

# University of California Irvine

# UCI Can't Implement

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$\mathbf{C}$	$\frac{\text{dontest}}{\text{contest}}$ (1)			
so	cal.sh	2 line		
	make problem directories a,, z r l in {az}; do mkdir \$1; cp temp.cpp \$1/\$1.cpp; done			
teı	mplate.cpp	12 line		
us: typ	nclude <bits stdc++.h=""> ing namespace std;  pedef int uci; efine int long long</bits>			

#define sz(x) ((int)x.size()) #define all(a) (a).begin(), (a).end() uci main() { cin.tie(0)->sync\_with\_stdio(0); cin.exceptions(cin.failbit);

#### bruh.pv

import sys

input = sys.stdin.readline

# Mathematics (2)

### 2.1 Equations

The extremum of  $ax^2 + bx + c = 0$  is given by x = -b/2a.

$$ax + by = e \Rightarrow x = \frac{ed - bf}{ad - bc}$$
$$cx + dy = f \Rightarrow y = \frac{af - ec}{ad - bc}$$

In general, given an equation Ax = b, the solution to a variable  $x_i$  is given by

$$x_i = \frac{\det A_i'}{\det A}$$

where  $A'_i$  is A with the i'th column replaced by b.

#### 2.2 Recurrences

If  $a_n = c_1 a_{n-1} + \cdots + c_k a_{n-k}$ , and  $r_1, \ldots, r_k$  are distinct roots of  $x^k - c_1 x^{k-1} - \cdots - c_k$ , there are  $d_1, \ldots, d_k$  s.t.

$$a_n = d_1 r_1^n + \dots + d_k r_k^n$$

Non-distinct roots r become polynomial factors, e.g.  $a_n = (d_1 n + d_2)r^n.$ 

### 2.3 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  $r = \sqrt{a^2 + b^2}$ ,  $\phi = \operatorname{atan2}(b, a)$ .

#### 2.4 Geometry

### 2.4.1 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:  $r = \frac{A}{n}$ 

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

Length of bisector (divides angles in two):

$$s_a = \sqrt{bc \left[ 1 - \left( \frac{a}{b+c} \right)^2 \right]}$$

Law of sines:  $\frac{\sin \alpha}{a} = \frac{\sin \beta}{b} = \frac{\sin \gamma}{c} = \frac{1}{2R}$ Law of cosines:  $a^2 = b^2 + c^2 - 2bc\cos\alpha$ Law of tangents:  $\frac{a+b}{a-b} = \tan\frac{\alpha+\beta}{2}/\tan\frac{\alpha-\beta}{2}$ 

#### 2.4.2 Quadrilaterals

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

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

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

#### 2.4.3 Spherical coordinates



$$\begin{aligned} x &= r \sin \theta \cos \phi & r &= \sqrt{x^2 + y^2 + z^2} \\ y &= r \sin \theta \sin \phi & \theta &= \arccos(z/\sqrt{x^2 + y^2 + z^2}) \\ z &= r \cos \theta & \phi &= \operatorname{atan2}(y, x) \end{aligned}$$

### 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:

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

#### 2.6Sums

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

$$1^{2} + 2^{2} + 3^{2} + \dots + n^{2} = \frac{n(2n+1)(n+1)}{6}$$

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

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

d41d8c, 50 lines

#### 2.7 Series

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

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

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

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

#### 2.8 Normal distribution

n	-4	-6	-8	-10
$\Phi(n)$	$3 \cdot 10^{-5}$	$10^{-9}$	$10^{-15}$	$10^{-23}$

# Data structures (3)

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

#include <bits/extc++.h>
using namespace \_\_gnu\_pbds;

template<class T>

#### HashMap.h

**Description:** Hash map with mostly the same API as unordered\_map, but  $\sim 3x$  faster. Uses 1.5x memory. Initial capacity must be a power of 2 (if provided).

#include <bits/extc++.h>
// To use most bits rather than just the lowest ones:
struct chash { // large odd number for C
 const uint64\_t C = l1(4e18 \* acos(0)) | 71;
 l1 operator()(l1 x) const { return \_\_builtin\_bswap64(x\*C); }
};
\_\_qnu\_pbds::qp\_hash\_table<ll,int,chash> h({},{},{},{},{},{1<<16});</pre>

#### IntSet.h

**Description:** set<int> implemented with bitsets. IntSet<(int)1e9> = 125 MB. 20 x faster than set<int> (50-200e6 ops = 1 s).

#### SegmentTree.h

**Description:** Zero-indexed max-tree. Can be changed by modifying iv, f. Low constant factor: 1e6 ops = 150 ms (2x faster than recursive).

```
Time: \mathcal{O}(\log N)
                                                     d41d8c, 19 lines
int iv = INT MIN;
int f(int a, int b) { return max(a, b); } // commutative
struct ST {
 vector<int> s; int n;
 ST(int n = 0) : s(2*n, iv), n(n) {}
 void update(int i, int v) { // a[i] = v
   for (s[i += n] = v; i /= 2;)
      s[i] = f(s[i * 2], s[i * 2 + 1]);
 int query(int b, int e) { // query [b, e)
   int r = iv;
    for (b += n, e += n; b < e; b /= 2, e /= 2) {
     if (b % 2) r = f(r, s[b++]);
     if (e % 2) r = f(s[--e], r);
    return r;
```

