



International Institute Of Information Technology - Hyderabad

# FLogic

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4	Contest (1)	
4	.bashrc	3 lines
5	alias c='g++ -Wall -Wconversion -Wfatal-errors -g -std=c++14 \	
6	-fsanitize=undefined,address'	
7	xmodmap -e 'clear lock' -e 'keycode 66=less greater' #caps = <>	
7	template.cpp	17 lines
7	#include <bits/stdc++.h>	
10	using namespace std;	
10	#define rep(i, a, b) for(int i = a; i < (b); ++i)	
11	#define all(x) begin(x), end(x)	
11	#define sz(x) (int)(x).size()	
11	#define pb push_back	
11	typedef long long ll;	
12	typedef pair<int, int> pii;	
12	typedef vector<int> vi;	
13	int main() {	
13	// freopen("sample.in", "r", stdin);	
13	// freopen("sample.out", "w", stdout);	
13	cin.tie(0)->sync_with_stdio(0);	
13	cin.exceptions(cin.failbit);	
13	}	
	Data structures (2)	
	OrderedSet.h	
	Description: A set (not multiset!) with support for finding the n'th element, and finding the index of an element. To get a map, change null_type.	
	Time: $\mathcal{O}(\log N)$	782797, 16 lines
	#include <bits/extc++.h>	
	using namespace __gnu_pbds;	
	template<class T>	
	using Tree = tree<T, null_type, less<T>, rb_tree_tag,	
	tree_order_statistics_node_update>;	
	void example() {	
	Tree<int> t, t2; t.insert(8);	
	auto it = t.insert(10).first;	
	assert(it == t.lower_bound(9));	
	assert(t.order_of_key(10) == 1);	
	assert(t.order_of_key(11) == 2);	
	assert(*t.find_by_order(0) == 8);	
	t.join(t2); // assuming T < T2 or T > T2, merge t2 into t	
	}	
	Dsu.h	
	Description: DSU with rollback	
	Time: $\mathcal{O}(\alpha(N))$	1262f5, 21 lines
	struct DSU {	
	int sets; vi p, s;	
	stack<pii> ss, sp;	
	DSU(int n) : p(n, -1), s(n, 1), sets(n) {}	
	bool IsSameSet(int a, int b) { return find(a) == find(b); }	
	int find(int x) {return p[x] == -1 ? x : p[x] = find(p[x]);}	
	bool join(int a, int b) {	

a = find(a), b = find(b);	
if (a == b) return false;	
if (s[a] < s[b]) swap(a, b);	
ss.push({a, s[a]}); sp.push({b, p[b]});	
sets--; s[a] += s[b]; p[b] = a; return true;	
}	
int time() {return sz(ss);}	
void rollback(int t) {	
while(time() > t) {	
p[sp.top().first] = sp.top().second; sp.pop();	
s[ss.top().first] = ss.top().second; ss.pop();	
}	
}	
};	
DsuBp.h	
Description: Graph, adding edges, checking bp color	
Time: $\mathcal{O}(\alpha(N))$	8e325a, 18 lines
struct DSU {	
int sets; vi p, s, l;	
DSU(int n) : p(n, -1), s(n, 1), l(n, 0), sets(n) {}	
bool IsSameColor(int a, int b) {	
find(a); find(b); return l[a] == l[b];	
}	
bool IsSameSet(int a, int b) { return find(a) == find(b); }	
int find(int x) {	
if(p[x] == -1) return x;	
int y = find(p[x]); l[x] ^= l[p[x]]; return p[x] = y;	
}	
void join(int a, int b) {	
int ca = a, cb = b; a = find(a), b = find(b);	
if (a == b) return;	
if (s[a] < s[b]) swap(a, b);	
sets--; s[a] += s[b]; l[b] = 1 ^ l[ca] ^ l[cb]; p[b] = a;	
}	
};	
MinQueue.h	
Description: Minimum Queue Applications	
Time: $\mathcal{O}(1)$ push pop etc.	68dd56, 24 lines
template<class T>	
struct MinQueue {	
deque<pair<T, T>> q;	
int ca = 0, cr = 0, plus = 0, sze = 0;	
void push(T x) {	
x -= plus;	
// change '>' to '<' and you get max-queue	
while (!q.empty() && q.back().first > x)	
q.pop_back();	
q.push_back({x, ca}); ca++; sze++;	
}	
T pop() {	
T re = 0;	
if (!q.empty() && q.front().second == cr) {	
re = q.front().first; q.pop_front();	
}	
cr++; sze--; return re + plus;	
}	
// Returns minimum in the queue	
T min() { return q.front().first + plus; }	
int size() { return sze; }	
// Adds x to every element in the queue	
void add(int x) { plus += x; }	
};	

Matrix.h

**Description:** Basic operations on square matrices.  
**Usage:** Matrix<int, 3> A;  
A.d = {{{{1,2,3}}, {{4,5,6}}, {{7,8,9}}}};  
vector<int> vec = {1,2,3};  
vec = (A^N) \* vec;

c43c7d, 26 lines

```
template<class T, int N> struct Matrix {
    typedef Matrix M;
    array<array<T, N>, N> d{};
    M operator*(const M& m) const {
        M a;
        rep(i,0,N) rep(j,0,N)
            rep(k,0,N) a.d[i][j] += d[i][k]*m.d[k][j];
        return a;
    }
    vector<T> operator*(const vector<T>& vec) const {
        vector<T> ret(N);
        rep(i,0,N) rep(j,0,N) ret[i] += d[i][j] * vec[j];
        return ret;
    }
    M operator^(ll p) const {
        assert(p >= 0);
        M a, b(*this);
        rep(i,0,N) a.d[i][i] = 1;
        while (p) {
            if (p&1) a = a*b;
            b = b*b;
            p >>= 1;
        }
        return a;
    }
};
```

SparseTable.h

**Description:** Range Minimum Queries on an array. Returns min(V[a], V[a + 1], ... V[b - 1]) in constant time.  
**Usage:** RMQ rmq(values);  
rmq.query(inclusive, exclusive);  
**Time:**  $\mathcal{O}(|V|\log|V| + Q)$

4a61f2, 21 lines

```
template<class T>
struct SparseTable {
    T (*op)(T, T);
    vi log2s; vector<vector<T>> st;
    SparseTable (const vector<T>& arr, T (*op)(T, T))
        : op(op), log2s(sz(arr)+1), st(sz(arr)) {
        rep(i,2,sz(log2s)) { log2s[i] = log2s[i/2] + 1; }
        rep(i,0,sz(arr)) {
            st[i].assign(log2s[sz(arr) - i] + 1);
            st[i][0] = arr[i];
        }
        rep(p, 1, log2s[sz(arr)] + 1) rep(i,0,sz(arr))
            if(i+(1<<p) <= sz(arr)) {
                st[i][p] = op(st[i][p-1], st[i+(1<<(p-1))][p-1]);
            }
    }
    T query (int l, int r) {
        int p = log2s[r-l+1];
        return op(st[l][p], st[r-(1<<p)+1][p]);
    }
};
```

FenwickTree.h

**Description:** Computes partial sums a[0] + a[1] + ... + a[pos - 1], and updates single elements a[i], taking the difference between the old and new value.  
**Time:** Both operations are  $\mathcal{O}(\log N)$ .

e62fac, 22 lines

```
struct FT {
```

```
    vector<ll> s;
    FT(int n) : s(n) {}
    void update(int pos, ll dif) { // a[pos] += dif
        for (; pos < sz(s); pos |= pos + 1) s[pos] += dif;
    }
    ll query(int pos) { // sum of values in [0, pos)
        ll res = 0;
        for (; pos > 0; pos &= pos - 1) res += s[pos-1];
        return res;
    }
    int lower_bound(ll sum) { // min pos st sum of [0, pos] >= sum
        // Returns n if no sum is >= sum, or -1 if empty sum is.
        if (sum <= 0) return -1;
        int pos = 0;
        for (int pw = 1 << 25; pw; pw >>= 1) {
            if (pos + pw <= sz(s) && s[pos + pw-1] < sum)
                pos += pw, sum -= s[pos-1];
        }
        return pos;
    }
};
```

FenwickTree2D.h

**Description:** Computes sums a[i,j] for all i<I, j<J, and increases single elements a[i,j]. Requires that the elements to be updated are known in advance (call fakeUpdate() before init()).  
**Time:**  $\mathcal{O}(\log^2 N)$ . (Use persistent segment trees for  $\mathcal{O}(\log N)$ .)

"FenwickTree.h"157f07, 22 lines

```
struct FT2 {
    vector<vi> ys; vector<FT> ft;
    FT2(int limx) : ys(limx) {}
    void fakeUpdate(int x, int y) {
        for (; x < sz(ys); x |= x + 1) ys[x].push_back(y);
    }
    void init() {
        for (vi& v : ys) sort(all(v)), ft.emplace_back(sz(v));
    }
    int ind(int x, int y) {
        return (int)(lower_bound(all(ys[x]), y) - ys[x].begin()); }
    void update(int x, int y, ll dif) {
        for (; x < sz(ys); x |= x + 1)
            ft[x].update(ind(x, y), dif);
    }
    ll query(int x, int y) {
        ll sum = 0;
        for (; x; x &= x - 1)
            sum += ft[x-1].query(ind(x-1, y));
        return sum;
    }
};
```

SegmentTree.h

**Description:** RMQ SegTree  
**Time:**  $\mathcal{O}(\log(N))$

f19f05, 42 lines

```
const ll INF = 1e18;
struct node {
    ll x;
};
```

```
template<class T>
struct SegmentTrees {
    vector<node> st, lazy;
    node def;
    SegmentTrees(int n) : st(4*n, {INF}), lazy(4*n, {INF}), def({
        INF}) {}
    inline node combine(node a, node b) {
        node ret; ret.x = min(a.x, b.x); return ret;
    }
};
```

```
void push(int pos) {
    if(lazy[pos].x != INF) {
        st[pos*2] = lazy[pos]; st[pos*2 + 1] = lazy[pos];
        lazy[pos*2] = lazy[pos]; lazy[pos*2+1] = lazy[pos];
        lazy[pos] = def;
    }
}
void update(int l,int r,T val,int left,int right,int pos=1) {
    if(l > r) return;
    if(l==left && r==right) {
        st[pos].x = val; lazy[pos] = {val};
    } else {
        push(pos);
        int mid = (left + right)/2;
        update(l, min(r,mid), val, left, mid, pos*2);
        update(max(l,mid+1), r, val, mid+1, right, pos*2+1);
        st[pos] = combine(st[pos*2], st[pos*2+1]);
    }
}
node query(int l,int r,int left,int right,int pos=1) {
    if(l>r) return def;
    if(l==left && r==right) return st[pos];
    else {
        push(pos); int mid = (left + right)/2;
        return combine(query(l, min(r,mid), val, left, mid, pos*2),
            query(max(l,mid+1), r, mid+1, right, pos*2+1));
    }
}
};
```

Treap.h

**Description:** cutting and moving array. everything is [l, r] 0 based indexing.  
**Usage:** Treap<int> tr(arr);  
**Time:**  $\mathcal{O}(\log N)$

419fcb, 166 lines

```
struct node {
    int prior, val, minl, lazy, size;
    bool rev;
    node *l, *r;
};
typedef node* pnode;

template<class T = int>
class Treap {
public:
    pnode root;
    pnode getnode(T val) {
        pnode t = new node;
        t->l = t->r = NULL;
        t->prior = rand(); t->size = 1; t->rev = false;
        t->lazy = 0; t->minl = t->val = val;
        return t;
    }
    inline int sz(pnode t) { return t ? t->size : 0;}
    // t may denote same node as l or r, so take care of that.
    void combine(pnode &t,pnode l,pnode r) {
        if(!l or !r) return void(t = (l ? l : r));
        t->size = sz(l) + sz(r); t->minl = min(l->minl, r->minl);
    }
    void operation(pnode t) {
        if(!t) return;
        // reset t;
        t->size = 1; t->minl = t->val;
        push(t->l); push(t->r);
        // combine
        combine(t, t->l, t); combine(t, t, t->r);
    }
    void push(pnode t) {
        if(!t) return;
```

```

    if(t->rev) {
        swap(t->r, t->l);
        if(t->r) t->r->rev = not t->r->rev;
        if(t->l) t->l->rev = not t->l->rev;
        t->rev = false;
    }
    if(t->lazy) {
        t->val += t->lazy;
        t->minl += t->lazy;
        if(t->r) t->r->lazy += t->lazy;
        if(t->l) t->l->lazy += t->lazy;
        t->lazy = 0;
    }
}
// l = [0, pos], r = rest
void split(pnode t,pnode &l,pnode &r,int pos,int add=0) {
    push(t);
    if(!t) return void(l=r=NULL);
    int curr_pos = add + sz(t->l);
    if(pos >= curr_pos) {
        split(t->r,t->r, r, pos, curr_pos + 1);
        l = t;
    } else {
        split(t->l, l, t->l, pos, add);
        r = t;
    }
    operation(t);
}
void merge(pnode &t,pnode l,pnode r) {
    push(l); push(r);
    if(!l or !r) return void(t = (l ? l : r));
    if(l->prior > r->prior) {
        merge(l->r, l->r, r);
        t = l;
    } else {
        merge(r->l, l, r->l);
        t = r;
    }
    operation(t);
}
void heapify(pnode t) {
    if(!t) return ;
    pnode max = t;
    if (t->l != NULL && t->l->prior > max->prior)
        max = t->l;
    if (t->r != NULL && t->r->prior > max->prior)
        max = t->r;
    if (max != t) {
        swap (t->prior, max->prior);
        heapify (max);
    }
}
// O(n) treap build given array is increasing
pnode build(T *arr,int n) {
    if(n==0) return NULL;
    int mid = n/2;
    pnode t = getnode(arr[mid]);
    t->l = build(arr, mid);
    t->r = build(arr + mid + 1, n - mid - 1);
    heapify(t); operation(t);
    return t;
}
Treap(vector<T> &arr) {
    root = NULL;
    for(int i=0;i<arr.size();i++) {
        T c = arr[i];
        merge(root, root, getnode(c));
    }
}
}

```

```

void add(int l,int r,T d) {
    if(l>r) return;
    pnode L, mid, R;
    split(root, L, mid, l-1); split(mid, mid, R, r-1);
    if(mid) {
        mid->lazy += d;
    }
    merge(L, L, mid); merge(root, L, R);
}
void reverse(int l,int r) {
    if(l>r) return;
    pnode L, mid, R;
    split(root, L, mid, l-1); split(mid, mid, R, r-1);
    if(mid) {
        mid->rev = not mid->rev;
    }
    merge(R, mid, R); merge(root, L, R);
}
void revolve(int l,int r,int cnt) {
    if(cnt<=0 or l>r) return;
    int len = r - l + 1;
    // cnt = len => no rotation;
    cnt %= len;
    if(cnt == 0) return;
    // pick cnt elements from the end // => (len - cnt) from front
    int mid = l + (len - cnt) - 1; pnode L, Range, R;
    split(root, L, Range, l-1); split(Range, Range, R, r - 1);
    pnode first, second;
    split(Range, first, second, (len-cnt-1));
    merge(Range, second, first);
    merge(L, L, Range); merge(root, L, R);
}
void insert(int after,T val) {
    pnode L, R; split(root, L, R, after);
    merge(L, L, getnode(val)); merge(root, L, R);
}
void del(int pos) {
    pnode L, mid, R;
    split(root, L, mid, pos-1); split(mid, mid, R, 0);
    if(mid) {
        delete mid;
    }
    merge(root, L, R);
}
T range_min(int l,int r) {
    pnode L, mid, R;
    split(root, L, mid, l-1); split(mid, mid, R, r-1);
    push(mid); T ans = mid->minl;
    merge(L, L, mid); merge(root, L, R);
    return ans;
}
void inorder(pnode curr) {
    push(curr); if(!curr) return;
    inorder(curr->l); cerr<<curr->val<<" "; inorder(curr->r);
}
int query(int pos) {
    pnode l, mid, r;
    split(root, l, mid, pos-1); split(mid, mid, r, 0);
    int ans = mid->val;
    merge(l, l, mid); merge(root, l, r);
    return ans;
}
}
};

```

## SQRT.h

**Description:** Square Root Decomposition

**Time:** Amul Knows

976251, 65 lines

```

const int N = 1e5 + 13, Q = 1e5 + 13, B = 500;
int S[N/B + 13][B + 13], len[N/B + 13], prv[N], nxt[N], st[N/B + 13], en[N/B + 13], A[N];
map<int,set<int>> pos; int n, q;

void add_link(int p,int val) {
    nxt[p] = val; prv[val] = p;
    if(p < 1 or p > n) return;
    int b = p / B;
    for(int i = st[b]; i <= en[b]; i++) {
        S[b][i - st[b] + 1] = nxt[i];
    }
    sort(S[b] + 1, S[b] + len[b] + 1);
}
// set A_x = y
void point_update(int x,int y) {
    // update the original link
    add_link(prv[x], nxt[x]); pos[A[x]].erase(x);
    // insert new links
    A[x] = y; pos[A[x]].insert(x);
    int pr = 0, nx = n + 1;
    if(*pos[A[x]].begin() != x) pr = *prev(pos[A[x]].find(x));
    if(*pos[A[x]].rbegin() != x) nx = *next(pos[A[x]].find(x));
    add_link(pr, x); add_link(x, nx);
}

int query_block(int s,int e,int k) {
    int ans = 0;
    for(int i = s; i <= e; i++)
        ans += ((S[i] + len[i] + 1) - upper_bound(S[i] + 1, S[i] + len[i] + 1, k));
    return ans;
}

int query_elements(int s,int e,int k) {
    int ans = 0;
    for(int i = s; i <= e; i++)
        ans += (nxt[i] > k);
    return ans;
}

int range_query(int l,int r) {
    int lb = l / B, rb = r / B;
    if(lb == rb) return query_elements(l, r, r);
    return query_elements(l, en[lb], r)
        + query_block(lb + 1, rb - 1, r)
        + query_elements(st[rb], r, r);
}
for(int i = 1; i <= n; i++) {
    nxt[i] = n + 1;
    if(!pos[A[i]].empty()) {
        prv[i] = *pos[A[i]].rbegin();
        nxt[prv[i]] = i;
    }
    pos[A[i]].insert(i);
}
for(int i = 1; i <= n; i++) {
    int b = i / B;
    if(!len[b])
        st[b] = i;
    en[b] = i;
    len[b]++;
    S[b][len[b]] = nxt[i];
}
for(int i = 0; i <= n/B; i++) {
    sort(S[i] + 1, S[i] + len[i] + 1);
}
}

```

LazyDynamicSegTree.h

**Description:** Segment Tree based on large [L, R] range (includes range updates)  
**Time:**  $\mathcal{O}(\log(R - L))$  in addition and deletion

391dcb, 31 lines

```
using T=ll; using U=ll; // exclusive right bounds
T t_id; U u_id; // t_id: total (normal), u_id: lazy (default)
T op(T a, T b){ return a+b; }
void join(U &a, U b){ a+=b; }
void apply(T &t, U u, int x){ t+=x*u; }
T part(T t, int r, int p){ return t/r*p; }
struct DynamicSegmentTree {
    struct Node { int l, r, lc, rc; T t; U u;
        Node(int l, int r):l(l),r(r),lc(-1),rc(-1),t(t_id),u(u_id){ }
    };
    vector<Node> tree;
    DynamicSegmentTree(int N) { tree.push_back({0,N}); }
    void push(Node &n, U u){ apply(n.t, u, n.r-n.l); join(n.u,u); }
    void push(Node &n){push(tree[n.lc],n,u);push(tree[n.rc],n,u);
        n.u=u_id;}
    T query(int l, int r, int i = 0) { auto &n = tree[i];
        if(r <= n.l || n.r <= l) return t_id;
        if(l <= n.l && n.r <= r) return n.t;
        if(n.lc < 0) return part(n.t, n.r-n.l, min(n.r,r)-max(n.l,l));
        return push(n), op(query(l,r,n.lc),query(l,r,n.rc));
    }
    void update(int l, int r, U u, int i = 0) { auto &n = tree[i];
        if(r <= n.l || n.r <= l) return;
        if(l <= n.l && n.r <= r) return push(n,u);
        if(n.lc < 0) { int m = (n.l + n.r) / 2;
            n.lc = tree.size(); n.rc = n.lc+1;
            tree.push_back({tree[i].l, m}); tree.push_back({m, tree[i].r});
        }
        push(tree[i]); update(l,r,u,tree[i].lc); update(l,r,u,tree[i].rc);
        tree[i].t = op(tree[tree[i].lc].t, tree[tree[i].rc].t);
    }
};
```

LineContainer.h

**Description:** Container where you can add lines of the form  $kx+m$ , and query maximum values at points  $x$ . Useful for dynamic programming (“convex hull trick”).  
**Time:**  $\mathcal{O}(\log N)$

8ec1c7, 30 lines

```
struct Line {
    mutable ll k, m, p;
    bool operator<(const Line& o) const { return k < o.k; }
    bool operator<(ll x) const { return p < x; }
};

struct LineContainer : multiset<Line, less<>> {
    // (for doubles, use inf = 1/.0, div(a,b) = a/b)
    static const ll inf = LLONG_MAX;
    ll div(ll a, ll b) { // floored division
        return a / b - ((a ^ b) < 0 && a % b); }
    bool isect(iterator x, iterator y) {
        if (y == end()) return x->p >= inf, 0;
        if (x->k == y->k) x->p = x->m > y->m ? inf : -inf;
        else x->p = div(y->m - x->m, x->k - y->k);
        return x->p >= y->p;
    }
    void add(ll k, ll m) {
        auto z = insert({k, m, 0}), y = z++, x = y;
        while (isect(y, z)) z = erase(z);
    }
};
```

LineContainer BellmanFord FloydWarshall Dinic

```
if (x != begin() && isect(--x, y)) isect(x, y = erase(y));
while ((y = x) != begin() && (--x)->p >= y->p)
    isect(x, erase(y));
}
ll query(ll x) {
    assert(!empty());
    auto l = *lower_bound(x);
    return l.k * x + l.m;
}
};
```

Graph (3)

3.1 Fundamentals

BellmanFord.h

**Description:** Calculates shortest paths from  $s$  in a graph that might have negative edge weights. Unreachable nodes get  $\text{dist} = \text{inf}$ ; nodes reachable through negative-weight cycles get  $\text{dist} = -\text{inf}$ . Assumes  $V^2 \max |w_i| < \sim 2^{63}$ . undirected graph with negative edge gets  $-\text{inf}$  for all vertices.  
**Time:**  $\mathcal{O}(VE)$

1101fe, 29 lines

```
const int INF = 1e9;
void bellmann_ford_extended(vector<vector<pii>> &adj,
    int source, vi &dist, vector<bool> &cyc) {
    dist.assign(adj.size(), INF);
    cyc.assign(adj.size(), false); // true when u is in a <0 cycle
    dist[source] = 0;
    for (int iter = 0; iter < adj.size() - 1; ++iter){
        bool relax = false;
        for (int u = 0; u < adj.size(); ++u)
            if (dist[u] == INF) continue;
            else for (auto &e : adj[u])
                if (dist[u]+e.second < dist[e.first])
                    dist[e.first] = dist[u]+e.second, relax = true;
            if(!relax) break;
        }
    bool ch = true;
    while (ch) { // keep going untill no more changes
        ch = false; // set dist to -INF when in cycle
        for (int u = 0; u < adj.size(); ++u)
            if (dist[u] == INF) continue;
            else for (auto &e : adj[u])
                if (dist[e.first] > dist[u] + e.second
                    && !cyc[e.first]) {
                    dist[e.first] = -INF;
                    ch = true; //return true for cycle detection only
                    cyc[e.first] = true;
                }
        }
    }
```

