

International Institute Of Information Technology - Hyderabad

FLogic

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int time() {return sz(ss);}

Contest (1)

```
.bashrc
alias c='q++ -Wall -Wconversion -Wfatal-errors -g -std=c++14 \
 -fsanitize=undefined,address
xmodmap -e 'clear lock' -e 'keycode 66=less greater' #caps = <>
template.cpp
                                                          17 lines
#include <bits/stdc++.h>
using namespace std;
#define rep(i, a, b) for(int i = a; i < (b); ++i)
#define all(x) begin(x), end(x)
#define sz(x) (int)(x).size()
#define pb push back
typedef long long ll;
typedef pair<int, int> pii;
typedef vector<int> vi;
int main() {
  // freopen("sample.in", "r", stdin);
  // freopen("sample.out", "w", stdout);
  cin.tie(0)->sync with stdio(0);
  cin.exceptions(cin.failbit);
```

Data structures (2)

```
OrderedSet.h
```

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

```
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;
```

```
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}\left(\alpha(N)\right)
                                                        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] \sim 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 \overline{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 {
    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 rmg(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
Time: Both operations are \mathcal{O}(\log N).
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;</pre>
  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;</pre>
    int pos = 0:
    for (int pw = 1 << 25; pw; pw >>= 1) {
      if (pos + pw \le 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 ele-
ments 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);
    for (vi& v : vs) 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 v. ll dif) {
    for (; x < sz(ys); x |= x + 1)
      ft[x].update(ind(x, y), dif);
  ll query(int x, int y) {
    11 \text{ 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 SeamentTrees
  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 guerv(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), 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. min1. lazv. 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 \rightarrow l = t \rightarrow r = NULL:
    t->prior = rand(); t->size = 1; t->rev = false;
    t \rightarrow lazv = 0: t \rightarrow min1 = t \rightarrow 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->min1 = min(l->min1, r->min1);
 void operation(pnode t) {
    if(!t) return;
    // reset t:
    t->size = 1; t->min1 = 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->lazv:
    t->min1 += 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++) {</pre>
   Tc = 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-l);
  if(mid) {
    mid->lazy += d;
```

SQRT LazyDynamicSegTree LineContainer

```
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-l);
    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
    int mid = l + (len - cnt) - 1; pnode L, Range, R;
    split(root, L, Range, l-1); split(Range, Range, R, r - l);
    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-l);
    push(mid); T ans = mid->min1;
    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++) {</pre>
        S[b][i - st[b] + 1] = nxt[i];
    sort(S[b] + 1, S[b] + len[b] + 1);
// set A \times = 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 guery block(int s,int e,int k) {
    int ans = 0;
    for(int i = s; i <= e; i++)</pre>
        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 l\bar{b} = l / B, rb = r / B:
    if(lb == rb) return query elements(l, r, r);
    return query elements(l, en[lb], r)
        + query \overline{b}lock(lb + 1, rb - 1, r)
        + query elements(st[rb], r, r);
for(int i = 1: i \le 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++) {</pre>
    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 \le 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 up-
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</pre>
    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);</pre>
    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]}
    push(tree[i]); update(l,r,u,tree[i].lc); update(l,r,u,tree[
    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 ("con-
vex hull trick").
Time: \mathcal{O}(\log N)
                                                       8ec1c7, 30 lines
struct Line {
  mutable ll k, m, p;
  bool operator<(const Line& 0) const { return k < 0.k; }</pre>
  bool operator<(ll x) const { return p < x; }</pre>
struct LineContainer : multiset<Line. less<>> 
  // (for doubles, use \inf = 1/.0, \operatorname{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);
    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 dist = inf; nodes reachable through negative-weight cycles get dist = -inf. Assumes $V^2 \max |w_i| < \sim 2^{63}$ undirected graph with negative edge gets -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
      cvcle
  dist[source] = 0;
  for (int iter = 0; iter < adj.size() - 1; ++iter){</pre>
   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])</pre>
          dist[e.first] = dist[u]+e.second, relax = true;
   if(!relax) break:
  bool ch = true:
  while (ch) {
                      // keep going untill no more changes
                      // set dist to -INF when in cycle
    ch = false;
    for (int u = 0; u < adj.size(); ++u)</pre>
     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;
```

FlovdWarshall.h

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

```
61c353, 12 lines
const ll inf = 1LL << 62;</pre>
void floydWarshALL(vector<vector<ll>>& m) {
  int n = sz(m):
  rep(i,0,n) m[i][i] = min(m[i][i], OLL);
  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) $O(V^2E)$: General (2) O(Flow E): General (3) $O(E\sqrt{V})$: when sum of edge capacities is O(n), we can treat edge with weight x as x edges with weight 1. (4) $O(EV \log(Flow))$: Dinics with scaling

