 cap, cost;
        };
    vector<edge> g;
    mcf() {}
    mcf(int _n, int _s,int _t) : n(_n), s(_s), t(_t), fl(0), cost(0) {
    queued.resize(n, 0);
        first.resize(n, -1);
        dist.resize(n);
        prev.resize(n);
        g.reserve(n*n);
    }:
    void addEdge(int u, int v, ll cap, ll cost) {
        g.pb(edge(v, cap, cost, first[u]));
        first[u] = sz(g) - 1;
g.pb(edge(u, 0, -cost, first[v]));
first[v] = sz(g) - 1;
    bool augment() {
        dist.assign(n, INF);
        dist[s] = 0;
        queued[s] = 1;
        queue<int> q;
        q.push(s);
        while(!q.empty()) {
            int u = q.front();
            q.pop();
            queued[u] = 0;
            for(int e = first[u]; e != -1; e = g[e].next) {
                int v = g[e].to;
ll ndist = dist[u] + g[e].cost;
                 if(g[e].cap > 0 \&\& ndist < dist[v]) {
                     dist[v] = ndist;
                     prev[v] = e;
                     if(!queued[v]) {
                         q.push(v);
                         queued[v] = 1;
                }
        return dist[t] < INF;</pre>
    pair<ll, ll> getFlow() {
        while(augment()) {
            11 cur = t, curf = INF;
            while(cur != s) {
                int e = prev[cur];
                curf = min(curf, g[e].cap);
                cur = g[e^1].to;
            fl += curf;
            cost += dist[t] * curf;
            cur = t;
            while(cur != s) {
                int e = prev[cur];
                g[e].cap -= curf;
                g[e^1].cap += curf;
                cur = g[e^1].to;
            }
        return {fl, cost};
};
```

Geometry

halfplane_intersection.cpp, convex_hull.cpp, point_integer.cpp, delaunay.cpp, line_integer.cpp, angular_sweep.cpp, polygon_double.cpp, line_double.cpp, circle.cpp, polygon_integer.cpp, shamos hoey.cpp, point double.cpp

halfplane_intersection.cpp

```
/* Half-plane intersection algorithm. The result of intersecting half-planes is either
 * empty or a convex polygon (maybe degenerated). This template depends on point double.cpp
  and line_double.cpp.
* h - (input) set of half-planes to be intersected. Each half-plane is described as a pair
* of points such that the half-plane is at the left of them.
  pol - the intersection of the half-planes as a vector of points. If not empty, these
  points describe the vertices of the resulting polygon in clock-wise order.
 ^{\star} WARNING: Some points of the polygon might be repeated. This may be undesirable in some
* cases but it's useful to distinguish between empty intersections and degenerated
 * polygons (such as a point, line, segment or half-line).
 * Time complexity: O(n logn)
struct halfplane: public line {
    ld ang;
    halfplane() {}
    halfplane(point _p, point _q) {
   p = _p; q = _q;
   point vec(q - p);
        ang = atan2(vec.y, vec.x);
    bool operator <(const halfplane& other) const {</pre>
        if (fabsl(ang - other.ang) < EPS) return right(p, q, other.p);</pre>
            return ang < other.ang;
    bool operator == (const halfplane& other) const {
        return fabsl(ang - other.ang) < EPS;
    bool out(point r) {
        return right(p, q, r);
};
vector<point> hp_intersect(vector<halfplane> h) {
    point box[4] = {\{-INF, -INF\}, {INF, -INF\}, {INF, INF\}}; for(int i = 0; i < 4; i++)
        h.pb(halfplane(box[i], box[(i+1) % 4]));
    sort(h.begin(), h.end());
    h.resize(unique(h.begin(), h.end()) - h.begin());
    deque<halfplane> dq;
    for(auto hp: h) {
        while (sz(dq) > 1 \&\& hp.out(intersect(dq.back(), dq[sz(dq) - 2])))
            dq.pop back();
        while (sz(dq) > 1 \&\& hp.out(intersect(dq[0], dq[1])))
            dq.pop front();
        dq.pb(hp);
    while (sz(dq) > 2 \& dq[0].out(intersect(dq.back(), dq[sz(dq) - 2])))
        dq.pop_back();
    while (sz(dq) > 2 \&\& dq.back().out(intersect(dq[0], dq[1])))
        dq.pop_front();
    if (sz(dq) < 3) return {};
    vector<point> pol(sz(dq));
for(int i = 0; i < sz(dq); i++) {</pre>
        pol[i] = intersect(dq[i], dq[(i+1) % sz(dq)]);
    return pol;
}
```

convex_hull.cpp

```
/* Finds the convex hull of a given set of points. This templates requires
* the struct point defined in point_integer.cpp or in point_double.cpp
*
```

```
* p - (input) vector of points for which the convex hull will be found. * ch - convex hull of \hat{p} in counter-clockwise order.
vector<point> convexHull(vector<point> p) {
    int n = sz(p);
    sort(p.begin(), p.end());
    vector<point> low, up;
    for (int i = 0; i < n; i++) {
        if(i \&\& p[i] == p[i - 1]) continue;
        while (sz(up) \ge 2 \&\& !right(up[sz(up)-2], up.back(), p[i]))
             up.pop_back();
        up.pb(p[i]);
        while (sz(low) \ge 2 \&\& !left(low[sz(low)-2], low.back(), p[i]))
            low.pop back();
        low.pb(p[i]);
    vector<point> ch;
    if(sz(low) == 1) return low;
    for (int i = 0; i < sz(low) - 1; i++)
        ch.pb(low[i]);
    for(int i = sz(up) - 1; i >= 1; i--)
        ch.pb(up[i]);
    return ch;
}
```

point integer.cpp

```
/* Basic structure of point and operations related with it. This template assumes
 * integer coordinates.
 * All operations' time complexity are O(1)
struct point {
    11 x, y;
    point(11 x, 11 y) : x(x), y(y) {}
    point() {}
    11 norm2()
       return *this * *this;
    bool operator==(const point& other) const {
       return x == other.x && y == other.y;
    point operator+(const point& other) const {
       return point(x + other.x, y + other.y);
    point operator-(const point& other) const {
       return point(x - other.x, y - other.y);
    point operator*(ll t) const {
       return point(x * t, y * t);
    point operator/(ll t) const {
       return point(x / t, y / t);
    11 operator*(const point& other) const {
        return x*other.x + y*other.y;
    11 operator^(const point& other) const { // cross product
       return x*other.y - y*other.x;
    bool operator<(const point& other) const { // for sweep line</pre>
       return x < other.x || (x == other.x && y < other.y);
    point rotate(point r) {
       return point(*this ^ r, *this * r);
};
point ccw90(1, 0);
point cw90(-1, 0);
11 dist2(point p, point q) { // squared distance
    return (p - q).norm2();
ll area2(point a, point b, point c) { // two times signed area of triangle abc
    return (b - a) ^ (c - a);
bool left(point a, point b, point c) {
    return area2(a, b, c) > 0;
```

```
bool right(point a, point b, point c) {
    return area2(a, b, c) < 0;
}

bool collinear(point a, point b, point c) {
    return abs(area2(a, b, c)) == 0;
}

// Returns 0 if vectors a and b are not parallel.

// If they are parallel, returns 1 if they have the same direction

// and returns -1 otherwise
int paral(point a, point b) {
    if((a ^ b) != 0) return 0;
    if((a.x > 0) == (b.x > 0) && (a.y > 0) == (b.y > 0))
        return 1;
    return -1;
}
```

delaunay.cpp

```
/* Builds a delaunay triangulation of a given set of points. This templates requires
 * the struct point defined in point integer.cpp
\mbox{\ensuremath{\star}} p - (input) vector of points for which the convex hull will be found.
* adj - the delaunay triangulation of `p` as an adjacency list in which every vertex
          has its original index in `p
struct quadEdge {
    point o;
    quadEdge *rot, *nxt;
    bool used;
    quadEdge(point o = point(INF, INF))
         : o(o), rot(nullptr), nxt(nullptr), used(false) {}
    quadEdge* rev() const {
         return rot->rot;
    quadEdge* lnext() const {
         return rot->rev()->nxt->rot;
    quadEdge* prev() const {
         return rot->nxt->rot;
    point dest() const {
         return rev()->o;
};
quadEdge* makeEdge(point from, point to) {
    vector<quadEdge*> e(4);
    e[0] = new quadEdge(from);
    e[1] = new quadEdge(to);
    e[2] = new quadEdge; e[3] = new quadEdge;
    \texttt{tie}(\texttt{e}[\texttt{0}] \texttt{-} \texttt{rot}, \ \texttt{e}[\texttt{1}] \texttt{-} \texttt{rot}, \ \texttt{e}[\texttt{2}] \texttt{-} \texttt{rot}, \ \texttt{e}[\texttt{3}] \texttt{-} \texttt{rot}) \ = \ \{\texttt{e}[\texttt{2}], \ \texttt{e}[\texttt{3}], \ \texttt{e}[\texttt{1}], \ \texttt{e}[\texttt{0}]\};
    tie(e[0]->nxt, e[1]->nxt, e[2]->nxt, e[3]->nxt) = {e[0], e[1], e[3], e[2]};
    return e[0];
}
void splice(quadEdge* a, quadEdge* b) {
    swap (a->nxt->rot->nxt, b->nxt->rot->nxt);
    swap(a->nxt, b->nxt);
}
void deleteEdge(quadEdge* &e, quadEdge* ne) {
    splice(e, e->prev());
splice(e->rev(), e->rev()->prev());
    delete e->rev()->rot;
    delete e->rev();
    delete e->rot;
    delete e;
    e = ne;
}
quadEdge* connect(quadEdge* a, quadEdge* b) {
    quadEdge* e = makeEdge(a->dest(), b->o);
splice(e, a->lnext());
    splice(e->rev(), b);
    return e;
}
```