#### SegmentTree2d.h

**Description:** 0-indexed sparse  $(O(N \log N) \text{ memory})$  2D segment tree. Computes max a[x],xr[y],yr] and sets single elements a[x][y]. Offline, so requires that the elements to be updated are known in advance. Before all "real" updates: fakeUpdate(all updates); init();. High constant factor: 1e5 ops = (1s, 80MB).

```
Time: \mathcal{O}\left(\log^2 N\right).
"SegmentTree.h"
                                                       d41d8c, 28 lines
struct ST2 {
  int nx; ve<ve<pii>>> pts; vector<ST> st;
  ST2(int nx) : nx(nx), pts(2*nx) {}
  void fakeUpdate(int x, int y) {
    for (int i = x+nx; i; i /= 2)
      pts[i].push_back({y, x});
  void init() {
    for (auto &v: pts) sort(all(v)), st.emplace_back(sz(v));
  int ind(int i, int y, int x) {
    return lower_bound(all(pts[i]),pii(y, x))-pts[i].begin();
  void update(int x, int y, int v) { // a[x]/y = v
    for (int i = x+nx; i; i \neq 2)
      st[i].update(ind(i, y, x), v);
  int query(int x1, int xr, int y1, int yr) { // [xl, xr)[yl, yr)
```

for (x1 += nx, xr += nx; x1 < xr; x1 /= 2, xr /= 2) {

st[xl].query(ind(xl, yl, 0), ind(xl, yr, 0))), xl++;

st[--xr].query(ind(xr, yl, 0), ind(xr, yr, 0)));

**if** (x1 % 2) r = f(r,

**if** (xr % 2) r = f(r,

return r;

```
};
```

#### LazySegmentTree.h

const int inf = 1e9;

**Description:** Segment tree with ability to add or set values of large intervals, and compute max of intervals. Can be changed to other things. Use with a bump allocator for better performance, and SmallPtr or implicit indices to save memory.

```
Usage: Node* tr = new Node(v, 0, sz(v)); 

Time: \mathcal{O}(\log N).

"../various/BumpAllocator.h"
```

```
struct Node {
 Node *1 = 0, *r = 0;
  int lo, hi, mset = inf, madd = 0, val = -inf;
 Node (int lo, int hi):lo(lo), hi(hi) {} // Large interval of -inf
 Node (vi& v, int lo, int hi) : lo(lo), hi(hi) {
    if (lo + 1 < hi) {
      int mid = lo + (hi - lo)/2;
      l = new Node(v, lo, mid); r = new Node(v, mid, hi);
      val = max(1->val, r->val);
    else val = v[lo];
 int query(int L, int R) {
    if (R <= lo || hi <= L) return -inf;</pre>
    if (L <= lo && hi <= R) return val;</pre>
    return max(l->query(L, R), r->query(L, R));
 void set(int L, int R, int x) {
    if (R <= lo || hi <= L) return;</pre>
    if (L <= lo && hi <= R) mset = val = x, madd = 0;</pre>
      push(), l\rightarrow set(L, R, x), r\rightarrow set(L, R, x);
      val = max(1->val, r->val);
 void add(int L, int R, int x) {
    if (R <= lo || hi <= L) return;</pre>
    if (L <= lo && hi <= R) {
      if (mset != inf) mset += x;
      else madd += x;
      val += x;
    else {
      push(), l\rightarrow add(L, R, x), r\rightarrow add(L, R, x);
      val = max(1->val, r->val);
 void push() {
      int mid = lo + (hi - lo)/2;
      l = new Node(lo, mid); r = new Node(mid, hi);
    if (mset != inf)
      1->set(lo,hi,mset), r->set(lo,hi,mset), mset = inf;
    else if (madd)
      1- add (lo, hi, madd), r- add (lo, hi, madd), madd = 0;
```

#### PersistSegmentTree.h

**Description:** Persistent, Implicit, and Lazy segment tree. Can add to large intervals and compute sum of intervals. Can be modified (e.g. set + max). iv is the default value, ip is default prop value. Bounds are inclusive on BOTH sides: [lq, rq]. High constant factor: 1e5 ops = (200ms, 300MB). Use bump allocator to 2x speed.

d41d8c, 13 lines

```
0. sz(v)-1):
Time: \mathcal{O}(\log N).
                                                      d41d8c, 31 lines
int iv = 0, ip = 0;
struct Node{
 Node *1, *r; int v, p=ip, le, ri;
 Node(int le, int ri, int v=iv, int p=ip, Node*l=0, Node*r=0)
   : l(l), r(r), v(v), p(p), le(le), ri(ri) {}
  Node (Node *1, Node *r)
   : l(1), r(r), v(1->v+r->v), le(1->le), ri(r->ri) {}
  void push() {
   int m = (le+ri)/2;
    if(!1) 1 = new Node(le, m), r = new Node(m+1, ri);
    if(p!=ip) l=l->update(le,m,p), r=r->update(m+1,ri,p), p=ip;
  Node* update(int lq, int rq, int x) { // [lq, rq]
    if(lq > ri || rq < le) return this;</pre>
    if(la <= le && ri <= ra)
      return new Node(le, ri, v+x*(ri-le+1), p+x, l, r);
    return new Node(l->update(lq,rq,x), r->update(lq,rq,x));
  int query(int lq, int rq){ // [lq, rq]
    if(lq > ri || rq < le) return iv;</pre>
    if(lq <= le && ri <= rq) return v;</pre>
   push();
    return 1->query(lq, rq) + r->query(lq, rq);
};
Node* build(vi &a, int 1, int r) {
 if(l == r) return new Node(l, l, a[l]);
  int m = (1+r)/2;
  return new Node(build(a, 1, m), build(a, m+1, r));
```