```
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 flow() { return max(oc - c, (T)0); } // if you need flows
 int N;
 vector<int> lvl, ptr, q; vector<vector<Edge>> adj;
 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++) {</pre>
      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]) {</pre>
       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++) {</pre>
     for(auto e : adj[i]) {
       int i = e.to:
       Flow[i][j] = e.flow();
};
```

MinCut.h

Description: After running max-flow, the left side of a min-cut from s to tis 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(\bar{i}t,0,n-ph) \{ // O(V^2) \rightarrow O(E \log V) \text{ 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;
```

MinCostMaxFlow.h

Description: Min-cost max-flow. cap[i][j] != 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 $\mathcal{O}(E^2)$

dist[s] = 0; ii di;

fe85cc, 81 lines

```
#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);
```

```
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) {
   II totflow = 0, totcost = 0;
    while (path(s), seen[t]) {
     II 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]);
      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])</pre>
            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(q, btoa);
Time: \mathcal{O}\left(\sqrt{V}E\right)
                                                       fd54bf, 43 lines
bool dfs(int a, int L, vector<vi>& q, 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) {

if (btoa[b] == -1 || dfs(btoa[b], L + 1, g, btoa, A, B))

```
return btoa[b] = a, 1;
 return 0;
int hopcroftKarp(vector<vi>& q, 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(q)) 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(q))
      res += dfs(a, 0, q, 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>& q, 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 : q[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 { // diikstra
      done[i0] = true;
      int i0 = p[j0], j1, delta = INT MAX;
      rep(j,1,m) if (!done[j]) {
        auto cur = a[i0 - 1][i - 1] - u[i0] - v[i];
        if (cur < dist[j]) dist[j] = cur, pre[j] = j0;
        if (dist[j] < delta) delta = dist[j], j1 = j;</pre>
      rep(j,0,m) {
        if (done[j]) u[p[j]] += delta, v[j] -= delta;
        else dist[i] -= delta;
      i0 = i1;
    } 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
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|||b)\&\&(!a|||c)\&\&(d|||!b)\&\&... becomes true, or reports that it is unsatis-
fiable. Negated variables are represented by bit-inversions (\sim X).
Usage: TwoSat ts(number of boolean variables):
ts.either(0, \sim3); // Var 0 is true or var 3 is false
ts.setValue(2); // Var 2 is true
```

```
Usage: TwoSat ts(number of boolean variables); ts.either(0, \sim3); // Var 0 is true or var 3 is false ts.setValue(2); // Var 2 is true ts.atMostOne(\{0, \sim 1, 2\}); // <= 1 of vars 0, \sim1 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.
```

```
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) {
  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;</pre>
    int cur = \simli[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;
};
```

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

```
780b64, 15 lines
vi eulerWalk(vector<vector<pii>>>& gr, int nedges, int src=0) {
  int n = sz(ar):
  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) = qr[x][it++];
   if (!eu[e]) {
     D[x]--, D[y]++;
     eu[e] = 1; s.push back(y);
  for (int x : D) if (x < 0 \mid \mid 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

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). ncomps will contain the number of components.

Time: $\mathcal{O}\left(E+V\right)$ f5e4c5, 53 lines

```
// 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]);
```

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.

```
// 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 : adi[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.ioin(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> adi(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];
     adi[u][e] = left;
     adj[left][e] = u;
     adj[right][e] = -1;
      free[right] = e;
    adj[u][d] = fan[i];
```

```
adi[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;
```

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 = \simB(), B X={}, B R={}) {
  if (!P.any()) { if (!X.any()) f(R); return; }
  auto q = (P | X). Find first();
  auto cands = P \& \neg eds[\overline{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.

```
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;</pre>
     q.push back(R.back().i);
     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[i++].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[i].i = i, T[i++].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.

3.7Trees

BinaryLifting.h

Description: Initialize par[N][logN] and depth[N] array. kthpar: logN, 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 (time[node], node) in a vector every time you iterate over an edge (node, x). lca(a, b) => rangeMin(time[a], 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 $O(\log N)$.

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 \land 1] = b ? x : this;
    y - > c[i \land 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();
```

DirectedMST HLD CentroidDecomposition

```
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}\left(E\log V\right)
"../data-structures/UnionFindRollback.h"
                                                       39e620, 60 lines
struct Edge { int a, b; ll w; };
struct Node {
 Edge key;
  Node *1, *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>& q) {
 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>>> cycs;
  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;
        cycs.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] : cycs) { // 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: O(logN * 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: adi = adi1: 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):
        sqt.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]) {
            sqt.update(st[head[x]], st[x], add, 0, t-1);
            x = par[head[x]];
        sqt.update(st[p], st[x], add, 0, t - 1);
    void range update(int u,int v,T add) {
        int \bar{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, sqt.query(st[p], st[x], 0, t - 1));
        return ans;
    T range min(int u,int v) {
        int = 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, log N = 17;
vi adj[N], sub(N), par(N,-\bar{1}), lvl(N), done(N), par adj(N);
vector<vi> dist(N, vi(logN, 0)), anc(N, vi(logN, \overline{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;
```

rep(x,0,n) {

return l:

vector<vi> child cntb(N), my(N);

if(par[y] >= 0)