FloydWarshall.h

**Description:** Calculates all-pairs shortest path in a directed graph that might have negative edge weights. Input is an distance matrix  $m$ , where  $m[i][j] = \text{inf}$  if  $i$  and  $j$  are not adjacent. As output,  $m[i][j]$  is set to the shortest distance between  $i$  and  $j$ ,  $\text{inf}$  if no path, or  $-\text{inf}$  if the path goes through a negative-weight cycle.  
**Time:**  $\mathcal{O}(N^3)$

61c353, 12 lines

```
const ll inf = 1LL << 62;
void floydWarshALL(vector<vector<ll>>& m) {
    int n = sz(m);
    rep(i,0,n) m[i][i] = min(m[i][i], 0LL);
    rep(k,0,n) rep(i,0,n) rep(j,0,n)
        if (m[i][k] != inf && m[k][j] != inf) {
            auto newDist = max(m[i][k] + m[k][j], -inf);
```

```
        m[i][j] = min(m[i][j], newDist);
    }
    rep(k,0,n) if (m[k][k] < 0) rep(i,0,n) rep(j,0,n)
        if (m[i][k] != inf && m[k][j] != inf) m[i][j] = -inf;
}
```

3.2 Network flow

Dinic.h

**Description:** Complexity: (1)  $\mathcal{O}(V^2E)$ : General (2)  $\mathcal{O}(\text{Flow}E)$ : General (3)  $\mathcal{O}(E\sqrt{V})$ : when sum of edge capacities is  $\mathcal{O}(n)$ , we can treat edge with weight  $x$  as  $x$  edges with weight 1. (4)  $\mathcal{O}(EV \log(\text{Flow}))$ : Dinics with scaling

allife, 61 lines

```
const int INF = 1e9 + 13;
template<class T = long long>
class Dinic {
    // {to: to, rev: reverse_edge_id, c: cap, oc: original cap}
    struct Edge {
        int to, rev;
        T c, oc;
        T flow() { return max(oc - c, (T)0); } // if you need flows
    };
    int N;
    vector<int> lvl, ptr, q; vector<vector<Edge>> adj;
public:
    vector<vector<T>> Flow;
    Dinic(int n) {
        N = n; Flow.assign(n, vector<T>(n, (T)0));
        lvl.resize(n); adj.resize(n); ptr.resize(n); q.resize(n);
    }
    // automatically adds a reversed edge
    void addEdge(int a, int b, T c, T rcap = 0) {
        adj[a].push_back({b, sz(adj[b]), c, c});
        adj[b].push_back({a, sz(adj[a]) - 1, rcap, rcap});
    }
    T dfs(int v, int t, T f) {
        if (v == t || !f) return f;
        for (int& i = ptr[v]; i < sz(adj[v]); i++) {
            Edge& e = adj[v][i];
            if (lvl[e.to] == lvl[v] + 1)
                if (T p = dfs(e.to, t, min(f, e.c))) {
                    e.c -= p, adj[e.to][e.rev].c += p;
                    return p;
                }
        }
        return 0;
    }
    T calc(int s, int t) {
        T flow = 0; q[0] = s;
        // bfs part, setting the lvl here
        for(int L = 0; L < 31; L++) do { // 'int L=30' maybe faster
            for random data
            lvl = ptr = vector<int>(sz(q));
            int qi = 0, qe = lvl[s] = 1;
            while (qi < qe && !lvl[t]) {
                int v = q[qi++];
                for (Edge e : adj[v])
                    if (!lvl[e.to] && e.c >> (30 - L))
                        q[qe++] = e.to, lvl[e.to] = lvl[v] + 1;
            }
            // dfs part, setting ptr and checking for a path.
            while (T p = dfs(s, t, INF)) flow += p;
        } while (lvl[t]);
        return flow;
    }
    bool leftOfMinCut(int a) { return lvl[a] != 0; }
    void buildFlow() {
        for(int i=0;i<N;i++) {
            for(auto e : adj[i]) {
                int j = e.to;
```

```
        Flow[i][j] = e.flow();
    }
}
};
```

MinCut.h

**Description:** After running max-flow, the left side of a min-cut from  $s$  to  $t$  is given by all vertices reachable from  $s$ , only traversing edges with positive residual capacity.

GlobalMinCut.h

**Description:** Find a global minimum cut in an undirected graph, as represented by an adjacency matrix.

**Time:**  $\mathcal{O}(V^3)$

```
pair<int, vi> globalMinCut(vector<vi> mat) {
    pair<int, vi> best = {INT_MAX, {}};
    int n = sz(mat);
    vector<vi> co(n);
    rep(i,0,n) co[i] = {i};
    rep(ph,1,n) {
        vi w = mat[0];
        size_t s = 0, t = 0;
        rep(it,0,n-ph) { // O(V^2) -> O(E log V) with prio. queue
            w[t] = INT_MIN;
            s = t, t = max_element(all(w)) - w.begin();
            rep(i,0,n) w[i] += mat[t][i];
        }
        best = min(best, {w[t] - mat[t][t], co[t]});
        co[s].insert(co[s].end(), all(co[t]));
        rep(i,0,n) mat[s][i] += mat[t][i];
        rep(i,0,n) mat[i][s] = mat[s][i];
        mat[0][t] = INT_MIN;
    }
    return best;
}
```

8b0e19, 21 lines

MinCostMaxFlow.h

**Description:** Min-cost max-flow.  $cap[i][j] \neq cap[j][i]$  is allowed; double edges are not. If costs can be negative, call setpi before maxflow, but note that negative cost cycles are not supported. To obtain the actual flow, look at positive values only.

**Time:** Approximately  $\mathcal{O}(E^2)$

#include <bits/extc++.h>

```
const ll INF = numeric_limits<ll>::max() / 4;
typedef vector<ll> VL;
```

```
struct MCMF {
    int N;
    vector<vi> ed, red;
    vector<VL> cap, flow, cost;
    vi seen;
    VL dist, pi;
    vector<pii> par;
```

```
MCMF(int N) :
    N(N), ed(N), red(N), cap(N, VL(N)), flow(cap), cost(cap),
    seen(N), dist(N), pi(N), par(N) {}
```

```
void addEdge(int from, int to, ll cap, ll cost) {
    this->cap[from][to] = cap;
    this->cost[from][to] = cost;
    ed[from].push_back(to);
    red[to].push_back(from);
}
```

```
void path(int s) {
    fill(all(seen), 0);
    fill(all(dist), INF);
    dist[s] = 0; ll di;
```

```
__gnu_pbds::priority_queue<pair<ll, int>> q;
vector<decltype(q)::point_iterator> its(N);
q.push({0, s});
```

```
auto relax = [&](int i, ll cap, ll cost, int dir) {
    ll val = di - pi[i] + cost;
    if (cap && val < dist[i]) {
        dist[i] = val;
        par[i] = {s, dir};
        if (its[i] == q.end()) its[i] = q.push({-dist[i], i});
        else q.modify(its[i], {-dist[i], i});
    }
};
```

```
while (!q.empty()) {
    s = q.top().second; q.pop();
    seen[s] = 1; di = dist[s] + pi[s];
    for (int i : ed[s]) if (!seen[i])
        relax(i, cap[s][i] - flow[s][i], cost[s][i], 1);
    for (int i : red[s]) if (!seen[i])
        relax(i, flow[i][s], -cost[i][s], 0);
}
rep(i,0,N) pi[i] = min(pi[i] + dist[i], INF);
}
```

```
pair<ll, ll> maxflow(int s, int t) {
    ll totflow = 0, totcost = 0;
    while (path(s), seen[t]) {
        ll fl = INF;
        for (int p,r,x = t; tie(p,r) = par[x], x != s; x = p)
            fl = min(fl, r ? cap[p][x] - flow[p][x] : flow[x][p]);
        totflow += fl;
        for (int p,r,x = t; tie(p,r) = par[x], x != s; x = p)
            if (r) flow[p][x] += fl;
            else flow[x][p] -= fl;
    }
    rep(i,0,N) rep(j,0,N) totcost += cost[i][j] * flow[i][j];
    return {totflow, totcost};
}
```

```
// If some costs can be negative, call this before maxflow:
void setpi(int s) { // (otherwise, leave this out)
    fill(all(pi), INF); pi[s] = 0;
    int it = N, ch = 1; ll v;
    while (ch-- && it--)
        rep(i,0,N) if (pi[i] != INF)
            for (int to : ed[i]) if (cap[i][to])
                if ((v = pi[i] + cost[i][to]) < pi[to])
                    pi[to] = v, ch = 1;
        assert(it >= 0); // negative cost cycle
    }
};
```

3.3 Matching

hopcroftKarp.h

**Description:** Fast bipartite matching algorithm. Graph  $g$  should be a list of neighbors of the left partition, and  $btoa$  should be a vector full of -1's of the same size as the right partition. Returns the size of the matching.  $btoa[i]$  will be the match for vertex  $i$  on the right side, or  $-1$  if it's not matched.

**Usage:** VI btoa(m, -1); hopcroftKarp(g, btoa);

**Time:**  $\mathcal{O}(\sqrt{VE})$

fd54bf, 43 lines

```
bool dfs(int a, int L, vector<vi>& g, vi& btoa, vi& A, vi& B) {
    if (A[a] != L) return 0;
    A[a] = -1;
    for (int b : g[a]) if (B[b] == L + 1) {
        B[b] = 0;
        if (btoa[b] == -1 || dfs(btoa[b], L + 1, g, btoa, A, B))
            return btoa[b] = a, 1;
    }
    return 0;
}
```

```
int hopcroftKarp(vector<vi>& g, vi& btoa) {
    int res = 0;
    vi A(g.size()), B(btoa.size()), cur, next;
    for (;;) {
        fill(all(A), 0);
        fill(all(B), 0);
        cur.clear();
        for (int a : btoa) if(a != -1) A[a] = -1;
        rep(a,0,sz(g)) if(A[a] == 0) cur.push_back(a);
        for (int lay = 1;; lay++) {
            bool islast = 0;
            next.clear();
            for (int a : cur) for (int b : g[a]) {
                if (btoa[b] == -1) {
                    B[b] = lay;
                    islast = 1;
                }
                else if (btoa[b] != a && !B[b]) {
                    B[b] = lay;
                    next.push_back(btoa[b]);
                }
            }
            if (islast) break;
            if (next.empty()) return res;
            for (int a : next) A[a] = lay;
            cur.swap(next);
        }
        rep(a,0,sz(g))
            res += dfs(a, 0, g, btoa, A, B);
    }
    return sz(btoa) - (int)count(all(btoa), -1);
}
```

MinimumVertexCover.h

**Description:** Finds a minimum vertex cover in a bipartite graph. The size is the same as the size of a maximum matching, and the complement is a maximum independent set.

"hopcroftKarp.h" bfb654, 20 lines

```
vi cover(vector<vi>& g, int n, int m) {
    vi match(m, -1);
    int res = hopcroftKarp(g, match);
    vector<bool> lfound(n, true), seen(m);
    for (int it : match) if (it != -1) lfound[it] = false;
    vi q, cover;
    rep(i,0,n) if (lfound[i]) q.push_back(i);
    while (!q.empty()) {
        int i = q.back(); q.pop_back();
        lfound[i] = 1;
        for (int e : g[i]) if (!seen[e] && match[e] != -1) {
            seen[e] = true;
            q.push_back(match[e]);
        }
    }
    rep(i,0,n) if (!lfound[i]) cover.push_back(i);
    rep(i,0,m) if (seen[i]) cover.push_back(n+i);
    assert(sz(cover) == res);
    return cover;
}
```

WeightedMatching.h

**Description:** Given a weighted bipartite graph, matches every node on the left with a node on the right such that no nodes are in two matchings and the sum of the edge weights is minimal. Takes cost[N][M], where cost[i][j] = cost for L[i] to be matched with R[j] and returns (min cost, match), where L[i] is matched with R[match[i]]. Negate costs for max cost.

**Time:**  $\mathcal{O}(N^2M)$

```
pair<int, vi> hungarian(const vector<vi> &a) {
    if (a.empty()) return {0, {}};
    int n = sz(a) + 1, m = sz(a[0]) + 1;
    vi u(n), v(m), p(m), ans(n - 1);
    rep(i, 1, n) {
        p[0] = i;
        int j0 = 0; // add "dummy" worker 0
        vi dist(m, INT_MAX), pre(m, -1);
        vector<bool> done(m + 1);
        do { // dijkstra
            done[j0] = true;
            int i0 = p[j0], j1, delta = INT_MAX;
            rep(j, 1, m) if (!done[j]) {
                auto cur = a[i0 - 1][j - 1] - u[i0] - v[j];
                if (cur < dist[j]) dist[j] = cur, pre[j] = j0;
                if (dist[j] < delta) delta = dist[j], j1 = j;
            }
            rep(j, 0, m) {
                if (done[j]) u[p[j]] += delta, v[j] -= delta;
                else dist[j] -= delta;
            }
            j0 = j1;
        } while (p[j0]);
        while (j0) { // update alternating path
            int j1 = pre[j0];
            p[j0] = p[j1], j0 = j1;
        }
    }
    rep(j, 1, m) if (p[j]) ans[p[j] - 1] = j - 1;
    return {-v[0], ans}; // min cost
}
```

1e0fe9, 31 lines

3.4 DFS algorithms

2sat.h

**Description:** Calculates a valid assignment to boolean variables a, b, c,... to a 2-SAT problem, so that an expression of the type  $(a \vee b) \wedge \neg (a \vee c) \wedge (d \vee b) \wedge \dots$  becomes true, or reports that it is unsatisfiable. Negated variables are represented by bit-inversions ( $\sim x$ ).

**Usage:** TwoSat ts(number of boolean variables);  
ts.either(0, ~3); // Var 0 is true or var 3 is false  
ts.setValue(2); // Var 2 is true  
ts.atMostOne({0, ~1, 2}); //  $\leq 1$  of vars 0, ~1 and 2 are true  
ts.solve(); // Returns true iff it is solvable  
ts.values[0..N-1] holds the assigned values to the vars

**Time:**  $\mathcal{O}(N + E)$ , where N is the number of boolean variables, and E is the number of clauses.

```
struct TwoSat {
    int N;
    vector<vi> gr;
    vi values; // 0 = false, 1 = true

    TwoSat(int n = 0) : N(n), gr(2*n) {}

    int addVar() { // (optional)
        gr.emplace_back();
        gr.emplace_back();
        return N++;
    }

    void either(int f, int j) {
```

5f9706, 56 lines

```
f = max(2*f, -1-2*f);
j = max(2*j, -1-2*j);
gr[f].push_back(j^1);
gr[j].push_back(f^1);
}

void setValue(int x) { either(x, x); }

void atMostOne(const vi& li) { // (optional)
    if (sz(li) <= 1) return;
    int cur = ~li[0];
    rep(i, 2, sz(li)) {
        int next = addVar();
        either(cur, ~li[i]);
        either(cur, next);
        either(~li[i], next);
        cur = ~next;
    }
    either(cur, ~li[1]);
}

vi val, comp, z; int time = 0;
int dfs(int i) {
    int low = val[i] = ++time, x; z.push_back(i);
    for(int e : gr[i]) if (!comp[e])
        low = min(low, val[e] ? dfs(e));
    if (low == val[i]) do {
        x = z.back(); z.pop_back();
        comp[x] = low;
        if (values[x>>1] == -1)
            values[x>>1] = x&1;
    } while (x != i);
    return val[i] = low;
}

bool solve() {
    values.assign(N, -1);
    val.assign(2*N, 0); comp = val;
    rep(i, 0, 2*N) if (!comp[i]) dfs(i);
    rep(i, 0, N) if (comp[2*i] == comp[2*i+1]) return 0;
    return 1;
}

};
```

780b64, 15 lines

EulerWalk.h

**Description:** Eulerian undirected/directed path/cycle algorithm. Input should be a vector of (dest, global edge index), where for undirected graphs, forward/backward edges have the same index. Returns a list of nodes in the Eulerian path/cycle with src at both start and end, or empty list if no cycle/path exists. To get edge indices back, add .second to s and ret.

**Time:**  $\mathcal{O}(V + E)$

```
vi eulerWalk(vector<vector<pii>>& gr, int nedges, int src=0) {
    int n = sz(gr);
    vi D(n), its(n), eu(nedges), ret, s = {src};
    D[src]++; // to allow Euler paths, not just cycles
    while (!s.empty()) {
        int x = s.back(), y, e, &it = its[x], end = sz(gr[x]);
        if (it == end){ ret.push_back(x); s.pop_back(); continue; }
        tie(y, e) = gr[x][it++];
        if (!eu[e]) {
            D[x]--, D[y]++;
            eu[e] = 1; s.push_back(y);
        }
    }
    for (int x : D) if (x < 0 || sz(ret) != nedges+1) return {};
    return {ret.rbegin(), ret.rend()};
}
```

CondensationGraph.h

**Description:** Finds strongly connected components in a directed graph. If vertices  $u, v$  belong to the same component, we can reach  $u$  from  $v$  and vice versa.

**Usage:** scc(graph, [&](VI& v) { ... }) visits all components in reverse topological order. comp[i] holds the component index of a node (a component only has edges to components with lower index). ncomp will contain the number of components.

**Time:**  $\mathcal{O}(E + V)$

```
// 0 based indexing
void condense(vector<vi> adj, vector<vi> &adj_scc,
              vector<vi> &comp, vi &root_of, int n) {
    vector<vi> rev_adj(n);
    rep(u, 0, n) {
        for(auto v : adj[u]) {
            rev_adj[v].push_back(u);
        }
    }
    vector<bool> vis(n, false); vi order, component, root_nodes;
    function<void(int)> dfs1 = [&](int x) {
        vis[x] = true;
        for(auto nx : adj[x]) {
            if(!vis[nx]) {
                dfs1(nx);
            }
            order.push_back(x);
        }
    };
    rep(i, 0, n) { if(!vis[i]) dfs1(i); }
    vis.clear(); vis.assign(n, false);
    // order is now kind of topologically sorted
    reverse(order.begin(), order.end());
    function<void(int)> dfs2 = [&](int x) {
        vis[x] = true;
        component.push_back(x);
        for(auto u : rev_adj[x]) {
            if(!vis[u]) {
                dfs2(u);
            }
        }
    };
    comp.clear(); comp.resize(n);
    root_of.clear(); root_of.resize(n);
    for(auto v : order) {
        if(!vis[v]) {
            dfs2(v);
            int root = component.front();
            for(auto u : component) root_of[u] = root;
            root_nodes.push_back(root);
            comp[root] = component;
            component.clear();
        }
    }
    adj_scc.clear(); adj_scc.resize(n);
    rep(u, 0, n) {
        for(auto v : adj[u]) {
            if(root_of[u] != root_of[v]) {
                adj_scc[root_of[u]].push_back(root_of[v]);
            }
        }
    }
}
```

f5e4c5, 53 lines

BridgeTree.h

**Description:** Finds all biconnected components in an undirected graph, and runs a callback for the edges in each. In a biconnected component there are at least two distinct paths between any two nodes. Note that a node can be in several components. An edge which is not in a component is a bridge, i.e., not part of any cycle.

```
5e0207, 31 lines
// 0 based indexing
int n, m, Tin;
vector<vii> adj, adjn;
vi vis, low;
vector<array<int, 3>> bridges;
Dsu<int> ds;

int dfs0(int x,int p=-1,int w=0) {
    vis[x] = 1; low[x] = Tin++;
    int crl = low[x];
    for(auto nx : adj[x]) {
        if(nx.ff == p) continue;
        else if (vis[nx.ff]) crl = min(crl, low[nx.ff]);
        else crl = min(crl, dfs0(nx.ff, x, nx.ss));
    }
    if(crl == low[x] and p != -1) bridges.pb({x, p, w});
    else if (p != -1) ds.join(x, p);
    return crl;
}

void build_bridgetree() {
    // CLEAR global variables
    ds.build(n); // INITIALIZE DSU HERE
    rep(i,0,n) if(!vis[i]) dfs0(i);
    for(auto arr : bridges) {
        int u = ds.find(arr[0]), v = ds.find(arr[1]), w = arr[2];
        if(u != v) {
            adjn[v].pb({u, w}); adjn[u].pb({v, w});
        }
    }
}
```

3.5 Coloring

EdgeColoring.h

**Description:** Given a simple, undirected graph with max degree  $D$ , computes a  $(D + 1)$ -coloring of the edges such that no neighboring edges share a color. ( $D$ -coloring is NP-hard, but can be done for bipartite graphs by repeated matchings of max-degree nodes.)

**Time:**  $\mathcal{O}(NM)$

```
e210e2, 31 lines
vi edgeColoring(int N, vector<pii> eds) {
    vi cc(N + 1), ret(sz(eds)), fan(N), free(N), loc;
    for (pii e : eds) ++cc[e.first], ++cc[e.second];
    int u, v, ncols = *max_element(all(cc)) + 1;
    vector<vi> adj(N, vi(ncols, -1));
    for (pii e : eds) {
        tie(u, v) = e;
        fan[0] = v;
        loc.assign(ncols, 0);
        int at = u, end = u, d, c = free[u], ind = 0, i = 0;
        while (d = free[v], !loc[d] && (v = adj[u][d]) != -1)
            loc[d] = ++ind, cc[ind] = d, fan[ind] = v;
        cc[loc[d]] = c;
        for (int cd = d; at != -1; cd ^= c ^ d, at = adj[at][cd])
            swap(adj[at][cd], adj[end = at][cd ^ c ^ d]);
        while (adj[fan[i]][d] != -1) {
            int left = fan[i], right = fan[++i], e = cc[i];
            adj[u][e] = left;
            adj[left][e] = u;
            adj[right][e] = -1;
            free[right] = e;
        }
        adj[u][d] = fan[i];
    }
```

```
adj[fan[i]][d] = u;
for (int y : {fan[0], u, end})
    for (int& z = free[y] = 0; adj[y][z] != -1; z++);
}
rep(i,0,sz(eds))
    for (tie(u, v) = eds[i]; adj[u][ret[i]] != v;) ++ret[i];
return ret;
}
```

3.6 Heuristics

MaximalCliques.h

**Description:** Runs a callback for all maximal cliques in a graph (given as a symmetric bitset matrix; self-edges not allowed). Callback is given a bitset representing the maximal clique.

**Time:**  $\mathcal{O}\left(3^{n/3}\right)$ , much faster for sparse graphs

```
b0d5b1, 12 lines
typedef bitset<128> B;
template<class F>
void cliques(vector<B>& eds, F f, B P = ~B(), B X={}, B R={}) {
    if (!P.any()) { if (!X.any()) f(R); return; }
    auto q = (P | X)._Find_first();
    auto cands = P & ~eds[q];
    rep(i,0,sz(eds)) if (cands[i]) {
        R[i] = 1;
        cliques(eds, f, P & eds[i], X & eds[i], R);
        R[i] = P[i] = 0; X[i] = 1;
    }
}
```

MaximumClique.h

**Description:** Quickly finds a maximum clique of a graph (given as symmetric bitset matrix; self-edges not allowed). Can be used to find a maximum independent set by finding a clique of the complement graph.