Usage: Node\* tr = new Node(0, sz(v)-1); or Node\* tr = build(v,

#### KineticSegmentTree.h

Description: Kinetic Segment Tree, intervals are [l, r) To change min to max, only change the min in rec\_get. Change all < to > in pull(). Lastly, change v1 to v2 and vice versa in the calculation of x. 2e5 ops is around 500

**Time:**  $\mathcal{O}(\log^2 N)$  for heaten,  $\mathcal{O}(\log N)$  for others

```
d41d8<u>c, 75 lines</u>
const int INFX = 2e12, INFY = 2e18;
struct Line {
  int k, b;
  Line (int k = 0, int b = 0): k(k), b(b) {}
  int operator()(int x) {return k * x + b;}
struct LazyKST {
  int n; vector<Line> tree; vector<int> melt, add;
  void pull(int v) {
   int v1 = v * 2, v2 = v * 2 + 1;
   if ((tree[v1].b==tree[v2].b && tree[v1].k<tree[v2].k)||</pre>
     tree[v1].b < tree[v2].b) swap(v1, v2);</pre>
    tree[v] = tree[v2]; melt[v] = min(melt[v1], melt[v2]);
    if (tree[v1].k < tree[v2].k) {</pre>
      int x=(tree[v1].b-tree[v2].b+tree[v2].k-tree[v1].k-1)
          / (tree[v2].k - tree[v1].k);
      melt[v] = min(melt[v], x);
  void push(int v) {
   int d = add[v]; add[v] = 0;
   if (v*2 < 4*n) {
      add[v*2]+=d, add[v*2+1]+=d, melt[v*2]-=d, melt[v*2+1]-=d;
      tree[v*2].b += tree[v*2].k * d;
      tree[v*2+1].b += tree[v*2+1].k * d;
  void build(int v, int vl, int vr, vector<Line>& arr) {
    add[v] = 0;
```

```
if (vr - vl == 1) {
     tree[v] = arr[vl], melt[v] = INFX; return; }
    int vm = (vl + vr) / 2;
   build(v*2,v1,vm,arr);build(v*2+1,vm,vr,arr); pull(v);
 LazyKST(vector<Line> arr = \{\}): n(arr.size()), tree(n*4),
   melt(n * 4), add(n * 4) {
   build(1, 0, n, arr);
 void rec upd(int v, int vl, int vr, int ind, Line val) {
   push(v);
    if (ind >= vr || ind < vl) return;</pre>
    if (vr - vl == 1) {
     tree[v] = val; return; }
    int vm = (vl + vr) / 2; rec_upd(v*2, vl, vm, ind, val);
   rec_upd(v*2+1, vm, vr, ind, val); pull(v);
 void upd(int ind, Line 1) {
   rec_upd(1, 0, n, ind, 1);
 void propagate(int v, int vl, int vr) {
   if (melt[v] > 0) return;
   int vm = (vl + vr) / 2; push(v);
   propagate(v*2, v1, vm); propagate(v*2+1, vm, vr); pull(v);
 void rec_heaten(int v, int vl, int vr, int l, int r, int d) {
   if (1 >= vr || r <= v1) return;</pre>
   if (1 <= v1 && r >= vr) {
     tree[v].b += tree[v].k \star d;
     if (vr - vl > 1) melt[v] -= d, add[v] += d;
     propagate(v, vl, vr); return; }
    int vm = (vl + vr) / 2; push(v);
    rec_heaten(v * 2, vl, vm, l, r, d);
    rec_heaten(v * 2 + 1, vm, vr, 1, r, d); pull(v);
 void heaten(int 1, int r, int d) {
   rec_heaten(1, 0, n, 1, r, d);
 int rec get(int v, int vl, int vr, int l, int r) {
   if (1 >= vr || r <= v1) return INFY;</pre>
   if (1 <= v1 && r >= vr) return tree[v].b;
    int vm = (vl + vr) / 2; push(v);
    return min(rec_qet(v*2,v1,vm,1,r),rec_qet(v*2+1,vm,vr,1,r))
 int get(int 1, int r) {
   return rec_get(1, 0, n, 1, r);
};
```