// number of nodes <= k in v.

auto cnt k = [&](vi &v, int k)

my[y].pb(dist[x][lvl[y]]);

rep(x,0,n) for(int y = x; y >= 0; y = par[y]) {

child cntb[y].pb(dist[x][lvl[par[y]]]);

sort(all(my[x])); sort(all(child cntb[x]));

int l = upper bound(all(v), k) - v.begin();

```
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][[vl[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(\overline{n}), max depth(n), dp(n), max up(n) {}
  void add edge(int s, int t) {
   adj[s].pb(t); adj[t].pb(s);
  int argmin(int i, int j) { return depth[i] < depth[j] ? i : j</pre>
  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); // =
    return i == j ? u : tour[argmin(table[h][i], table[h][i]
         -(1<<h)1)1;
 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++){</pre>
     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
    for(int i = 1; i < (int)start times.size(); i++){</pre>
      while(pos end[tour[start times[st.top()]]] < start times[</pre>
          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);
    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;
}</pre>
```

3.8 Math

Number of Spanning Trees Create an $N \times N$ matrix mat, and for each edge $a \to b \in G$, do mat[a][b]--, mat[b][b]++ (and mat[b][a]--, mat[a][a]++ if G is undirected). Remove the ith 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.)

ee09e2, 12 lines

Hashing Kmp Manacher MinRotation SuffixArray Z

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 \mathbb{Z}_2 which can be done easily as $\operatorname{Permanent}(A) = \operatorname{Determinent}(A)$.

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>

```
int x; B b; A(int x=0) : x(x), b(x) {}
 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; }</pre>
typedef A<1000000007, A<1000000009, 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: 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

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

Description: For each position in a string, computes p[0][i] = half length of longest even palindrome around pos i, <math>p[1][i] = longest odd (half rounded down).

```
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++;</pre>
```

SuffixArrav.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: $O(n \log n)$

```
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) = 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)</pre>
      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;
```

DvnamicAhoCorasik.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); }
    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++) {</pre>
 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 LIM < mod and that mod is a prime. 6f684f, 3 lines

```
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

b83e45, 8 lines

```
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: $\mathcal{O}(\sqrt{m})$ c040b8, 11 lines

```
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 \le 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 (\overline{A}.count(e = e * f % m))
      return n * i - A[e];
  return -1;
```

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

Time: $\log(m)$, with a large constant.

```
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);
Il 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 \mod c (or a^b \mod c) for 0 \le a, b \le c \le 7.2 \cdot 10^{18}.
Time: \mathcal{O}(1) for modmul, \mathcal{O}(\log b) for modpow
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 \neq 2)
    if (e & 1) ans = modmul(ans, b, mod);
```

ModSart.h

Description: Tonelli-Shanks algorithm for modular square roots. Finds xs.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;
  II 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:
    II qs = modpow(q, 1LL \ll (r - m - 1), p);
    q = qs * qs % p;
    x = x * qs % p;
    b = b * q % p;
```

```
Linear Diophantine.h
Description: Solving linear diophantine eqns (ax + by = c). 87b1ec, 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,
    g = gcd(abs(a), abs(b), x0, y0);
    if (c % g)
        return false;
    x0 *= c / a:
    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 /= q;
    b /= q;
    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)</pre>
        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)</pre>
        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]) {</pre>
    cp.push back(\{i, i * i / 2\});
    for (int j = i * i; j \le 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;</pre>
    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 ex-
tend A randomly.
Time: 7 times the complexity of a^b \mod 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}\left(n^{1/4}\right), less for numbers with small factors.
"ModMulLL.h", "MillerRabin.h"
ull pollard(ull n) {
  auto f = [n](uli 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}.
ll euclid(ll a, ll b, ll &x, ll &y) {
 if (!b) return x = 1, y = 0, a;
  ll 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 \le x < \operatorname{lcm}(m, n). Assumes mn < 2^{62}
Time: \log(n)
"euclid.h"
ll crt(ll a, ll m, ll b, ll n) {
 if (n > m) swap(a, b), swap(m, n);
 ll 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/q : x;
```