**Time:** Runs in about 1s for  $n=155$  and worst case random graphs ( $p=.90$ ). Runs faster for sparse graphs.

```
f7c0bc, 49 lines
typedef vector<bitset<200>> vb;
struct Maxclique {
    double limit=0.025, pk=0;
    struct Vertex { int i, d=0; };
    typedef vector<Vertex> vv;
    vb e;
    vv V;
    vector<vi> C;
    vi qmax, q, S, old;
    void init(vv& r) {
        for (auto& v : r) v.d = 0;
        for (auto& v : r) for (auto j : r) v.d += e[v.i][j.i];
        sort(all(r), [](auto a, auto b) { return a.d > b.d; });
        int mxD = r[0].d;
        rep(i,0,sz(r)) r[i].d = min(i, mxD) + 1;
    }
    void expand(vv& R, int lev = 1) {
        S[lev] += S[lev - 1] - old[lev];
        old[lev] = S[lev - 1];
        while (sz(R)) {
            if (sz(q) + R.back().d <= sz(qmax)) return;
            q.push_back(R.back().l);
            vv T;
            for(auto v:R) if (e[R.back().i][v.i]) T.push_back({v.i});
            if (sz(T)) {
                if (S[lev]++ / ++pk < limit) init(T);
                int j = 0, mxk = 1, mnk = max(sz(qmax) - sz(q) + 1, 1);
                C[1].clear(), C[2].clear();
                for (auto v : T) {
                    int k = 1;
                    auto f = [&](int i) { return e[v.i][i]; };
                    while (any_of(all(C[k]), f)) k++;
                }
```

```
if (k > mxk) mxk = k, C[mxk + 1].clear();
if (k < mnk) T[j++].i = v.i;
C[k].push_back(v.i);
}
if (j > 0) T[j - 1].d = 0;
rep(k,mnk,mxk + 1) for (int i : C[k])
    T[j].i = i, T[j++].d = k;
expand(T, lev + 1);
} else if (sz(q) > sz(qmax)) qmax = q;
q.pop_back(), R.pop_back();
}
}
vi maxClique() { init(V), expand(V); return qmax; }
Maxclique(vb conn) : e(conn), C(sz(e)+1), S(sz(C)), old(S) {
    rep(i,0,sz(e)) V.push_back({i});
}
};
```

MaximumIndependentSet.h

**Description:** To obtain a maximum independent set of a graph, find a max clique of the complement. If the graph is bipartite, see MinimumVertexCover.

3.7 Trees

BinaryLifting.h

**Description:** Initialize  $\text{par}[N][\log N]$  and  $\text{depth}[N]$  array.  $\text{kthpar}$ :  $\log N$ ,  $\text{lca}$ :

```
5e6ba9, 25 lines
void dfs(int x,int pr=0,int d=0) {
    depth[x] = d; par[x][0] = pr;
    rep(i,1,logN) par[x][i]=par[par[x][i-1]][i-1];
    for(auto next : adj[x])
        if(next != pr)
            dfs(next, x, d+1);
}

int kthpar(int x,int k) {
    int ret = x;
    rep(i, 0, logN) if((k>>i)&1) ret = par[ret][i];
    return ret;
}

int lca(int u,int v) {
    if(depth[u] > depth[v]) swap(u,v);
    v = kthpar(v, depth[v] - depth[u]);
    if(u == v) return u;
    for(int i=logN-1;i>=0;i--) {
        if(par[u][i] != par[v][i]) {
            u=par[u][i]; v=par[v][i];
        }
    }
    return par[u][0];
}
```

LCA.h

**Description:** Push  $(\text{time}[\text{node}], \text{node})$  in a vector every time you iterate over an edge  $(\text{node}, x)$ .  $\text{lca}(a, b) \Rightarrow \text{rangeMin}(\text{time}[a], \text{time}[b])$

**Time:**  $\mathcal{O}(N \log N + Q)$

LinkCutTree.h

**Description:** Represents a forest of unrooted trees. You can add and remove edges (as long as the result is still a forest), and check whether two nodes are in the same tree.

**Time:** All operations take amortized  $\mathcal{O}(\log N)$ .

```
5909e2, 90 lines
struct Node { // Splay tree. Root's pp contains tree's parent.
    Node *p = 0, *pp = 0, *c[2];
};
```



```

bool flip = 0;
Node() { c[0] = c[1] = 0; fix(); }
void fix() {
    if (c[0]) c[0]->p = this;
    if (c[1]) c[1]->p = this;
    // (+ update sum of subtree elements etc. if wanted)
}
void pushFlip() {
    if (!flip) return;
    flip = 0; swap(c[0], c[1]);
    if (c[0]) c[0]->flip ^= 1;
    if (c[1]) c[1]->flip ^= 1;
}
int up() { return p ? p->c[1] == this : -1; }
void rot(int i, int b) {
    int h = i ^ b;
    Node *x = c[i], *y = b == 2 ? x : x->c[h], *z = b ? y : x;
    if ((y->p = p)) p->c[up()] = y;
    c[i] = z->c[i ^ 1];
    if (b < 2) {
        x->c[h] = y->c[h ^ 1];
        z->c[h ^ 1] = b ? x : this;
    }
    y->c[i ^ 1] = b ? this : x;
    fix(); x->fix(); y->fix();
    if (p) p->fix();
    swap(pp, y->pp);
}
void splay() {
    for (pushFlip(); p; ) {
        if (p->p) p->p->pushFlip();
        p->pushFlip(); pushFlip();
        int c1 = up(), c2 = p->up();
        if (c2 == -1) p->rot(c1, 2);
        else p->p->rot(c2, c1 != c2);
    }
}
Node* first() {
    pushFlip();
    return c[0] ? c[0]->first() : (splay(), this);
}
};

struct LinkCut {
    vector<Node> node;
    LinkCut(int N) : node(N) {}

    void link(int u, int v) { // add an edge (u, v)
        assert(!connected(u, v));
        makeRoot(&node[u]);
        node[u].pp = &node[v];
    }
    void cut(int u, int v) { // remove an edge (u, v)
        Node *x = &node[u], *top = &node[v];
        makeRoot(top); x->splay();
        assert(top == (x->pp ? x->c[0]));
        if (x->pp) x->pp = 0;
        else {
            x->c[0] = top->p = 0;
            x->fix();
        }
    }
    bool connected(int u, int v) { // are u, v in the same tree?
        Node* nu = access(&node[u])->first();
        return nu == access(&node[v])->first();
    }
}
void makeRoot(Node* u) {
    access(u);
    u->splay();
}

```

```

    if(u->c[0]) {
        u->c[0]->p = 0;
        u->c[0]->flip ^= 1;
        u->c[0]->pp = u;
        u->c[0] = 0;
        u->fix();
    }
}
Node* access(Node* u) {
    u->splay();
    while (Node* pp = u->pp) {
        pp->splay(); u->pp = 0;
        if (pp->c[1]) {
            pp->c[1]->p = 0; pp->c[1]->pp = pp; }
        pp->c[1] = u; pp->fix(); u = pp;
    }
    return u;
}
};

```

### DirectedMST.h

**Description:** Finds a minimum spanning tree/arborescence of a directed graph, given a root node. If no MST exists, returns -1.

**Time:**  $\mathcal{O}(E \log V)$

```

"/data-structures/UnionFindRollback.h" 39e620, 60 lines

struct Edge { int a, b; ll w; };
struct Node {
    Edge key;
    Node *l, *r;
    ll delta;
    void prop() {
        key.w += delta;
        if (l) l->delta += delta;
        if (r) r->delta += delta;
        delta = 0;
    }
    Edge top() { prop(); return key; }
};
Node *merge(Node *a, Node *b) {
    if (!a || !b) return a ? b;
    a->prop(), b->prop();
    if (a->key.w > b->key.w) swap(a, b);
    swap(a->l, (a->r = merge(b, a->r)));
    return a;
}
void pop(Node*& a) { a->prop(); a = merge(a->l, a->r); }

pair<ll, vi> dmst(int n, int r, vector<Edge>& g) {
    RollbackUF uf(n);
    vector<Node*> heap(n);
    for (Edge e : g) heap[e.b] = merge(heap[e.b], new Node{e});
    ll res = 0;
    vi seen(n, -1), path(n, par(n));
    seen[r] = r;
    vector<Edge> Q(n), in(n, {-1,-1}), comp;
    deque<tuple<int, int, vector<Edge>>> cycts;
    rep(s,0,n) {
        int u = s, qi = 0, w;
        while (seen[u] < 0) {
            if (!heap[u]) return {-1,{};};
            Edge e = heap[u]->top();
            heap[u]->delta -= e.w, pop(heap[u]);
            Q[qi] = e, path[qi++] = u, seen[u] = s;
            res += e.w, u = uf.find(e.a);
            if (seen[u] == s) {
                Node* cyc = 0;
                int end = qi, time = uf.time();
                do cyc = merge(cyc, heap[w = path[--qi]]);
                while (uf.join(u, w));
            }
        }
    }
}

```

```

        u = uf.find(u), heap[u] = cyc, seen[u] = -1;
        cycts.push_front({u, time, {&Q[qi], &Q[end]}});
    }
}
rep(i,0,qi) in[uf.find(Q[i].b)] = Q[i];
}

for (auto& [u,t,comp] : cycts) { // restore sol (optional)
    uf.rollback(t);
    Edge inEdge = in[u];
    for (auto& e : comp) in[uf.find(e.b)] = e;
    in[uf.find(inEdge.b)] = inEdge;
}
rep(i,0,n) par[i] = in[i].a;
return {res, par};
}

```

### HLD.h

**Description:** Heavy Light Decomposition

**Time:**  $\mathcal{O}(\log N * \text{TimetakenbyRangeQueryDS})$

3a0a27, 74 lines

```

// requires a segment tree with init function
class HLD {
    SegmentTrees sgt; vector<vi> adj;
    vi sz, par, head, sc, st, ed;
    int t, n;
public:
    HLD(vector<vector<int>>& adj1,int n1): sz(n1+1), par(n1+1),
        head(n1+1), sc(n1+1), st(n1+1), ed(n1+1) {
        n = n1; adj = adj1; t = 0;
    }
    void dfs_sz(int x,int p = 0) {
        sz[x] = 1; par[x] = p; head[x] = x;
        for(auto nx : adj[x]) {
            if(nx == p) continue;
            dfs_sz(nx, x);
            sz[x] += sz[nx];
            if(sz[nx] > sz[sc[x]]) sc[x] = nx;
        }
    }
    void dfs_hld(int x,int p = 0) {
        st[x] = t++;
        if(sc[x]) {
            head[sc[x]] = head[x];
            dfs_hld(sc[x], x);
        }
        for(auto nx : adj[x]) {
            if(nx == p or nx == sc[x]) continue;
            dfs_hld(nx, x);
        }
        ed[x] = t - 1;
    }
    void build(int base = 1) {
        dfs_sz(base);
        dfs_hld(base);
        sgt.init(t);
    }
    bool anc(int x,int y) {
        if(x == 0) return true; if(y == 0) return false;
        return (st[x] <= st[y] and ed[x] >= ed[y]);
    }
    int lca(int x,int y) {
        if(anc(x, y)) return x; if(anc(y, x)) return y;
        while(!anc(par[head[x]], y)) x = par[head[x]];
        while(!anc(par[head[y]], x)) y = par[head[y]];
        x = par[head[x]]; y = par[head[y]];
        // one will overshoot the lca and the other will reach lca.
        return anc(x, y) ? y : x;
    }
}

```

```

}
void update_up(int x,int p,ll add) {
    while(head[x] != head[p]) {
        sgt.update(st[head[x]], st[x], add, 0, t-1);
        x = par[head[x]];
    }
    sgt.update(st[p], st[x], add, 0, t - 1);
}
void range_update(int u,int v,T add) {
    int l = lca(u, v);
    update_up(u, l, add); update_up(v, l, add);
    update_up(l, l, -add);
}
T query_up(int x,int p) {
    T ans = 0;
    while(head[x] != head[p]) {
        ans = min(ans, sgt.query(st[head[x]], st[x], 0, t - 1));
        x = par[head[x]];
    }
    ans = min(ans, sgt.query(st[p], st[x], 0, t - 1));
    return ans;
}
T range_min(int u,int v) {
    int l = lca(u, v);
    return min(query_up(u, l), query_up(v, l));
}
};

```

## CentroidDecomposition.h

Time:  $\mathcal{O}(N \log N + Q)$

ca0a79, 59 lines

```

const int N = 5e4 + 13, logN = 17;
vi adj[N], sub(N), par(N,-1), lvl(N), done(N), par_adj(N);
vector<vi> dist(N, vi(logN, 0)), anc(N, vi(logN, 0));
int nn = 0, root;
void dfs_size(int x,int p) {
    nn++; sub[x] = 1;
    for(auto nx : adj[x]) if(!done[nx] and nx != p) {
        dfs_size(nx, x); sub[x] += sub[nx];
    }
}
int find_ct(int x,int p) {
    for(auto nx : adj[x]) if(!done[nx] and nx != p and sub[nx] > nn/2)
        return find_ct(nx, x);
    return x;
}
void dfs(int x,int p,int ct) {
    anc[x][lvl[ct]] = ct;
    for(auto nx : adj[x]) if(!done[nx] and nx != p) {
        dist[nx][lvl[ct]] = 1 + dist[x][lvl[ct]];
        dfs(nx, x, ct);
    }
}
// par_adj[ct] = adjacent vertex to parent of ct in OT in subtree of ct.
int decompose(int x,int p=-1) {
    nn = 0; dfs_size(x, x);
    int ct = find_ct(x, x);
    if(p) lvl[ct] = 1 + lvl[p];
    done[ct] = 1; par[ct] = p;
    dfs(ct, ct, ct);
    for(auto nx : adj[ct]) if(!done[nx]) {
        int nct = decompose(nx, ct);
        par_adj[nct] = nx;
    }
    return ct;
}

```

```

vector<vi> child_cntb(N), my(N);
rep(x,0,n) for(int y = x; y >= 0; y = par[y]) {
    my[y].pb(dist[x][lvl[y]]);
    if(par[y] >= 0)
        child_cntb[y].pb(dist[x][lvl[par[y]]]);
}
rep(x,0,n) {
    sort(all(my[x])); sort(all(child_cntb[x]));
}
// number of nodes <= k in v.
auto cnt_k = [&](vi &v,int k) {
    int l = upper_bound(all(v), k) - v.begin();
    return l;
};
auto k_dists = [&](int x,int k) {
    int ans = cnt_k(my[x], k);
    int ch = x, q = x; x = par[x];
    while(x >= 0) {
        ans += (cnt_k(my[x], k - dist[q][lvl[x]]));
        ans -= (cnt_k(child_cntb[ch], k - dist[q][lvl[x]]));
        ch = x; x = par[x];
    }
    return ans;
};

```

## AuxiliaryTrees.h

Description: Creates a auxiliary tree of  $k$  nodes.

Time:  $\mathcal{O}(k)$

1b4c5b, 62 lines

```

using vvi = vector<vector<int>>
struct Tree {
    int n;
    vvi adj;
    vi pos, tour, depth, pos_end, max_depth, dp, max_up;
    Tree(int n) : n(n), adj(n), max_depth(n), dp(n), max_up(n) {}
    void add_edge(int s, int t) {
        adj[s].pb(t); adj[t].pb(s);
    }
    vvi table;
    int argmin(int i, int j) { return depth[i] < depth[j] ? i : j; }
    void rootify(int r) {
        pos.resize(n); pos_end.resize(n);
        function<void (int,int,int)> dfs = [&](int u, int p, int d)
        {
            pos[u] = pos_end[u] = depth.size();
            tour.pb(u); depth.pb(d);
            for (int v: adj[u]) {
                if (v != p) {
                    dfs(v, u, d+1);
                    pos_end[u] = depth.size();
                    tour.pb(u);
                    depth.pb(d);
                }
            }
        }; dfs(r, r, 0);
        int logn = sizeof(int)*__CHAR_BIT__-1-__builtin_clz(tour.size()); // log2
        table.resize(logn+1, vi(tour.size()));
        iota(all(table[0]), 0);
        for (int h = 0; h < logn; ++h)
            for (int i = 0; i+(1<<h) < tour.size(); ++i)
                table[h+1][i] = argmin(table[h][i], table[h][i+(1<<h)])
                ;
    }
    int lca(int u, int v) {
        int i = pos[u], j = pos[v]; if (i > j) swap(i, j);
        int h = sizeof(int)*__CHAR_BIT__-1-__builtin_clz(j-i); // = log2
    }
}

```

```

    return i == j ? u : tour[argmin(table[h][i], table[h][j] - (1<<h))];
}
int getDepth(int u){
    return depth[pos[u]];
}
void aux_Tree(vi nodes, vvi &adj_aux, vi &start_times){
    // adj_aux stores the children
    for(int x : nodes) start_times.pb(pos[x]);
    sort(all(start_times));
    for(int i = 1; i < (int) nodes.size(); i++){
        start_times.pb(pos[lca(tour[start_times[i]], tour[start_times[i - 1]])]);
    }
    sort(all(start_times));
    start_times.erase(unique(start_times.begin(), start_times.end()), start_times.end());
    adj_aux.resize(start_times.size());
    stack<int> st;
    // nodes now indexed according to start_times
    st.push(0);
    for(int i = 1; i < (int) start_times.size(); i++){
        while(pos_end[tour[start_times[st.top()]]] < start_times[i]){
            st.pop();
        }
        adj_aux[st.top()].pb(i);
        st.push(i);
    }
}
};

```

## Blossom.h

Description: Blossom Algorithm

1b2a6f, 52 lines

```

vector<int> Blossom(vector<vector<int>>& graph) {
    int n = graph.size(), timer = -1;
    vector<int> mate(n, -1), label(n), parent(n), orig(n), aux(n, -1), q;
    auto lca = [&](int x, int y) {
        for (timer++; ; swap(x, y)) {
            if (x == -1) continue;
            if (aux[x] == timer) return x;
            aux[x] = timer;
            x = (mate[x] == -1 ? -1 : orig[parent[mate[x]]]);
        }
    };
    auto blossom = [&](int v, int w, int a) {
        while (orig[v] != a) {
            parent[v] = w; w = mate[v];
            if (label[w] == 1) label[w] = 0, q.push_back(w);
            orig[v] = orig[w] = a; v = parent[w];
        }
    };
    auto augment = [&](int v) {
        while (v != -1) {
            int pv = parent[v], nv = mate[pv];
            mate[v] = pv; mate[pv] = v; v = nv;
        }
    };
    auto bfs = [&](int root) {
        fill(label.begin(), label.end(), -1);
        iota(orig.begin(), orig.end(), 0);
        q.clear();
        label[root] = 0; q.push_back(root);
        for (int i = 0; i < (int)q.size(); ++i) {
            int v = q[i];
            for (auto x : graph[v]) {
                if (label[x] == -1) {

```

```
label[x] = 1; parent[x] = v;
if (mate[x] == -1)
    return augment(x), 1;
label[mate[x]] = 0; q.push_back(mate[x]);
} else if (label[x] == 0 && orig[v] != orig[x]) {
    int a = lca(orig[v], orig[x]);
    blossom(x, v, a); blossom(v, x, a);
}
}
}
return 0;
};
// Time halves if you start with (any) maximal matching.
for (int i = 0; i < n; i++)
    if (mate[i] == -1)
        bfs(i);
return mate;
}
```

3.8 Math

**Number of Spanning Trees** Create an  $N \times N$  matrix  $\text{mat}$ , and for each edge  $a \rightarrow b \in G$ , do  $\text{mat}[a][b]--$ ,  $\text{mat}[b][b]++$  (and  $\text{mat}[b][a]--$ ,  $\text{mat}[a][a]++$  if  $G$  is undirected). Remove the  $i$ th row and column and take the determinant; this yields the number of directed spanning trees rooted at  $i$  (if  $G$  is undirected, remove any row/column).

**Mirsky’s Theorem** Max length chain is equal to min partitioning into antichains. Max chain is height of poset.

**Dilworth’s Theorem** Min partition into chains is equal to max length antichain. From poset create bipartite graph. Any edge from  $v_i - v_j$  implies  $LV_i - RV_j$ . Let  $A$  be the set of vertices such that neither  $LV_i$  nor  $RV_i$  are in vertex cover.  $A$  is an antichain of size  $n$ -max matching. To get min partition into chains, take a vertex from left side, keep taking vertices till a matching exist. Consider this as a chain. Its size is  $n$  - max matching.

**Matrix-tree Theorem** Let matrix  $T = [t_{ij}]$ , where  $t_{ij}$  is negative of the number of multiedges between  $i$  and  $j$ , for  $i \neq j$ , and  $t_{ii} = \deg_i$ . Number of spanning trees of a graph is equal to the determinant of a matrix obtained by deleting any  $k$ -th row and  $k$ -th column from  $T$ . If  $G$  is a multigraph and  $e$  is an edge of  $G$ , then the number  $\tau(G)$  of spanning trees of  $G$  satisfies recurrence  $\tau(G) = \tau(G - e) + \tau(G/e)$ , when  $G - e$  is the multigraph obtained by deleting  $e$ , and  $G/e$  is the contraction of  $G$  by  $e$  (multiple edges arising from the contraction are preserved.)

Hashing Kmp Manacher

**Cycle Spaces** The (binary) cycle space of an undirected graph is the set of its Eulerian subgraphs. This set of subgraphs can be described algebraically as a vector space over the two-element finite field. One way of constructing a cycle basis is to form a spanning forest of the graph, and then for each edge  $e$  that does not belong to the forest, form a cycle  $C \ C_e$  consisting of  $e$  together with the path in the forest connecting the endpoints of  $e$ . The set of cycles  $C_e$  formed in this way are linearly independent (each one contains an edge  $e$  that does not belong to any of the other cycles) and has the correct size  $m - n + c$  to be a basis, so it necessarily is a basis. This is fundamental cycle basis.

**Cut Spaces** The family of all cut sets of an undirected graph is known as the cut space of the graph. It forms a vector space over the two-element finite field of arithmetic modulo two, with the symmetric difference of two cut sets as the vector addition operation, and is the orthogonal complement of the cycle space. To compute the basis vector for the cut space, consider any spanning tree of the graph. For every edge  $e$  in the spanning tree, remove the edge and consider the cut formed. Thus dimension of the basis vector for cut space is  $n-1$ .

**Number of perfect matchings of a bipartite graph** is equal to the permanent of the adjacency matrix obtained. To check the parity of the number of perfect matchings, we can evaluate the permanent of the matrix in  $Z_2$  which can be done easily as  $\text{Permanent}(A) = \text{Determinant}(A)$ .

**Tutte Matrix.** For a simple undirected graph  $G$ , Let  $M$  be a matrix with entries  $A_{i,j} = 0$  if  $(i,j) \notin E$  and  $A_{i,j} = -A_{j,i} = X$  if  $(i,j) \in E$ .  $X$  could be any random value. If the determinants are non-zero, then a perfect matching exists, while other direction might not hold for very small probability.

**Kirchhoff’s Theorem.** For a multigraph  $G$  with no loops, define Laplacian matrix as  $L = D - A$ .  $D$  is a diagonal matrix with  $D_{i,i} = \deg(i)$ , and  $A$  is an adjacency matrix. If you remove any row and column of  $L$ , the determinant gives a number of spanning trees.

**Brook’s Theorem** If a graph is not a complete graph or an odd cycle then it can be coloured with max degree  $\#$  of colours.