#### LiChao.h

Description: Extended Li Chao tree, a lazy segment tree that maintains a sequence  $c_0, c_1, \ldots c_{n-1}$ . Can point query  $c_r$ . Can range update  $c(x) = (\max, \operatorname{add})(c(x), f(x)), f$  is linear. However, can change to any F that is a family of functions closed under function addition, and  $f \neq q \implies$ the functions intersect at one point. d41d8c, 34 lines

```
struct LiChao {
 struct Func { 11 a, b; // a*x + b
   11 operator()(11 x) const { return a*x+b; }
   Func operator+(Func r) const { return {a+r.a, b+r.b}; } };
 static constexpr Func ID_ADD{0, 0}, ID_MAX{0, 11(-1e9)};
 ve<Func> v, p; int n;
 LiChao(int N = 0) { for (n = 1; n < N; n *= 2);
   v.resize(n*2, ID_MAX); p.resize(n*2, ID_ADD); }
 void push(int i) {
   if (i < n) rep(j, 0, 2) p[i*2+j] = p[i*2+j] + p[i],
     v[i*2+j] = v[i*2+j] + p[i]; p[i] = ID_ADD; }
 // For each x in [l,r), set c[x] = max(c[x], f(x))
```

```
void max(int 1, int r, Func f, int i=1, int b=0, int e=-1) {
    if (e<0) e=n; if (1 >= e || b >= r || i >= n*2) return;
    int m = (b+e) / 2; push(i);
    if (b >= 1 && e <= r) { auto& q = v[i];</pre>
     if (g(m) < f(m)) swap(g, f);
      if (g(b) < f(b)) \max(1, r, f, i*2, b, m);
      else max(1, r, f, i*2+1, m, e); }
    else max(1, r, f, i*2, b, m), max(1, r, f, i*2+1, m, e); }
  // For each x in [l,r), set c[x] = c[x] + f(x)
  // O(\log^2 n), O(1) \text{ if } [l,r) = [0,n)
  void add(int 1, int r, Func f, int i=1, int b=0, int e=-1) {
    if (e < 0) e = n; if (1 >= e || b >= r) return;
    if (b >= 1 && e <= r) p[i] = p[i] + f, v[i] = v[i] + f;
    else { int m = (b+e) / 2; push(i);
     max(b, m, v[i], i*2, b, m);
      max(m, e, v[i], i*2+1, m, e); v[i] = ID_MAX;
      add(1, r, f, i*2, b, m); add(1, r, f, i*2+1, m, e); }
  // O(\log n), get value of c[x]
  auto query(int x) { int i = x+n; auto ret = v[i](x);
    while (i \neq 2) ret = ::max(ret+p[i](x), v[i](x));
    return ret; } };
UnionFindRollback.h
Description: Disjoint-set data structure with undo. If undo is not needed,
skip st, time() and rollback().
Usage: int t = uf.time(); ...; uf.rollback(t);
Time: \mathcal{O}(\log(N))
                                                     d41d8c, 21 lines
struct RollbackUF {
 vi e; vector<pii> st;
 RollbackUF(int n) : e(n, -1) {}
  int size(int x) { return -e[find(x)]; }
  int find(int x) { return e[x] < 0 ? x : find(e[x]); }
  int time() { return sz(st); }
  void rollback(int t) {
    for (int i = time(); i --> t;)
      e[st[i].first] = st[i].second;
    st.resize(t);
 bool join(int a, int b) {
    a = find(a), b = find(b);
    if (a == b) return false;
    if (e[a] > e[b]) swap(a, b);
    st.push back({a, e[a]});
    st.push_back({b, e[b]});
    e[a] += e[b]; e[b] = a;
    return true;
```

 $// O(\log^2 n), O(\log n) \text{ if } [l,r) = [0,n]$ 

#### SubMatrix.h

};

Description: Calculate submatrix sums quickly, given upper-left and lowerright corners (half-open).

```
Usage: SubMatrix<int> m (matrix);
m.sum(0, 0, 2, 2); // top left 4 elements
Time: \mathcal{O}(N^2+Q)
```

```
template<class T>
struct SubMatrix {
  vector<vector<T>> p;
  SubMatrix(vector<vector<T>>& v) {
    int R = sz(v), C = sz(v[0]);
    p.assign(R+1, vector<T>(C+1));
    rep(r, 0, R) rep(c, 0, C)
      p[r+1][c+1] = v[r][c] + p[r][c+1] + p[r+1][c] - p[r][c];
  T sum(int u, int 1, int d, int r) {
    return p[d][r] - p[d][l] - p[u][r] + p[u][l];
```

```
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;
```

d41d8c, 26 lines

d41d8c, 30 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];
  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;
};
```

#### LineContainer.h

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

```
struct Line {
  mutable 11 k, m, p;
 bool operator<(const Line& o)const{return k </*>*/ o.k;}
 bool operator<(11 x) const { return p < x; }</pre>
struct LineContainer : multiset<Line, less<>>> {
  // (for doubles, use inf = 1/.0, div(a,b) = a/b)
  static const 11 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 \rightarrow 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) {
```

**if** (x != begin() && isect(--x, y)) isect(x, y = erase(y));

**auto**  $z = insert(\{k, m, 0\}), y = z++, x = y;$ 

**while** ((y = x) != begin() && (--x)->p >= y->p)

while (isect(y, z)) z = erase(z);

isect(x, erase(y));

11 query(11 x) {

assert(!empty());

```
auto 1 = *lower_bound(x);
   return 1.k * x + 1.m;
};
```

#### Treap.h

**Description:** A short self-balancing tree. It acts as a sequential container with log-time splits/merges. Can also insert, move and reverse ranges. Currently lazy. Remove push() and uncomment "..new Node.." for just persistent. For both, add method t->update(prop) { return new Node(t + prop) } and use it to make push() create new nodes. 1e5 (split/merge) = 25ms. 1e5 (move/reverse) = 150ms. 1e5 (lazy+persistent) = (500ms, 600MB).