Bézout's identity

For $a \neq b \neq 0$, then d = qcd(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 ϕ function is defined as $\phi(n) := \#$ of positive integers $\leq n$ that are coprime with n. $\phi(1) = 1$, p prime $\Rightarrow \phi(p^k) = (p-1)p^{k-1}$, $m, n \text{ coprime } \Rightarrow \phi(mn) = \phi(m)\phi(n).$ If $n = p_1^{k_1} p_2^{k_2} \dots p_r^{k_r}$ then $\phi(n) = p_1^{k_1} p_2^{k_2} \dots p_r^{k_r}$ $(p_1-1)p_1^{k_1-1}...(p_r-1)p_r^{k_r-1}.$ $\phi(n)=n\cdot\prod_{p\mid n}(1-1/p).$ $\sum_{d|n} \phi(d) = n, \sum_{1 \le k \le n, \gcd(k,n)=1} k = n\phi(n)/2, n > 1$ **Euler's thm**: a, n coprime $\Rightarrow a^{\phi(n)} \equiv 1 \pmod{n}$. **Fermat's little thm**: $p \text{ prime } \Rightarrow a^{p-1} \equiv 1 \pmod{p} \ \forall a.$ 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)
```

ContinuedFractions FracBinarySearch IntPerm

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

5.4 Fractions

ContinuedFractions.h

Description: Given N and a real number $x \ge 0$, finds the closest rational approximation p/q with $p, q \le N$. It will obey $|p/q - x| \le 1/qN$.

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 ∞ ; 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 \sim 1e7; long double for N \sim 1e9
pair<ll, ll> approximate(d x, ll N) {
 II LP = 0, LQ = 1, P = 1, Q = 0, inf = LLONG MAX; dy = x;
  for (;;) {
   II \lim = \min(P ? (N-LP) / P : \inf. 0 ? (N-L0) / 0 : \inf).
      a = (ll)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: $\mathcal{O}\left(\log(N)\right)$ 27ab3e, 25 lines

```
struct Frac { ll p, q; };
template<class F>
Frac fracBS(F f, ll 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) {
   ll 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

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

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 $\mathbb{Z}_{>a}^{\times}$ 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:

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

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

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

Combinatorics (6)

6.1 Permutations

Factorial

	n	1 2	3 4	5	6	7	8	3	9	10	
	n!	1 2	6 24	120	0 720	504	0 403	320 36	32880	3628800	
	n	11		12	13	1	4	15	16	3628800 17 3 3.6e14 171	
_	n!	4.0e	7 4.	8e8	6.2e	9 8.7	e10 1	.3e12	2.1e1	3.6e14	
	n	20	2	5	30	40	50	100	150	171	
_	n!	2e18	3 2e	25 3	e32	8e47	3e64	9e15'	7 6e26	2 >dbl_m	AX

IntPerm.h

Description: Permutation -> integer conversion. (Not order preserving.) Integer -> permutation can use a lookup table.

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.

$$p(0) = 1, \ p(n) = \sum_{k \in \mathbb{Z} \setminus \{0\}} (-1)^{k+1} p(n - k(3k - 1)/2)$$
$$p(n) \sim 0.145/n \cdot \exp(2.56\sqrt{n})$$
$$\frac{n \quad | \ 0 \ 1 \ 2 \ 3 \ 4 \ 5 \ 6 \ 7 \ 8 \ 9 \ 20 \quad 50 \quad 100}{p(n) \quad | \ 1 \ 1 \ 2 \ 3 \ 5 \ 7 \ 11 \ 15 \ 22 \ 30 \ 627 \sim 2e5 \sim 2e8}$$

Lucas' Theorem

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

multinomial Polynomial PolyRoots PolyInterpolate

Binomials

multinomial.h

Description: Computes
$$\binom{k_1+\cdots+k_n}{k_1,k_2,\ldots,k_n} = \frac{(\sum k_i)!}{k_1!k_2!\ldots k_n!}$$
.

ll multinomial(vi& v) { ll 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:

General purpose numbers

Bernoulli numbers

EGF of Bernoulli numbers is $B(t) = \frac{t}{e^{t-1}}$ (FFT-able). $B[0,\ldots] = [1,-\frac{1}{2},\frac{1}{6},0,-\frac{1}{20},0,\frac{1}{42},\ldots]$

Sums of powers:

$$\sum_{i=1}^{n} n^{m} = \frac{1}{m+1} \sum_{k=0}^{m} {m+1 \choose k} B_{k} \cdot (n+1)^{m+1-k}$$

Euler-Maclaurin formula for infinite sums:

$$\sum_{i=m}^{\infty} f(i) = \int_{m}^{\infty} f(x)dx - \sum_{k=1}^{\infty} \frac{B_{k}}{k!} f^{(k-1)}(m)$$

$$\sum_{i=m}^{\infty} f(m) = f'(m) = f'''(m)$$

$$\approx \int_{m}^{\infty} f(x)dx + \frac{f(m)}{2} - \frac{f'(m)}{12} + \frac{f'''(m)}{720} + O(f^{(5)}(m))$$

Stirling numbers of the first kind

Number of permutations on n items with k cycles.

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

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. $\pi(j) \ge j$, k j:s s.t. $\pi(j) > j$.