Strings (4)

```
Hashing.h
Description: Various self-explanatory methods for string hashing. Use on
Codeforces, which lacks 64-bit support and where solutions can be hacked.
<sys/time.h> eb5e9e, 36 lines
typedef uint64_t ull;
static int C; // initialized below

// Arithmetic mod two primes and 2^32 simultaneously.
// "typedef uint64_t H;" instead if Thue-Morse does not apply.
template<int M, class B>
```

```
struct A {
    int x; B b; A(int x=0) : x(x), b(b) {}
    A(int x, B b) : x(x), b(b) {}
    A operator+(A o){int y = x+o.x; return{y - (y>=M)*M, b+o.b};}
    A operator-(A o){int y = x-o.x; return{y + (y< 0)*M, b-o.b};}
    A operator*(A o) { return {(int)(1LL*x*o.x % M), b*o.b}; }
    explicit operator ull() { return x ^ (ull) b << 21; }
};
typedef A<10000000007, A<10000000009, unsigned>> H;

struct HashInterval {
    vector<H> ha, pw;
    HashInterval(string& str) : ha(SZ(str)+1), pw(ha) {
        pw[0] = 1;
        rep(i,0,sz(str))
            ha[i+1] = ha[i] * C + str[i],
            pw[i+1] = pw[i] * C;
    }
    H hashInterval(int a, int b) { // hash [a, b)
        return ha[b] - ha[a] * pw[b - a];
    }
};

int main() {
    timeval tp;
    gettimeofday(&tp, 0);
    C = (int)tp.tv_usec; // (less than modulo)
    assert((ull)(H(1)*2+1-3) == 0);
    // ...
}
```

Kmp.h

**Description:**  $\text{pi}[x]$  computes the length of the longest prefix of  $s$  that ends at  $x$ , other than  $s[0...x]$  itself (abacaba -> 0010123). Can be used to find all occurrences of a string.  
**Time:**  $\mathcal{O}(n)$

```
vi pi(const string& s) {
    vi p(sz(s));
    rep(i,1,sz(s)) {
        int g = p[i-1];
        while (g && s[i] != s[g]) g = p[g-1];
        p[i] = g + (s[i] == s[g]);
    }
    return p;
}
```

```
vi match(const string& s, const string& pat) {
    vi p = pi(pat + '\0' + s), res;
    rep(i,sz(p)-sz(s),sz(p))
        if (p[i] == sz(pat)) res.push_back(i - 2 * sz(pat));
    return res;
}
```

Manacher.h

**Description:** For each position in a string, computes  $\text{p}[0][i] = \text{half length of longest even palindrome around pos } i$ ,  $\text{p}[1][i] = \text{longest odd (half rounded down)}$ .  
**Time:**  $\mathcal{O}(N)$

```
array<vi, 2> manacher(const string& s) {
    int n = sz(s);
    array<vi,2> p = {vi(n+1), vi(n)};
    rep(z,0,2) for (int i=0,l=0,r=0; i < n; i++) {
        int t = r-i!z;
        if (i<r) p[z][i] = min(t, p[z][l+t]);
        int L = i-p[z][i], R = i+p[z][i]-!z;
        while (L>=1 && R+1<n && s[L-1] == s[R+1])
            p[z][i]++, L--, R++;
    }
```

```
    if (R>r) l=L, r=r;
  }
  return p;
}
```

MinRotation.h

**Description:** Finds the lexicographically smallest rotation of a string.  
**Usage:** rotate(v.begin(), v.begin()+minRotation(v), v.end());  
**Time:**  $\mathcal{O}(N)$

```
int minRotation(string s) {
    int a=0, N=sz(s); s += s;
    rep(b,0,N) rep(k,0,N) {
        if (a+k == b || s[a+k] < s[b+k]) {b += max(0, k-1); break;}
        if (s[a+k] > s[b+k]) { a = b; break; }
    }
    return a;
}
```

SuffixArray.h

**Description:** Builds suffix array for a string. sa[i] is the starting index of the suffix which is *i*'th in the sorted suffix array. The returned vector is of size  $n + 1$ , and sa[0] = n. The lcp array contains longest common prefixes for neighbouring strings in the suffix array: lcp[i] = lcp(sa[i], sa[i-1]), lcp[0] = 0. The input string must not contain any zero bytes.  
**Time:**  $\mathcal{O}(n \log n)$

```
struct SuffixArray {
    vi sa, lcp;
    SuffixArray(string& s, int lim=256) { // or basic_string<int>
        int n = sz(s) + 1, k = 0, a, b;
        vi x(all(s)+1), y(n), ws(max(n, lim)), rank(n);
        sa = lcp = y, iota(all(sa), 0);
        for (int j = 0, p = 0; p < n; j = max(1, j * 2), lim = p) {
            p = j, iota(all(y), n - j);
            rep(i,0,n) if (sa[i] >= j) y[p++] = sa[i] - j;
            fill(all(ws), 0);
            rep(i,0,n) ws[x[i]]++;
            rep(i,1,lim) ws[i] += ws[i - 1];
            for (int i = n; i--;) sa[--ws[x[y[i]]]] = y[i];
            swap(x, y), p = 1, x[sa[0]] = 0;
            rep(i,1,n) a = sa[i - 1], b = sa[i], x[b] =
                (y[a] == y[b] && y[a + j] == y[b + j]) ? p - 1 : p++;
        }
        rep(i,1,n) rank[sa[i]] = i;
        for (int i = 0, j; i < n - 1; lcp[rank[i++]] = k)
            for (k && k--, j = sa[rank[i] - 1];
                s[i + k] == s[j + k]; k++);
    }
};
```

Z.h

**Description:** z[x] computes the length of the longest common prefix of s[i:] and s, except z[0] = 0. (abacaba -> 0010301)  
**Time:**  $\mathcal{O}(n)$

```
vi Z(const string& S) {
    vi z(sz(S));
    int l = -1, r = -1;
    rep(i,1,sz(S)) {
        z[i] = i >= r ? 0 : min(r - i, z[i - l]);
        while (i + z[i] < sz(S) && S[i + z[i]] == S[z[i]])
            z[i]++;
        if (i + z[i] > r)
            l = i, r = i + z[i];
    }
    return z;
}
```

DynamicAhoCorasik.h

**Description:** Deletion happens by creating another aho corasik. Aho-Corasick automaton, used for multiple pattern matching. Initialize with AhoCorasick ac(patterns); the automaton start node will be at index 0. find(word) returns for each position the index of the longest word that ends there, or -1 if none. findAll(−, word) finds all words (up to  $N\sqrt{N}$  many if no duplicate patterns) that start at each position (shortest first). Duplicate patterns are allowed; empty patterns are not. To find the longest words that start at each position, reverse all input. For large alphabets, split each symbol into chunks, with sentinel bits for symbol boundaries.  
**Time:** construction takes  $\mathcal{O}(26N)$ , where  $N$  = sum of length of patterns. find(x) is  $\mathcal{O}(N)$ , where  $N$  = length of x. findAll is  $\mathcal{O}(NM)$ .

```
struct AhoCorasick {
    enum {alpha = 26, first = 'A'}; // change this!
    struct Node {
        // (nmatches is optional)
        int back, next[alpha], start = -1, end = -1, nmatches = 0;
        Node(int v) { memset(next, v, sizeof(next)); }
    };
    vector<Node> N;
    vi backp;
    void insert(string& s, int j) {
        assert(!s.empty());
        int n = 0;
        for (char c : s) {
            int& m = N[n].next[c - first];
            if (m == -1) { n = m = sz(N); N.emplace_back(-1); }
            else n = m;
        }
        if (N[n].end == -1) N[n].start = j;
        backp.push_back(N[n].end);
        N[n].end = j;
        N[n].nmatches++;
    }
    AhoCorasick(vector<string>& pat) : N(1, -1) {
        rep(i,0,sz(pat)) insert(pat[i], i);
        N[0].back = sz(N);
        N.emplace_back(0);

        queue<int> q;
        for (q.push(0); !q.empty(); q.pop()) {
            int n = q.front(), prev = N[n].back;
            rep(i,0,alpha) {
                int &ed = N[n].next[i], y = N[prev].next[i];
                if (ed == -1) ed = y;
                else {
                    N[ed].back = y;
                    (N[ed].end == -1 ? N[ed].end : backp[N[ed].start])
                        = N[y].end;
                    N[ed].nmatches += N[y].nmatches;
                    q.push(ed);
                }
            }
        }
    }
    vi find(string word) {
        int n = 0;
        vi res; // ll count = 0;
        for (char c : word) {
            n = N[n].next[c - first];
            res.push_back(N[n].end);
            // count += N[n].nmatches;
        }
        return res;
    }
    vector<vi> findAll(vector<string>& pat, string word) {
        vi r = find(word);
        vector<vi> res(sz(word));
```

```
        rep(i,0,sz(word)) {
            int ind = r[i];
            while (ind != -1) {
                res[i - sz(pat[ind]) + 1].push_back(ind);
                ind = backp[ind];
            }
            return res;
        }
    };

    vector<string> vc;
    vc.push_back(s);
    for(int i=0;i<LIM;i++) {
        if(ad[0][i].vs.size()>0) {
            for(auto x: ad[0][i].vs) {
                vc.push_back(x);
            }
            ad[0][i]=Aho();
        }
        else {
            for(auto x: vc) {
                ad[0][i].add(x);
            }
            ad[0][i].build_aho();
            break;
        }
    }
}
```

Number theory (5)

5.1 Modular arithmetic

ModInverse.h

**Description:** Pre-computation of modular inverses. Assumes  $\text{LIM} \leq \text{mod}$  and that mod is a prime.

```
const ll mod = 1000000007, LIM = 200000;
ll* inv = new ll[LIM] - 1; inv[1] = 1;
rep(i,2,LIM) inv[i] = mod - (mod / i) * inv[mod % i] % mod;
```

ModPow.h

```
const ll mod = 1000000007; // faster if const

ll modpow(ll b, ll e) {
    ll ans = 1;
    for (; e; b = b * b % mod, e /= 2)
        if (e & 1) ans = ans * b % mod;
    return ans;
}
```

ModLog.h

**Description:** Returns the smallest  $x > 0$  s.t.  $a^x = b \pmod m$ , or  $-1$  if no such  $x$  exists. modLog(a,1,m) can be used to calculate the order of  $a$ .  
**Time:**  $\mathcal{O}(\sqrt{m})$

```
ll modLog(ll a, ll b, ll m) {
    ll n = (ll) sqrt(m) + 1, e = 1, f = 1, j = 1;
    unordered_map<ll, ll> A;
    while (j <= n && (e = f = e * a % m) != b % m)
        A[e * b % m] = j++;
    if (e == b % m) return j;
    if (__gcd(m, e) == __gcd(m, b))
        rep(i,2,n+2) if (A.count(e = e * f % m))
            return n * i - A[e];
    return -1;
}
```

ModSum.h

**Description:** Sums of mod'ed arithmetic progressions.  
modsum(to, c, k, m) =  $\sum_{i=0}^{to-1} (ki+c)\%m$ . divsum is similar but for floored division.  
**Time:**  $\log(m)$ , with a large constant.

5c5bc5, 16 lines

```
typedef unsigned long long ull;
ull sumsq(ull to) { return to / 2 * ((to-1) | 1); }

ull divsum(ull to, ull c, ull k, ull m) {
    ull res = k / m * sumsq(to) + c / m * to;
    k %= m; c %= m;
    if (!k) return res;
    ull to2 = (to * k + c) / m;
    return res + (to - 1) * to2 - divsum(to2, m-1 - c, m, k);
}

ll modsum(ull to, ll c, ll k, ll m) {
    c = ((c % m) + m) % m;
    k = ((k % m) + m) % m;
    return to * c + k * sumsq(to) - m * divsum(to, c, k, m);
}
```

ModMulLL.h

**Description:** Calculate  $a \cdot b \bmod c$  (or  $a^b \bmod c$ ) for  $0 \leq a, b \leq c \leq 7.2 \cdot 10^{18}$ .  
**Time:**  $\mathcal{O}(1)$  for modmul,  $\mathcal{O}(\log b)$  for modpow

bbbd8f, 11 lines

```
typedef unsigned long long ull;
ull modmul(ull a, ull b, ull M) {
    ll ret = a * b - M * ull(1.L / M * a * b);
    return ret + M * (ret < 0) - M * (ret >= (ll)M);
}

ull modpow(ull b, ull e, ull mod) {
    ull ans = 1;
    for (; e; b = modmul(b, b, mod), e /= 2)
        if (e & 1) ans = modmul(ans, b, mod);
    return ans;
}
```

ModSqrt.h

**Description:** Tonelli-Shanks algorithm for modular square roots. Finds  $x$  s.t.  $x^2 = a \pmod p$  ( $-x$  gives the other solution).  
**Time:**  $\mathcal{O}(\log^2 p)$  worst case,  $\mathcal{O}(\log p)$  for most  $p$

"ModPow.h" 19a793, 24 lines

```
ll sqrt(ll a, ll p) {
    a %= p; if (a < 0) a += p;
    if (a == 0) return 0;
    assert(modpow(a, (p-1)/2, p) == 1); // else no solution
    if (p % 4 == 3) return modpow(a, (p+1)/4, p);
    // a^(n+3)/8 or 2^(n+3)/8 * 2^(n-1)/4 works if p % 8 == 5
    ll s = p - 1, n = 2;
    int r = 0, m;
    while (s % 2 == 0)
        ++r, s /= 2;
    while (modpow(n, (p - 1) / 2, p) != p - 1) ++n;
    ll x = modpow(a, (s + 1) / 2, p);
    ll b = modpow(a, s, p), g = modpow(n, s, p);
    for (; r = m) {
        ll t = b;
        for (m = 0; m < r && t != 1; ++m)
            t = t * t % p;
        if (m == 0) return x;
        ll gs = modpow(g, 1LL << (r - m - 1), p);
        g = gs * gs % p;
        x = x * gs % p;
        b = b * g % p;
    }
}
```

LinearDiophantine.h

**Description:** Solving linear diophantine eqns  $(ax + by = c)$ . 87blec, 74 lines

```
int gcd(int a, int b, int& x, int& y) {
    if (b == 0) {
        x = 1;
        y = 0;
        return a;
    }
    int x1, y1;
    int d = gcd(b, a % b, x1, y1);
    x = y1;
    y = x1 - y1 * (a / b);
    return d;
}

bool find_any_solution(int a, int b, int c, int &x0, int &y0,
    int &g) {
    g = gcd(abs(a), abs(b), x0, y0);
    if (c % g) {
        return false;
    }
    x0 *= c / g;
    y0 *= c / g;
    if (a < 0) x0 = -x0;
    if (b < 0) y0 = -y0;
    return true;
}

void shift_solution(int &x, int &y, int a, int b, int cnt) {
    x += cnt * b;
    y -= cnt * a;
}

int find_all_solutions(int a, int b, int c, int minx, int maxx,
    int miny, int maxy) {
    int x, y, g;
    if (!find_any_solution(a, b, c, x, y, g))
        return 0;
    a /= g;
    b /= g;

    int sign_a = a > 0 ? +1 : -1;
    int sign_b = b > 0 ? +1 : -1;

    shift_solution(x, y, a, b, (minx - x) / b);
    if (x < minx)
        shift_solution(x, y, a, b, sign_b);
    if (x > maxx)
        return 0;
    int lx1 = x;

    shift_solution(x, y, a, b, (maxx - x) / b);
    if (x > maxx)
        shift_solution(x, y, a, b, -sign_b);
    int rx1 = x;

    shift_solution(x, y, a, b, -(miny - y) / a);
    if (y < miny)
        shift_solution(x, y, a, b, -sign_a);
    if (y > maxy)
        return 0;
    int lx2 = x;

    shift_solution(x, y, a, b, -(maxy - y) / a);
    if (y > maxy)
        shift_solution(x, y, a, b, sign_a);
    int rx2 = x;
}
```

```
if (lx2 > rx2)
    swap(lx2, rx2);
int lx = max(lx1, lx2);
int rx = min(rx1, rx2);

if (lx > rx)
    return 0;
return (rx - lx) / abs(b) + 1;
}
```

5.2 Primality

FastEratosthenes.h

**Description:** Prime sieve for generating all primes smaller than LIM.  
**Time:** LIM=1e9  $\approx$  1.5s

6b2912, 20 lines

```
const int LIM = 1e6;
bitset<LIM> isPrime;
vi eratosthenes() {
    const int S = (int)round(sqrt(LIM)), R = LIM / 2;
    vi pr = {2}, sieve(S+1); pr.reserve((int)(LIM/log(LIM)*1.1));
    vector<pii> cp;
    for (int i = 3; i <= S; i += 2) if (!sieve[i]) {
        cp.push_back({i, i * i / 2});
        for (int j = i * i; j <= S; j += 2 * i) sieve[j] = 1;
    }
    for (int L = 1; L <= R; L += S) {
        array<bool, S> block{};
        for (auto &[p, idx] : cp)
            for (int i=idx; i < S+L; idx = (i+=p)) block[i-L] = 1;
        rep(i,0,min(S, R - L))
            if (!block[i]) pr.push_back((L + i) * 2 + 1);
    }
    for (int i : pr) isPrime[i] = 1;
    return pr;
}
```

MillerRabin.h

**Description:** Deterministic Miller-Rabin primality test. Guaranteed to work for numbers up to  $7 \cdot 10^{18}$ ; for larger numbers, use Python and extend A randomly.  
**Time:** 7 times the complexity of  $a^b \bmod c$ .

"ModMulLL.h" 60dcd1, 12 lines

```
bool isPrime(ull n) {
    if (n < 2 || n % 6 % 4 != 1) return (n | 1) == 3;
    ull A[] = {2, 325, 9375, 28178, 450775, 9780504, 1795265022},
        s = __builtin_ctzll(n-1), d = n >> s;
    for (ull a : A) { // ^ count trailing zeroes
        ull p = modpow(a%n, d, n), i = s;
        while (p != 1 && p != n - 1 && a % n && i--)
            p = modmul(p, p, n);
        if (p != n-1 && i != s) return 0;
    }
    return 1;
}
```

Factor.h

**Description:** Pollard-rho randomized factorization algorithm. Returns prime factors of a number, in arbitrary order (e.g. 2299 -> {11, 19, 11}).

**Time:**  $\mathcal{O}(n^{1/4})$ , less for numbers with small factors.

"ModMulLL.h", "MillerRabin.h" a33cf6, 18 lines

```
ull pollard(ull n) {
    auto f = [n](ull x) { return modmul(x, x, n) + 1; };
    ull x = 0, y = 0, t = 30, prd = 2, i = 1, q;
    while (t++ % 40 || __gcd(prd, n) == 1) {
        if (x == y) x = ++i, y = f(x);
        if ((q = modmul(prd, max(x,y) - min(x,y), n))) prd = q;
        x = f(x), y = f(f(y));
    }
}
```

```
    }
    return __gcd(prd, n);
}
vector<ull> factor(ull n) {
    if (n == 1) return {};
    if (isPrime(n)) return {n};
    ull x = pollard(n);
    auto l = factor(x), r = factor(n / x);
    l.insert(l.end(), all(r));
    return l;
}
```

### 5.3 Divisibility

**euclid.h**  
**Description:** Finds two integers  $x$  and  $y$ , such that  $ax + by = \gcd(a, b)$ . If you just need gcd, use the built in `__gcd` instead. If  $a$  and  $b$  are coprime, then  $x$  is the inverse of  $a \pmod b$ .

```
33ba8f, 5 lines
11 euclid(11 a, 11 b, 11 &x, 11 &y) {
    if (!b) return x = 1, y = 0, a;
    11 d = euclid(b, a % b, y, x);
    return y -= a/b * x, d;
}
```

**CRT.h**  
**Description:** Chinese Remainder Theorem.  
`crt(a, m, b, n)` computes  $x$  such that  $x \equiv a \pmod m, x \equiv b \pmod n$ . If  $|a| < m$  and  $|b| < n$ ,  $x$  will obey  $0 \leq x < \text{lcm}(m, n)$ . Assumes  $mn < 2^{62}$ .  
**Time:**  $\log(n)$

```
04d93a, 7 lines
11 crt(11 a, 11 m, 11 b, 11 n) {
    if (n > m) swap(a, b), swap(m, n);
    11 x, y, g = euclid(m, n, x, y);
    assert((a - b) % g == 0); // else no solution
    x = (b - a) % n * x % n / g * m + a;
    return x < 0 ? x + m*n/g : x;
}
```

### Bézout’s identity

For  $a \neq 0, b \neq 0$ , then  $d = \gcd(a, b)$  is the smallest positive integer for which there are integer solutions to

$$ax + by = d$$

If  $(x, y)$  is one solution, then all solutions are given by

$$\left(x + \frac{kb}{\gcd(a,b)}, y - \frac{ka}{\gcd(a,b)}\right), \quad k \in \mathbb{Z}$$

**phiFunction.h**  
**Description:** *Euler’s*  $\phi$  function is defined as  $\phi(n) := \#$  of positive integers  $\leq n$  that are coprime with  $n$ .  $\phi(1) = 1, p$  prime  $\Rightarrow \phi(p^k) = (p - 1)p^{k-1}$ ,  $m, n$  coprime  $\Rightarrow \phi(mn) = \phi(m)\phi(n)$ . If  $n = p_1^{k_1}p_2^{k_2}...p_r^{k_r}$  then  $\phi(n) = (p_1 - 1)p_1^{k_1-1}...(p_r - 1)p_r^{k_r-1}$ .  $\phi(n) = n \cdot \prod_{p|n} (1 - 1/p)$ .  
 $\sum_{d|n} \phi(d) = n, \sum_{1 \leq k \leq n, \gcd(k,n)=1} k = n\phi(n)/2, n > 1$   
**Euler’s thm:**  $a, n$  coprime  $\Rightarrow a^{\phi(n)} \equiv 1 \pmod n$ .  
**Fermat’s little thm:**  $p$  prime  $\Rightarrow a^{p-1} \equiv 1 \pmod p \ \forall a$ .

```
cf7d6d, 8 lines
const int LIM = 5000000;
int phi[LIM];

void calculatePhi() {
    rep(i, 0, LIM) phi[i] = i&1 ? i : i/2;
    for (int i = 3; i < LIM; i += 2) if(phi[i] == i)
```

```
    for (int j = i; j < LIM; j += i) phi[j] -= phi[j] / i;
}
```

### 5.4 Fractions

**ContinuedFractions.h**  
**Description:** Given  $N$  and a real number  $x \geq 0$ , finds the closest rational approximation  $p/q$  with  $p, q \leq N$ . It will obey  $|p/q - x| \leq 1/qN$ .  
For consecutive convergents,  $p_{k+1}q_k - q_{k+1}p_k = (-1)^k$ . ( $p_k/q_k$  alternates between  $> x$  and  $< x$ .) If  $x$  is rational,  $y$  eventually becomes  $\infty$ ; if  $x$  is the root of a degree 2 polynomial the  $a$ ’s eventually become cyclic.  
**Time:**  $\mathcal{O}(\log N)$

```
dd6c5e, 21 lines
typedef double d; // for N ~ 1e7; long double for N ~ 1e9
pair<11, 11> approximate(d x, 11 N) {
    11 LP = 0, LQ = 1, P = 1, Q = 0, inf = LLONG_MAX; d y = x;
    for (;;) {
        11 lim = min(P ? (N-LP) / P : inf, Q ? (N-LQ) / Q : inf),
           a = (11)floor(y), b = min(a, lim),
           NP = b*P + LP, NQ = b*Q + LQ;
        if (a > b) {
            // If b > a/2, we have a semi-convergent that gives us a
            // better approximation; if b = a/2, we *may* have one.
            // Return {P, Q} here for a more canonical approximation.
            return (abs(x - (d)NP / (d)NQ) < abs(x - (d)P / (d)Q)) ?
                make_pair(NP, NQ) : make_pair(P, Q);
        }
        if (abs(y = 1/(y - (d)a)) > 3*N) {
            return {NP, NQ};
        }
        LP = P; P = NP;
        LQ = Q; Q = NQ;
    }
}
```

**FracBinarySearch.h**  
**Description:** Given  $f$  and  $N$ , finds the smallest fraction  $p/q \in [0, 1]$  such that  $f(p/q)$  is true, and  $p, q \leq N$ . You may want to throw an exception from  $f$  if it finds an exact solution, in which case  $N$  can be removed.  
**Usage:** `fracBS({}(Frac f) { return f.p>=3*f.q; }, 10);` //  $\{1, 3\}$   
**Time:**  $\mathcal{O}(\log(N))$

```
27ab3e, 25 lines
struct Frac { 11 p, q; };

template<class F>
Frac fracBS(F f, 11 N) {
    bool dir = 1, A = 1, B = 1;
    Frac lo{0, 1}, hi{1, 1}; // Set hi to 1/0 to search (0, N]
    if (f(lo)) return lo;
    assert(f(hi));
    while (A || B) {
        11 adv = 0, step = 1; // move hi if dir, else lo
        for (int si = 0; step; (step *= 2) >= si) {
            adv += step;
            Frac mid{lo.p * adv + hi.p, lo.q * adv + hi.q};
            if (abs(mid.p) > N || mid.q > N || dir == !f(mid)) {
                adv -= step; si = 2;
            }
        }
        hi.p += lo.p * adv;
        hi.q += lo.q * adv;
        dir = !dir;
        swap(lo, hi);
        A = B; B = !!adv;
    }
    return dir ? hi : lo;
}
```

### 5.5 Pythagorean Triples

The Pythagorean triples are uniquely generated by

$$a = k \cdot (m^2 - n^2), \quad b = k \cdot (2mn), \quad c = k \cdot (m^2 + n^2),$$

with  $m > n > 0, k > 0, m \perp n$ , and either  $m$  or  $n$  even.