```
Time: \mathcal{O}(\log N)
#define cnt(n) ((n) ? (n) ->c : 0)
struct Node {
    Node *1 = 0, *r = 0;
    int v, y, c = 1, rev = 0;
    Node (int v) : v(v), y(rand()) {}
    Node (Node *a) { *this = *a; }
    void pull() { c = cnt(l) + cnt(r) + 1; }
    void push() {
        if (rev) { swap(l, r);
            if (1) 1->rev ^= 1; if (r) r->rev ^= 1; rev = 0;
// l = [0, i), r = [i, n]
void split(Node* t, Node*& 1, Node*& r, int i) {
    if (!t) { 1 = r = 0; return; }
    t \rightarrow push(); // t = new Node(t);
    // Replace with comments for lower_bound(i)
    if (i <= cnt(t->1) /*x->v*/) split(t->1, 1, t->1, i), r = t;
    else split(t->r, t->r, r, i - cnt(t->l) - 1 /*i*/, l = t;
    t->pull();
```

```
void merge(Node*& t, Node* 1, Node* r) {
    if (!1 || !r) t = 1 ? 1 : r;
    else if (1->v < r->v) {
        r\rightarrow push(); // r = new Node(r);
        merge (r->1, 1, r->1); t = r;
    else {
        1-push(); // l = new Node(l);
        merge(1->r, 1->r, r); t = 1;
    t->pull();
void insert(Node*& t, Node* n, int i) { // (0 indexed)
    Node *1, *r;
    split(t, 1, r, i), merge(l, l, n), merge(t, l, r);
// Move [l, r] to index k (0 indexed)
void move(Node*& t, int 1, int r, int k) {
    Node *a, *b, *c;
    split(t, a, c, r+1), split(a, a, b, 1), merge(t, a, c);
    if (k<=1) insert(t, b, k);
    else insert(t, b, k - (r - 1 + 1) /*k*/);
// Reverse [l, r] (0 indexed)
void rev(Node*& t, int 1, int r) {
    Node *a, *b, *c;
    split(t, a, c, r+1), split(a, a, b, 1);
```

b->rev ^= 1;

merge(a, a, b), merge(t, a, c);

**Description:** Computes partial sums a[0, i) and adds to single elements a[i]. Time:  $\mathcal{O}(\log N)$ .

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

#### FenwickTree2d.h

**Description:** 0-indexed sparse  $(O(N \log N) \text{ memory})$  2D fenwick tree. Computes sums a[i,j] for all i<I, j<J, and increases single elements a[i,j]. Offline, so requires that the elements to be updated are known in advance. Before all "real" updates: fakeUpdate(all updates); init();. Low constant factor: 1e5 ops = (200 ms, 35 MB)

**Time:**  $\mathcal{O}(\log^2 N)$ . (Use persistent segment trees for  $\mathcal{O}(\log N)$ .) "FenwickTree.h"

```
struct FT2 {
  vector<vi> ys; vector<FT> ft;
  FT2(int limx) : ys(limx) {}
  void fakeUpdate(int x, int y) {
    for (; x < sz(ys); x = x + 1) ys[x].push_back(y);
  void init() {
    for (vi& v : ys) sort(all(v)), ft.emplace_back(sz(v));
  int ind(int x, int y) {
    return (int) (lower_bound(all(ys[x]), y) - ys[x].begin()); }
  void update(int x, int y, ll dif) {
    for (; x < sz(ys); x | = x + 1)
      ft[x].update(ind(x, y), dif);
  11 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:
};
```

#### RMQ.h

**Description:** Range Minimum Queries on an array. Returns min(V[a], V[a +1], ... V[b - 1]) in constant time. Usage: RMQ rmq(values);

d41d8c, 16 lines

```
rmq.query(inclusive, exclusive);
Time: \mathcal{O}(|V|\log|V|+Q)
```

template < class T> struct RMO { vector<vector<T>> jmp;

RMQ(const vector<T>& V) : jmp(1, V) {

```
for (int pw = 1, k = 1; pw * 2 <= sz(V); pw *= 2, ++k) {
      jmp.emplace_back(sz(V) - pw * 2 + 1);
      rep(j,0,sz(jmp[k]))
        jmp[k][j] = min(jmp[k - 1][j], jmp[k - 1][j + pw]);
  T query(int a, int b) {
    assert (a < b); // or return inf if a == b
    int dep = 31 - __builtin_clz(b - a);
    return min(jmp[dep][a], jmp[dep][b - (1 << dep)]);</pre>
};
```

#### MoQueries.h

Description: Answer interval or tree path queries by finding an approximate TSP through the queries, and moving from one query to the next by adding/removing points at the ends. If values are on tree edges, change step to add/remove the edge (a, c) and remove the initial add call (but keep in). Time:  $\mathcal{O}(N\sqrt{Q})$ 