```
__int128 det3(point a, point b, point c) {
    vector<__int128> len = {a.norm2(), b.norm2(), c.norm2()};
return a.x * (b.y * len[2] - c.y * len[1])
             -a.y * (b.x * len[2] - c.x * len[1])
             + len[0] * (b ^ c);
}
bool inCircle(point a, point b, point c, point d) {
      int128 det = -det3(b, c, d);
    \overline{det} += det3(a, c, d);
    det -= det3(a, b, d);
    det += det3(a, b, c);
    return det > 0;
pair<quadEdge*, quadEdge*> buildTr(int 1, int r, vector<point>& p) { if(r - 1 <= 3) {
        quadEdge* a = makeEdge(p[1], p[1 + 1]), *b = makeEdge(p[1 + 1], p[r - 1]);
        if (r - 1 == 2) return mp(a, a->rev());
        splice(a->rev(), b);
        ll sg = area2(p[1], p[1 + 1], p[1 + 2]);
        quadEdge* c = sg ? connect(b, a) : 0;
        if (sg \ge 0) return mp(a, b \ge vev());
        else return mp(c->rev(), c);
    int m = (1 + r) >> 1;
quadEdge *ldo, *ldi, *rdo, *rdi;
    tie(ldo, ldi) = buildTr(l, m, p);
    tie(rdi, rdo) = buildTr(m, r, p);
    while(1) {
        if(left(rdi->o, ldi->o, ldi->dest()))
             ldi = ldi->lnext();
        else if(right(ldi->o, rdi->o, rdi->dest()))
            rdi = rdi->rev()->nxt;
        else break;
    quadEdge* basel = connect(rdi->rev(), ldi);
    auto valid = [&](quadEdge* e) {
        return right(e->dest(), basel->o, basel->dest());
    if(ldi->o == ldo->o) ldo = basel->rev();
    if(rdi->o == rdo->o) rdo = basel;
    while(1) {
        quadEdge *lcand = basel->rev()->nxt;
        if(valid(lcand)) {
             while(inCircle(basel->dest(), basel->o, lcand->dest(), lcand->nxt->dest()))
                 deleteEdge(lcand, lcand->nxt);
        quadEdge *rcand = basel->prev();
        if(valid(rcand)) {
             while(inCircle(basel->dest(), basel->o, rcand->dest(), rcand->prev()->dest()))
                 deleteEdge(rcand, rcand->prev());
        if(!valid(lcand) && !valid(rcand))
             break;
        if(!valid(lcand)
             || (valid(rcand) && inCircle(lcand->dest(), lcand->o, rcand->o, rcand->dest())))
             basel = connect(rcand, basel->rev());
        else basel = connect(basel->rev(), lcand->rev());
    return mp(ldo, rdo);
}
void delaunay(vector<point> p, vector<vector<int>>& adj) {
    vector<point> temp = p;
    map<point, int> m;
    for (int i = 0; i < sz(p); i++) m[p[i]] = i;
    sort(p.begin(), p.end());
    adj.resize(sz(p));
    auto add edge = [&](point a, point b) {
            adj[m[a]].pb(m[b]);
    };
    bool col = 1;
    for(int i = 2; i < sz(p); i++) col &= collinear(p[0], p[1], p[i]);
        for(int i = 0; i + 1 < sz(p); i++) {
            add_edge(p[i], p[i + 1]);
add_edge(p[i + 1], p[i]);
        quadEdge* e = buildTr(0, sz(p), p).first;
        vector<quadEdge*> edges = {e};
        for(int i = 0; i < sz(edges); e = edges[i++]) {
             for(quadEdge* at = e; !at->used; at = at->nxt) {
                 at->used = 1;
```

line integer.cpp

```
/* Basic structure of line defined by two 2D points which the line goes through.
   ^{\star} Some of the functions assume that the line is in fact a segment with the two
       points as its endpoints. Those operations are preceded by a comment stating
   * that this is the case.
  * This template depends on point_integer.cpp and works only with integers.
  * All operations' time complexity are O(1)
struct line {
           point p, q;
           line(point p, point q) : p(p), q(q) {}
           line() {}
           bool has(const point& r) const {
                     return paral((r - p), (q - p));
           bool operator == (const line& other) const { // assumes that direction does not matter
                     return has(other.p) && has(other.q);
           bool isVert() {
                     return p.x == q.x;
           // the following operations are for segments only
           bool segHas(point r) {
                     && (\min(p.y, q.y) \le r.y \le x.y 
           line rotate(point r) { // rotates segment pivoted in p
                     return line(p, p + (q - p).rotate(r));
           bool operator<(const line& other) const { // for Shamos-Hoey</pre>
                     // the case when q == other.p is such that we cosider this == other.
                      // that might give wrong answer, so be careful
                      if(p == other.p) return left(p, q, other.q);
                      if(!isVert() && (other.isVert() || p.x < other.p.x))</pre>
                                return left(p, q, other.p);
                     return left(p, other.q, other.p);
           }
};
int paraline(line a, line b) {
           return paral(a.q - a.p, b.q - b.p);
// the following functions are for segments only
bool checkInter(line a, line b) {
   if(a.segHas(b.p) || a.segHas(b.q) || b.segHas(a.p) || b.segHas(a.q))
                     return 1;
           return left(a.p, a.q, b.p) != left(a.p, a.q, b.q)
                                && left(b.p, b.q, a.p) != left(b.p, b.q, a.q);
}
```

angular_sweep.cpp

```
/* Version of point_integer with < operator for angular sweep.
 * The angle range is ]-\pi, \pi] and the template considers that the angle of the point
 * (0, 0) is 0. This template only contains a the operators needed for the for applying
 * the angular sort.
 */

struct point {
    ll x, y;
    point(ll x, ll y) : x(x), y(y) {}
    point() {}
    ll operator*(const point& other) const {
        return x*other.x + y*other.y;
    }
    ll operator^(const point& other) const { // cross product</pre>
```

```
return x*other.y - y*other.x;
}
int side() const {
    return y > 0 || (y == 0 && x < 0);
}
bool operator==(const point& other) const {
    return x == other.x && y == other.y;
}
bool operator<(const point& other) const { // for angular sweep
    int this_side = side(), other_side = other.side();
    if(this_side != other_side) return this_side < other_side;
    if(*this == point(0, 0)) return 0;
    if(other == point(0, 0)) return 1;
    return (*this ^ other) > 0;
}
};
```

polygon_double.cpp

```
/* Basic structure of polygon.
* This template depends on point double.cpp since it can work with double coordinates.
 * All operations' time complexity are O(1) unless stated otherwise.
struct polygon {
   vector<point> p;
    int n;
    polygon() : n(0) {}
    polygon(vector<point> _p) {
       p = _p;
        n = \overline{sz}(p);
    void add(point q) {
        p.pb(q);
        n++;
    // If positive, the polygon is in ccw order. It is in cw order otherwise.
    ld orientation() { // O(n)
        1d acum = 0;
        for(int i = 0; i < n; i++)
            acum += p[i] ^ p[(i + 1) % n];
        return acum;
    ld area() { // O(n)
        return abs(orientation()) / 2.0;
    void turnCcw() { // O(n)
        if(orientation() < -EPS)</pre>
            reverse(p.begin(), p.end());
    bool has(point q) { // O(log n). The polygon must be convex and in ccw order.
        if(right(p[0], p[1], q) || left(p[0], p[n-1], q)) return 0;
        int lo = 1, hi = n;
while(lo + 1 < hi) {
            int mid = (lo + hi) >> 1;
            if(!right(p[0], p[mid], q)) lo = mid;
            else hi = mid;
        return hi != n ? !right(p[lo], p[hi], q) : dist(p[0], q) < dist2(p[0], p[n-1]) + EPS;
    ld calipers() { // O(n). The polygon must be convex and in ccw order.
        ld ans = 0;
        for(int i = 0, j = 1; i < n; i++) {
            point vec_i = p[(i+1) n] - p[i];
while((vec_i ^ (p[(j+1) n] - p[j])) > EPS)
                j = (j + 1) % n;
            ans = max(ans, dist(p[i], p[j])); // Example with polygon diameter
        return ans;
    int extreme(const function<bool(point, point)> &cmp) {
        auto isExtreme = [&](int i, bool& curDir) -> bool {
            curDir = cmp(p[(i + 1) % n], p[i]);
            return !cmp(p[(i + n - 1) % n], p[i]) && !curDir;
        bool lastDir, curDir;
        if(isExtreme(0, lastDir)) return 0;
        int lo = 0, hi = n;
        while (lo + 1 < hi) {
```

```
int m = (lo + hi) >> 1;
           if (isExtreme (m, curDir)) return m;
           bool relDir = cmp(p[m], p[lo]);
           if((!lastDir && curDir) || (lastDir == curDir && relDir == curDir)) {
               lo = m;
               lastDir = curDir;
           } else hi = m;
       return lo;
   pair<int, int> tangent(point q) { // O(log n) for convex polygon in ccw orientation
       // Finds the indices of the two tangents to an external point q
       auto leftTangent = [&](point r, point s) -> bool {
           return right(q, r, s);
       auto rightTangent = [&](point r, point s) -> bool {
           return left(q, r, s);
       return {extreme(leftTangent), extreme(rightTangent)};
   int maximize(point v) { // O(log n) for convex polygon in ccw orientation
       // Finds the extreme point in the direction of the vector
       return extreme([&](point p, point q) {return p * v > q * v + EPS;});
   void normalize() { // p[0] becomes the lowest leftmost point
       rotate(p.begin(), min element(p.begin(), p.end()), p.end());
   polygon operator+(polygon& other) { // Minkowsky sum
       vector<point> sum;
       normalize();
       other.normalize();
       ld dir;
       for(int i = 0, j = 0; i < n \mid \mid j < other.n; i += dir > -EPS, j = dir < EPS) {
           return polygon(sum);
   }
};
```

line double.cpp

```
/* Basic structure of line defined by two 2D points which the line goes through.
^{\star} Some of the functions assume that the line is in fact a segment with the two
  points as its endpoints. Those operations are preceded by a comment stating
* that this is the case.
* This template depends on point double.cpp and hence the coordinates don't need
* to be integers.
* All operations' time complexity are O(1)
struct line {
   point p, q;
   line(point p, point q) : p(p), q(q) {}
   line() {}
   bool has(point r) const {
       return paral((r - p), (q - p));
   bool operator == (const line& other) const { // assumes that direction does not matter
       return has (other.p) && has (other.q);
   bool isVert() {
       return abs(p.x - q.x) <= EPS;
   point proj(point r) {
       point q vec = q - p, r vec = r - p;
        return p + q_vec * (q_vec * r_vec / q vec.norm2());
    ld dist(point r) {
        return (r - proj(r)).norm();
    // the following operations are for segments only
   bool segHas(point r) {
        return collinear(p, q, r)
            && (\min(p.x, q.x) < r.x + EPS && r.x < \max(p.x, q.x) + EPS)
            && (\min(p.y, q.y) < r.y + EPS && r.y < \max(p.y, q.y) + EPS);
    line rotate(point r) { // rotates segment pivoted in p
```

circle.cpp

```
/* Basic structure of circle and operations related with it. This template works
* only with double numbers since most of the operations of a circle can't be
* done with only integers. Therefore, this template depends on point_double.cpp.
* All operations' time complexity are O(1)
const ld PI = acos(-1);
struct circle {
   point o; ld r;
    circle() {}
    circle(point o, ld r) : o(o), r(r) {}
    bool has(point p) {
        return (o - p).norm2() < r*r + EPS;
    vector<point> operator/(circle c) { // Intersection of circles.
        vector<point> inter;
                                                   // The points in the output are in ccw order.
        1d d = (o - c.o).norm();
        if(r + c.r < d - EPS | | d + min(r, c.r) < max(r, c.r) - EPS)
            return {};
        1d x = (r*r - c.r*c.r + d*d) / (2*d);
        ld y = sqrt(r*r - x*x);
point v = (c.o - o) / d;
        inter.pb(o + v*x + v.rotate(cw90)*y);
        if (y > EPS) inter.pb(o + v*x + v.rotate(ccw90)*y);
        return inter;
    vector<point> tang(point p) {
        ld d = sqrt((p - o).norm2() - r*r);
return *this / circle(p, d);
    bool in(circle c){ // non strictly inside
        ld d = (o - c.o).norm();
return d + r < c.r + EPS;</pre>
};
```

polygon_integer.cpp

```
/* Basic structure of polygon.

*
 * This template depends on point_integer.cpp since it only works for integer coordinates.

* All operations' time complexity are O(1) unless stated otherwise.