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

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

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

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 \cdots n_k n^{k-2}
# with degrees d_i: (n-2)!/((d_1-1)!\cdots(d_n-1)!)
```

Catalan numbers

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

$$C_0 = 1, \ C_{n+1} = \frac{2(2n+1)}{n+2} C_n, \ C_{n+1} = \sum_{n=1}^{\infty} C_i C_{n-n}$$

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

- 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.
- \bullet 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**

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

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

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

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

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

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

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

Generating Functions Blog 1 and 2.

Polynomials and recurrences

```
Polynomial.h
                                                        c9b7b0, 17 lines
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: 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:
  Polv 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 ^{\wedge} (p(h) > 0))
      rep(it, 0, 60)  { // while (h - l > 1e-8)
        double m = (l + h) / 2, f = p(m);
        if ((f \le 0) \land 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 + ... + a[n-1] * x^{n-1}. For
numerical precision, pick x[k] = c * \cos(k/(n-1) * \pi), k = 0 \dots n-1.
Time: 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;

BerlekampMassev.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"
                                                      96548b, 20 lines
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;
   II 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 $S[0\ldots \ge n-1]$ and $tr[0\ldots n-1]$. Faster than matrix multiplication. Useful together with Berlekamp–Massey.

Usage: linearRec({0, 1}, {1, 1}, k) // k'th Fibonacci number Time: $\mathcal{O}\left(n^2\log k\right)$

```
typedef vector<ll> Poly;
ll linearRec(Poly S, Poly tr, ll k) {
 int n = sz(tr);
  auto combine = [&](Poly a, Poly b) {
   Polv 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.

```
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_{8eeeaf, 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 = le9; jmp > le-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.

4756fc. 7 lines

```
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

```
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^Tx 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^Tx 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*\#pivots), where a pivot may be e.g. an edge relaxation.

\mathcal{O}(2^n) in the general case.

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 \mid \mid 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;</pre>
```

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;</pre>
      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: $\mathcal{O}(N^3)$

```
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) \operatorname{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.

```
Time: \mathcal{O}(N^3)
                                                      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: $\mathcal{O}\left(n^2m\right)$

```
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[i] -= 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)</pre>
SolveLinear2.h
```

Description: To get all uniquely determined values of x back from Solve-Linear, make the following changes:

```
"SolveLinear.h"
rep(j,0,n) if (j != i) // instead of <math>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));</pre>
 vi col(m); iota(all(col), 0);
 rep(i,0,n) {
    for (br=i; br<n; ++br) if (A[br].any()) break;</pre>
```

```
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)</pre>
```

MatrixInverse.h

return n;

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 mod p, and k is doubled in each step.

```
Time: \mathcal{O}\left(n^3\right)
                                                        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 = i, c = k;
    if (fabs(A[r][c]) < 1e-12) return i;</pre>
    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[i][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];
```

Tridiagonal.h

Description: x = tridiagonal(d, p, q, b) solves the equation system

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

This is useful for solving problems on the type

```
a_i = b_i a_{i-1} + c_i a_{i+1} + d_i, \ 1 \le i \le 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, ..., -1, 1\}, \{0, c_1, c_2, ..., c_n\},\
                           \{b_1, b_2, \ldots, b_n, 0\}, \{a_0, d_1, d_2, \ldots, 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;
     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;
```

Fourier transforms

FastFourierTransform.h

Description: 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: conv(a, b) = c, where $c[x] = \sum a[i]b[x-i]$. 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:** $O(N \log N)$ with N = |A| + |B| (~1s for $N = 2^{22}$)

```
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) {
      Cz = 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 \ll L;
 vector<C> in(\overline{n}), out(\overline{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, mod).
Time: \mathcal{O}(N \log N), where N = |A| + |B| (twice as slow as NTT or FFT)
"FastFourierTransform.h"
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));</pre>
 vector<C > L(n), R(\overline{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)) {
    II av = II(real(outl[i])+.5), cv = II(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 $g = \sum_{x} a[x]g^{xk}$ $root^{(mod-1)/N}$. 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 $2^{\hat{a}}$. For arbitrary modulo, see FFTMod. 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"
```

00ced6, 35 lines

const ll mod = (119 << 23) + 1, root = 62; // = 998244353 // For p < 2^3 0 there is also e.g. 5 << 25, 7 << 26, 479 << 21

```
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]]);</pre>
  for (int k = 1; k < n; k *= 2)
    for (int i = 0; i < n; i += 2 * k) rep(j,0,k) {
      Il 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 %
  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 u} a[x] \cdot b[y], \text{ where } \oplus \text{ is one of AND, OR, XOR.} The size
of a must be a power of two.
Time: \mathcal{O}(N \log N)
void FST(vi& a, bool inv) {
 for (int n = sz(a), step = 1; step < n; step *= 2) {</pre>
    for (int i = 0: i < n: i += 2 * step) rep(i,i,i+step) {
      int &u = a[i], &v = a[i + 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);
  if (inv) for (int& x : a) x \neq sz(a); // XOR only
```

// and 483 << 21 (same root). The last two are > 10^9.