```
void add(int ind, int end) { ... } // add a[ind] (end = 0 or 1)
void del(int ind, int end) { ... } // remove a[ind]
int calc() { ... } // compute current answer
vi mo(vector<pii> Q) {
 int L = 0, R = 0, blk = 350; // \sim N/sqrt(Q)
  vi s(sz(Q)), res = s;
#define K(x) pii(x.first/blk, x.second ^ -(x.first/blk & 1))
  iota(all(s), 0);
  sort(all(s), [&](int s, int t) { return K(Q[s]) < K(Q[t]); });
  for (int qi : s) {
   pii q = O[qi];
   while (L > q.first) add(--L, 0);
   while (R < q.second) add(R++, 1);</pre>
    while (L < q.first) del(L++, 0);
    while (R > q.second) del(--R, 1);
    res[gi] = calc();
  return res:
vi moTree(vector<array<int, 2>> 0, vector<vi>& ed, int root=0) {
 int N = sz(ed), pos[2] = {}, blk = 350; // \sim N/sqrt(Q)
  vi s(sz(Q)), res = s, I(N), L(N), R(N), in(N), par(N);
  add(0, 0), in[0] = 1;
  auto dfs = [\&] (int x, int p, int dep, auto\& f) -> void {
   par[x] = p;
   L[x] = N;
   if (dep) I[x] = N++;
   for (int y : ed[x]) if (y != p) f(y, x, !dep, f);
   if (!dep) I[x] = N++;
   R[x] = N;
 dfs(root, -1, 0, dfs);
#define K(x) pii(I[x[0]] / blk, I[x[1]] ^ -(I[x[0]] / blk & 1))
  iota(all(s), 0);
  sort(all(s), [\&](int s, int t){ return K(Q[s]) < K(Q[t]); });
  for (int qi : s) rep(end, 0, 2) {
   int &a = pos[end], b = Q[qi][end], i = 0;
#define step(c) { if (in[c]) { del(a, end); in[a] = 0; } \
                  else { add(c, end); in[c] = 1; } a = c; }
    while (!(L[b] <= L[a] && R[a] <= R[b]))</pre>
     I[i++] = b, b = par[b];
    while (a != b) step(par[a]);
   while (i--) step(I[i]);
   if (end) res[qi] = calc();
  return res;
```

#### Wavelet.h

Description: Counts num elements in range equal to q, kth smallest, or range count. All indicies are 0 indexed and inclusive on both ides

```
Time: Operations are \mathcal{O}(logalpha).
struct wavelet{
 struct node{
   int 1, r; vector<int> val, mr; node *le{}, *ri{};
   node(int lv, int rv, vector<int> &a):l(lv), r(rv){
     for(int i : a) val.push_back(i);}
   node(int lv, int rv):1(lv), r(rv){}};
 node *root;
  // l and r are min and max values
 wavelet(vector<int> &a, int 1, int r) {
    root = new node(l, r, a); build(root);}
 void build(node *no) {
   if (no->1 == no->r) return;
   int m = no->1 + (no->r-no->1)/2, run{};
    no->le=new node(no->l, m); no->ri=new node(m+1, no->r);
    for(int i : no->val){
     if(i <= m) no->le->val.push_back(i);
     else no->ri->val.push back(i), run++;
     no->mr.push back(run); }
   build(no->le); build(no->ri);
 int count(int r, int q){
   node *no = root;
    while (no->1 != no->r) {
     if(r < 0) return 0;
     int m = no->1 + (no->r-no->1)/2;
     if (q \le m) r = r-no->mr[r], no = no->le;
     else r = no->mr[r]-1, no = no->ri;
   } return r+1;
 int count(int 1, int r, int q) {
   return count (r, q) - count (1-1, q);
 // gets kth smallest element in a[l, r]
  // return -1 if there is none
 int kth(int 1, int r, int k){
   if (k <= 0 || k > r-1+1) return -1;
     node *no = root;
     while (no->1 != no->r) {
       int c=r-no->mr[r]+1-(l==0?0:l-1-no->mr[l-1]+1);
         1 = 1 == 0 ? 0 : 1-1-no->mr[1-1]+1;
         r = r-no->mr[r]; no = no->le;
       }else{
         1 = (1 == 0 ? 0 : no->mr[1-1]);
          r = no->mr[r]-1; k -= c; no = no->ri;
        }} return no->1;
 // counts l \le i \le r and x \le a \lceil i \rceil \le y
 int rc(int 1, int r, int x, int y, node *o=nullptr) {
   if(!o) return rc(l, r, x, y, root);
   if(1 > r || o->1 > y || o->r < x) return 0;</pre>
   if(x <= o->1 && o->r <= y) return r-1+1;</pre>
    return rc(l==0?0:1-1-o->mr[1-1]+1,r-o->mr[r],x,y,o->le)
      +rc(l==0?0:o->mr[l-1],o->mr[r]-1,x,y,o->ri);
 void swap(int i, node *no = nullptr) {
   if(!no){swap(i, root); return;}
   if (no->l==no->r || i==(int)no->mr.size()-1) return;
   int m = no->1 + (no->r-no->1)/2;
   std::swap(no->val[i], no->val[i+1]);
   if(no->mr[i]-(i == 0 ? 0 : no->mr[i-1])){
     if(no->mr[i+1]-no->mr[i]) swap(no->mr[i]-1, no->ri);
     else no->mr[i]--;
    }else{
```

```
if (no->mr[i+1] - no->mr[i])no->mr[i]++;
      else swap(i-no->mr[i], no->le);}
};
```

# Numerical (4)

#### 4.1 Polynomials and recurrences

#### Polynomial.h