*/

struct polygon {
    vector<point> p;
    int n;
    polygon() : n(0) {}
```

```
polygon(vector<point> _p) {
   p = _p;
n = sz(p);
void add(point q) {
    p.pb(q);
    n++;
// If positive, the polygon is in ccw order. It is in cw order otherwise.
11 orientation() { // O(n)
    11 \text{ acum} = 0;
    for(int i = 0; i < n; i++)
       acum += p[i] ^ p[(i + 1) % n];
    return acum;
ll area2() { // O(n)
    return abs(orientation());
void turnCcw() { // O(n)
    if(orientation() < 0)</pre>
        reverse(p.begin(), p.end());
bool has(point q) { // O(log n). The polygon must be convex and in ccw order
    if(right(p[0], p[1], q) || left(p[0], p[n-1], q)) return 0;
    int lo = 1, hi = n;
while(lo + 1 < hi) {
        int mid = (lo + hi) >> 1;
        if(!right(p[0], p[mid], q)) lo = mid;
        else hi = mid;
    return hi != n ? !right(p[lo], p[hi], q) : dist2(p[0], q) <= dist2(p[0], p[n-1]);
ll calipers() { // O(n). The polygon must be convex and in ccw order.
    11 \text{ ans} = 0;
    for(int i = 0, j = 1; i < n; i++) {
        ans = \max(ans, dist2(p[i], p[j])); // Example with polygon diameter squared
    return ans;
int extreme(const function<bool(point, point)> &cmp) {
   auto isExtreme = [&](int i, bool& curDir) -> bool {
    curDir = cmp(p[(i + 1) % n], p[i]);
        return !cmp(p[(i + n - 1) % n], p[i]) && !curDir;
    };
    bool lastDir, curDir;
    if(isExtreme(0, lastDir)) return 0;
    int lo = 0, hi = n;
    while (lo + 1 < hi) {
        int m = (lo + hi) >> 1;
        if(isExtreme(m, curDir)) return m;
        bool relDir = cmp(p[m], p[lo]);
        if((!lastDir && curDir) || (lastDir == curDir && relDir == curDir)) {
            lo = m;
            lastDir = curDir;
        } else hi = m;
    }
    return lo;
pair<int, int> tangent(point q) { // O(log n) for convex polygon in ccw orientation
    // Finds the indices of the two tangents to an external point q
    auto leftTangent = [&](point r, point s) -> bool {
        return right(q, r, s);
    auto rightTangent = [&](point r, point s) -> bool {
        return left(q, r, s);
    };
    return {extreme(leftTangent), extreme(rightTangent)};
int maximize(point v) { // O(log n) for convex polygon in ccw orientation
    \ensuremath{//} Finds the extreme point in the direction of the vector
    return extreme([&](point p, point q) {return p * v > q * v;});
void normalize() { // p[0] becomes the lowest leftmost point
    rotate(p.begin(), min_element(p.begin(), p.end()), p.end());
polygon operator+(polygon& other) { // Minkowsky sum
    vector<point> sum;
    normalize();
    other.normalize();
    for(ll i = 0, j = 0, dir; i < n || j < other.n; i += dir >= 0, j += dir <= 0) {
        sum.pb(p[i % n] + other.p[j % other.n]);
        dir = (p[(i + 1) % n] - p[i % n])
```

shamos_hoey.cpp

```
/* Shamos-Hoey algorithm for checking wether a collection of segments have an intersection.
* This template depends on point_integer.cpp and line_integer.cpp.
  seg - (input) collection of segments.
 * Time complexity: O(n logn)
bool shamos hoey(vector<line> seg) {
    // create sweep line events {x, type, seg_id}
    vector < array < 11, 3 > ev;
    for(int i = 0; i < sz(seg); i++) {
    if(seg[i].q < seg[i].p) swap(seg[i].p, seg[i].q);</pre>
        ev.pb({seg[i].p.x, 0, i});
ev.pb({seg[i].q.x, 1, i});
    sort(ev.begin(), ev.end());
    set<line> s;
    for(auto e: ev) {
        line at = seg[e[2]];
        if(!e[1]) {
            auto nxt = s.lower_bound(at);
            if((nxt != s.end() && checkInter(*nxt, at))
                || (nxt != s.begin() && checkInter(*(--nxt), at)))
                     return 1;
            s.insert(at);
        } else {
            auto nxt = s.upper bound(at), cur = nxt, prev = --cur;
            if(nxt != s.end() && prev != s.begin()
                && checkInter(*nxt, *(--prev))) return 1;
            s.erase(cur);
    return 0;
```

point_double.cpp

```
/* Basic structure of point and operations related with it. This template works
* with double coordinates.
* All operations' time complexity are O(1)
typedef long double ld;
const ld EPS = 1e-9;
struct point {
   ld x, y;
   point(ld x, ld y) : x(x), y(y) {}
   point() {}
   ld norm2()
       return *this * *this;
   ld norm() {
       return sqrt(norm2());
   bool operator==(const point& other) const {
       return abs(x - other.x) < EPS && abs(y - other.y) < EPS;
   point operator+(const point& other) const {
       return point(x + other.x, y + other.y);
   point operator-(const point& other) const {
       return point(x - other.x, y - other.y);
   point operator*(ld t) const {
       return point(x * t, y * t);
```

```
point operator/(ld t) const {
        return point(x / t, y / t);
    ld operator*(const point& other) const {
        return x*other.x + y*other.y;
    ld operator^(const point& other) const { // cross product
       return x*other.y - y*other.x;
    bool operator<(const point& other) const { // for sweep line
        return x < other.x - EPS || (abs(x - other.x) < EPS && y < other.y - EPS);
    point rotate(point r) {
        return point(*this ^ r, *this * r);
    point rotate(ld ang) {
        return rotate(point(sin(ang), cos(ang)));
    id angle(point& other) { // only works for angles in the range [0, PI]
    ld cos_val = min(1.0L, max(-1.0L, *this * other / (norm() * other.norm())));
        return acos(cos val);
};
point ccw90(1, 0);
point cw90(-1, 0);
ld dist2(point p, point q) { // squared distance
   return (p - q).norm2();
ld dist(point p, point q) {
    return sqrt(dist2(p, q));
ld area2(point a, point b, point c) { // two times signed area of triangle abc
   return (b - a) ^ (c - a);
bool left(point a, point b, point c) {
    return area2(a, b, c) > EPS;
bool right(point a, point b, point c) {
   return area2(a, b, c) < -EPS;
bool collinear(point a, point b, point c) {
   return abs(area2(a, b, c)) < EPS;
// Returns 0 if vectors a and b are not parallel.
// If they are parallel, returns 1 if they have the same direction
// and returns -1 otherwise
int paral(point a, point b)
    if((a ^ b) != 0) return 0;
    if((a.x > EPS) == (b.x > EPS) && (a.y > EPS) == (b.y > EPS))
        return 1;
    return -1;
}
```

Graphs

manhattan_mst.cpp, lca.cpp, block_cut_tree.cpp, articulation_points.cpp, bridges.cpp, two_sat.cpp,
tarjan scc.cpp

manhattan_mst.cpp

```
/* Computes O(n) edges that contains all the edges of the Manhattan MST
* of a set of points int the 2D plane. Must apply Kruskal to these edges
^{\star} to obtain the actual MST.
* xs - (input) coordinates in the x-axis
  ys - (input) coordinates in the y-axis
* E - superset of the manhattan MST edges
* Complexity: O(n logn)
vector<pair<ll, pair<int, int>> > E;
void manhattan mst(vector<ll> xs, vector<ll> ys) {
    int n = sz(xs);
    vector<int> ord(n);
    for (int i = 0; i < n; i++)
        ord[i]=i;
    for (int s = 0; s < 2; s++) {
        for(int t = 0; t < 2; t++) {
            auto cmp = [&](int i, int j) -> bool {
   if(xs[i] + ys[i] == xs[j] + ys[j]) return ys[i] < ys[j];</pre>
                 return xs[i] + ys[i] < xs[j] + ys[j];
             sort(ord.begin(), ord.end(), cmp);
            map<int, int> id;
for(int i: ord) {
                 for(auto it = id.lower_bound(-ys[i]); it != id.end(); it = id.erase(it)) {
                     int j = it -> ss;
                     if(xs[j] - ys[j] > xs[i] - ys[i]) break;
                     E.pb(mp(abs(xs[i] - xs[j]) + abs(ys[i] - ys[j]), mp(i, j)));
                 id[-ys[i]] = i;
             swap(xs,ys);
        for (int i = 0; i < n; i++)
            xs[i] *= -1;
    }
}
```

lca.cpp

```
/* Binary lifting algorithm to compute the lowest common ancestor of two nodes in a tree
*
* To use this template, first you need to add the (undirected) edges using `addEdge` and,
* after all the edges has been added, call `eulerTour(root, root)` where `root` is the
* root of the tree. Then you can use `lca(u, v)` to fin the lowest common ancestor of two
* nodes `u` and `v`.

* LOGN - logarithm in base 2 of N
* anc[u][j] - (output) 2^j-th ancestor o u
*
* Time complexity: O(n log n) for precomputing and O(log n) per query.
* Space complexity: O(n log n)
*/
const int LOGN = 20;
int anc[N][LOGN], tam[N], tin[N];
vector<int> adj[N];
int node_id;

void addEdge(int u, int v) {
   adj[u].pb(v);
   adj[v].pb(u);
}
```

```
// Call this with eulerToor(root, root)
int eulerTour(int u, int p) {
    anc[u][0] = p;
    tam[u] = 1;
tin[u] = node_id++;
    for (int i = 1; i < LOGN; i++)
         anc[u][i] = anc[anc[u][i-1]][i-1];
     for (int v: adj[u]) if (v != p)
         tam[u] += eulerTour(v, u);
    return tam[u];
}
// Check if u is ancestor of v in O(1)
bool isAnc(int u, int v) {
    return tin[u] <= tin[v] && tin[v] < tin[u] + tam[u];</pre>
int lca(int u, int v) {
    if(isAnc(u, v)) return u;
if(isAnc(v, u)) return v;
     for(int i = LOGN - 1; i >= 0; i--)
         if(!isAnc(anc[u][i], v)) u = anc[u][i];
     return anc[u][0];
}
```

block_cut_tree.cpp

```
/* Builds the block-cut tree of a given graph. The block-cut tree is a graph
  decomposition in biconnected components (blocks) and articulation points (cuts).
 * Every block is uniquely defined by a set of edges. This template considers
^{\star} the the articulation point only belongs to the cut.
* Usage: add undirected edges with addEdge(u, v) and then call build(n), when n is
* the total number of vertices in the graph. WARNING: make sure to do not add
  parallel edges. Also, notice that you have to clear adj, adj_bct and edge_cont
 * by yourself between different test cases.
* n - (input) number of vertices.
 * belong - belong[i] is the component in which the vertex i is in, either a block
            or a cut.
 * edge belong - edge belong[i] is the block in which the edge i is in.
 * ap - ap[i] is true if the vertex i is an articulation point. WARNING: this template
        considers isolated vertices as articulation points.
 * adj bct - the list of adjacency of the block-cut tree
 ^{\star} Complexity: O(n) to find all biconnected components and articulation points.
               O(n log n) to build the block-cut tree.
int edge_cont, comp, temp, stp;
pair<int, int> edge[N];
int belong[N], low[N], disc[N], st[N], edge belong[N];
bool ap[N];
vector<int> adj[N], adj bct[N];
void bcc(int v, int p=-1) {
    low[v] = disc[v] = ++temp;
    int child = 0;
    if(sz(adj[v]) == 0) ap[v] = 1;
    for(int ed: adj[v]) {
        int u = edge[ed].first ^ edge[ed].second ^ v;
        if(!disc[u]) {
            child++;
            st[stp++] = ed;
            bcc(u, v);
            low[v] = min(low[v], low[u]);
            if(low[u] >= disc[v]) {
                if (p != -1 || child > 1) ap[v] = 1;
                while(1) {
                    int e = st[--stp];
                    edge_belong[e] = comp;
                    belong[edge[e].first] = comp;
                    belong[edge[e].second] = comp;
                    if(e == ed) break;
                }
                comp++;
        } else if(u != p) {
            low[v] = min(low[v], disc[u]);
            if(disc[u] < disc[v])</pre>
                st[stp++] = ed;
    }
```