```
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;
```

```
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
```

WalshHadamard.h

```
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(i, 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 \neq sz(P);
Description: Given B_1, \ldots 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}\left(N\log^2 N\right)
                                                             c0e86b, 18 lines
void online(const Polv &B, CD a1, int n, Polv &A) {
  const int m = SZ(B) - 1;
  A.assign(n + 1, 0); A[1] = a1;
  auto bst = B.begin(), ast = A.begin();
  REP(i, 1, n) {
    A[i + 1] += A[i] * B[1];
    if (i + 2 \le 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 \le 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
$$r = \sqrt{a^2 + b^2}$$
, $\phi = \operatorname{atan2}(b, a)$.

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): $m_a = \frac{1}{2}\sqrt{2b^2 + 2c^2 - a^2}$

OnlineFFT Point

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 sines: $\frac{\sin\alpha}{a} = \frac{\sin\beta}{b} = \frac{\sin\gamma}{c} = \frac{1}{2R}$ Law of cosines: $a^2 = b^2 + c^2 - 2bc\cos\alpha$ Quadrilaterals $\tan\frac{\alpha + \beta}{b}$ With fiden lengths a,b,c,d, diagonals e,f, diagonals angle θ , area A and magic flux $F = bb^2 + d^2n \frac{\alpha}{2} = \frac{\alpha^2\beta}{2} = c^2$:

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

Spherical coordinates

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



$$x = r \sin \theta \cos \phi \qquad r = \sqrt{x^2 + y^2 + z^2}$$

$$y = r \sin \theta \sin \phi \qquad \theta = a\cos(z/\sqrt{x^2 + y^2 + z^2})$$

$$z = r \cos \theta \qquad \phi = a\tan(y, x)$$

Derivatives/Integrals

$$\frac{d}{dx}\arcsin x = \frac{1}{\sqrt{1-x^2}} \qquad \frac{d}{dx}\arccos x = -\frac{1}{\sqrt{1-x^2}}$$

$$\frac{d}{dx}\tan x = 1 + \tan^2 x \qquad \frac{d}{dx}\arctan x = \frac{1}{1+x^2}$$

$$\int \tan ax = -\frac{\ln|\cos ax|}{a} \qquad \int x\sin ax = \frac{\sin ax - ax\cos ax}{a^2}$$

$$\int e^{-x^2} = \frac{\sqrt{\pi}}{2}\operatorname{erf}(x) \qquad \int xe^{ax}dx = \frac{e^{ax}}{a^2}(ax-1)$$

Integration by parts:

$$\int_{a}^{b} f(x)g(x)dx = [F(x)g(x)]_{a}^{b} - \int_{a}^{b} F(x)g'(x)dx$$

Geometric primitives

Point.h

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

```
<br/><br/>bits/stdc++.h>
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) {</pre>
```

return os << "(" << p.x << "," << p.y << ")";

84d6d3, 11 lines

```
// 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 /S on the result of the cross product. "Point h"

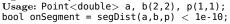


```
f6bf6b, 4 lines
template<class P>
double lineDist(const P& a, const P& b, const P& p) {
 return (double)(b-a).cross(p-a)/(b-a).dist();
```

SegmentDistance.h

Description:

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



"Point.h"

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<|l> 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 = seqInter(s1,e1,s2,e2);
if (sz(inter)==1)
```

cout << "segments intersect at " << inter[0] << endl;</pre> "Point.h", "OnSegment.h" 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, e^2\}
(0,0)} is returned. The wrong position will be returned if P
is Point<ll> and the intersection point does not have inte-
ger coordinates. Products of three coordinates are used in 1
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;</pre>
"Point.h"
template<class P>
pair<int, P> lineInter(P s1, P e1, P s2, P e2) {
```

return $\{-(s1.cross(e1, s2) == 0), P(0, 0)\};$

auto p = s2.cross(e1, e2), q = s2.cross(e2, s1);

sideOf.h

Description: Returns where p is as seen from s towards e. $1/0/-1 \Leftrightarrow \text{left/on}$ 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"
```

return (a > l) - (a < -l);

auto d = (e1 - s1).cross(e2 - s2);

return $\{1, (s1 * p + e1 * q) / d\};$

if (d == 0) // if parallel

```
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;
```

OnSegment.h

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

```
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.