```
d41d8c, 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" d41d8c, 23 lines

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

```
typedef vector<double> vd;
vd interpolate(vd x, vd y, int n) {
 vd res(n), temp(n);
 rep(k, 0, n-1) rep(i, k+1, n)
```

```
y[i] = (y[i] - y[k]) / (x[i] - x[k]);
double last = 0; temp[0] = 1;
rep(k, 0, n) rep(i, 0, n) {
 res[i] += y[k] * temp[i];
  swap(last, temp[i]);
 temp[i] -= last * x[k];
return res;
```

#### BerlekampMassey.h

**Description:** Recovers any n-order linear recurrence relation from the first 2n terms of the recurrence. Useful for guessing linear recurrences after bruteforcing the first terms. Should work on any field, but numerical stability for floats is not guaranteed. Output will have size  $\leq n$ .

```
Usage: berlekampMassey({0, 1, 1, 3, 5, 11}) // {1, 2}
Time: \mathcal{O}(N^2)
```

```
"../number-theory/ModPow.h"
vector<ll> berlekampMassey(vector<ll> s) {
 int n = sz(s), L = 0, m = 0;
  vector<ll> C(n), B(n), T;
 C[0] = B[0] = 1;
 11 b = 1;
  rep(i, 0, n) \{ ++m;
   11 d = s[i] % mod;
   rep(j, 1, L+1) d = (d + C[j] * s[i - j]) % mod;
   if (!d) continue;
   T = C; 11 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 (11& 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  $S[0... \ge n-1]$  and tr[0...n-1]. Faster than matrix multiplication. Useful together with Berlekamp-Massey. Usage: linearRec({0, 1}, {1, 1}, k) // k'th Fibonacci number Time:  $\mathcal{O}\left(n^2 \log k\right)$ d41d8c, 26 lines

```
typedef vector<11> Polv;
ll linearRec(Poly S, Poly tr, ll k) {
 int n = sz(tr);
  auto combine = [&](Poly a, Poly b) {
   Poly res(n \star 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);
```

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

Usage: double func(double x) { return 4+x+.3\*x\*x; }

#### Optimization 4.2

#### 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 and no local maximum. The maximum error in the result is eps. Works equally well for maximization with a small change in the code. See Ternary-Search.h in the Various chapter for a discrete version.

```
double xmin = qss(-1000,1000,func);
Time: \mathcal{O}(\log((b-a)/\epsilon))
                                                        d41d8c, 14 lines
double gss(double a, double b, double (*f)(double)) {
 double r = (sgrt(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);
   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<sub>d41d8c, 14 lines</sub>

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

#### IntegrateAdaptive.h

**Description:** Fast integration using an adaptive Simpson's rule. Make sure (b-a)/eps < 1e-13 due to double's precision.

```
Usage: double sphereVolume = quad(-1, 1, [](double x) {
return quad(-1, 1, [&] (double y)
return quad(-1, 1, [&](double z)
return x*x + y*y + z*z < 1; {);});});
```

```
Time: \mathcal{O}\left((b-a)/\sqrt[4]{\epsilon}\right)
                                                          d41d8c, 15 lines
typedef double d;
#define S(a,b) (f(a) + 4*f((a+b) / 2) + f(b)) * (b-a) / 6
template <class F>
d rec(F& f, d a, d b, d eps, d S) {
 dc = (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)</pre>
    return T + (T - S) / 15;
  return rec(f, a, c, eps / 2, S1) + rec(f, c, b, eps / 2, S2);
```

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

#### Simplex.h

**Description:** Solves a general linear maximization problem: maximize  $c^T x$ subject to Ax < b, x > 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);
```

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

```
Time: \mathcal{O}(NM * \#pivots), where a pivot may be e.g. an edge relaxation.
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 | MP(X[j], N[j]) < MP(X[s], N[s])) s=j
struct LPSolver {
  int m, n;
  vi N, B;
  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 \mid | MP(D[i][n+1] / D[i][s], B[i])
                      < MP(D[r][n+1] / D[r][s], B[r])) r = i;
      if (r == -1) return false;
      pivot(r, s);
```

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

#### 4.3 Matrices

#### Determinant.h

**Description:** Calculates determinant of a matrix. Destroys the matrix. **Time:**  $\mathcal{O}(N^3)$ 

d41d8c, 15 lines

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

#### IntDeterminant.h

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

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

d41d8c, 18 lines

```
const l1 mod = 12345;
l1 det(vector<vector<l1>>& a) {
  int n = sz(a); l1 ans = 1;
  rep(i,0,n) {
    rep(j,i+1,n) {
    while (a[j][i] != 0) { // gcd step
        l1 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;
}
```

#### ${ m Solve Linear.h}$

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

d41d8c, 38 lines