```
}
void addEdge(int u, int v) {
    adj[u].pb(edge_cont);
adj[v].pb(edge_cont);
    edge[edge_cont++] = \{u, v\};
void build(int n) {
    for(int i = 0; i < n; i++) {</pre>
         low[i] = disc[i] = ap[i] = 0;
    temp = comp = stp = 0;
    for (int i = 0; i < n; i++)
         if(!disc[i])
             bcc(i);
    for(int u = 0; u < n; i++) if(ap[u]) {
         for(int e: adj[u]) {
             adj_bct[comp].pb(edge_belong[e]);
             adj_bct[edge_belong[e]].pb(comp);
         belong[u] = comp;
         comp++;
    for(int i = 0; i < comp; i++) {
         sort(all(adj bct[i]));
         adj bct[i].resize(unique(all(adj bct[i])) - adj bct[i].begin());
}
```

articulation points.cpp

```
/* Finds the articulation points of an undirected graph. The edges might be parallel or
 * self-loops without problems.
* Usage: first add the edges with `addEdge(u, v)`, and then call `findAps(n)`.
* n - (input) number of vertices.
* ap - ap[i] is true if the vertex i is an articulation point.
* Complexity: O(V + E) time and space
int low[N], disc[N], t_in;
bool ap[N];
vector<int> adj[N];
void compute(int v, int root, int p=-1) {
    low[v] = disc[v] = ++t in;
    int childs = 0;
    for(int u : adj[v]) {
        if(!disc[u]) {
            childs++;
            compute(u, root, v);
            low[v] = min(low[v], low[u]);
            if(low[u] >= disc[v] && (v != root || childs > 1))
                ap[v] = 1;
        } else if(u != p) {
            low[v] = min(low[v], disc[u]);
    }
}
void addEdge(int u, int v) {
   adj[u].pb(v);
    adj[v].pb(u);
void findAps(int n) {
   for (int i = 0; i < n; i++) {
        low[i] = disc[i] = ap[i] = 0;
    t in = 0;
    for(int i = 0; i < n; i++) {
       if(!disc[i]) compute(i, i);
}
// Must be call between testcases
void clearGraph(int n) {
    for (int i = 0; i < n; i++)
```

```
adj[i].clear();
}
```

bridges.cpp

```
/* Finds the bridges of an undirected graph. The edges might be parallel or self-loops
* without problems.
* Usage: first add the edges with `addEdge(u, v)`, and then call `findBridges(n)`
* n - (input) number of vertices.
* bridge - bridge[i] is true if the i-th edge is a bridge.
* Complexity: O(V + E) time and space
vector<pair<int, int>> edge;
int low[N], disc[N], t in;
bool bridge[N];
vector<int> adj[N];
void compute(int v, int p=-1) {
    low[v] = disc[v] = ++t_in;
    for(int e : adj[v]) {
        int u = edge[e].first ^ edge[e].second ^ v;
        if(!disc[u]) {
            compute(u, e);
            low[v] = min(low[v], low[u]);
            if(low[u] > disc[v]) bridge[e] = 1;
        } else if(e != p) {
            low[v] = min(low[v], disc[u]);
}
void addEdge(int u, int v) {
    adj[u].pb(sz(edge));
    adj[v].pb(sz(edge));
    edge.pb({u, v});
void findBridges(int n) {
    for (int i = 0; i < n; i++) {
       low[i] = disc[i] = 0;
    for(int i = 0; i < sz(edge); i++) {
       bridge[i] = 0;
    t in = 0;
    for(int i = 0; i < n; i++) {
       if(!disc[i]) compute(i);
}
// Must be call between testcases
void clearGraph(int n) {
    for (int i = 0; i < n; i++)
       adj[i].clear();
    edge.clear();
}
```

two_sat.cpp

```
/* Returns true if a given set of conditions of the form a v b is
  * satisfiable. NOTE: you must clear adj and adjt between testcases.

* n - (input) number of atomic propositions
  * assign - one valid assignation of truth values

* Complexity: O(n)
  */

// Useful when negation of atomic expression is given as a negative number int conv(int x) {
    if(x > 0) return (x - 1) << 1;
    else return (-x - 1) << 1 | 1;
}

vector<int> adj[N], adjt[N];
```

```
void addEdge(int u, int v) {
    adj[u].pb(v);
    adjt[v].pb(u);
}
// Add the condition a v\ b
void addProp(int a, int b) {
    a = conv(a), b = conv(b);
    addEdge(a^1, b);
    addEdge(b^1, a);
bool vis[N];
int comp[N];
stack<int> order;
void toposort(int v) {
    vis[v] = 1;
    for(int u: adj[v]) if(!vis[u])
        toposort(u);
    order.push(v);
void mark(int v, int id) {
    comp[v] = id;
    for (int u: adjt[v]) if (comp[u] == -1)
        mark(u, id);
}
bool assign[N];
// Call this function after adding al conditions with addProp
bool twoSat(int n) {
    for(int i = 0; i < (n << 1); i++) {
        comp[i] = -1;
        vis[i] = 0;
    for(int i = 0; i < (n << 1); i++) if(!vis[i])
        toposort(i);
    int cont = 0;
    for(int i = 0; i < (n << 1); i++) {
        int v = order.top();
        order.pop();
        if(comp[v] == -1)
            mark(v, cont++);
    for (int i = 0; i < n; i++) {
        if(comp[i << 1] == comp[i << 1 | 1])
             return 0;
        assign[i] = comp[i << 1] > comp[i << 1 | 1];
    return 1;
}
```

tarjan_scc.cpp

```
/* Computes the Strongly Connected Components of a graph
 * n - (input) number of nodes of the graph
 * adj - (input) vector of adjacency of the graph
* cont - number of SCCs of the graph

* SCC[i] - list of nodes inside the i-th SCC
 ^{\star} belong[i] - the index of the SCC in which the node i is in
 ^{\star} adjSCC - vector of adjacency of the compressed graph
 * Complexity: O(n) to compute cont, SCC and belong
                 O(n log(n)) to compute adjSCC
 * /
int n, temp, cont;
vector<int> adj[N], SCC[N], adjSCC[N];
bool vis[N];
int low[N], disc[N], belong[N];
stack<int> S;
void dfs(int v) {
    disc[v] = low[v] = ++temp;
    S.push(v);
    vis[v] = 1;
    for(int u: adj[v]) {
         if(!disc[u])
```

```
dfs(u);
          if(vis[u])
               low[v] = min(low[u], low[v]);
     if(disc[v] == low[v]) {
          while (1) {
   int u = S.top();
               S.pop();
vis[u] = 0;
               belong[u] = cont;
               SCC[cont].pb(u);
               if(u == v) break;
          cont++;
     }
}
void tarjan() {
   for(int i = 0; i < n; i++)
        if(!disc[i])</pre>
     dfs(i);
for(int i = 0; i < cont; i++) {</pre>
          for(int u: SCC[i])
               for(int v: adj[u]) if(belong[v] != i) {
   adjSCC[i].pb(belong[v]);
          sort(adjSCC[i].begin(), adjSCC[i].end());
          adjSCC[i].resize(unique(adjSCC[i].begin(), adjSCC[i].end()) - adjSCC[i].begin());
}
```

Math

factorize.cpp, mod_exp.cpp, mod_inv_any.cpp, miller_rabin.cpp, extended_gcd.cpp, phi.cpp, divisors.cpp,
mod_inv.cpp, polynomials_operations_full.cpp, gcd.cpp, polynomials_operations.cpp, gauss.cpp,
inverseln.cpp, gauss_mod.cpp, gauss_mod2.cpp, matrix_exp.cpp

factorize.cpp

mod_exp.cpp

mod_inv_any.cpp

```
/* Returns the modular multiplicative inverse of a number or -1 if gcd(a, m) != 1

* a, m - positive integers

* Complexity: O(log(m))

*/
ll gcdExt(ll a, ll b, ll &x, ll &y) {
    if (b == 0) {
        x = 1;
        y = 0;
        return a;
    }
    ll x1, y1;
    ll res = gcdExt(b, a % b, x1, y1);
    x = y1;
    y = x1 - y1 * (a / b);
    return res;
}
ll inv(ll a, ll m = MOD) {
    ll x, y;
```

```
ll g = gcdExt(a, m, x, y);
if (g != 1) return -1;
return (m + x % m) % m;
}
```

miller_rabin.cpp

```
11 \mod \exp(11 b, 11 e, 11 m = MOD) {
    11 \text{ res} = 1;
    b %= m;
    while (e > 0) {
   if (e & 1) res = (res * b) % m;
   b = (b * b) % m;
        e /= 2;
    return res;
}
    Returns true if n is composite
    n - positive integer
    a - fermat base 2 <= a <= n - 2
    d_{r} s - greatest s such that = 2^s * d
bool checkComposite(ll n, ll a, ll d, int s) {
    11 x = modexp(a, d, n);
    if (x == 1 \mid \mid x == n - 1) return false;
    for (int i = 1; i < s; i++) {
        x = (x * x) % n;
        if (x == n - 1) return false;
    return true;
}
/\star Returns true if a number is prime. Deterministic for 64-bit integers.
 * n - positive integer
 * Complexity: O(\log(n)) the constant is at least 12
bool millerRabin(ll n) {
   if (n < 2) return false;
    int r = 0;
    11 d = n - 1;
    while ((d \& 1) == 0) {
        d >>= 1;
        r++;
    for (int a: {2, 3, 5, 7, 11, 13, 17, 19, 23, 29, 31, 37}) {
        if (n == a) return true;
        if (checkComposite(n, a, d, r)) return false;
    return true;
}
```

extended_gcd.cpp

```
/* Extended Euclidean algorithm. Returns gcd(a, b) and set the parameters
  * x and y to numbers such that ax + by = gcd(a, b).
  *
  * Time complexity: O(log(min(a, b)))
  */
  ll gcdExt(ll a, ll b, ll &x, ll &y) {
    if (b == 0) {
        x = 1;
        y = 0;
        return a;
    }
    ll x1, y1;
    ll res = gcdExt(b, a % b, x1, y1);
    x = y1;
    y = x1 - y1 * (a / b);
    return res;
}
```

phi.cpp

divisors.cpp

```
/* Get divisors of a number

* a - number

* d - output param to store the divisors (needs to be empty)

* 
* Complexity: O(n)

*/

void getDivisors(1l a, vector<1l> &d) {
    for (1l i = 1; i * i <= a; i++) {
        if (a % i == 0) {
            d.push_back(i);
            if (i * i != a) d.push_back(a / i);
        }
    }
}</pre>
```

mod_inv.cpp

```
/* Returns the modular multiplicative inverse of a number

* a - integer
* m - prime module

* Complexity: O(log(m))
*/
11 modexp(11 b, 11 e, 11 m = MOD) {
    11 res = 1;
    b % = m;
    while (e > 0) {
        if (e & 1) res = (res * b) % m;
        b = (b * b) % m;
        e /= 2;
    }
    return res;
}

11 inv(11 a, 11 m = MOD) {
    return modexp(a, m-2, m);
}
```

polynomials_operations_full.cpp