### 5.6 Primes

$p = 962592769$  is such that  $2^{21} \mid p - 1$ , which may be useful. For hashing use 970592641 (31-bit number), 31443539979727 (45-bit), 3006703054056749 (52-bit). There are 78498 primes less than 1 000 000.

Primitive roots exist modulo any prime power  $p^a$ , except for  $p = 2, a > 2$ , and there are  $\phi(\phi(p^a))$  many. For  $p = 2, a > 2$ , the group  $\mathbb{Z}_{2^a}^\times$  is instead isomorphic to  $\mathbb{Z}_2 \times \mathbb{Z}_{2^{a-2}}$ .

### 5.7 Estimates

$$\sum_{d|n} d = O(n \log \log n).$$

The number of divisors of  $n$  is at most around 100 for  $n < 5e4$ , 500 for  $n < 1e7$ , 2000 for  $n < 1e10$ , 200 000 for  $n < 1e19$ .

### 5.8 Mobius Function

$$\mu(n) = \begin{cases} 0 & n \text{ is not square free} \\ 1 & n \text{ has even number of prime factors} \\ -1 & n \text{ has odd number of prime factors} \end{cases}$$

Mobius Inversion:

$$g(n) = \sum_{d|n} f(d) \Leftrightarrow f(n) = \sum_{d|n} \mu(d)g(n/d)$$

Other useful formulas/forms:

$$\sum_{d|n} \mu(d) = [n = 1] \text{ (very useful)}$$

$$g(n) = \sum_{n|d} f(d) \Leftrightarrow f(n) = \sum_{n|d} \mu(d/n)g(d)$$

$$g(n) = \sum_{1 \leq m \leq n} f(\lfloor \frac{n}{m} \rfloor) \Leftrightarrow f(n) = \sum_{1 \leq m \leq n} \mu(m)g(\lfloor \frac{n}{m} \rfloor)$$

## Combinatorics (6)

### 6.1 Permutations

$n$	1	2	3	4	5	6	7	8	9	10
$n!$	1	2	6	24	120	720	5040	40320	362880	3628800
$n$	11	12	13	14	15	16	17			
$n!$	4.0e7	4.8e8	6.2e9	8.7e10	1.3e12	2.1e13	3.6e14			
$n$	20	25	30	40	50	100	150	171		
$n!$	2e18	2e25	3e32	8e47	3e64	9e157	6e262	>DBL_MAX		

IntPerm.h  
**Description:** Permutation -> integer conversion. (Not order preserving.) Integer -> permutation can use a lookup table.  
**Time:**  $\mathcal{O}(n)$

```
044568, 6 lines
int permToInt(vi& v) {
    int use = 0, i = 0, r = 0;
    for(int x:v) r = r * ++i + __builtin_popcount(use & ~(1<<x)),
        use |= 1 << x; // (note: minus, not ~!)
    return r;
}
```

**Cycles**  
Let  $g_S(n)$  be the number of  $n$ -permutations whose cycle lengths all belong to the set  $S$ . Then

$$\sum_{n=0}^\infty g_S(n) \frac{x^n}{n!} = \exp\left(\sum_{n \in S} \frac{x^n}{n}\right)$$

**Derangements**  
Permutations of a set such that none of the elements appear in their original position.

$$D(n) = (n-1)(D(n-1) + D(n-2)) = nD(n-1) + (-1)^n = \left\lfloor \frac{n!}{e} \right\rfloor$$

**Burnside’s lemma**  
Given a group  $G$  of symmetries and a set  $X$ , the number of elements of  $X$  up to symmetry equals

$$\frac{1}{|G|} \sum_{g \in G} |X^g|,$$

where  $X^g$  are the elements fixed by  $g$  ( $g.x = x$ ).

If  $f(n)$  counts “configurations” (of some sort) of length  $n$ , we can ignore rotational symmetry using  $G = \mathbb{Z}_n$  to get

$$g(n) = \frac{1}{n} \sum_{k=0}^{n-1} f(\gcd(n, k)) = \frac{1}{n} \sum_{k|n} f(k) \phi(n/k).$$

6.2 Partitions and subsets

**Partition function**  
Number of ways of writing  $n$  as a sum of positive integers, disregarding the order of the summands.

$$p(0) = 1, \quad p(n) = \sum_{k \in \mathbb{Z} \setminus \{0\}} (-1)^{k+1} p(n - k(3k - 1)/2)$$

$$p(n) \sim 0.145/n \cdot \exp(2.56\sqrt{n})$$

$n$	0	1	2	3	4	5	6	7	8	9	20	50	100
$p(n)$	1	1	2	3	5	7	11	15	22	30	627	$\sim 2e5$	$\sim 2e8$

multinomial

**Lucas’ Theorem**  
Let  $n, m$  be non-negative integers and  $p$  a prime. Write  $n = n_k p^k + \dots + n_1 p + n_0$  and  $m = m_k p^k + \dots + m_1 p + m_0$ . Then  $\binom{n}{m} \equiv \prod_{i=0}^k \binom{n_i}{m_i} \pmod{p}$ .

**Binomials**  
multinomial.h  
**Description:** Computes  $\binom{k_1 + \dots + k_n}{k_1, k_2, \dots, k_n} = \frac{(\sum k_i)!}{k_1! k_2! \dots k_n!}$ .

```
a0a312, 6 lines
11 multinomial(vi& v) {
    11 c = 1, m = v.empty() ? 1 : v[0];
    rep(i, 1, sz(v)) rep(j, 0, v[i])
        c = c * ++m / (j+1);
    return c;
}
```

6.3 General purpose numbers

**Bernoulli numbers**  
EGF of Bernoulli numbers is  $B(t) = \frac{t}{e^t - 1}$  (FFT-able).  
 $B[0, \dots] = [1, -\frac{1}{2}, \frac{1}{6}, 0, -\frac{1}{30}, 0, \frac{1}{42}, \dots]$

Sums of powers:  
$$\sum_{i=1}^n n^m = \frac{1}{m+1} \sum_{k=0}^m \binom{m+1}{k} B_k \cdot (n+1)^{m+1-k}$$

Euler-Maclaurin formula for infinite sums:

$$\begin{aligned} \sum_{i=m}^\infty f(i) &= \int_m^\infty f(x) dx - \sum_{k=1}^\infty \frac{B_k}{k!} f^{(k-1)}(m) \\ &\approx \int_m^\infty f(x) dx + \frac{f(m)}{2} - \frac{f'(m)}{12} + \frac{f'''(m)}{720} + O(f^{(5)}(m)) \end{aligned}$$

**Stirling numbers of the first kind**  
Number of permutations on  $n$  items with  $k$  cycles.

$$c(n, k) = c(n-1, k-1) + (n-1)c(n-1, k), \quad c(0, 0) = 1$$
$$\sum_{k=0}^n c(n, k) x^k = x(x+1) \dots (x+n-1)$$

$$c(8, k) = 8, 0, 5040, 13068, 13132, 6769, 1960, 322, 28, 1$$
$$c(n, 2) = 0, 0, 1, 3, 11, 50, 274, 1764, 13068, 109584, \dots$$

**Eulerian numbers**  
Number of permutations  $\pi \in S_n$  in which exactly  $k$  elements are greater than the previous element.  $k$   $j$ :s s.t.  $\pi(j) > \pi(j+1)$ ,  $k+1$   $j$ :s s.t.  $\pi(j) \geq j$ ,  $k$   $j$ :s s.t.  $\pi(j) > j$ .

$$E(n, k) = (n-k)E(n-1, k-1) + (k+1)E(n-1, k)$$

$$E(n, 0) = E(n, n-1) = 1$$

$$E(n, k) = \sum_{j=0}^k (-1)^j \binom{n+1}{j} (k+1-j)^n$$

**Stirling numbers of the second kind**  
Partitions of  $n$  distinct elements into exactly  $k$  groups.

$$S(n, k) = S(n-1, k-1) + kS(n-1, k)$$

$$S(n, 1) = S(n, n) = 1$$

$$S(n, k) = \frac{1}{k!} \sum_{j=0}^k (-1)^{k-j} \binom{k}{j} j^n$$

**Bell numbers**  
Total number of partitions of  $n$  distinct elements.  $B(n) = 1, 1, 2, 5, 15, 52, 203, 877, 4140, 21147, \dots$  For  $p$  prime,

$$B(p^m + n) \equiv mB(n) + B(n+1) \pmod{p}$$

**Labeled unrooted trees**  
# on  $n$  vertices:  $n^{n-2}$   
# on  $k$  existing trees of size  $n_i$ :  $n_1 n_2 \dots n_k n^{k-2}$   
# with degrees  $d_i$ :  $(n-2)! / ((d_1-1)! \dots (d_n-1)!)$

**Catalan numbers**  
$$C_n = \frac{1}{n+1} \binom{2n}{n} = \binom{2n}{n} - \binom{2n}{n+1} = \frac{(2n)!}{(n+1)!n!}$$

$$C_0 = 1, \quad C_{n+1} = \frac{2(2n+1)}{n+2} C_n, \quad C_{n+1} = \sum C_i C_{n-i}$$

$$C_n = 1, 1, 2, 5, 14, 42, 132, 429, 1430, 4862, 16796, 58786, \dots$$

- sub-diagonal monotone paths in an  $n \times n$  grid.
- strings with  $n$  pairs of parenthesis, correctly nested.
- binary trees with with  $n+1$  leaves (0 or 2 children).
- ordered trees with  $n+1$  vertices.
- ways a convex polygon with  $n+2$  sides can be cut into triangles by connecting vertices with straight lines.
- permutations of  $[n]$  with no 3-term increasing subseq.

6.4 Probability theory  
$$V[X] = E[(X - E[X])^2] = E[X^2] - E[X]^2$$

Algebra (7)

7.1 Equations and Generating Functions

$$c^a + c^{a+1} + \dots + c^b = \frac{c^{b+1} - c^a}{c - 1}, c \neq 1$$

$$e^x = 1 + x + \frac{x^2}{2!} + \frac{x^3}{3!} + \dots, \quad (-\infty < x < \infty)$$

$$\ln(1+x) = x - \frac{x^2}{2} + \frac{x^3}{3} - \frac{x^4}{4} + \dots, \quad (-1 < x \leq 1)$$

$$\sqrt{1+x} = 1 + \frac{x}{2} - \frac{x^2}{8} + \frac{2x^3}{32} - \frac{5x^4}{128} + \dots, (-1 \leq x \leq 1)$$

$$\sin x = x - \frac{x^3}{3!} + \frac{x^5}{5!} - \frac{x^7}{7!} + \dots, (-\infty < x < \infty)$$

$$\cos x = 1 - \frac{x^2}{2!} + \frac{x^4}{4!} - \frac{x^6}{6!} + \dots, (-\infty < x < \infty)$$

$$1^4 + 2^4 + 3^4 + \dots + n^4 = \frac{n(n+1)(2n+1)(3n^2+3n-1)}{30}$$

Generating Functions Blog 1 and 2.

7.2 Polynomials and recurrences

Polynomial.h

```
struct Poly {
    vector<double> a;
    double operator()(double x) const {
        double val = 0;
        for (int i = sz(a); i--;) (val *= x) += a[i];
        return val;
    }
    void diff() {
        rep(i,1,sz(a)) a[i-1] = i*a[i];
        a.pop_back();
    }
    void divroot(double x0) {
        double b = a.back(), c; a.back() = 0;
        for(int i=sz(a)-1; i--;) c = a[i], a[i] = a[i+1]*x0+b, b=c;
        a.pop_back();
    }
};
```

PolyRoots.h

**Description:** Finds the real roots to a polynomial.  
**Usage:** polyRoots({{2,-3,1}},-1e9,1e9) // solve x^2-3x+2 = 0  
**Time:**  $\mathcal{O}(n^2 \log(1/\epsilon))$

```
"Polynomial.h"
vector<double> polyRoots(Poly p, double xmin, double xmax) {
    if (sz(p.a) == 2) { return {-p.a[0]/p.a[1]}; }
    vector<double> ret;
    Poly der = p;
    der.diff();
    auto dr = polyRoots(der, xmin, xmax);
    dr.push_back(xmin-1);
    dr.push_back(xmax+1);
    sort(all(dr));
    rep(i,0,sz(dr)-1) {
        double l = dr[i], h = dr[i+1];
        bool sign = p(l) > 0;
        if (sign ^ (p(h) > 0)) {
            rep(it,0,60) { // while (h - l > 1e-8)
                double m = (l + h) / 2, f = p(m);
                if ((f <= 0) ^ sign) l = m;
                else h = m;
            }
            ret.push_back((l + h) / 2);
        }
    }
    return ret;
}
```

PolyInterpolate.h

**Description:** Given  $n$  points  $(x[i], y[i])$ , computes an  $n-1$ -degree polynomial  $p$  that passes through them:  $p(x) = a[0] * x^0 + \dots + a[n-1] * x^{n-1}$ . For numerical precision, pick  $x[k] = c * \cos(k/(n-1) * \pi), k = 0 \dots n-1$ .

```
Time:  $\mathcal{O}(n^2)$ 
typedef vector<double> vd;
vd interpolate(vd x, vd y, int n) {
    vd res(n), temp(n);
    rep(k,0,n-1) rep(i,k+1,n)
        y[i] = (y[i] - y[k]) / (x[i] - x[k]);
    double last = 0; temp[0] = 1;
    rep(k,0,n) rep(i,0,n) {
        res[i] += y[k] * temp[i];
        swap(last, temp[i]);
        temp[i] -= last * x[k];
    }
    return res;
}
```

BerlekampMassey.h

**Description:** Recovers any  $n$ -order linear recurrence relation from the first  $2n$  terms of the recurrence. Useful for guessing linear recurrences after brute-forcing the first terms. Should work on any field, but numerical stability for floats is not guaranteed. Output will have size  $\leq n$ .  
**Usage:** berlekampMassey({0, 1, 1, 3, 5, 11}) // {1, 2}  
**Time:**  $\mathcal{O}(N^2)$

```
"../number-theory/ModPow.h"
vector<ll> berlekampMassey(vector<ll> s) {
    int n = sz(s), L = 0, m = 0;
    vector<ll> C(n), B(n), T;
    C[0] = B[0] = 1;

    ll b = 1;
    rep(i,0,n) { ++m;
        ll d = s[i] % mod;
        rep(j,1,L+1) d = (d + C[j] * s[i - j]) % mod;
        if (!d) continue;
        T = C; ll coef = d * modpow(b, mod-2) % mod;
        rep(j,m,n) C[j] = (C[j] - coef * B[j - m]) % mod;
        if (2 * L > i) continue;
        L = i + 1 - L; B = T; b = d; m = 0;
    }

    C.resize(L + 1); C.erase(C.begin());
    for (ll& x : C) x = (mod - x) % mod;
    return C;
}
```

LinearRecurrence.h

**Description:** Generates the  $k$ 'th term of an  $n$ -order linear recurrence  $S[i] = \sum_j S[i-j-1]tr[j]$ , given  $S[0 \dots \geq n-1]$  and  $tr[0 \dots n-1]$ . Faster than matrix multiplication. Useful together with Berlekamp-Massey.  
**Usage:** linearRec({0, 1}, {1, 1}, k) //  $k$ 'th Fibonacci number  
**Time:**  $\mathcal{O}(n^2 \log k)$

```
typedef vector<ll> Poly;
ll linearRec(Poly S, Poly tr, ll k) {
    int n = sz(tr);

    auto combine = [&](Poly a, Poly b) {
        Poly res(n * 2 + 1);
        rep(i,0,n+1) rep(j,0,n+1)
            res[i + j] = (res[i + j] + a[i] * b[j]) % mod;
        for (int i = 2 * n; i > n; --i) rep(j,0,n)
            res[i - 1 - j] = (res[i - 1 - j] + res[i] * tr[j]) % mod;
        res.resize(n + 1);
        return res;
    };
}
```

```
Poly pol(n + 1), e(pol);
pol[0] = e[1] = 1;

for (++k; k; k /= 2) {
    if (k % 2) pol = combine(pol, e);
    e = combine(e, e);
}

ll res = 0;
rep(i,0,n) res = (res + pol[i + 1] * S[i]) % mod;
return res;
}
```

7.3 Optimization

GoldenSectionSearch.h

**Description:** Finds the argument minimizing the function  $f$  in the interval  $[a,b]$  assuming  $f$  is unimodal on the interval, i.e. has only one local minimum. The maximum error in the result is  $eps$ . Works equally well for maximization with a small change in the code. See TernarySearch.h in the Various chapter for a discrete version.  
**Usage:** double func(double x) { return 4\*x+.3\*x\*x; }  
double xmin = gss(-1000,1000,func);  
**Time:**  $\mathcal{O}(\log((b-a)/\epsilon))$

```
double gss(double a, double b, double (*f)(double)) {
    double r = (sqrt(5)-1)/2, eps = 1e-7;
    double x1 = b - r*(b-a), x2 = a + r*(b-a);
    double f1 = f(x1), f2 = f(x2);
    while (b-a > eps)
        if (f1 < f2) { //change to > to find maximum
            b = x2; x2 = x1; f2 = f1;
            x1 = b - r*(b-a); f1 = f(x1);
        } else {
            a = x1; x1 = x2; f1 = f2;
            x2 = a + r*(b-a); f2 = f(x2);
        }
    return a;
}
```

HillClimbing.h

**Description:** Poor man's optimization for unimodal functions

```
"Seecaf, 14 lines"
typedef array<double, 2> P;

template<class F> pair<double, P> hillClimb(P start, F f) {
    pair<double, P> cur(f(start), start);
    for (double jmp = 1e9; jmp > 1e-20; jmp /= 2) {
        rep(j,0,100) rep(dx,-1,2) rep(dy,-1,2) {
            P p = cur.second;
            p[0] += dx*jmp;
            p[1] += dy*jmp;
            cur = min(cur, make_pair(f(p), p));
        }
    }
    return cur;
}
```

Integrate.h

**Description:** Simple integration of a function over an interval using Simpson's rule. The error should be proportional to  $h^4$ , although in practice you will want to verify that the result is stable to desired precision when epsilon changes.

```
template<class F>
double quad(double a, double b, F f, const int n = 1000) {
    double h = (b - a) / 2 / n, v = f(a) + f(b);
    rep(i,1,n*2)
        v += f(a + i*h) * (i&1 ? 4 : 2);
    return v * h / 3;
}
```



```
}

IntegrateAdaptive.h
Description: Fast integration using an adaptive Simpson's rule.
Usage: double sphereVolume = quad(-1, 1, [](double x) {
return quad(-1, 1, [&](double y) {
return quad(-1, 1, [&](double z) {
return x*x + y*y + z*z < 1; });});});});
92dd79, 15 lines
```

```
typedef double d;
#define S(a,b) (f(a) + 4*f((a+b) / 2) + f(b)) * (b-a) / 6

template <class F>
d rec(F& f, d a, d b, d eps, d S) {
d c = (a + b) / 2;
d S1 = S(a, c), S2 = S(c, b), T = S1 + S2;
if (abs(T - S) <= 15 * eps || b - a < 1e-10)
return T + (T - S) / 15;
return rec(f, a, c, eps / 2, S1) + rec(f, c, b, eps / 2, S2);
}

template<class F>
d quad(d a, d b, F f, d eps = 1e-8) {
return rec(f, a, b, eps, S(a, b));
}
```

Simplex.h

**Description:** Solves a general linear maximization problem: maximize  $c^T x$  subject to  $Ax \leq b, x \geq 0$ . Returns -inf if there is no solution, inf if there are arbitrarily good solutions, or the maximum value of  $c^T x$  otherwise. The input vector is set to an optimal  $x$  (or in the unbounded case, an arbitrary solution fulfilling the constraints). Numerical stability is not guaranteed. For better performance, define variables such that  $x = 0$  is viable.

**Usage:** vvd A = {{1,-1}, {-1,1}, {-1,-2}};

vd b = {1,1,-4}, c = {-1,-1}, x;

T val = LPSolver(A, b, c).solve(x);

**Time:**  $\mathcal{O}(NM * \text{\#pivots})$ , where a pivot may be e.g. an edge relaxation.  $\mathcal{O}(2^n)$  in the general case.

```
aa8530, 68 lines

typedef double T; // long double, Rational, double + mod<P>...
typedef vector<T> vd;
typedef vector<vd> vvd;

const T eps = 1e-8, inf = 1/.0;
#define MP make_pair
#define ltj(X) if(s == -1 || MP(X[j],N[j]) < MP(X[s],N[s])) s=j

struct LPSolver {
int m, n;
vi N, B;
vvd D;

LPSolver(const vvd& A, const vd& b, const vd& c) :
m(sz(b)), n(sz(c)), N(n+1), B(m), D(m+2, vd(n+2)) {
rep(i,0,m) rep(j,0,n) D[i][j] = A[i][j];
rep(i,0,m) { B[i] = n+i; D[i][n] = -1; D[i][n+1] = b[i]; }
rep(j,0,n) { N[j] = j; D[m][j] = -c[j]; }
N[n] = -1; D[m+1][n] = 1;
}

void pivot(int r, int s) {
T *a = D[r].data(), inv = 1 / a[s];
rep(i,0,m+2) if (i != r && abs(D[i][s]) > eps) {
T *b = D[i].data(), inv2 = b[s] * inv;
rep(j,0,n+2) b[j] -= a[j] * inv2;
b[s] = a[s] * inv2;
}
rep(j,0,n+2) if (j != s) D[r][j] *= inv;
rep(i,0,m+2) if (i != r) D[i][s] *= -inv;
D[r][s] = inv;
}
```

```
swap(B[r], N[s]);
}

bool simplex(int phase) {
int x = m + phase - 1;
for (;;) {
int s = -1;
rep(j,0,n+1) if (N[j] != -phase) ltj(D[x]);
if (D[x][s] >= -eps) return true;
int r = -1;
rep(i,0,m) {
if (D[i][s] <= eps) continue;
if (r == -1 || MP(D[i][n+1] / D[i][s], B[i])
< MP(D[r][n+1] / D[r][s], B[r])) r = i;
}
if (r == -1) return false;
pivot(r, s);
}
}

T solve(vd &x) {
int r = 0;
rep(i,1,m) if (D[i][n+1] < D[r][n+1]) r = i;
if (D[r][n+1] < -eps) {
pivot(r, n);
if (!simplex(2) || D[m+1][n+1] < -eps) return -inf;
rep(i,0,m) if (B[i] == -1) {
int s = 0;
rep(j,1,n+1) ltj(D[i]);
pivot(i, s);
}
}
bool ok = simplex(1); x = vd(n);
rep(i,0,m) if (B[i] < n) x[B[i]] = D[i][n+1];
return ok ? D[m][n+1] : inf;
}
}
```

7.4 Matrices

Determinant.h

**Description:** Calculates determinant of a matrix. Destroys the matrix.

**Time:**  $\mathcal{O}(N^3)$

```
bd5cec, 15 lines

double det(vector<vector<double>>& a) {
int n = sz(a); double res = 1;
rep(i,0,n) {
int b = i;
rep(j,i+1,n) if (fabs(a[j][i]) > fabs(a[b][i])) b = j;
if (i != b) swap(a[i], a[b]), res *= -1;
res *= a[i][i];
if (res == 0) return 0;
rep(j,i+1,n) {
double v = a[j][i] / a[i][i];
if (v != 0) rep(k,i+1,n) a[j][k] -= v * a[i][k];
}
}
return res;
}
```

IntDeterminant.h

**Description:** Calculates determinant using modular arithmetics. Modulos can also be removed to get a pure-integer version.

```
3313dc, 18 lines

const ll mod = 12345;
ll det(vector<vector<ll>>& a) {
int n = sz(a); ll ans = 1;
rep(i,0,n) {
rep(j,i+1,n) {
```

```
while (a[j][i] != 0) { // gcd step
ll t = a[i][i] / a[j][i];
if (t) rep(k,i,n)
a[i][k] = (a[i][k] - a[j][k] * t) % mod;
swap(a[i], a[j]);
ans *= -1;
}
}
ans = ans * a[i][i] % mod;
if (!ans) return 0;
}
return (ans + mod) % mod;
}
```

SolveLinear.h

**Description:** Solves  $A * x = b$ . If there are multiple solutions, an arbitrary one is returned. Returns rank, or -1 if no solutions. Data in  $A$  and  $b$  is lost.