```
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

```
a number of rotations around the origin). Useful for rotational sweeping.
Sometimes also represents points or vectors.
Usage: vector<Ângle> 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
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 \mid | (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) {</pre>
  // add a.dist2() and b.dist2() to also compare distances
  return make tuple(a.t, a.half(), a.y * (ll)b.x) <</pre>
         make_tuple(b.t, b.half(), a.x * (ll)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)\};
```

Description: A class for ordering angles (as represented by int points and

8.3 Circles

"Point.h"

3af81c, 9 lines

CircleIntersection.h

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

```
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 \text{ 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 \text{ mid} = a + \text{vec*p}, \text{ per} = \text{vec.perp()} * \text{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 {};</pre>
 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: $\mathcal{O}(n)$

"../../content/geometry/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 = [\&](Pp, Pq) {
    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;</pre>
   auto s = max(0., -a-sqrt(det)), t = min(1., -a+sqrt(det));
   if (t < 0 | | 1 \le s) return arg(p, q) * r2;
   Pu = p + d * s, v = p + d * t;
    return arg(p,u) * r2 + u.cross(v)/2 + arg(v,g) * 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 circumcirle 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"



```
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 \circ = 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(v, P{3, 3}, false);
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;</pre>
    cnt ^= ((a.y<p[i].y) - (a.y<q.y)) * a.cross(p[i], q) > 0;
 return cnt:
```

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"

```
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
tvpedef 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;</pre>
    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 counterclockwise order. Points on the edge of the hull between two other points are not considered part of the hull.



```
"Point.h"
typedef Point<ll> P;
```

```
vector<P> convexHull(vector<P> pts) {
  if (sz(pts) <= 1) return pts;</pre>
  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.i)
    for (;; j = (j + 1) % n) {
      res = \max(\text{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)</pre>
    return false;
  while (abs(a - b) > 1) {
   int c = (a + b) / 2;
    (sideOf(l[0], l[c], p) > 0 ? b : a) = c;
  return sqn(l[a].cross(l[b], p)) < r;</pre>
```

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)
                                                     7cf45b, 39 lines
"Point.h"
#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) sqn(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 \mid | 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; });</pre>
  pair<ll, pair<P, P>> ret{LLONG MAX, {P(), P()}};
  int j = 0;
  for (Pp: 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; }</pre>
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 \times 0 = INF, \times 1 = -INF, y = INF, y = -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 {
  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)</pre>
      best = min(best, search(s, p));
    return best:
 // find nearest point to a point, and its squared distance
 // (requires an arbitrary operator< for Point)</pre>
 pair<T, P> nearest(const P& p) {
    return search(root, p);
};
```

FastDelaunav.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], $t[0][1], t[0][2], t[1][0], \dots$, all counter-clockwise.

```
Time: \mathcal{O}(n \log n)
"Point.h"
typedef Point<ll> P;
typedef struct Quad* Q;
typedef int128 t lll; // (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: }
  0& r() { return rot->rot; }
  0 prev() { return rot->o->rot: }
 0 next() { return r()->prev(); }
bool circ(P p, P a, P b, P c) { // is p in the circumcircle?
 lll 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;
0 makeEdge(P orig. P dest) {
  0 r = H ? H : new Quad{new Quad{new Quad{new Quad{0}}}};
 H = r -> 0: r -> r() -> r() = r:
  rep(i,0,4) r = r->rot, r->p = arb, r->o = i & 1 ? r : r->r();
  r \rightarrow p = orig; r \rightarrow 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 = 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) \le 3) {
    Q a = makeEdge(s[0], s[1]), b = makeEdge(s[1], s.back());
    if (sz(s) == 2) return { a, a->r() };
```

Point3D.h

template<class T> struct Point3D {

typedef Point3D P;

typedef const P& R;

```
splice(a->r(), b);
    auto side = s[0].cross(s[1], s[2]);
   0 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 \rightarrow 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);
  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())) { \
     0 t = e->dir; \
      splice(e, e->prev()); \
      splice(e->r(), e->r()->prev()); \
      e -> 0 = H; H = e; e = t; \setminus
  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());
     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 {}:
  0 e = rec(pts).first;
  vector<0> q = {e};
  int ai = 0:
  while (e->o->F().cross(e->F(), e->p) < 0) e = e->o;
#define ADD { Q c = e; do { c \rightarrow mark = 1; pts.push back(c \rightarrow p); \
  q.push back(c->r()); c = c->next(); \} while (c \overline{!}= 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& trilist) {
  double v = 0:
  for (auto i : trilist) v += p[i.a].cross(p[i.b]).dot(p[i.c]);
  return v / 6:
```

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

```
explicit Point3D(T x=0, T y=0, T z=0) : x(x), y(y), z(z) {}
  bool operator<(R p) const {</pre>
    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: \mathcal{O}\left(n^2\right)
"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[i];
      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;
};</pre>
```

sphericalDistance.h

Description: Returns the shortest distance on the sphere with radius radius between the points with azimuthal angles (longitude) f1 (θ_1) and f2 (θ_2) from x axis and zenith angles (latitude) t1 (θ_1) and t2 (θ_2) from z axis (0 = 1) 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.