```
/* Useful polynomials operations of competitive programming
  * From https://github.com/e-maxx-eng/e-maxx-eng-aux/blob/master/src/polynomial.cpp
  *
  * Complexity: varies
  */
#include <bits/stdc++.h>
```

```
using namespace std;
namespace algebra {
   const int maxn = 1 << 21;
    const int magic = 250; // threshold for sizes to run the naive algo
    mt19937 rng(chrono::steady_clock::now().time_since_epoch().count());
    template<typename T>
    T bpow(T x, size_t n) {
        if(n == 0) {
           return T(1);
        } else {
            auto t = bpow(x, n / 2);
            t *= t;
            return n % 2 ? x * t : t;
        }
    }
    template<int m>
    struct modular {
        // https://en.wikipedia.org/wiki/Berlekamp-Rabin algorithm
        // solves x^2 = y \pmod{m} assuming m is prime in O(\log m). // returns nullopt if no sol.
        optional<modular> sqrt() const {
            static modular y;
            y = *this;
            if(r == 0)
                return 0;
            } else if(bpow(y, (m - 1) / 2) != modular(1)) {
                return nullopt;
            } else {
                while(true) {
                    modular z = rng();
                    if(z * z == *this) {
                        return z;
                    struct lin {
                        modular a, b;
                        lin(modular a, modular b): a(a), b(b) {}
                        lin(modular a): a(a), b(0) {}
                        lin operator * (const lin& t) {
                            return {
                                a * t.a + b * t.b * y,
                                 a * t.b + b * t.a
                    x(z, 1); // z + x
                    x = bpow(x, (m - 1) / 2);
                    if(x.b != modular(0)) {
                        return x.b.inv();
                }
            }
        int64 t r;
        modular() : r(0) {}
        \label{eq:modular(int64_trr): r(rr) {if(abs(r) >= m) r %= m; if(r < 0) r += m;}
        modular inv() const {return bpow(*this, m - 2);}
        modular operator - () const {return r ? m - r : 0;}
        modular operator * (const modular &t) const {return r * t.r % m;}
        modular operator / (const modular &t) const {return *this * t.inv();}
        modular operator += (const modular &t) {r += t.r; if(r >= m) r -= m; return *this;}
        modular operator -= (const modular &t) {r -= t.r; if(r < 0) r += m; return *this;}
        modular operator + (const modular &t) const {return modular(*this) += t;}
        modular operator - (const modular &t) const {return modular(*this) -= t;}
        modular operator *= (const modular &t) {return *this = *this * t;}
        modular operator /= (const modular &t) {return *this = *this / t;}
        bool operator == (const modular &t) const {return r == t.r;}
        bool operator != (const modular &t) const {return r != t.r;}
        operator int() const {return r;}
        int64 t rem() const {return 2 * r > m ? r - m : r;}
    } ;
    template<int T>
    istream& operator >> (istream &in, modular<T> &x) {
       return in >> x.r;
    template<typename T>
    T fact(int n) {
```

```
static T F[maxn];
    return F[n] ? F[n] : F[n] = n ? fact<T>(n - 1) * T(n) : T(1);
template<typename T>
T rfact(int n) {
    static T RF[maxn];
    return RF[n] ? RF[n] : RF[n] = T(1) / fact<T>(n);
namespace fft {
    typedef double ftype;
    typedef complex<ftype> point;
    point w[maxn];
    const ftype pi = acos(-1);
    bool initiated = 0;
    void init() {
        if(!initiated) {
    for(int i = 1; i < maxn; i *= 2) {</pre>
                 for (int j = 0; j < i; j++)
                     w[i + j] = polar(ftype(1), pi * j / i);
             initiated = 1;
        }
    }
    void fft(auto *in, point *out, int n, int k = 1) {
        if(n == 1) {
             *out = *in;
         } else {
            n /= 2;
             fft(in, out, n, 2 * k);
             fft(in + k, out + n, n, 2 * k);
             for(int i = 0; i < n; i++) {
                 auto t = out[i + n] * w[i + n];
                 out[i + n] = out[i] - t;
                 out[i] += t;
             }
        }
    void mul_slow(vector<auto> &a, const vector<auto> &b) {
        if(a.empty() || b.empty()) {
             a.clear();
         } else {
             a.resize(a.size() + b.size() - 1);
             for(int k = a.size() - 1; k >= 0; k--) {
                 a[k] *= b[0];
                 for(int j = 1; j < min(k + 1, (int)b.size()); j++) {
                     a[k] += a[k - j] * b[j];
             }
        }
    }
    template<typename T>
    void mul(vector<T> &a, vector<T> b) {
         if(min(a.size(), b.size()) < magic) {</pre>
            mul slow(a, b);
             return;
        init();
        static const T split = 1 << 15;</pre>
        size t n = a.size() + b.size() - 1;
        while( builtin popcount(n) != 1) {
             n++;
        a.resize(n);
        b.resize(n);
        static point A[maxn], B[maxn];
         static point C[maxn], D[maxn];
        for(size_t i = 0; i < n; i++) {
             A[i] = point(a[i].rem() % split, a[i].rem() / split);
             B[i] = point(b[i].rem() % split, b[i].rem() / split);
        fft(A, C, n); fft(B, D, n);
        for(size_t i = 0; i < n; i++) {
    A[i] = C[i] * (D[i] + conj(D[(n - i) % n]));
    B[i] = C[i] * (D[i] - conj(D[(n - i) % n]));</pre>
        fft(A, C, n); fft(B, D, n);
        reverse(C + 1, C + n);
reverse(D + 1, D + n);
```

```
int t = 2 * n;
        for(size_t i = 0; i < n; i++) {
            T AO = llround(real(C[i]) / t);
            T A1 = llround(imag(C[i]) / t + imag(D[i]) / t);
T A2 = llround(real(D[i]) / t);
            a[i] = A0 + A1 * split - A2 * split * split;
    }
template<typename T>
struct poly {
    vector<T> a;
    void normalize() { // get rid of leading zeroes
        while (!a.empty() && a.back() == T(0)) {
           a.pop_back();
    }
    poly(){}
    poly(T a0) : a{a0}{normalize();}
    poly(const vector<T> &t) : a(t) {normalize();}
    poly operator -() const {
        auto t = *this;
        for(auto &it: t.a) {
            it = -it;
        return t;
    }
    poly operator += (const poly &t) {
        a.resize(max(a.size(), t.a.size()));
        for(size_t i = 0; i < t.a.size(); i++) {</pre>
            a[i] += t.a[i];
        normalize();
        return *this;
    poly operator -= (const poly &t) {
        a.resize(max(a.size(), t.a.size()));
for(size_t i = 0; i < t.a.size(); i++) {
            a[i] -= t.a[i];
        normalize();
        return *this;
    poly operator + (const poly &t) const {return poly(*this) += t;}
    poly operator - (const poly &t) const {return poly(*this) -= t;}
    poly mod_xk(size_t k) const { // get first k coefficients
        return vector<T>(begin(a), begin(a) + min(k, a.size()));
    poly mul_xk(size_t k) const { // multiply by x^k
        auto res = a;
        res.insert(begin(res), k, 0);
        return res;
    poly div_xk(size_t k) const { // drop first k coefficients
        return vector<T>(begin(a) + min(k, a.size()), end(a));
    poly substr(size t l, size t r) const { // return mod xk(r).div xk(l)
        return vector<T>(
            begin(a) + min(l, a.size()),
            begin(a) + min(r, a.size())
        );
    }
    poly inv(size_t n) const { // get inverse series mod x^n
        assert((*this)[0] != T(0));
        poly ans = T(1) / a[0];
        size t a = 1;
        while (a < n) {
            poly C = (ans * mod_xk(2 * a)).substr(a, 2 * a);
            ans -= (ans * C).mod xk(a).mul xk(a);
            a *= 2;
        return ans.mod xk(n);
    }
```

```
poly operator *= (const poly &t) {fft::mul(a, t.a); normalize(); return *this;}
        poly operator * (const poly &t) const {return poly(*this) *= t;}
        poly reverse(size_t n) const { // computes x^n A(x^{-1})
             auto res = a;
             res.resize(max(n, res.size()));
             return vector<T>(res.rbegin(), res.rbegin() + n);
        poly reverse() const {
             return reverse(deg() + 1);
        pair<poly, poly> divmod slow(const poly &b) const { // when divisor or quotient is small
             vector<T> A(a);
             vector<T> res;
             while(A.size() >= b.a.size()) {
                 res.push back(A.back() / b.a.back());
                 if(res.back() != T(0)) {
                     for(size_t i = 0; i < b.a.size(); i++) {
    A[A.size() - i - 1] -= res.back() * b.a[b.a.size() - i - 1];
                 A.pop_back();
             std::reverse(begin(res), end(res));
             return {res, A};
        pair<poly, poly> divmod(const poly &b) const { // returns quotiend and remainder of a mod b
             assert(!b.is zero());
             if(deg() < b.deg()) {
                 return {poly{0}, *this};
             int d = deg() - b.deg();
             if(min(d, b.deg()) < magic) {</pre>
                 return divmod slow(b);
             poly D = (reverse().mod_xk(d + 1) * b.reverse().inv(d + 1)).mod_xk(d + 1).reverse(d + 1)
1);
             return {D, *this - D * b};
        poly operator / (const poly &t) const {return divmod(t).first;}
        poly operator % (const poly &t) const {return divmod(t).second;}
        poly operator /= (const poly &t) {return *this = divmod(t).first;}
        poly operator %= (const poly &t) {return *this = divmod(t).second;}
        poly operator *= (const T &x) {
             for(auto &it: a) {
                 it *= x;
             normalize();
             return *this;
        poly operator /= (const T &x) {
             for(auto &it: a) {
                 it /= x;
             normalize();
             return *this;
        poly operator * (const T &x) const {return poly(*this) *= x;} poly operator / (const T &x) const {return poly(*this) /= x;}
        poly conj() const \{ // A(x) \rightarrow A(-x) \}
             auto res = *this;
             for(int i = 1; i \le deg(); i += 2) {
                 res[i] = -res[i];
             return res;
         }
         void print(int n) const {
            for (int i = 0; i < n; i++) {
                 cout << (*this)[i] << ' ';
             cout << "\n";
         }
        void print() const {
            print(deg() + 1);
        T eval(T x) const \{ // \text{ evaluates in single point } x \}
             T res(0);
```

```
for(int i = deg(); i >= 0; i--) {
        res *= x;
        res += a[i];
    return res;
}
T lead() const { // leading coefficient
    assert(!is zero());
    return a.back();
}
int deg() const { // degree, -1 for P(x) = 0
    return (int)a.size() - 1;
bool is_zero() const {
   return a.empty();
T operator [] (int idx) const {
    return idx < 0 \mid \mid idx > deg() ? T(0) : a[idx];
T& coef(size t idx) { // mutable reference at coefficient
    return a[idx];
bool operator == (const poly &t) const {return a == t.a;}
bool operator != (const poly &t) const {return a != t.a;}
poly deriv() { // calculate derivative
    vector<T> res(deg());
    for(int i = 1; i <= deg(); i++) {
    res[i - 1] = T(i) * a[i];
    return res;
poly integr() { // calculate integral with C = 0
    vector<T> res(deg() + 2);
    for(int i = 0; i \le deg(); i++) {
        res[i + 1] = a[i] / T(i + 1);
    return res;
}
size t trailing xk() const \{ // \text{ Let } p(x) = x^k * t(x), \text{ return } k \}
    if(is_zero()) {
        return -1;
    int res = 0;
    while (a[res] == T(0)) {
       res++;
    return res;
}
poly log(size_t n) { // calculate log p(x) mod x^n }
    assert(a[0] == T(1));
    return (deriv().mod xk(n) * inv(n)).integr().mod xk(n);
poly exp(size_t n) { // calculate exp p(x) mod x^n
    if(is zero()) {
        return T(1);
    assert(a[0] == T(0));
    poly ans = T(1);
    size_t a = 1;
    while (a < n) {
        poly C = ans.log(2 * a).div xk(a) - substr(a, 2 * a);
        ans -= (ans * C).mod_xk(a).mul_xk(a);
        a *= 2;
    return ans.mod xk(n);
}
poly pow_bin(int64_t k, size_t n) { // O(n log n log k)
    if(k == 0) {
        return poly(1).mod_xk(n);
    } else {
        auto t = pow(k / 2, n);
        t *= t;
        return (k % 2 ? *this * t : t).mod xk(n);
```