**Time:**  $\mathcal{O}(n^2m)$

```
44c9ab, 38 lines

typedef vector<double> vd;
const double eps = 1e-12;

int solveLinear(vector<vd>& A, vd& b, vd& x) {
int n = sz(A), m = sz(x), rank = 0, br, bc;
if (n) assert(sz(A[0]) == m);
vi col(m); iota(all(col), 0);

rep(i,0,n) {
double v, bv = 0;
rep(r,i,n) rep(c,i,m)
if ((v = fabs(A[r][c])) > bv)
br = r, bc = c, bv = v;
if (bv <= eps) {
rep(j,i,n) if (fabs(b[j]) > eps) return -1;
break;
}
swap(A[i], A[br]);
swap(b[i], b[br]);
swap(col[i], col[bc]);
rep(j,0,n) swap(A[j][i], A[j][bc]);
bv = 1/A[i][i];
rep(j,i+1,n) {
double fac = A[j][i] * bv;
b[j] -= fac * b[i];
rep(k,i+1,m) A[j][k] -= fac*A[i][k];
}
rank++;
}

x.assign(m, 0);
for (int i = rank; i--;) {
b[i] /= A[i][i];
x[col[i]] = b[i];
rep(j,0,i) b[j] -= A[j][i] * b[i];
}
return rank; // (multiple solutions if rank < m)
}
```

SolveLinear2.h

**Description:** To get all uniquely determined values of  $x$  back from SolveLinear, make the following changes:

```
08e495, 7 lines

"SolveLinear.h"

rep(j,0,n) if (j != i) // instead of rep(j,i+1,n)
// ... then at the end:
x.assign(m, undefined);
rep(i,0,rank) {
rep(j,rank,m) if (fabs(A[i][j]) > eps) goto fail;
x[col[i]] = b[i] / A[i][i];
fail;;
}
```

SolveLinearBinary.h

**Description:** Solves  $Ax = b$  over  $\mathbb{F}_2$ . If there are multiple solutions, one is returned arbitrarily. Returns rank, or -1 if no solutions. Destroys  $A$  and  $b$ .  
**Time:**  $\mathcal{O}(n^2m)$

fa2d7a, 34 lines

**typedef** bitset<1000> bs;

```
int solveLinear(vector<bs>& A, vi& b, bs& x, int m) {
    int n = sz(A), rank = 0, br;
    assert(m <= sz(x));
    vi col(m); iota(all(col), 0);
    rep(i,0,n) {
        for (br=i; br<n; ++br) if (A[br].any()) break;
        if (br == n) {
            rep(j,i,n) if(b[j]) return -1;
            break;
        }
        int bc = (int)A[br]._Find_next(i-1);
        swap(A[i], A[br]);
        swap(b[i], b[br]);
        swap(col[i], col[bc]);
        rep(j,0,n) if (A[j][i] != A[j][bc]) {
            A[j].flip(i); A[j].flip(bc);
        }
        rep(j,i+1,n) if (A[j][i]) {
            b[j] ^= b[i];
            A[j] ^= A[i];
        }
        rank++;
    }
}
```

```
x = bs();
for (int i = rank; i--;) {
    if (!b[i]) continue;
    x[col[i]] = 1;
    rep(j,0,i) b[j] ^= A[j][i];
}
return rank; // (multiple solutions if rank < m)
```

MatrixInverse.h

**Description:** Invert matrix  $A$ . Returns rank; result is stored in  $A$  unless singular (rank < n). Can easily be extended to prime moduli; for prime powers, repeatedly set  $A^{-1} = A^{-1}(2I - AA^{-1}) \pmod{p^k}$  where  $A^{-1}$  starts as the inverse of  $A \pmod p$ , and  $k$  is doubled in each step.  
**Time:**  $\mathcal{O}(n^3)$

ebfff6, 35 lines

```
int matInv(vector<vector<double>>& A) {
    int n = sz(A); vi col(n);
    vector<vector<double>> tmp(n, vector<double>(n));
    rep(i,0,n) tmp[i][i] = 1, col[i] = i;

    rep(i,0,n) {
        int r = i, c = i;
        rep(j,i,n) rep(k,i,n)
            if (fabs(A[j][k]) > fabs(A[r][c]))
                r = j, c = k;
        if (fabs(A[r][c]) < 1e-12) return i;
        A[i].swap(A[r]); tmp[i].swap(tmp[r]);
        rep(j,0,n)
            swap(A[j][i], A[j][c]), swap(tmp[j][i], tmp[j][c]);
        swap(col[i], col[c]);
        double v = A[i][i];
        rep(j,i+1,n) {
            double f = A[j][i] / v;
            A[j][i] = 0;
            rep(k,i+1,n) A[j][k] -= f*A[i][k];
            rep(k,0,n) tmp[j][k] -= f*tmp[i][k];
        }
    }
}
```

```
rep(j,i+1,n) A[i][j] /= v;
rep(j,0,n) tmp[i][j] /= v;
A[i][i] = 1;
}

for (int i = n-1; i > 0; --i) rep(j,0,i) {
    double v = A[j][i];
    rep(k,0,n) tmp[j][k] -= v*tmp[i][k];
}

rep(i,0,n) rep(j,0,n) A[col[i]][col[j]] = tmp[i][j];
return n;
}
```

Tridiagonal.h

**Description:**  $x = \text{tridiagonal}(d,p,q,b)$  solves the equation system

$$\begin{pmatrix} b_0 \\ b_1 \\ b_2 \\ b_3 \\ \vdots \\ b_{n-1} \end{pmatrix} = \begin{pmatrix} d_0 & p_0 & 0 & 0 & \cdots & 0 \\ q_0 & d_1 & p_1 & 0 & \cdots & 0 \\ 0 & q_1 & d_2 & p_2 & \cdots & 0 \\ \vdots & \vdots & \vdots & \ddots & \ddots & \vdots \\ 0 & 0 & \cdots & q_{n-3} & d_{n-2} & p_{n-2} \\ 0 & 0 & \cdots & 0 & q_{n-2} & d_{n-1} \end{pmatrix} \begin{pmatrix} x_0 \\ x_1 \\ x_2 \\ x_3 \\ \vdots \\ x_{n-1} \end{pmatrix}.$$

This is useful for solving problems on the type

$$a_i = b_i a_{i-1} + c_i a_{i+1} + d_i, 1 \leq i \leq n,$$

where  $a_0, a_{n+1}, b_i, c_i$  and  $d_i$  are known.  $a$  can then be obtained from

$$\{a_i\} = \text{tridiagonal}(\{1, -1, -1, \dots, -1, 1\}, \{0, c_1, c_2, \dots, c_n\}, \{b_1, b_2, \dots, b_n, 0\}, \{a_0, d_1, d_2, \dots, d_n, a_{n+1}\}).$$

Fails if the solution is not unique.  
If  $|d_i| > |p_i| + |q_{i-1}|$  for all  $i$ , or  $|d_i| > |p_{i-1}| + |q_i|$ , or the matrix is positive definite, the algorithm is numerically stable and neither tr nor the check for diag[i] == 0 is needed.  
**Time:**  $\mathcal{O}(N)$

8f9fa8, 26 lines

```
typedef double T;
vector<T> tridiagonal(vector<T> diag, const vector<T>& super,
    const vector<T>& sub, vector<T> b) {
    int n = sz(b); vi tr(n);
    rep(i,0,n-1) {
        if (abs(diag[i]) < 1e-9 * abs(super[i])) { // diag[i] == 0
            b[i+1] -= b[i] * diag[i+1] / super[i];
            if (i+2 < n) b[i+2] -= b[i] * sub[i+1] / super[i];
            diag[i+1] = sub[i]; tr[++i] = 1;
        } else {
            diag[i+1] -= super[i]*sub[i]/diag[i];
            b[i+1] -= b[i]*sub[i]/diag[i];
        }
    }
    for (int i = n; i--;) {
        if (tr[i]) {
            swap(b[i], b[i-1]);
            diag[i-1] = diag[i];
            b[i] /= super[i-1];
        } else {
            b[i] /= diag[i];
            if (i) b[i-1] -= b[i]*super[i-1];
        }
    }
    return b;
}
```

7.5 Fourier transforms

FastFourierTransform.h

**Description:**  $\text{fft}(a)$  computes  $\hat{f}(k) = \sum_x a[x] \exp(2\pi i \cdot kx/N)$  for all  $k$ .  $N$  must be a power of 2. Useful for convolution:  $\text{conv}(a, b) = c$ , where  $c[i] = \sum a[i]b[x-i]$ . For convolution of complex numbers or more than two vectors: FFT, multiply pointwise, divide by  $n$ , reverse(start+1, end), FFT back. Rounding is safe if  $(\sum a_i^2 + \sum b_i^2) \log_2 N < 9 \cdot 10^{14}$  (in practice  $10^{16}$ ; higher for random inputs). Otherwise, use NTT/FFTMod.  
**Time:**  $\mathcal{O}(N \log N)$  with  $N = |A| + |B|$  ( $\sim 1s$  for  $N = 2^{22}$ )

00ced6, 35 lines

```
typedef complex<double> C;
typedef vector<double> vd;
void fft(vector<C>& a) {
    int n = sz(a), L = 31 - __builtin_clz(n);
    static vector<complex<long double>> R(2, 1);
    static vector<C> rt(2, 1); // (^ 10% faster if double)
    for (static int k = 2; k < n; k *= 2) {
        R.resize(n); rt.resize(n);
        auto x = polar(1.0L, acos(-1.0L) / k);
        rep(i,k,2*k) rt[i] = R[i] = i&1 ? R[i/2] * x : R[i/2];
    }
    vi rev(n);
    rep(i,0,n) rev[i] = (rev[i / 2] | (i & 1) << L) / 2;
    rep(i,0,n) if (i < rev[i]) swap(a[i], a[rev[i]]);
    for (int k = 1; k < n; k *= 2)
        for (int i = 0; i < n; i += 2 * k) rep(j,0,k) {
            C z = rt[j+k] * a[i+j+k]; // (25% faster if hand-rolled)
            a[i + j + k] = a[i + j] - z;
            a[i + j] += z;
        }
}
vd conv(const vd& a, const vd& b) {
    if (a.empty() || b.empty()) return {};
    vd res(sz(a) + sz(b) - 1);
    int L = 32 - __builtin_clz(sz(res)), n = 1 << L;
    vector<C> in(n), out(n);
    copy(all(a), begin(in));
    rep(i,0,sz(b)) in[i].imag(b[i]);
    fft(in);
    for (C& x : in) x *= x;
    rep(i,0,n) out[i] = in[-i & (n - 1)] - conj(in[i]);
    fft(out);
    rep(i,0,sz(res)) res[i] = imag(out[i]) / (4 * n);
    return res;
}
```

FastFourierTransformMod.h

**Description:** Higher precision FFT, can be used for convolutions modulo arbitrary integers as long as  $N \log_2 N \cdot \text{mod} < 8.6 \cdot 10^{14}$  (in practice  $10^{16}$  or higher). Inputs must be in  $[0, \text{mod})$ .  
**Time:**  $\mathcal{O}(N \log N)$ , where  $N = |A| + |B|$  (twice as slow as NTT or FFT)  
**"FastFourierTransform.h"**

b82773, 22 lines

```
typedef vector<ll> vl;
template<int M> vl convMod(const vl &a, const vl &b) {
    if (a.empty() || b.empty()) return {};
    vl res(sz(a) + sz(b) - 1);
    int B=32-__builtin_clz(sz(res)), n=1<<B, cut=int(sqrt(M));
    vector<C> L(n), R(n), outs(n), outl(n);
    rep(i,0,sz(a)) L[i] = C((int)a[i] / cut, (int)a[i] % cut);
    rep(i,0,sz(b)) R[i] = C((int)b[i] / cut, (int)b[i] % cut);
    fft(L), fft(R);
    rep(i,0,n) {
        int j = -i & (n - 1);
        outl[j] = (L[i] + conj(L[j])) * R[i] / (2.0 * n);
        outs[j] = (L[i] - conj(L[j])) * R[i] / (2.0 * n) / 1i;
    }
    fft(outl), fft(outs);
    rep(i,0,sz(res)) {
        ll av = ll(real(outl[i])+.5), cv = ll(imag(outs[i])+.5);
        ll bv = ll(imag(outl[i])+.5) + ll(real(outs[i])+.5);
        res[i] = ((av % M * cut + bv) % M * cut + cv) % M;
    }
}
```

```
    }
    return res;
}
```

NumberTheoreticTransform.h

**Description:** ntt(a) computes  $\hat{f}(k) = \sum_x a[x]g^{xk}$  for all  $k$ , where  $g = \text{root}^{(\text{mod}-1)/N}$ .  $N$  must be a power of 2. Useful for convolution modulo specific nice primes of the form  $2^a b + 1$ , where the convolution result has size at most  $2^a$ . For arbitrary modulo, see FFTMod.  $\text{conv}(a, b) = c$ , where  $c[x] = \sum a[i]b[x-i]$ . For manual convolution: NTT the inputs, multiply pointwise, divide by n, reverse(start+1, end), NTT back. Inputs must be in [0, mod).

```
Time:  $\mathcal{O}(N \log N)$ 
"../number-theory/ModPow.h"
const ll mod = (119 << 23) + 1, root = 62; // = 998244353
// For p < 2^30 there is also e.g. 5 << 25, 7 << 26, 479 << 21
// and 483 << 21 (same root). The last two are > 10^9.
typedef vector<ll> vl;
void ntt(vl &a) {
    int n = sz(a), L = 31 - __builtin_clz(n);
    static vl rt(2, 1);
    for (static int k = 2, s = 2; k < n; k *= 2, s++) {
        rt.resize(n);
        ll z[] = {1, modpow(root, mod >> s)};
        rep(i,k,2*k) rt[i] = rt[i / 2] * z[i & 1] % mod;
    }
    vi rev(n);
    rep(i,0,n) rev[i] = (rev[i / 2] | (i & 1) << L) / 2;
    rep(i,0,n) if (i < rev[i]) swap(a[i], a[rev[i]]);
    for (int k = 1; k < n; k *= 2)
        for (int i = 0; i < n; i += 2 * k) rep(j,0,k) {
            ll z = rt[j + k] * a[i + j + k] % mod, &ai = a[i + j];
            a[i + j + k] = ai - z + (z > ai ? mod : 0);
            ai += (ai + z >= mod ? z - mod : z);
        }
}
vl conv(const vl &a, const vl &b) {
    if (a.empty() || b.empty()) return {};
    int s = sz(a) + sz(b) - 1, B = 32 - __builtin_clz(s), n = 1
        << B;
    int inv = modpow(n, mod - 2);
    vl L(a), R(b), out(n);
    L.resize(n), R.resize(n);
    ntt(L), ntt(R);
    rep(i,0,n) out[-i & (n - 1)] = (ll)L[i] * R[i] % mod * inv %
        mod;
    ntt(out);
    return {out.begin(), out.begin() + s};
}
```

FastSubsetTransform.h

**Description:** Transform to a basis with fast convolutions of the form  $c[z] = \sum_{z=x \oplus y} a[x] \cdot b[y]$ , where  $\oplus$  is one of AND, OR, XOR. The size of  $a$  must be a power of two.

```
Time:  $\mathcal{O}(N \log N)$ 
464cf3, 16 lines
void FST(vi& a, bool inv) {
    for (int n = sz(a), step = 1; step < n; step *= 2) {
        for (int i = 0; i < n; i += 2 * step) rep(j,i,i+step) {
            int &u = a[j], &v = a[j + step]; tie(u, v) =
                inv ? pii(v - u, u) : pii(v, u + v); // AND
                inv ? pii(v, u - v) : pii(u + v, u); // OR
                pii(u + v, u - v); // XOR
        }
    }
    if (inv) for (int& x : a) x /= sz(a); // XOR only
}
vi conv(vi a, vi b) {
```

```
FST(a, 0); FST(b, 0);
rep(i,0,sz(a)) a[i] *= b[i];
FST(a, 1); return a;
}

WalshHadamard.h
Description:  $C_k = \sum_{i \otimes j = k} A_i B_j$ 
Usage: Apply the transform, point multiply and invert
Time:  $\mathcal{O}(N \log N)$ 
922b72, 11 lines
void WalshHadamard(Poly &P, bool invert) {
    for (int len = 1; 2 * len <= sz(P); len <= 1) {
        for (int i = 0; i < sz(P); i += 2 * len) {
            rep(j, 0, len) {
                auto u = P[i + j], v = P[i + len + j];
                P[i + j] = u + v, P[i + len + j] = u - v; // XOR
            }
        }
    }
    if (invert) for (auto &x : P) x /= sz(P);
}
```

OnlineFFT.h

```
Description: Given  $B_1, \dots B_m$ , compute  $A_i = \sum_{j=1}^{i-1} A_j * B_{i-j}$ 
Usage: 1-indexed, pad B[i] = 0 for i > m
Time:  $\mathcal{O}(N \log^2 N)$ 
c0e86b, 18 lines
void online(const Poly &B, CD al, int n, Poly &A) {
    const int m = SZ(B) - 1;
    A.assign(n + 1, 0); A[1] = al;
    auto bst = B.begin(), ast = A.begin();
    REP(i, 1, n) {
        A[i + 1] += A[i] * B[1];
        if (i + 2 <= n) A[i + 2] += A[i] * B[2];
        for (int pw = 2; i % pw == 0 && pw + 1 <= m; pw <= 1) {
            Poly blockA(ast + i - pw, ast + i);
            Poly blockB(bst + pw + 1, bst + min(pw * 2, m) + 1);
            Poly prod = conv(blockA, blockB);
            REP(j, 0, sz(prod)) {
                if (i + 1 + j <= n)
                    A[i + 1 + j] += prod[j];
            }
        }
    }
}
```

Geometry (8)

8.1 Basics Trigonometry

$\sin(v + w) = \sin v \cos w + \cos v \sin w$   
 $\cos(v + w) = \cos v \cos w - \sin v \sin w$

$\tan(v + w) = \frac{\tan v + \tan w}{1 - \tan v \tan w}$   
 $\sin v + \sin w = 2 \sin \frac{v + w}{2} \cos \frac{v - w}{2}$   
 $\cos v + \cos w = 2 \cos \frac{v + w}{2} \cos \frac{v - w}{2}$

$(V + W) \tan(v - w)/2 = (V - W) \tan(v + w)/2$

where  $V, W$  are lengths of sides opposite angles  $v, w$ .

$a \cos x + b \sin x = r \cos(x - \phi)$   
 $a \sin x + b \cos x = r \sin(x + \phi)$

where  $r = \sqrt{a^2 + b^2}, \phi = \text{atan2}(b, a)$ .

Triangles

Side lengths:  $a, b, c$   
Semiperimeter:  $p = \frac{a + b + c}{2}$   
Area:  $A = \sqrt{p(p - a)(p - b)(p - c)}$   
Circumradius:  $R = \frac{abc}{4A}$   
Inradius:  $r = \frac{A}{p}$

Length of median (divides triangle into two equal-area triangles):  
 $m_a = \frac{1}{2} \sqrt{2b^2 + 2c^2 - a^2}$   
Length of bisector (divides angles in two):

$s_a = \sqrt{bc \left[ 1 - \left( \frac{a}{b + c} \right)^2 \right]}$   
Law of sines:  $\frac{\sin \alpha}{a} = \frac{\sin \beta}{b} = \frac{\sin \gamma}{c} = \frac{1}{2R}$   
Law of cosines:  $a^2 = b^2 + c^2 - 2bc \cos \alpha$   
Law of tangents:  $\frac{a + b}{a - b} = \frac{\tan \frac{\alpha + \beta}{2}}{\tan \frac{\alpha - \beta}{2}}$

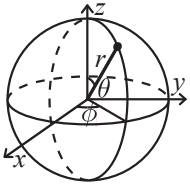
Quadrilaterals

With side lengths  $a, b, c, d$ , diagonals  $e, f$ , diagonals angle  $\theta$ , area  $A$  and magic flux  $F = b^2 + d^2 - a^2 - c^2$ :

$4A = 2ef \cdot \sin \theta = F \tan \theta = \sqrt{4e^2 f^2 - F^2}$

For cyclic quadrilaterals the sum of opposite angles is  $180^\circ$ ,  $ef = ac + bd$ , and  $A = \sqrt{(p - a)(p - b)(p - c)(p - d)}$ .

Spherical coordinates



$x = r \sin \theta \cos \phi$   
 $y = r \sin \theta \sin \phi$   
 $z = r \cos \theta$   
 $r = \sqrt{x^2 + y^2 + z^2}$   
 $\theta = \text{acos}(z / \sqrt{x^2 + y^2 + z^2})$   
 $\phi = \text{atan2}(y, x)$

Derivatives/Integrals

$$\frac{d}{dx} \arcsin x = \frac{1}{\sqrt{1-x^2}}$$
$$\frac{d}{dx} \tan x = 1 + \tan^2 x$$
$$\int \tan ax = -\frac{\ln|\cos ax|}{a}$$
$$\int e^{-x^2} = \frac{\sqrt{\pi}}{2}\text{erf}(x)$$

$$\frac{d}{dx} \arccos x = -\frac{1}{\sqrt{1-x^2}}$$
$$\frac{d}{dx} \arctan x = \frac{1}{1+x^2}$$
$$\int x \sin ax = \frac{\sin ax - ax \cos ax}{a^2}$$
$$\int xe^{ax}dx = \frac{e^{ax}}{a^2}(ax-1)$$

Integration by parts:

$$\int_a^b f(x)g(x)dx = [F(x)g(x)]_a^b - \int_a^b F(x)g'(x)dx$$

8.2 Geometric primitives

Point.h

Description: Class to handle points in the plane. T can be e.g. double or long long. (Avoid int.)