```
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)

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);
```

mint() : value() {}

```
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)
                                                         9e9d8d, 19 lines
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]; });</pre>
  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]));
    if (mx.second == -1) return {};
    cur = mx.first:
    R.push back(mx.second);
  return R;
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}\left(k\log\frac{n}{k}\right)
                                                         753a4c, 19 lines
template<class F. class G. class T>
void rec(int from, int to, F& f, G& q, int& i, T& p, T q) {
  if (p == q) return;
  if (from == to) {
    q(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 q) {
  if (to <= from) return;</pre>
  int i = from; auto p = f(i), q = f(to-1);
  rec(from, to-1, f, g, i, p, q);
  g(i, to, q);
Mint.h
Description: Modular Arithmetic (Integer) for Aditya and Ario 17/10/22, 27 lines
template <int32 t MOD>
struct mint {
  int32 t value;
```

mint(int64 t value) : value(value < 0 ? value % MOD + MOD

inline mint<MOD> operator + (mint<MOD> other) const { return

inline mint<MOD> operator - (mint<MOD> other) const { return

: value >= MOD ? value % MOD : value) {}

explicit operator bool() const { return value; }

mint<MOD>(*this) += other; }

mint<MOD>(*this) -= other; }

mint(int32 t value , std::nullptr t) : value(value) {}

```
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 > \overline{n}) { return mint < MOD > (value) - n: }
template <int32 t MOD> mint<MOD> operator * (int64 t value,
     mint < MOD > \overline{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</pre>
      & 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) < \ldots < f(i) \ge \cdots \ge 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))
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;
                                                                       void rec(int L, int R, int LO, int HI) {
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 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)) {
    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[i]] = max(v[x-w[i]], i);
  for (a = t; v[a+m-t] < 0; a--);
  return a;
KnuthDP.h
Description: When doing DP on intervals: a[i][j] = \min_{i < k < j} (a[i][k] + a[i][j])
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) < f(a,d) + f(b,c) for all a < b < c < 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 \leq hi(i)} (f(i, k)) where the (minimal)
optimal k increases with i, computes \bar{a}[i] for i = L..R - 1.
Time: \mathcal{O}((N + (hi - lo)) \log N)
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); }
```

if (L >= R) return;

int mid = $(L + R) \gg 1$;

pair<ll, int> best(LLONG_MAX, L0);

rec(L, mid, L0, best.second+1);

rec(mid+1, R, best.second, HI);

store(mid, best.second, best.first);

rep(k, max(L0,lo(mid)), min(HI,hi(mid)))

best = min(best, make pair(f(mid, k), k));

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 \land (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
// 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)))
                                                                       vmi res;
// 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(i,0,N) rep(mask,0,1<< N) if((mask>>i)&1) {
                                                                     9.1
  fhat[i][mask] += fhat[i][mask ^ (1 << j)];</pre>
  qhat[i][mask] += qhat[i][mask ^ (1 << j)];
// Do the convolution and store into h[][] = \{0\}
rep(mask, 0, (1 << N)) rep(i, 0, N+1) rep(i, 0, i+1)
            h[i][mask] += fhat[j][mask] * qhat[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];
                                                                     Bitset
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}\left(N * log(A[i])\right)
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(
    if (!basis[i]) { // If there is no basis vector with the i'
         th bit set, then insert this vector into the basis
                                                                     main.pv
      basis[i] = mask; ++sz; return;
    mask ^= basis[i]; // Otherwise subtract the basis vector
         from this vector
```

```
PolyModPoly.h
Description: Poly Mod Poly
                                                     c40f13, 33 lines
#define rsz resize
poly RSZ(poly p, int x) { p.rsz(x); return p; }
poly rev(poly p) { reverse(all(p)); return p; }
poly inv(poly A, int n) { // Q-(1/Q-A)/(-Q^{-2})
  poly B{1/A[0]};
  while (sz(B) < n) {
    int x = 2*sz(B);
    B = RSZ(2*B-conv(RSZ(A,x),conv(B,B)),x); } // fft
  return RSZ(B,n);
pair<poly, poly> divi(const poly& f, const poly& g) {
  if (sz(f) < sz(g)) return {{}, f};
  auto g = mul(inv(rev(g),sz(f)-sz(g)+1),rev(f));
  q = rev(RSZ(q,sz(f)-sz(q)+1));
  auto r = RSZ(f-mul(q,g),sz(g)-1); return \{q,r\};
typedef vector<mi> vmi; // mi = modular int
struct MultipointEval {
  poly stor[1<<18];
  void prep(vmi v, int ind = 1) { // v -> places to evaluate at
    if (sz(v) == 1) { stor[ind] = {-v[0],1}; return; }
    int m = sz(v)/2:
    prep(vmi(begin(v),begin(v)+m),2*ind);
    prep(vmi(m+all(v)), 2*ind+1);
    stor[ind] = conv(stor[2*ind],stor[2*ind+1]);
  void eval(vmi v, int ind = 1) {
    v = divi(v,stor[ind]).s;
   if (sz(stor[ind]) == 2) { res.pb(v[0]); return; }
    eval(v,2*ind); eval(v,2*ind+1);
       Bit hacks
   • x & -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\&-x, r = x+c; (((r^x) >> 2)/c) | r is the
     next number after X with the same number of bits set.
   • 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)
9.2 Python
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
```