```
}
}
// O(d * n) with the derivative trick from
// https://codeforces.com/blog/entry/73947?#comment-581173
poly pow_dn(int64_t k, size_t n) {
    vector<T> Q(n);
    Q[0] = bpow(a[0], k);
    for(int i = 1; i < (int)n; i++) {
   for(int j = 0; j <= min(deg(), i + 1); j++) {
      Q[i] += a[j] * Q[i - j] * (T(k + 1) * T(j) - T(i));
        Q[i + 1] /= i * a[0];
    return 0;
}
// calculate p^k(n) mod x^n in O(n log n)
// might be guite slow due to high constant
poly pow(int64_t k, size_t n) {
    if(is zero())
        return *this;
    int i = trailing xk();
    if(i > 0) {
        return i * k \ge (int64 t)n ? poly(0) : div xk(i).pow(k, n - i * k).mul xk(i * k);
    if(min(deg(), (int)n) <= magic) {</pre>
        return pow_dn(k, n);
    if(k <= magic)
        return pow bin(k, n);
    T j = a[i];
poly t = *this / j;
return bpow(j, k) * (t.log(n) * T(k)).exp(n).mod_xk(n);
// returns nullopt if undefined
optional<poly> sqrt(size_t n) const {
    if(is_zero())
        return *this;
    int i = trailing_xk();
if(i % 2) {
        return nullopt;
    } else if(i > 0)
        auto ans = div xk(i).sgrt(n - i / 2);
        return ans ? ans->mul xk(i / 2) : ans;
    auto st = (*this)[0].sqrt();
    if(st) {
        poly ans = *st;
         size_t a = 1;
        while (a < n) {
            a *= 2;
             ans -= (ans - mod_xk(a) * ans.inv(a)).mod_xk(a) / 2;
        return ans.mod xk(n);
    return nullopt;
}
poly mulx(T a) { // component-wise multiplication with a^k
    T cur = 1;
    poly res(*this);
    for(int i = 0; i <= deg(); i++) {
        res.coef(i) *= cur;
        cur *= a;
    return res;
}
poly mulx_sq(T a) { // component-wise multiplication with a^{k^2}
    T cur = a;
    T \text{ total} = 1;
    T aa = a * a:
    poly res(*this);
    for(int i = 0; i \le deg(); i++) {
        res.coef(i) *= total;
        total *= cur;
        cur *= aa;
    return res;
```

```
vector<T> chirpz even(T z, int n) { // P(1), P(z^2), P(z^4), ..., P(z^2(n-1))
                          int m = deg(\bar{)};
                          if(is zero()) {
                                  return vector<T>(n, 0);
                          vector<T> vv(m + n);
                          T zi = T(1) / z;
                          T zz = zi * zi;
                          T cur = zi;
                          T \text{ total} = 1;
                          for (int i = 0; i \le max(n - 1, m); i++) {
                                  if(i <= m) {vv[m - i] = total;}
                                   if(i < n) \{vv[m + i] = total;\}
                                   total *= cur;
                                  cur *= zz;
                          poly w = (mulx sq(z) * vv).substr(m, m + n).mulx sq(z);
                          vector<T> res(n);
for(int i = 0; i < n; i++) {</pre>
                                  res[i] = w[i];
                          return res;
                 }
                 vector<T> chirpz(T z, int n) { // P(1), P(z), P(z^2), ..., P(z^(n-1)) auto even = chirpz_even(z, (n + 1) / 2);
                          auto odd = mulx(z).chirpz even(z, n / 2);
                          vector<T> ans(n);
                          for (int i = 0; i < n / 2; i++) {
                                  ans[2 * i] = even[i];
                                   ans[2 * i + 1] = odd[i];
                          if(n % 2 == 1) {
                                  ans[n-1] = even.back();
                          return ans;
                 \verb|vector<T>| eval(vector<poly> & tree, int v, auto l, auto r) | { // auxiliary evaluation function } | { // auxiliary evaluation func
                         if(r - 1 == 1) {
                                  return {eval(*1)};
                          } else {
                                  auto m = 1 + (r - 1) / 2;
auto A = (*this % tree[2 * v]).eval(tree, 2 * v, 1, m);
auto B = (*this % tree[2 * v + 1]).eval(tree, 2 * v + 1, m, r);
                                   A.insert(end(A), begin(B), end(B));
                                  return A;
                         }
                 }
                 vector<T> eval(vector<T> x) { // evaluate polynomial in (x1, ..., xn)}
                          int n = x.size();
                          if(is zero()) {
                                  return vector<T>(n, T(0));
                          vector<poly> tree(4 * n);
                          build(tree, 1, begin(x), end(x));
                          return eval(tree, 1, begin(x), end(x));
                 poly inter(vector<poly> &tree, int v, auto l, auto r, auto ly, auto ry) { // auxiliary
interpolation function
                         if(r - 1 == 1) {
                                   return {*ly / a[0]};
                          } else {
                                   auto m = 1 + (r - 1) / 2;
                                   auto my = ly + (ry - ly) / 2;
                                   auto A = (*this \% tree[2 * v]).inter(tree, 2 * v, 1, m, ly, my);
auto B = (*this \% tree[2 * v + 1]).inter(tree, 2 * v + 1, m, r, my, ry);
                                  return A * tree[2 * v + 1] + B * tree[2 * v];
                 }
                 static auto resultant (poly a, poly b) \{\ //\ \text{computes resultant of a and b}
                         if(b.is zero()) {
                                  return 0:
                          } else if(b.deg() == 0) {
                                  return bpow(b.lead(), a.deg());
                           } else {
                                  int pw = a.deg();
                                   a \% = b:
                                   pw -= a.deg();
                                   auto mul = bpow(b.lead(), pw) * T((b.deg() & a.deg() & 1) ? -1 : 1);
                                   auto ans = resultant(b, a);
```

```
return ans * mul;
        }
        static poly kmul(auto L, auto R) { // computes (x-a1)(x-a2)...(x-an) without building tree
            if(R - L == 1) {
                return vector<T>{-*L, 1};
            } else {
                auto M = L + (R - L) / 2;
                return kmul(L, M) * kmul(M, R);
        }
        static poly build(vector<poly> &res, int v, auto L, auto R) { // builds evaluation tree for
(x-a1)(x-a2)...(x-an)
            if(R - L == 1) {
                return res[v] = vector<T>{-*L, 1};
            } else {
                auto M = L + (R - L) / 2;
                return res[v] = build(res, 2 * v, L, M) * build(res, 2 * v + 1, M, R);
        }
        static auto inter(vector<T> x, vector<T> y) { // interpolates minimum polynomial from (xi,
yi) pairs
            int n = x.size();
            vector<poly> tree(4 * n);
            return build(tree, 1, begin(x), end(x)).deriv().inter(tree, 1, begin(x), end(x),
begin(y), end(y));
        static poly xk(size t n) { // P(x) = x^n}
            return poly(T(1)).mul_xk(n);
        static poly ones(size t n) { // P(x) = 1 + x + ... + x^{n-1}}
            return vector\langle T \rangle (n, 1);
        static poly expx(size_t n) { // P(x) = e^x (mod x^n)
           return ones(n).borel();
        // [x^k] (a corr b) = sum_{i+j=k} ai*b{m-j}
                             = sum_{i-j=k-m} ai*bj
        static poly corr(poly a, poly b) { // cross-correlation
            return a * b.reverse();
        poly invborel() const { // ak *= k!
            auto res = *this;
for(int i = 0; i <= deg(); i++) {
                res.coef(i) *= fact<T>(i);
            return res;
        }
        poly borel() const { // ak /= k!
            auto res = *this;
            for (int i = 0; i \le deg(); i++) {
                res.coef(i) *= rfact<T>(i);
            return res;
        poly shift(T a) const \{ // P(x + a) \}
            return (expx(deg() + 1).mulx(a).reverse() * invborel()).div xk(deg()).borel();
    static auto operator * (const auto& a, const poly<auto>& b) {
        return b * a;
};
using namespace algebra;
const int mod = 998244353;
typedef modular<mod> base;
typedef poly<br/>base> polyn;
void solve() {
    int n, m;
cin >> n >> m;
```

```
vector<base> a(n);
    copy_n(istream_iterator<base>(cin), n, begin(a));
    polyn(a).pow(m, n).print(n);
}

signed main() {
    //freopen("input.txt", "r", stdin);
    ios::sync_with_stdio(0);
    cin.tie(0);
    int t;
    t = 1;// cin >> t;
    while(t--) {
        solve();
    }
}
```

gcd.cpp

```
/* Greatest common divisor

*
 * a, b - non-negative integers

*
 * Complexity: O(log(min(a, b)))
 */
ll gcd(ll a, ll b) {
    return b ? gcd(b, a % b) : a;
}
```

polynomials_operations.cpp

```
/* Useful polynomials operation of competitive programming
 * From https://github.com/e-maxx-eng/e-maxx-eng-aux/blob/master/src/polynomial.cpp
 * The main operations are kept in this file.
* Complexity: varies
struct Poly {
    vector<ll> p;
    Poly(int deg) { // for deg < 0, p(x) = 0
         p.resize(max(0, deg+1));
    Poly(vector<ll> &coefs) : p(coefs) {
   cout << coefs[0] << " " << coefs[1] << endl;</pre>
    int deg() { // degree is -1 for p(x) = 0
        return int(p.size()) - 1;
    Poly operator*(ll x) { // O(deg())
         Poly b = *this;
         for (ll &c : b.p) c *= x;
         return b;
    Poly operator*=(ll x) {
        return *this = *this * x;
    Poly mul slow(Poly &a) { // O(deg() * a.deg())
         Poly b(deg() + a.deg());
         for (int i = 0; i <= deg(); i++) {
    for (int j = 0; j <= a.deg(); j++) {
        b.p[i+j] += p[i] * a.p[j];
         return b;
    }
    void print() { // debug
         for (int i = deg(); i >= 0; i--) {
             if (p[i] < 0) cout << "- ";
             else cout << "+ ";
```

```
cout << abs(p[i]);
    if (i != 0) cout << "x^" << i << " ";
}
cout << endl;
}
};</pre>
```

gauss.cpp