<bits/stdc++.h>e8d121, 44 lines

template <class T> int sgn(T x) { return (x > T(0)) - (x < T(0)); }

template <class T> struct Point { typedef Point P; T x, y; explicit Point(T x = 0, T y = 0) : x(x), y(y) {} bool operator<(P p) const { return tie(x, y) < tie(p.x, p.y); }

bool operator==(P p) const { return tie(x, y) == tie(p.x, p.y); }

P operator+(P p) const { return P(x + p.x, y + p.y); }

P operator-(P p) const { return P(x - p.x, y - p.y); }

P operator\*(T d) const { return P(x \* d, y \* d); }

P operator/(T d) const { return P(x / d, y / d); }

T dot(P p) const { return x \* p.x + y \* p.y; }

T cross(P p) const { return x \* p.y - y \* p.x; }

T cross(P a, P b) const { return (a - \*this).cross(b - \*this); }

T dist2() const { return x \* x + y \* y; }

// abs() == dist()

double dist() const { return sqrt((double)dist2()); }

// angle to x-axis in interval [-pi, pi]

double angle() const { return atan2(y, x); }

P unit() const { return \*this / dist(); }

// makes dist()==1

P perp() const { return P(-y, x); }

// rotates +90 degrees

P normal() const { return perp().unit(); }

P translate(P v) { return P(x + v.x, y + v.y); }

// scale an object by a certain ratio alpha around a

// center c, we need to shorten or lengthen the vector

// from c to every point by a factor alpha, while

// conserving the direction

P scale(P c, double factor) { return c + (\*this - c) \* factor; }

// returns point rotated 'a' radians ccw around the origin

P rotate(double a) const { return P(x \* cos(a) - y \* sin(a), x \* sin(a) + y \* cos(a)); }

friend ostream &operator<<(ostream &os, P p) { return os << "(" << p.x << ", " << p.y << ")"; }

// Additional random shit

bool isPerp(P p) { return P(x, y).dot(p) == 0; }

double angle(P p) { double costheta = P(x, y).dot(p) / (\*this).dist() / p.dist(); return acos(fmax(-1.0, fmin(1.0, costheta))); }

T orient(P b, P c) { return (b - \*this).cross(c - \*this); }

};

lineDistance.h

Description: Returns the signed distance between point p and the line containing points a and b. Positive value on left side and negative on right as seen from a towards b. a==b gives nan. P is supposed to be Point<T> or Point3D<T> where T is e.g. double or long long. It uses products in intermediate steps so watch out for overflow if using int or long long. Using Point3D will always give a non-negative distance. For Point3D, call .dist on the result of the cross product.

f6bf6b, 4 lines

template<class P> double lineDist(const P& a, const P& b, const P& p) { return (double) (b-a).cross(p-a) / (b-a).dist(); }

};

SegmentDistance.h

Description: Returns the shortest distance between point p and the line segment from point s to e.

Usage: Point<double> a, b(2,2), p(1,1); bool onSegment = segDist(a,b,p) < 1e-10;

5c88f4, 6 lines

typedef Point<double> P; double segDist(P& s, P& e, P& p) { if (s==e) return (p-s).dist(); auto d = (e-s).dist2(), t = min(d,max(.0, (p-s).dot(e-s))); return ((p-s)\*d-(e-s)\*t).dist()/d; }

};

SegmentIntersection.h

Description: If a unique intersection point between the line segments going from s1 to e1 and from s2 to e2 exists then it is returned. If no intersection point exists an empty vector is returned. If infinitely many exist a vector with 2 elements is returned, containing the endpoints of the common line segment. The wrong position will be returned if P is Point<ll> and the intersection point does not have integer coordinates. Products of three coordinates are used in intermediate steps so watch out for overflow if using int or long long.

Usage: vector<P> inter = segInter(s1,e1,s2,e2); if (sz(inter)==1) cout << "segments intersect at " << inter[0] << endl;

9d57f2, 13 lines

template<class P> vector<P> segInter(P a, P b, P c, P d) { auto oa = c.cross(d, a), ob = c.cross(d, b), oc = a.cross(b, c), od = a.cross(b, d); // Checks if intersection is single non-endpoint point. if (sgn(oa) \* sgn(ob) < 0 && sgn(oc) \* sgn(od) < 0) return {(a \* ob - b \* oa) / (ob - oa)}; set<P> s; if (onSegment(c, d, a)) s.insert(a); if (onSegment(c, d, b)) s.insert(b); if (onSegment(a, b, c)) s.insert(c); if (onSegment(a, b, d)) s.insert(d); return {all(s)}; }

lineIntersection.h

Description: If a unique intersection point of the lines going through s1,e1 and s2,e2 exists {1, point} is returned. If no intersection point exists {0, (0,0)} is returned and if infinitely many exists {-1, (0,0)} is returned. The wrong position will be returned if P is Point<ll> and the intersection point does not have integer coordinates. Products of three coordinates are used in intermediate steps so watch out for overflow if using int or ll.

Usage: auto res = lineInter(s1,e1,s2,e2); if (res.first == 1) cout << "intersection point at " << res.second << endl;

a01f81, 8 lines

template<class P> pair<int, P> lineInter(P s1, P e1, P s2, P e2) { auto d = (e1 - s1).cross(e2 - s2); if (d == 0) // if parallel return {-(s1.cross(e1, s2) == 0), P(0, 0)}; auto p = s2.cross(e1, e2), q = s2.cross(e2, s1); return {1, (s1 \* p + e1 \* q) / d}; }

sideOf.h

Description: Returns where p is as seen from s towards e. 1/0/-1 ⇔ left/on line/right. If the optional argument eps is given 0 is returned if p is within distance eps from the line. P is supposed to be Point<T> where T is e.g. double or long long. It uses products in intermediate steps so watch out for overflow if using int or long long.

Usage: bool left = sideOf(p1,p2,q)==1;

Point.h"3af81c, 9 lines

template<class P> int sideOf(P s, P e, P p) { return sgn(s.cross(e, p)); }

template<class P> int sideOf(const P& s, const P& e, const P& p, double eps) { auto a = (e-s).cross(p-s); double l = (e-s).dist()\*eps; return (a > l) - (a < -l); }

OnSegment.h

Description: Returns true iff p lies on the line segment from s to e. Use (segDist(s,e,p)<=epsilon) instead when using Point<double>.

Point.h" c597e8, 3 lines

template<class P> bool onSegment(P s, P e, P p) { return p.cross(s, e) == 0 && (s - p).dot(e - p) <= 0; }

linearTransformation.h

Description: Apply the linear transformation (translation, rotation and scaling) which takes line p0-p1 to line q0-q1 to point r.

Point.h"03a306, 6 lines

typedef Point<double> P; P linearTransformation(const P& p0, const P& p1, const P& q0, const P& q1, const P& r) { P dp = p1-p0, dq = q1-q0, num(dp.cross(dq), dp.dot(dq)); return q0 + P((r-p0).cross(num), (r-p0).dot(num))/dp.dist2(); }

Angle.h

Description: A class for ordering angles (as represented by int points and a number of rotations around the origin). Useful for rotational sweeping. Sometimes also represents points or vectors.

**Usage:** vector<Angle> v = {w[0], w[0].t360() ...}; // sorted

int j = 0; rep(i,0,n) { while (v[j] < v[i].t180()) ++j; }

// sweeps j such that (j-i) represents the number of positively oriented triangles with vertices at 0 and i

0f0602, 35 lines

```
struct Angle {
    int x, y;
    int t;
    Angle(int x, int y, int t=0) : x(x), y(y), t(t) {}
    Angle operator-(Angle b) const { return {x-b.x, y-b.y, t}; }
    int half() const {
        assert(x || y);
        return y < 0 || (y == 0 && x < 0);
    }
    Angle t90() const { return {-y, x, t + (half() && x >= 0)}; }
    Angle t180() const { return {-x, -y, t + half()}; }
    Angle t360() const { return {x, y, t + 1}; }
};

bool operator<(Angle a, Angle b) {
    // add a.dist2() and b.dist2() to also compare distances
    return make_tuple(a.t, a.half(), a.y * (11)b.x) <
           make_tuple(b.t, b.half(), a.x * (11)b.y);
}

// Given two points, this calculates the smallest angle between
// them, i.e., the angle that covers the defined line segment.
pair<Angle, Angle> segmentAngles(Angle a, Angle b) {
    if (b < a) swap(a, b);
    return (b < a.t180() ?
           make_pair(a, b) : make_pair(b, a.t360()));
}

Angle operator+(Angle a, Angle b) { // point a + vector b
    Angle r(a.x + b.x, a.y + b.y, a.t);
    if (a.t180() < r) r.t--;
    return r.t180() < a ? r.t360() : r;
}

Angle angleDiff(Angle a, Angle b) { // angle b - angle a
    int tu = b.t - a.t; a.t = b.t;
    return {a.x*b.x + a.y*b.y, a.x*b.y - a.y*b.x, tu - (b < a)};
}
```

CircleIntersection.h

**Description:** Computes the pair of points at which two circles intersect. Returns false in case of no intersection.

Point.h"84d6d3, 11 lines

```
typedef Point<double> P;
bool circleInter(P a,P b,double r1,double r2,pair<P, P>* out) {
    if (a == b) { assert(r1 != r2); return false; }
    P vec = b - a;
    double d2 = vec.dist2(), sum = r1+r2, dif = r1-r2,
           p = (d2 + r1*r1 - r2*r2)/(d2*2), h2 = r1*r1 - p*p*d2;
    if (sum*sum < d2 || dif*dif > d2) return false;
    P mid = a + vec*p, per = vec.perp() * sqrt(fmax(0, h2) / d2);
    *out = {mid + per, mid - per};
    return true;
}
```

CircleTangents.h

**Description:** Finds the external tangents of two circles, or internal if r2 is negated. Can return 0, 1, or 2 tangents – 0 if one circle contains the other (or overlaps it, in the internal case, or if the circles are the same); 1 if the circles are tangent to each other (in which case .first = .second and the tangent line is perpendicular to the line between the centers). .first and .second give the tangency points at circle 1 and 2 respectively. To find the tangents of a circle with a point set r2 to 0.

Point.h"b0153d, 13 lines

```
template<class P>
```

```
vector<pair<P, P>> tangents(P c1, double r1, P c2, double r2) {
    P d = c2 - c1;
    double dr = r1 - r2, d2 = d.dist2(), h2 = d2 - dr * dr;
    if (d2 == 0 || h2 < 0) return {};
    vector<pair<P, P>> out;
    for (double sign : {-1, 1}) {
        P v = (d * dr + d.perp() * sqrt(h2) * sign) / d2;
        out.push_back({c1 + v * r1, c2 + v * r2});
    }
    if (h2 == 0) out.pop_back();
    return out;
}
```

CirclePolygonIntersection.h

**Description:** Returns the area of the intersection of a circle with a ccw polygon.

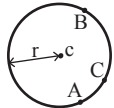
**Time:**  $\mathcal{O}(n)$

Point.h"alee63, 19 lines

```
typedef Point<double> P;
#define arg(p, q) atan2(p.cross(q), p.dot(q))
double circlePoly(P c, double r, vector<P> ps) {
    auto tri = [&](P p, P q) {
        auto r2 = r * r / 2;
        P d = q - p;
        auto a = d.dot(p)/d.dist2(), b = (p.dist2()-r*r)/d.dist2();
        auto det = a * a - b;
        if (det <= 0) return arg(p, q) * r2;
        auto s = max(0., -a-sqrt(det)), t = min(1., -a+sqrt(det));
        if (t < 0 || 1 <= s) return arg(p, q) * r2;
        P u = p + d * s, v = p + d * t;
        return arg(p,u) * r2 + u.cross(v)/2 + arg(v,q) * r2;
    };
    auto sum = 0.0;
    rep(i,0,sz(ps))
        sum += tri(ps[i] - c, ps[(i + 1) % sz(ps)] - c);
    return sum;
}
```

circumcircle.h

**Description:**



The circumcircle of a triangle is the circle intersecting all three vertices. ccRadius returns the radius of the circle going through points A, B and C and ccCenter returns the center of the same circle.

Point.h"1caa3a, 9 lines

```
typedef Point<double> P;
double ccRadius(const P& A, const P& B, const P& C) {
    return (B-A).dist()*(C-B).dist()*(A-C).dist()/
           abs((B-A).cross(C-A))/2;
}

P ccCenter(const P& A, const P& B, const P& C) {
    P b = C-A, c = B-A;
    return A + (b*c.dist2()-c*b.dist2()).perp()/b.cross(c)/2;
}
```

MinimumEnclosingCircle.h

**Description:** Computes the minimum circle that encloses a set of points.

**Time:** expected  $\mathcal{O}(n)$

circumcircle.h"09dd0a, 17 lines

```
pair<P, double> mec(vector<P> ps) {
    shuffle(all(ps), mt19937(time(0)));
    P o = ps[0];
    double r = 0, EPS = 1 + 1e-8;
    rep(i,0,sz(ps)) if ((o - ps[i]).dist() > r * EPS) {
        o = ps[i], r = 0;
        rep(j,0,i) if ((o - ps[j]).dist() > r * EPS) {
            o = (ps[i] + ps[j]) / 2;

```

```
        r = (o - ps[i]).dist();
        rep(k,0,j) if ((o - ps[k]).dist() > r * EPS) {
            o = ccCenter(ps[i], ps[j], ps[k]);
            r = (o - ps[i]).dist();
        }
    }
    return {o, r};
}
```

## 8.4 Polygons

InsidePolygon.h

**Description:** Returns true if p lies within the polygon. If strict is true, it returns false for points on the boundary. The algorithm uses products in intermediate steps so watch out for overflow.

**Usage:** vector<P> v = {P{4,4}, P{1,2}, P{2,1}};

bool in = inPolygon(ps[i], ps[j], ps[k]);

**Time:**  $\mathcal{O}(n)$

Point.h", "OnSegment.h", "SegmentDistance.h"2bf504, 11 lines

```
template<class P>
bool inPolygon(vector<P> &p, P a, bool strict = true) {
    int cnt = 0, n = sz(p);
    rep(i,0,n) {
        P q = p[(i + 1) % n];
        if (onSegment(p[i], q, a)) return !strict;
        //or: if (segDist(p[i], q, a) <= eps) return !strict;
        cnt ^= ((a.y<p[i].y) - (a.y<q.y)) * a.cross(p[i], q) > 0;
    }
    return cnt;
}
```

PolygonArea.h

**Description:** Returns twice the signed area of a polygon. Clockwise enumeration gives negative area. Watch out for overflow if using int as T!

Point.h"f12300, 6 lines

```
template<class T>
T polygonArea2(vector<Point<T>&& v) {
    T a = v.back().cross(v[0]);
    rep(i,0,sz(v)-1) a += v[i].cross(v[i+1]);
    return a;
}
```

PolygonCenter.h

**Description:** Returns the center of mass for a polygon.

**Time:**  $\mathcal{O}(n)$

Point.h"9706dc, 9 lines

```
typedef Point<double> P;
P polygonCenter(const vector<P>& v) {
    P res(0, 0); double A = 0;
    for (int i = 0, j = sz(v) - 1; i < sz(v); j = i++) {
        res = res + (v[i] + v[j]) * v[j].cross(v[i]);
        A += v[j].cross(v[i]);
    }
    return res / A / 3;
}
```

PolygonCut.h

**Description:**

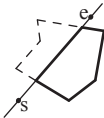
Returns a vector with the vertices of a polygon with everything to the left of the line going from s to e cut away.

**Usage:** vector<P> p = ...;

p = polygonCut(p, P(0,0), P(1,0));

Point.h", "LineIntersection.h"f2b7d4, 13 lines

```
typedef Point<double> P;
vector<P> polygonCut(const vector<P>& poly, P s, P e) {
    vector<P> res;
```



```
rep(i,0,sz(poly)) {
    P cur = poly[i], prev = i ? poly[i-1] : poly.back();
    bool side = s.cross(e, cur) < 0;
    if (side != (s.cross(e, prev) < 0))
        res.push_back(lineInter(s, e, cur, prev).second);
    if (side)
        res.push_back(cur);
}
return res;
}
```

ConvexHull.h

**Description:** Returns a vector of the points of the convex hull in counter-clockwise order. Points on the edge of the hull between two other points are not considered part of the hull.  
**Time:**  $\mathcal{O}(n \log n)$



```
"Point.h" 310954, 13 lines
typedef Point<ll> P;
vector<P> convexHull(vector<P> pts) {
    if (sz(pts) <= 1) return pts;
    sort(all(pts));
    vector<P> h(sz(pts)+1);
    int s = 0, t = 0;
    for (int it = 2; it--; s = --t, reverse(all(pts)))
        for (P p : pts) {
            while (t >= s + 2 && h[t-2].cross(h[t-1], p) <= 0) t--;
            h[t++] = p;
        }
    return {h.begin(), h.begin() + t - (t == 2 && h[0] == h[1])};
}
```

HullDiameter.h

**Description:** Returns the two points with max distance on a convex hull (ccw, no duplicate/collinear points).  
**Time:**  $\mathcal{O}(n)$

```
"Point.h" c571b8, 12 lines
typedef Point<ll> P;
array<P, 2> hullDiameter(vector<P> S) {
    int n = sz(S), j = n < 2 ? 0 : 1;
    pair<ll, array<P, 2>> res({0, {S[0], S[0]}});
    rep(i,0,j)
        for (; j = (j + 1) % n) {
            res = max(res, {(S[i] - S[j]).dist2(), {S[i], S[j]}});
            if ((S[(j + 1) % n] - S[j]).cross(S[i + 1] - S[i]) >= 0)
                break;
        }
    return res.second;
}
```

PointInsideHull.h

**Description:** Determine whether a point t lies inside a convex hull (CCW order, with no collinear points). Returns true if point lies within the hull. If strict is true, points on the boundary aren't included.  
**Time:**  $\mathcal{O}(\log N)$

```
"Point.h", "sideOf.h", "OnSegment.h" 71446b, 14 lines
typedef Point<ll> P;

bool inHull(const vector<P>& l, P p, bool strict = true) {
    int a = 1, b = sz(l) - 1, r = !strict;
    if (sz(l) < 3) return r && onSegment(l[0], l.back(), p);
    if (sideOf(l[0], l[a], l[b]) > 0) swap(a, b);
    if (sideOf(l[0], l[a], p) >= r || sideOf(l[0], l[b], p) <= -r)
        return false;
    while (abs(a - b) > 1) {
        int c = (a + b) / 2;
        (sideOf(l[0], l[c], p) > 0 ? b : a) = c;
    }
}
```

```
return sgn(l[a].cross(l[b], p)) < r;
}
```

LineHullIntersection.h

**Description:** Line-convex polygon intersection. The polygon must be ccw and have no collinear points. lineHull(line, poly) returns a pair describing the intersection of a line with the polygon:  $\bullet(-1, -1)$  if no collision,  $\bullet(i, -1)$  if touching the corner  $i$ ,  $\bullet(i, i)$  if along side  $(i, i + 1)$ ,  $\bullet(i, j)$  if crossing sides  $(i, i + 1)$  and  $(j, j + 1)$ . In the last case, if a corner  $i$  is crossed, this is treated as happening on side  $(i, i + 1)$ . The points are returned in the same order as the line hits the polygon. extrVertex returns the point of a hull with the max projection onto a line.  
**Time:**  $\mathcal{O}(\log n)$

```
"Point.h" 7cf45b, 39 lines
#define cmp(i, j) sgn(dir.perp().cross(poly[(i)%n]-poly[(j)%n]))
#define extr(i) cmp(i + 1, i) >= 0 && cmp(i, i - 1 + n) < 0
template <class P> int extrVertex(vector<P>& poly, P dir) {
    int n = sz(poly), lo = 0, hi = n;
    if (extr(0)) return 0;
    while (lo + 1 < hi) {
        int m = (lo + hi) / 2;
        if (extr(m)) return m;
        int ls = cmp(lo + 1, lo), ms = cmp(m + 1, m);
        (ls < ms || (ls == ms && ls == cmp(lo, m)) ? hi : lo) = m;
    }
    return lo;
}
```

```
#define cmpL(i) sgn(a.cross(poly[i], b))
template <class P>
array<int, 2> lineHull(P a, P b, vector<P>& poly) {
    int endA = extrVertex(poly, (a - b).perp());
    int endB = extrVertex(poly, (b - a).perp());
    if (cmpL(endA) < 0 || cmpL(endB) > 0)
        return {-1, -1};
    array<int, 2> res;
    rep(i,0,2) {
        int lo = endB, hi = endA, n = sz(poly);
        while ((lo + 1) % n != hi) {
            int m = ((lo + hi + (lo < hi ? 0 : n)) / 2) % n;
            (cmpL(m) == cmpL(endB) ? lo : hi) = m;
        }
        res[i] = (lo + !cmpL(hi)) % n;
        swap(endA, endB);
    }
    if (res[0] == res[1]) return {res[0], -1};
    if (!cmpL(res[0]) && !cmpL(res[1]))
        switch ((res[0] - res[1] + sz(poly) + 1) % sz(poly)) {
            case 0: return {res[0], res[0]};
            case 2: return {res[1], res[1]};
        }
    return res;
}
```

8.5 Misc. Point Set Problems

ClosestPair.h

**Description:** Finds the closest pair of points.  
**Time:**  $\mathcal{O}(n \log n)$

```
"Point.h" ac41a6, 17 lines
typedef Point<ll> P;
pair<P, P> closest(vector<P> v) {
    assert(sz(v) > 1);
    set<P> S;
    sort(all(v), [](P a, P b) { return a.y < b.y; });
    pair<ll, pair<P, P>> ret{LLONG_MAX, {P(), P()}};
    int j = 0;
    for (P p : v) {
        P d{1 + (ll)sqrt(ret.first), 0};

```

```
while (v[j].y <= p.y - d.x) S.erase(v[j++]);
auto lo = S.lower_bound(p - d), hi = S.upper_bound(p + d);
for (; lo != hi; ++lo)
    ret = min(ret, {(lo - p).dist2(), {lo, p}});
S.insert(p);
}
return ret.second;
}
```

kdTree.h

**Description:** KD-tree (2d, can be extended to 3d)

```
"Point.h" bac5b0, 63 lines
typedef long long T;
typedef Point<T> P;
const T INF = numeric_limits<T>::max();
```

```
bool on_x(const P& a, const P& b) { return a.x < b.x; }
bool on_y(const P& a, const P& b) { return a.y < b.y; }
```

```
struct Node {
    P pt; // if this is a leaf, the single point in it
    T x0 = INF, x1 = -INF, y0 = INF, y1 = -INF; // bounds
    Node *first = 0, *second = 0;

    T distance(const P& p) { // min squared distance to a point
        T x = (p.x < x0 ? x0 : p.x > x1 ? x1 : p.x);
        T y = (p.y < y0 ? y0 : p.y > y1 ? y1 : p.y);
        return (P(x,y) - p).dist2();
    }
}
```

```
Node(vector<P>&& vp) : pt(vp[0]) {
    for (P p : vp) {
        x0 = min(x0, p.x); x1 = max(x1, p.x);
        y0 = min(y0, p.y); y1 = max(y1, p.y);
    }
    if (vp.size() > 1) {
        // split on x if width >= height (not ideal...)
        sort(all(vp), x1 - x0 >= y1 - y0 ? on_x : on_y);
        // divide by taking half the array for each child (not
        // best performance with many duplicates in the middle)
        int half = sz(vp)/2;
        first = new Node({vp.begin(), vp.begin() + half});
        second = new Node({vp.begin() + half, vp.end()});
    }
};
```

```
struct KDTree {
    Node* root;
    KDTree(const vector<P>& vp) : root(new Node({all(vp)})) {}
}
```

```
pair<T, P> search(Node *node, const P& p) {
    if (!node->first) {
        // uncomment if we should not find the point itself:
        // if (p == node->pt) return {INF, P()};
        return make_pair((p - node->pt).dist2(), node->pt);
    }
}
```

```
Node *f = node->first, *s = node->second;
T bfirst = f->distance(p), bsec = s->distance(p);
if (bfirst > bsec) swap(bsec, bfirst), swap(f, s);
```

```
// search closest side first, other side if needed
auto best = search(f, p);
if (bsec < best.first)
    best = min(best, search(s, p));
return best;
}
```



```
// find nearest point to a point, and its squared distance
// (requires an arbitrary operator< for Point)
pair<T, P> nearest(const P& p) {
    return search(root, p);
}
};
```