```
/* Solves a system of linear equations
 * Complexity: O(n^3)
struct Gauss {
 int n, m;
  vector<int> pos;
 int rank = 0;
 vector<vector<double>> a;
 const double EPS = 1e-9;
  // n equations, m-1 variables, last column is for coefficients
 Gauss (int n, int m, vector<vector<double>> &a) : n(n), m(m), a(a) {
   pos.assign(m, -1);
  /* if a solution exists, it will be stored in ans
   0 - no solution
    1 - unique solution
    2 - infinite number of solutions */
  int solve(vector<double> &ans) {
    for (int col = 0, row = 0; col < m && row < n; col++) {
      int sel = row;
      for (int i = row+1; i < n; i++) {
       if (abs(a[i][col]) > abs(a[sel][col])) sel = i;
      if (abs(a[sel][col]) < EPS) continue;</pre>
      swap(a[sel], a[row]);
      pos[col] = row;
      for (int i = 0; i < n; i++) {
        if (i != row) {
          double mult = a[i][col] / a[row][col];
          a[i][col] = 0.0;
          for (int j = col+1; j <= m; j++) {
            a[i][j] -= a[row][j] * mult;
        }
      ++row, ++rank;
    ans.assign(m, 0);
    bool multiple = false;
    for (int i = 0; i < m; i++) {
     if (pos[i] != -1) ans[i] = a[pos[i]][m] / a[pos[i]][i];
      else multiple = true;
    for (int i = 0; i < n; i++) {
      double sum = 0.0;
      for (int j = 0; j < m; j++) {
  sum += ans[j] * a[i][j];
      if (abs(sum - a[i][m]) > EPS) return 0;
    if (multiple) return 2;
    return 1;
};
```

inverse1n.cpp

gauss_mod.cpp

```
/* Solves a system of linear equations in Z mod
 * Complexity: O(n^3)
ll gcdExt(ll a, ll b, ll &x, ll &y) {
    if (b == 0) {
       x = 1;
       y = 0;
       return a;
    11 x1, y1;
    ll res = gcdExt(b, a % b, x1, y1);
    x = y1;
    y = x1 - y1 * (a / b);
    return res;
ll inv(ll a, ll m = MOD) {
    11 x, y;
    ll g = gcdExt(a, m, x, y);
    if (q != 1) return -1;
    return (m + x % m) % m;
struct Gauss {
 int n. m;
 vector<int> pos;
 int rank = 0;
 11 mod;
  vector<vector<ll>> a;
  // n equations, m-1 variables, last column is for coefficients
  Gauss (int n, int m, 11 mod, vector<vector<11>> &a) : n(n), m(m), mod(mod), a(a) {
   pos.assign(m, -1);
  /\ast if a solution exists, it will be stored in ans
   0 - no solution
    1 - unique solution
    2 - infinite number of solutions */
  int solve(vector<11> &ans) {
    for (int col = 0, row = 0; col < m && row < n; col++) \{
      int sel = row;
      for (int i = row+1; i < n; i++) {
       if (a[i][col] > 0) {
          sel = i;
          break;
      if (a[sel][col] == 0) continue;
      swap(a[sel], a[row]);
      pos[col] = row;
      for (int i = 0; i < n; i++) {
        if (i != row) {
          11 mult = a[i][col] * inv(a[row][col], mod) % mod;
          a[i][col] = 0;
          for (int j = col+1; j <= m; j++) {
    a[i][j] = (a[i][j] + mod - a[row][j] * mult % mod) % mod;
```

```
++row, ++rank;
    ans.assign(m, 0);
    bool multiple = false;
    for (int i = 0; i < m; i++) {
  if (pos[i] != -1) ans[i] = a[pos[i]][m] * inv(a[pos[i]][i], mod) % mod;</pre>
      else multiple = true;
    for (int i = 0; i < n; i++) {
      11 \text{ sum} = 0.0;
       for (int j = 0; j < m; j++) {
         sum = (sum + ans[j] * a[i][j] % mod) % mod;
      if (sum != a[i][m]) return 0;
    if (multiple) return 2;
    return 1;
};
```

gauss_mod2.cpp

```
/* Solves a system of linear equations in Z 2, it is faster than gauss mod
* by the use of bitset.
* Complexity: O(n^3)
template<int M>
struct Gauss {
 int n, m;
 array<int, M> pos;
 int rank = 0;
 vector<bitset<M>>> a;
  \ensuremath{//} n equations, m-1 variables, last column is for coefficients
 Gauss(int n, int m, vector<br/>
(m), (m), (m), (m), (m)
   pos.fill(-1);
  int solve(bitset<N> &ans) {
    for (int col = 0, row = 0; col < m && row < n; col++) {
      int one = -1;
      for (int i = row; i < n; i++) {
        if (a[i][col]) {
          one = i;
          break;
        }
      if (one == -1) { continue; }
      swap(a[one], a[row]);
      pos[col] = row;
      for (int i = row + 1; i < n; i++) {
        if (a[i][col])
          a[i] ^= a[row];
      ++row, ++rank;
    ans.reset();
    for (int i = m - 1; i >= 0; i--) {
      if (pos[i] == -1) ans[i] = true;
      else {
        int k = pos[i];
        for (int j = i + 1; j < m; j++) if (a[k][j]) ans[i] = ans[i] ^ ans[j]; ans[i] = ans[i] ^ a[k][m];
```

```
for (int i = rank; i < n; i++) if (a[i][m]) return 0;
    return 1;
};</pre>
```

matrix_exp.cpp

```
/\!\!\!\!\!\!^{\star} Returns the matrix exponentiation
* base - d x d matrix
 \,^{\star} n - non-negative integer for the exponent
* res - output param
 * Complexity: O(d^2 * log(n))
void matProd(vector<vector<ll>> &a, vector<vector<ll>> &b, vector<vector<ll>> &c) {
    int n = a.size();
    int m = a[0].size();
int p = b[0].size();
    vector<vector<ll>> res(n, vector<ll>(p, 0));
    for (int i = 0; i < n; i++) {
        for (int j = 0; j < p; j++) {
    for (int k = 0; k < m; k++) {
                 res[i][j] += a[i][k] * b[k][j];
         }
    c = res;
void matExp(vector<vector<ll>>> &base, ll n, vector<vector<ll>>> &res) {
    int d = base.size();
    for (int i = 0; i < d; i++) {
         for (int j = 0; j < d; j++) {
            res[i][j] = ll(i == j);
    while (n > 0) {
        if (n \% 2 == 1) matProd(res, base, res);
         matProd(base, base, base);
        n /= 2;
}
```

Numerical methods

roots newton.cpp, ternary search.cpp, fast double bs.cpp, stable sum.cpp, simpson integration.cpp

roots_newton.cpp

```
/* Find some root of a function f with derivative df.

* Complexity: The precision doubles for each iteration, in ideal conditions.

* For roots with multiplicity greater than one, the precision increases linearly.

*/
double f(double x);
double df(double x);

double findRoot(double x0=0.0) {
    double x = x0;
    double fx = f(x);
    while (abs(fx) > EPS) {
        x -= fx / df(x);
        fx = f(x);
    }
    return x;
}
```

ternary_search.cpp

```
/* Find the minimum of a function that first strictly decreases, then has its
    * minimum and finally strictly increases
    *
    * Complexity: O(log(n))
    */

double ternary_search(double 1, double r) {
    while (r - 1 > EPS) { // TODO set EPS}
        double m1 = 1 + (r - 1) / 3;
        double m2 = r - (r - 1) / 3;
        if (f(m1) > f(m2)) 1 = m1; // < for maximum
        else r = m2;
    }
    return 1;
}</pre>
```

fast_double_bs.cpp

```
/* Binary search with doubles considering that small numbers are
* closer to each other, works only if 1 > 0
*/
double bs(double 1, double r) {
   while (r - 1 > EPS) {
        double m;
        if (r > 2*1) m = sqrt(a*b);
        else m = (a + b) / 2;
   }
}
```

stable_sum.cpp

```
/* From Handbook of geometry for competitive programmers - Victor Lecomte
* This is an algorithm to make sums of positive numbers more precise
* Complexity is O(n) amortized and the relative error of the sum is
* 2 log_2(n) eps, down from n * eps precision of the serial sum
* eps is the machine precision.
*/

struct stableSum {
  int cnt = 0;
    vector<double> v, pref{0};
```

```
// add number a to the sum
void operator+=(double a) {
    assert(a >= 0);
    int s = ++cnt;
    while (s % 2 == 0) {
        a += v.back();
        v.pop_back();
        pref.pop_back();
        s /= 2;
    }
    v.push_back(a);
    pref.push_back(pref.back() + 1);
}

// return the sum value
double val() {
    return pref.back();
}
```

simpson_integration.cpp

```
/* Returns the integral of some function f in the interval [a, b].

* Complexity: O(2n)
*/
double f(double x);

double integral(double a, double b, int n) {
    n *= 2;
    double h = (b - a) / n;
    double s = f(a) + f(b);
    for (int i = 1; i < n; i++) {
        s += f(a + h*i) * (1 + (i & 1)) * 2;
    }
    return s * h / 3;
}</pre>
```

Strings

 $z_{\rm function.cpp}$, manacher.cpp, rolling_hash.cpp, kmp.cpp, suffix_automata.cpp, kmp_automata.cpp, suffix array.cpp, min cyclic string.cpp

z_function.cpp

```
/* Computes z function, where z[i] is the length of the longest prefix of s[i, n)
 * that is also a prefix of s[0, n). For convention, this template assumes z[0] = 0
 * Complexity: O(n)
vector<int> z_function(string s) {
    int n = sz(s);
    vector<int> z(n);
    for (int i = 1, l = 0, r = 0; i < n; i++) {
        if(i \le r)
            z[i] = min(r - i + 1, z[i - 1]);
        while(i + z[i] < n && s[z[i]] == s[i + z[i]])
            z[i]++;
        if(i + z[i] - 1 > r) {
            1 = i;
            r = i + z[i] - 1;
    return z;
}
```

manacher.cpp

```
/* Computes the following for every position of the string s:
 * best[0][i] - length of the longest palindrome that ends in s[i]
* best[1][i] - length of the longest palindrome that starts in s[i]
* d1[i] - number of odd length palindromes with center in s[i]
* d2[i] - number of even length palindromes with center in s[i]
           (we consider the center of an even length palindrome as
           the rightmost of the two characters in the center)
* Complexity: O(n)
int best[2][N];
int d1[N], d2[N];
void manacher(string &s) {
   n = sz(s);
    for (int i = 0; i < n; i++)
       best[0][i] = best[1][i] = 1;
    for (int i = 0, l = 0, r = -1; i < n; i++) {
        int k = (i > r) ? 1 : min(d1[1 + r - i], r - i + 1);
        while (0 \le i - k \&\& i + k \le n \&\& s[i - k] == s[i + k]) {
            best[1][i-k] = max(best[1][i-k], k << 1 | 1);
            best[0][i+k] = max(best[0][i+k], k << 1 | 1);
        d1[i] = k--;
        if^{(i+k>r)} {
            l = i - k;
            r = i + k;
    for (int i = 0, l = 0, r = -1; i < n; i++) {
        int k = (i > r) ? 0 : min(d2[1 + r - i + 1], r - i + 1);
        while (0 \le i - k - 1 \&\& i + k \le n \&\& s[i - k - 1] == s[i + k]) {
            best[1][i-k-1] = max(best[1][i-k-1], k*2 + 2);
            best[0][i+k] = max(best[0][i+k], k*2 + 2);
            k++;
        d2[i] = k--;
        if(i + k > r)  {
            1 = i - k - 1;
            r = i + k;
```

```
}
```

rolling_hash.cpp

```
/st Computes the suffixes of a string s. Can answer queries of the
* any substring of s.
* START - first character of the alphabet
* K - number of prime bases
* P - array of prime bases
* Complexity: O(n) precomputation, O(1) queries
const 11 MOD = 1e9 + 7;
const 11 K = 2;
const 11 P[K] = \{29, 31\};
const char START = 'a';
array<11, K> pot[N];
ll mul(ll a, ll b) {
   return a * b % MOD;
11 sum(11 a, 11 b) {
   a += b;
    if(a >= MOD) a -= MOD;
   return a;
}
// Must be called once in the beginning of the program to compute
// powers of the prime bases
void prec() {
   for (int k = 0; k < K; k++)
        pot[0][k] = 1;
    for (int i = 1; i < N; i++)
       for (int k = 0; k < K; k++)
            pot[i][k] = mul(pot[i-1][k], P[k]);
}
struct hsh {
    vector<array<11, K> > suf;
    hsh() {}
   hsh(string \&s) : suf(sz(s) + 1) {
        for (int k = 0; k < K; k++) suf[sz(s)][k] = 0;
        for(int i = sz(s) - 1; i >= 0; i--)
            for (int k = 0; k < K; k++)
                suf[i][k] = sum(mul(suf[i + 1][k], P[k]), s[i] - START + 1);
    // Queries the hashing of the substring s[l, r)
    array<ll, K> que(int 1, int r) {
        array<11, K> cur;
        for (int k = 0; k < K; k++)
            cur[k] = sum(suf[l][k], MOD - mul(suf[r][k], pot[r - l][k]));
        return cur;
};
```

kmp.cpp

suffix automata.cpp

```
/* Online algorithm that computes the suffix automata of a string.
 * Every state of the suffix automata represents a set of end positions of some substrings
* and every path on the suffix automata represents a different substring.
^{\star} You can use the extend method to add character by character to the suffix automata
* or use the constructor with a string to build the suffix automara for the whole
* string at once.
 ^{\star} Complexity: O(n log ALPH) time for building the suffix automata for n characters.
               O(n) space
struct suffixAutomata {
    struct state {
        int len, link;
        map<char, int> next;
        state() : len(0), link(-1) {}
state(int len) : len(len) {}
        state(int len, int link, map<char, int> next)
            : len(len), link(link), next(next) {}
    };
    vector<state> st;
    int last:
    suffixAutomata() : last(0) {
        st.pb({});
    suffixAutomata(string &s) : last(0) {
        st.pb({});
        for(auto c: s)
            extend(c);
    void extend(char c) {
        st.pb({st[last].len + 1});
        int cur = sz(st) - 1;
        int p = last;
        while (p != -1 \&\& !st[p].next.count(c)) {
            st[p].next[c] = cur;
            p = st[p].link;
        if (p == -1) st [cur]. link = 0;
        else {
            int q = st[p].next[c];
            if(st[p].len + 1 == st[q].len) st[cur].link = q;
                int clone = sz(st);
                 st.pb({st[p].len + 1, st[q].link, st[q].next});
                 while (p != -1 \&\& st[p].next[c] == q) {
                     st[p].next[c] = clone;
                     p = st[p].link;
                 st[q].link = st[cur].link = clone;
             }
        last = cur;
    }
} ;
```

kmp_automata.cpp

```
/* Computes the kmp automata. The i-th state of this automata corresponds to any string
* which its longest suffix that is a prefix of string has length i. The entry kmp[i][j]
* corresponds to the resulting state of transitioning from state i with character
* START + j.
* Remember that the i-th entry of the pi function does not correspond to the i-th state
* of the automata.

*
* START - first character of the alphabet
* ALPH - alphabet size

* Complexity: O(n * ALPH)
*/

const int ALPH = 26;
const char START = 'a';

struct kmpAutomata {
    vector<int> pi;
    vector<array<int, ALPH> > kmp;
```

```
string s:
    int go(int i, int j){
        if(kmp[i][j] != -1) return kmp[i][j];
        int ans;
        if(s[i] == j + START) ans = i + 1;
        else if(i == 0) ans = 0;
        else ans = go(pi[i-1], j);
        return kmp[i][j] = ans;
    kmpAutomata(string _s) : s(_s) {
        s += '#';
        int n = sz(s);
        pi.resize(n);
        pi[0] = 0;
        int k = 0;
        for(int i = 1; i < n; i++) {
            while (k \&\& s[i] != s[k])
                k = pi[k-1];
            if(s[i] == s[k]) k++;
            pi[i] = k;
        kmp.resize(n);
        for(int i = 0; i < n; i++)
            for(int j = 0; j < ALPH; j++)
kmp[i][j] = -1;
        for (int i = 0; i < n; i++)
            for (int j = 0; j < ALPH; j++)
                go(i, j);
};
```

suffix_array.cpp

```
/* Computes the Suffix Array of a set of strings.
\,^\star v - (input) vector of strings for which the suffix array will be built.
* s - concatenation of the strings in v divided by the character '$'
* n - suffix array size.
* sa - suffix array. sa[i] means that the suffix s[sa[i], n) is the lexicografical i-th
        suffix of s
\star lcp - array of longest common prefixes. Must call computeLcp to calculate the entries
        of this array. lcp[i] means that the longest common prefix between s[sa[i-1], n)
        and s[sa[i], n) has length lcp[i]
* Complexity: O(n logn) time for building the suffix array.
                O(n) time for computing the lcp array.
                O(n) memory.
* /
struct suffixArray {
   string s;
    vector<int> sa, lcp;
    int n;
    void cSort(int k, vector<int>& ra) {
        int maxi = max(300, n);
        vector<int> c(maxi, 0), temp sa(n);
        for (int i = 0; i < n; i++)
            c[i + k < n ? ra[i + k] : 0]++;
        for(int i = 1; i < maxi; i++)
    c[i] += c[i - 1];</pre>
        for(int i = n - 1; i >= 0; i--)
            temp sa[--c[sa[i] + k < n ? ra[sa[i] + k] : 0]] = sa[i];
        sa.swap(temp sa);
    suffixArray() {}
    suffixArray(vector<string>& v) {
        for(auto str: v) {
            s += str:
            s += '$';
        n = sz(s);
        sa.resize(n);
        vector<int> ra(n), temp_ra(n);
        for (int i = 0; i < n; i++) {
            ra[i] = s[i];
            sa[i] = i;
        for (int k = 1; k < n; k <<= 1) {
            cSort(k, ra);
            cSort(0, ra);
            int r = temp_ra[sa[0]] = 0;
for(int i = 1; i < n; i++)</pre>
```

```
ra.swap(temp ra);
            if (r == n - 1) break;
        }
    void computeLcp() {
       vector<int> phi(n);
        int k = 0;
        phi[sa[0]] = -1;
        for(int i = 1; i < n; i++)
    phi[sa[i]] = sa[i-1];
        vector<int> plcp(n);
        for(int i = 0; i < n - 1; i++) {
   if(phi[i] == -1) {
                plcp[i] = 0;
                continue;
            while (s[i + k] == s[phi[i] + k]) k++;
            plcp[i] = k;
            k = \max(k - 1, 0);
        lcp.resize(n);
        for(int i = 0; i < n; i++)
    lcp[i] = plcp[sa[i]];</pre>
};
```

min_cyclic_string.cpp

```
/* Returns the start index of the minimum cyclic string
* Uses Lyndon Factorization
* Complexity: O(n)
int min_cyclic_string(string s) {
    s += s;
    int n = s.size();
    int i = 0, res = 0;
    while (i < n / 2) {
        res = i;
int j = i + 1, k = i;
        while (j < n \&\& s[k] \le s[j]) {
            if (s[k] < s[j]) {
                 k = i;
            } else {
                k++;
            j++;
        while (i \leq k) {
            i += j - k;
    return res;
}
```

Utils

random.cpp, k-mino.cpp

random.cpp

```
/* Random number generator

*
    *Complexity: O(1)
    */
mt19937 rng(chrono::steady_clock::now().time_since_epoch().count());
int x = rng() % n; // to generante random number

// to generate a uniform random number in the range [a, b]
11 random(11 a, 11 b) {
    return uniform_int_distribution<11> (a, b) (rng);
}
```

k-mino.cpp

```
/\star Example code to generate k-minos
 * Complexity: not efficient
int dx[4] = \{0, 0, -1, 1\};
int dy[4] = \{-1, 1, 0, 0\};
struct Tab {
    int w, h;
    int tab[10][10];
    Tab() {
         w = 0;
         h = 0;
         for (int i = 0; i < 10; i++) {
    for (int j = 0; j < 10; j++) {
                  tab[i][j] = 0;
         }
    }
    Tab (const Tab &t) {
         w = t.w;
         h = t.h;
         for (int i = 0; i < 10; i++) {
             for (int j = 0; j < 10; j++) {
    tab[i][j] = t.tab[i][j];
             }
         }
    }
    void shiftDown() {
         for (int i = h; i > 0; i--) {
             for (int j = 0; j < w; j++) {
                  tab[i][j] = tab[i-1][j];
         for (int j = 0; j < w; j++) tab[0][j] = 0;
         h++;
    }
    void shiftRight() {
         for (int j = w; j > 0; j--) {
    for (int i = 0; i < h; i++) {
                  tab[i][j] = tab[i][j-1];
         for (int i = 0; i < h; i++) tab[i][0] = 0;
         w++;
    pair<ll, 11> getHash() {
         ll h[2];
         for (int k = 0; k < 2; k++) {
             h[k] = 0;
```

```
for (int i = 0; i < 5; i++) {
                 for (int j = 0; j < 10; j++) {
   h[k] *= 2;
                     h[k] += tab[5*k + i][j];
            }
        }
        return {h[0], h[1]};
    }
    void print() {
        for (int i = 0; i < h; i++) {
             for (int j = 0; j < w; j++) {
    cout << (tab[i][j] ? '0' : ' ');
             cout << "\n";
        cout << "\n";
    }
};
vector<Tab> ps[11];
void rec(int k) {
    if (k == 1) {
        Tab t;
        t.w = 1;
        t.h = 1;
        t.tab[0][0] = 1;
        ps[k].push_back(t);
        return;
    set<pair<11, 11>> hashes;
    rec(k-1);
    for (auto &t : ps[k-1]) {
        for (int i = -1; i <= t.h; i++) {
    for (int j = -1; j <= t.w; j++) {
                 if (i >= 0 && i < 10 && j >= 0 && j < 10 && t.tab[i][j] == 1) continue;
                 bool any = false;
                 for (int a = 0; a < 4; a++) {
                     int ai = i + dx[a];
                     int aj = j + dy[a];
                     if (ai < 0 || ai >= 10 || aj < 0 || aj >= 10 || t.tab[ai][aj] == 0) continue;
                      any = true;
                      break;
                 if (any) {
                     Tab nt(t);
                      if (i == -1) {
                          i++;
                         nt.shiftDown();
                      } else if (i == t.h) {
                          nt.h++;
                      if (j == -1) {
                         j++;
                          nt.shiftRight();
                      } else if (j == t.w) {
                         nt.w++;
                     nt.tab[i][j] = 1;
                      pair<11, 11> h = nt.getHash();
                      if (hashes.find(h) == hashes.end()) {
                          ps[k].push back(nt);
                          hashes.insert(h);
                     }
                 }
           }
       }
   }
}
int cx[51][51];
void solve() {
    int n, m;
    cin >> n >> m;
    for (int i = 0; i < n; i++) {
        for (int j = 0; j < n; j++) {
    cin >> cx[i][j];
```