FastDelaunay.h

**Description:** Fast Delaunay triangulation. Each circumcircle contains none of the input points. There must be no duplicate points. If all points are on a line, no triangles will be returned. Should work for doubles as well, though there may be precision issues in 'circ'. Returns triangles in order {t[0][0], t[0][1], t[0][2], t[1][0], ... }, all counter-clockwise.  
**Time:**  $O(n \log n)$

"Point.h"	eefdf5, 88 lines
-----------	------------------

```
typedef Point<ll> P;
typedef struct Quad* Q;
typedef __int128_t ll1; // (can be ll if coords are < 2e4)
P arb(LLONG_MAX,LLONG_MAX); // not equal to any other point
```

```
struct Quad {
    Q rot, o; P p = arb; bool mark;
    P& F() { return r()->p; }
    Q& r() { return rot->rot; }
    Q prev() { return rot->o->rot; }
    Q next() { return r()->prev(); }
} *H;
```

```
bool circ(P p, P a, P b, P c) { // is p in the circumcircle?
    ll1 p2 = p.dist2(), A = a.dist2()-p2,
        B = b.dist2()-p2, C = c.dist2()-p2;
    return p.cross(a,b)*C + p.cross(b,c)*A + p.cross(c,a)*B > 0;
}
```

```
Q makeEdge(P orig, P dest) {
    Q r = H ? H : new Quad{new Quad{new Quad{new Quad{0}}}};
    H = r->o; r->r()->r() = r;
    rep(i,0,4) r = r->rot, r->p = arb, r->o = i & 1 ? r : r->r();
    r->p = orig; r->F() = dest;
    return r;
}
```

```
void splice(Q a, Q b) {
    swap(a->o->rot->o, b->o->rot->o); swap(a->o, b->o);
}
```

```
Q connect(Q a, Q b) {
    Q q = makeEdge(a->F(), b->p);
    splice(q, a->next());
    splice(q->r(), b);
    return q;
}
```

```
pair<Q,Q> rec(const vector<P>& s) {
    if (sz(s) <= 3) {
        Q a = makeEdge(s[0], s[1]), b = makeEdge(s[1], s.back());
        if (sz(s) == 2) return { a, a->r() };
        splice(a->r(), b);
        auto side = s[0].cross(s[1], s[2]);
        Q c = side ? connect(b, a) : 0;
        return {side < 0 ? c->r() : a, side < 0 ? c : b->r() };
    }
}
```

```
#define H(e) e->F(), e->p
#define valid(e) (e->F().cross(H(base)) > 0)
Q A, B, ra, rb;
int half = sz(s) / 2;
tie(ra, A) = rec({all(s) - half});
tie(B, rb) = rec({sz(s) - half + all(s)});
while ((B->p.cross(H(A)) < 0 && (A = A->next())) ||
        (A->p.cross(H(B)) > 0 && (B = B->r()->o)));
Q base = connect(B->r(), A);
```

FastDelaunay PolyhedronVolume Point3D 3dHull

```
if (A->p == ra->p) ra = base->r();
if (B->p == rb->p) rb = base;

#define DEL(e, init, dir) Q e = init->dir; if (valid(e)) \
    while (circ(e->dir->F(), H(base), e->F())) { \
        Q t = e->dir; \
        splice(e, e->prev()); \
        splice(e->r(), e->r()->prev()); \
        e->o = H; H = e; e = t; \
    }
for (;;) {
    DEL(LC, base->r(), o); DEL(RC, base, prev());
    if (!valid(LC) && !valid(RC)) break;
    if (!valid(LC) || (valid(RC) && circ(H(RC), H(LC))))
        base = connect(RC, base->r());
    else
        base = connect(base->r(), LC->r());
}
return { ra, rb };
```

```
vector<P> triangulate(vector<P> pts) {
    sort(all(pts)); assert(unique(all(pts)) == pts.end());
    if (sz(pts) < 2) return {};
    Q e = rec(pts).first;
    vector<Q> q = {e};
    int qi = 0;
    while (e->o->F().cross(e->F(), e->p) < 0) e = e->o;
    #define ADD { Q c = e; do { c->mark = 1; pts.push_back(c->p); \
        q.push_back(c->r()); c = c->next(); } while (c != e); }
    ADD; pts.clear();
    while (qi < sz(q)) if (!(e = q[qi++])->mark) ADD;
    return pts;
}
```

8.6 3D

PolyhedronVolume.h

**Description:** Magic formula for the volume of a polyhedron. Faces should point outwards.

	3058c3, 6 lines
--	-----------------

```
template<class V, class L>
double signedPolyVolume(const V& p, const L& trilst) {
    double v = 0;
    for (auto i : trilst) v += p[i.a].cross(p[i.b]).dot(p[i.c]);
    return v / 6;
}
```

Point3D.h

**Description:** Class to handle points in 3D space. T can be e.g. double or long long.

	8058ae, 32 lines
--	------------------

```
template<class T> struct Point3D {
    typedef Point3D P;
    typedef const P& R;
    T x, y, z;
    explicit Point3D(T x=0, T y=0, T z=0) : x(x), y(y), z(z) {}
    bool operator<(R p) const {
        return tie(x, y, z) < tie(p.x, p.y, p.z); }
    bool operator==(R p) const {
        return tie(x, y, z) == tie(p.x, p.y, p.z); }
    P operator+(R p) const { return P(x+p.x, y+p.y, z+p.z); }
    P operator-(R p) const { return P(x-p.x, y-p.y, z-p.z); }
    P operator*(T d) const { return P(x*d, y*d, z*d); }
    P operator/(T d) const { return P(x/d, y/d, z/d); }
    T dot(R p) const { return x*p.x + y*p.y + z*p.z; }
    P cross(R p) const {
        return P(y*p.z - z*p.y, z*p.x - x*p.z, x*p.y - y*p.x);
    }
    T dist2() const { return x*x + y*y + z*z; }
```

```
double dist() const { return sqrt((double)dist2()); }
//Azimuthal angle (longitude) to x-axis in interval [-pi, pi]
double phi() const { return atan2(y, x); }
//Zenith angle (latitude) to the z-axis in interval [0, pi]
double theta() const { return atan2(sqrt(x*x+y*y),z); }
P unit() const { return *this/(T)dist(); } //makes dist()==1
//returns unit vector normal to *this and p
P normal(P p) const { return cross(p).unit(); }
//returns point rotated 'angle' radians ccw around axis
P rotate(double angle, P axis) const {
    double s = sin(angle), c = cos(angle); P u = axis.unit();
    return u*dot(u)*(1-c) + (*this)*c - cross(u)*s;
}
};
```

3dHull.h

**Description:** Computes all faces of the 3-dimension hull of a point set. \*No four points must be coplanar\*, or else random results will be returned. All faces will point outwards.

**Time:**  $O(n^2)$

"Point3D.h"	5b45fc, 49 lines
-------------	------------------

```
typedef Point3D<double> P3;
```

```
struct PR {
    void ins(int x) { (a == -1 ? a : b) = x; }
    void rem(int x) { (a == x ? a : b) = -1; }
    int cnt() { return (a != -1) + (b != -1); }
    int a, b;
};
```

```
struct F { P3 q; int a, b, c; };
```

```
vector<F> hull3d(const vector<P3>& A) {
    assert(sz(A) >= 4);
    vector<vector<PR>> E(sz(A), vector<PR>(sz(A), {-1, -1}));
    #define E(x,y) E[f.x][f.y]
    vector<F> FS;
    auto mf = [&](int i, int j, int k, int l) {
        P3 q = (A[j] - A[i]).cross((A[k] - A[i]));
        if (q.dot(A[l]) > q.dot(A[i]))
            q = q * -1;
        F f{q, i, j, k};
        E(a,b).ins(k); E(a,c).ins(j); E(b,c).ins(i);
        FS.push_back(f);
    };
    rep(i,0,4) rep(j,i+1,4) rep(k,j+1,4)
        mf(i, j, k, 6 - i - j - k);
```

```
rep(i,4,sz(A)) {
    rep(j,0,sz(FS)) {
        F f = FS[j];
        if (f.q.dot(A[i]) > f.q.dot(A[f.a])) {
            E(a,b).rem(f.c);
            E(a,c).rem(f.b);
            E(b,c).rem(f.a);
            swap(FS[j--], FS.back());
            FS.pop_back();
        }
        int nw = sz(FS);
        rep(j,0,nw) {
            F f = FS[j];
        }
        #define C(a, b, c) if (E(a,b).cnt() != 2) mf(f.a, f.b, i, f.c);
        C(a, b, c); C(a, c, b); C(b, c, a);
    }
}
for (F& it : FS) if ((A[it.b] - A[it.a]).cross(
    A[it.c] - A[it.a]).dot(it.q) <= 0) swap(it.c, it.b);
return FS;
```

```
};
```

sphericalDistance.h

**Description:** Returns the shortest distance on the sphere with radius radius between the points with azimuthal angles (longitude) f1 ( $\phi_1$ ) and f2 ( $\phi_2$ ) from x axis and zenith angles (latitude) t1 ( $\theta_1$ ) and t2 ( $\theta_2$ ) from z axis (0 = north pole). All angles measured in radians. The algorithm starts by converting the spherical coordinates to cartesian coordinates so if that is what you have you can use only the two last rows. dx\*radius is then the difference between the two points in the x direction and d\*radius is the total distance between the points.

611f07, 8 lines

```
double sphericalDistance(double f1, double t1,
    double f2, double t2, double radius) {
    double dx = sin(t2)*cos(f2) - sin(t1)*cos(f1);
    double dy = sin(t2)*sin(f2) - sin(t1)*sin(f1);
    double dz = cos(t2) - cos(t1);
    double d = sqrt(dx*dx + dy*dy + dz*dz);
    return radius*2*asin(d/2);
}
```

Various (9)

9.1 Intervals

IntervalContainer.h

**Description:** Add and remove intervals from a set of disjoint intervals. Will merge the added interval with any overlapping intervals in the set when adding. Intervals are [inclusive, exclusive).

**Time:**  $\mathcal{O}(\log N)$

edce47, 23 lines

```
set<pii>::iterator addInterval(set<pii>& is, int L, int R) {
    if (L == R) return is.end();
    auto it = is.lower_bound({L, R}), before = it;
    while (it != is.end() && it->first <= R) {
        R = max(R, it->second);
        before = it = is.erase(it);
    }
    if (it != is.begin() && (--it)->second >= L) {
        L = min(L, it->first);
        R = max(R, it->second);
        is.erase(it);
    }
    return is.insert(before, {L,R});
}
```

```
void removeInterval(set<pii>& is, int L, int R) {
    if (L == R) return;
    auto it = addInterval(is, L, R);
    auto r2 = it->second;
    if (it->first == L) is.erase(it);
    else (int&)it->second = L;
    if (R != r2) is.emplace(R, r2);
}
```

9e9d8d, 19 lines

IntervalCover.h

**Description:** Compute indices of smallest set of intervals covering another interval. Intervals should be [inclusive, exclusive). To support [inclusive, inclusive], change (A) to add || R.empty(). Returns empty set on failure (or if G is empty).

**Time:**  $\mathcal{O}(N \log N)$

```
template<class T>
vi cover(pair<T, T> G, vector<pair<T, T>> I) {
    vi S(sz(I)), R;
    iota(all(S), 0);
    sort(all(S), [&](int a, int b) { return I[a] < I[b]; });
    T cur = G.first;
```

```
int at = 0;
while (cur < G.second) { // (A)
    pair<T, int> mx = make_pair(cur, -1);
    while (at < sz(I) && I[S[at]].first <= cur) {
        mx = max(mx, make_pair(I[S[at]].second, S[at]));
        at++;
    }
    if (mx.second == -1) return {};
    cur = mx.first;
    R.push_back(mx.second);
}
return R;
```

753a4c, 19 lines

ConstantIntervals.h

**Description:** Split a monotone function on [from, to) into a minimal set of half-open intervals on which it has the same value. Runs a callback g for each such interval.

**Usage:** constantIntervals(0, sz(v), [&](int x){return v[x];}, [&](int lo, int hi, T val){...});

**Time:**  $\mathcal{O}(k \log \frac{n}{k})$

```
template<class F, class G, class T>
void rec(int from, int to, F& f, G& g, int& i, T& p, T q) {
    if (p == q) return;
    if (from == to) {
        g(i, to, p);
        i = to; p = q;
    } else {
        int mid = (from + to) >> 1;
        rec(from, mid, f, g, i, p, f(mid));
        rec(mid+1, to, f, g, i, p, q);
    }
}

template<class F, class G>
void constantIntervals(int from, int to, F f, G g) {
    if (to <= from) return;
    int i = from; auto p = f(i), q = f(to-1);
    rec(from, to-1, f, g, i, p, q);
    g(i, to, q);
}
```

pt7ea2, 27 lines

9.2 Misc. algorithms

Mint.h

**Description:** Modular Arithmetic (Integer) for Aditya and Ario

```
template <int32_t MOD>
struct mint {
    int32_t value;
    mint() : value() {}
    mint(int64_t value_) : value(value_ < 0 ? value_ % MOD + MOD
        : value_ >= MOD ? value_ % MOD : value_) {}
    mint(int32_t value_, std::nullptr_t) : value(value_) {}
    explicit operator bool() const { return value; }
    inline mint<MOD> operator + (mint<MOD> other) const { return
        mint<MOD>(*this) += other; }
    inline mint<MOD> operator - (mint<MOD> other) const { return
        mint<MOD>(*this) -= other; }
    inline mint<MOD> operator * (mint<MOD> other) const { return
        mint<MOD>(*this) *= other; }
    inline mint<MOD> & operator += (mint<MOD> other) { this->
        value += other.value; if (this->value >= MOD) this->
        value -= MOD; return *this; }
    inline mint<MOD> & operator -= (mint<MOD> other) { this->
        value -= other.value; if (this->value < 0) this->
        value += MOD; return *this; }
    inline mint<MOD> & operator *= (mint<MOD> other) { this->
        value = (uint_fast64_t)this->value * other.value % MOD;
        return *this; }
```

```
inline mint<MOD> operator - () const { return mint<MOD>(this
->value ? MOD - this->value : 0, nullptr); }
inline bool operator == (mint<MOD> other) const { return
    value == other.value; }
inline bool operator != (mint<MOD> other) const { return
    value != other.value; }
inline mint<MOD> pow(uint64_t k) const { return mint<MOD>(
    powa(value, k, MOD), nullptr); }
inline mint<MOD> inv() const { return mint<MOD>(powa(value,
    MOD-2, MOD), nullptr); }
inline mint<MOD> operator / (mint<MOD> other) const { return
    *this * other.inv(); }
inline mint<MOD> & operator /= (mint<MOD> other) { return *
    this *= other.inv(); }

};
template <int32_t MOD> mint<MOD> operator + (int64_t value,
    mint<MOD> n) { return mint<MOD>(value) + n; }
template <int32_t MOD> mint<MOD> operator - (int64_t value,
    mint<MOD> n) { return mint<MOD>(value) - n; }
template <int32_t MOD> mint<MOD> operator * (int64_t value,
    mint<MOD> n) { return mint<MOD>(value) * n; }
template <int32_t MOD> mint<MOD> operator / (int64_t value,
    mint<MOD> n) { return mint<MOD>(value) / n; }
template <int32_t MOD> std::istream & operator >> (std::istream
    & in, mint<MOD> & n) { int64_t value; in >> value; n =
    value; return in; }
template <int32_t MOD> std::ostream & operator << (std::ostream
    & out, mint<MOD> n) { return out << n.value; }
```

TernarySearch.h

**Description:** Find the smallest i in  $[a, b]$  that maximizes  $f(i)$ , assuming that  $f(a) < \dots < f(i) \geq \dots \geq f(b)$ . To reverse which of the sides allows non-strict inequalities, change the < marked with (A) to <=, and reverse the loop at (B). To minimize f, change it to >, also at (B).

**Usage:** int ind = ternSearch(0,n-1,&)(int i){return a[i];});

**Time:**  $\mathcal{O}(\log(b-a))$

9155b4, 11 lines

```
template<class F>
int ternSearch(int a, int b, F f) {
    assert(a <= b);
    while (b - a >= 5) {
        int mid = (a + b) / 2;
        if (f(mid) < f(mid+1)) a = mid; // (A)
        else b = mid+1;
    }
    rep(i,a+1,b+1) if (f(a) < f(i)) a = i; // (B)
    return a;
}
```

LIS.h

**Description:** Compute indices for the longest increasing subsequence.

**Time:**  $\mathcal{O}(N \log N)$

2932a0, 17 lines

```
template<class I> vi lis(const vector<I>& S) {
    if (S.empty()) return {};
    vi prev(sz(S));
    typedef pair<I, int> p;
    vector<p> res;
    rep(i,0,sz(S)) {
        // change 0 -> i for longest non-decreasing subsequence
        auto it = lower_bound(all(res), p{S[i], 0});
        if (it == res.end()) res.emplace_back(), it = res.end()-1;
        *it = {S[i], i};
        prev[i] = it == res.begin() ? 0 : (it-1)->second;
    }
    int L = sz(res), cur = res.back().second;
    vi ans(L);
    while (L--) ans[L] = cur, cur = prev[cur];
    return ans;
```



```
}

```

FastKnapsack.h

**Description:** Given N non-negative integer weights w and a non-negative target t, computes the maximum S <= t such that S is the sum of some subset of the weights.

**Time:**  $\mathcal{O}(N \max(w_i))$

b20ccc, 16 lines

```
int knapsack(vi w, int t) {
    int a = 0, b = 0, x;
    while (b < sz(w) && a + w[b] <= t) a += w[b++];
    if (b == sz(w)) return a;
    int m = *max_element(all(w));
    vi u, v(2*m, -1);
    v[a+m-t] = b;
    rep(i,b,sz(w)) {
        u = v;
        rep(x,0,m) v[x+w[i]] = max(v[x+w[i]], u[x]);
        for (x = 2*m; --x > m;) rep(j, max(0,u[x]), v[x])
            v[x-w[j]] = max(v[x-w[j]], j);
    }
    for (a = t; v[a+m-t] < 0; a--);
    return a;
}
```

9.3 Dynamic programming

KnuthDP.h

**Description:** When doing DP on intervals:  $a[i][j] = \min_{i < k < j} (a[i][k] + a[k][j]) + f(i, j)$ , where the (minimal) optimal  $k$  increases with both  $i$  and  $j$ , one can solve intervals in increasing order of length, and search  $k = p[i][j]$  for  $a[i][j]$  only between  $p[i][j - 1]$  and  $p[i + 1][j]$ . This is known as Knuth DP. Sufficient criteria for this are if  $f(b, c) \leq f(a, d)$  and  $f(a, c) + f(b, d) \leq f(a, d) + f(b, c)$  for all  $a \leq b \leq c \leq d$ . Consider also: LineContainer (ch. Data structures), monotone queues, ternary search.

**Time:**  $\mathcal{O}(N^2)$

DivideAndConquerDP.h

**Description:** Given  $a[i] = \min_{lo(i) \leq k < hi(i)} (f(i, k))$  where the (minimal) optimal  $k$  increases with  $i$ , computes  $a[i]$  for  $i = L..R - 1$ .

**Time:**  $\mathcal{O}((N + (hi - lo)) \log N)$

d38d2b, 18 lines

```
struct DP { // Modify at will:
    int lo(int ind) { return 0; }
    int hi(int ind) { return ind; }
    ll f(int ind, int k) { return dp[ind][k]; }
    void store(int ind, int k, ll v) { res[ind] = pii(k, v); }

    void rec(int L, int R, int LO, int HI) {
        if (L >= R) return;
        int mid = (L + R) >> 1;
        pair<ll, int> best(LLONG_MAX, LO);
        rep(k, max(LO, lo(mid)), min(HI, hi(mid)))
            best = min(best, make_pair(f(mid, k), k));
        store(mid, best.second, best.first);
        rec(L, mid, LO, best.second+1);
        rec(mid+1, R, best.second, HI);
    }
    void solve(int L, int R) { rec(L, R, INT_MIN, INT_MAX); }
};
```

Convolution.h

**Description:** Getting different convolutiona

**Time:**  $\mathcal{O}(n * 2^n)$

c8b662, 41 lines

```
// Zeta/SOS, N*2^N
rep(i,0,M)
    for(int mask = (1<<M) - 1; mask >= 0; mask--)
        if((mask>>i)&1)
```

```
F[mask] += F[mask ^ (1 << i)];
// Rev mask loop and invert bit condition for superset sum

// Base from SOS
for(int i = M - 1; i >= 0; i--)
    for(int mask = (1 << M) - 1; mask >= 0; mask--)
        if((mask >> i)&1)
            F[mask] -= F[mask ^ (1 << i)];
// Rev mask loop and invert condition for base from Sum over superset
```

```
// Mobius, F[s] = SUM(-1^{s/s'}) * F[s']), N*2^N
// F[1011] = F[1011] - F[0011] - F[1001] - F[1010] + F[1000]
...
rep(i,0,M) rep(mask, 0, 1<<M) if((mask>>i)&1)
    F[mask] -= F[mask ^ (1 << i)];
```

```
// sos(mu(f(x))) = f(x) = mu(sos(f(x)))

// fog[s] = SUM(F[s']*g[s/s']), N^2 * 2^N
// Make fhat[][] = {0} and ghat[][] = {0}
rep(mask,0,1<<N) {
    fhat[__builtin_popcount(mask)][mask] = f[mask];
    ghat[__builtin_popcount(mask)][mask] = g[mask];
}
// Apply zeta transform on fhat[][] and ghat[][]
rep(i,0,N+1) rep(j,0,N) rep(mask,0,1<<N) if((mask>>j)&1) {
    fhat[i][mask] += fhat[i][mask ^ (1 << j)];
    ghat[i][mask] += ghat[i][mask ^ (1 << j)];
}
// Do the convolution and store into h[][] = {0}
rep(mask,0,(1<<N)) rep(i,0,N+1) rep(j,0,i+1)
    h[i][mask] += fhat[j][mask] * ghat[i - j][mask];
// Apply inverse SOS dp on h[][]
rep(i,0,N+1) rep(j,0,N) rep(mask,0,1<<N) if((mask>>j)&1)
    h[i][mask] -= h[i][mask ^ (1 << j)];
```

```
rep(mask,0,1<<N) fog[mask] = h[__builtin_popcount(mask)][mask];
```

9.4 XOR Basis

XorBasis.h

**Description:** Representing each number in 2D vector space, finding basis of that vector space. total elements =  $2^{sz}$ . ways to represent  $x$  is  $2^{n-sz}$ , unique basis combination for every subset.

**Time:**  $\mathcal{O}(N * \log(A[i]))$

8bb3b7, 10 lines

```
int basis[d], sz; // basis[i] keeps the mask of the vector
whose f value is i
void insertVector(int mask) {
    for (int i = 0; i < d; i++) {
        if ((mask & 1 << i) == 0) continue; // continue if i != f(
            mask)
        if (!basis[i]) { // If there is no basis vector with the i'
            th bit set, then insert this vector into the basis
            basis[i] = mask; ++sz; return;
        }
        mask ^= basis[i]; // Otherwise subtract the basis vector
            from this vector
    }
}
```

9.5 Bit hacks

- $x \& \neg x$  is the least bit in  $x$ .
- for (int x = m; x; x=(x-1)&m; {...} loops over all subset masks of m (except m itself).

- $c = x\&\neg x, r = x+c; (((r^x) >> 2)/c) \mid r$  is the next number after  $x$  with the same number of bits set.

Bitset

- `bitset<100> b(5)` or `bitset<100> b; b = 23;`.
- `b.any()`, `b.all()`, check if any or all bits set to true
- `b[i]`, `b.count()`
- `for(int i = BS.Find_first(); i < BS.size(); i = BS.Find_next(i))`

9.6 Python

main.py

**Description:** Python Basics

7 lines

```
from math import comb
from collections import defaultdict
from itertools import accumulate
from functools import lru_cache
@lru_cache(maxsize=None)
def f(n):
    pass
```