```
typedef vector<double> vd;
```

```
const double eps = 1e-12;
int solveLinear(vector<vd>& A, vd& b, vd& x) {
 int n = sz(A), m = sz(x), rank = 0, br, bc;
 if (n) assert(sz(A[0]) == m);
 vi col(m); iota(all(col), 0);
 rep(i,0,n) {
   double v, bv = 0;
   rep(r,i,n) rep(c,i,m)
     if ((v = fabs(A[r][c])) > bv)
       br = r, bc = c, bv = v;
   if (bv <= eps) {
     rep(j,i,n) if (fabs(b[j]) > eps) return -1;
     break;
   swap(A[i], A[br]);
   swap(b[i], b[br]);
   swap(col[i], col[bc]);
   rep(j,0,n) swap(A[j][i], A[j][bc]);
   bv = 1/A[i][i];
   rep(j,i+1,n) {
     double fac = A[j][i] * bv;
     b[j] = fac * b[i];
     rep(k,i+1,m) A[j][k] = fac*A[i][k];
   rank++;
 x.assign(m, 0);
 for (int i = rank; i--;) {
   b[i] /= A[i][i];
   x[col[i]] = b[i];
   rep(j, 0, i) b[j] -= A[j][i] * b[i];
 return rank; // (multiple solutions if rank < m)
```

#### SolveLinear2.h

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

```
"SolveLinear.h" d41d8c, 7 lines rep(j,0,n) if (j != i) // instead of rep(j,i+1,n) // ... then at the end: x.assign(m, undefined); rep(j,0,rank) { rep(j,rank,m) if (fabs(A[i][j]) > eps) goto fail; x[col[i]] = b[i] / A[i][i]; fail:; }
```

#### SolveLinearBinarv.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:**  $\mathcal{O}\left(n^2m\right)$ 

```
typedef bitset<1000> bs;
int solveLinear(vector<bs>& A, vi& b, bs& x, int m) {
  int n = sz(A), rank = 0, br;
  assert (m <= sz(x));
  vi col(m); iota(all(col), 0);
  rep(i,0,n) {
    for (br=i; br<n; ++br) if (A[br].any()) break;
    if (br == n) {
      rep(j,i,n) if(b[j]) return -1;
      break;
    }
  int bc = (int)A[br]._Find_next(i-1);
  swap(A[i], A[br]);</pre>
```

```
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

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

```
d41d8c, 35 lines
int matInv(vector<vector<double>>& A) {
 int n = sz(A); vi col(n);
 vector<vector<double>> tmp(n, vector<double>(n));
 rep(i, 0, n) tmp[i][i] = 1, col[i] = i;
 rep(i,0,n) {
   int r = i, c = i;
   rep(j,i,n) rep(k,i,n)
     if (fabs(A[j][k]) > fabs(A[r][c]))
       r = j, c = k;
    if (fabs(A[r][c]) < 1e-12) return i;</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[j][i] = 0;
      rep(k, i+1, n) A[j][k] -= f*A[i][k];
     rep(k, 0, n) tmp[j][k] -= f*tmp[i][k];
    rep(j,i+1,n) A[i][j] /= v;
    rep(j,0,n) tmp[i][j] /= v;
   A[i][i] = 1;
 for (int i = n-1; i > 0; --i) rep(j,0,i) {
    double v = A[j][i];
    rep(k, 0, n) tmp[j][k] -= v*tmp[i][k];
 rep(i,0,n) rep(j,0,n) A[col[i]][col[j]] = tmp[i][j];
```

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

d41d8c, 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];</pre>
     diag[i+1] = sub[i]; tr[++i] = 1;
    } else {
     diag[i+1] -= super[i]*sub[i]/diag[i];
     b[i+1] -= b[i] * sub[i] / diag[i];
  for (int i = n; i--;) {
   if (tr[i]) {
     swap(b[i], b[i-1]);
     diag[i-1] = diag[i];
     b[i] /= super[i-1];
    } else {
     b[i] /= diag[i];
     if (i) b[i-1] -= b[i] * super[i-1];
 return b;
```

#### 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  $N = |A| + |B| \ (\sim 1s \text{ for } N = 2^{22})$ d41d8c, 35 lines

```
typedef complex<double> C;
typedef vector<double> vd;
void fft(vector<C>& a) {
  int n = sz(a), L = 31 - _builtin_clz(n);
  static vector<complex<long double>> R(2, 1);
  static vector<C> rt(2, 1); // (^ 10% faster if double)
  for (static int k = 2; k < n; k \neq 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 << L;</pre>
  vector<C> in(n), out(n);
  copy(all(a), begin(in));
  rep(i,0,sz(b)) in[i].imag(b[i]);
 fft(in);
  for (C& x : in) x *= x;
  rep(i, 0, n) out[i] = in[-i & (n - 1)] - conj(in[i]);
  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> v1;
template<int M> v1 convMod(const v1 &a, const v1 &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(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)) {
    11 \text{ av} = 11(\text{real}(\text{outl}[i]) + .5), \text{ cv} = 11(\text{imag}(\text{outs}[i]) + .5);
    11 \text{ bv} = 11(\text{imag(outl[i])} + .5) + 11(\text{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  $g = \sum_{x} a[x]g^{xk}$  $\operatorname{root}^{(mod-1)/N}$ . 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  $2^a$ . 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"
const 11 mod = (119 << 23) + 1, root = 62; // = 998244353
// For p < 2^30 there is also e.g. 5 << 25, 7 << 26, 479 << 21
^{\prime\prime}// and ^{\prime}483 << 21 (same root). The last two are > 10^{\circ}9.
```

```
typedef vector<ll> v1;
void ntt(vl &a) {
 int n = sz(a), L = 31 - __builtin_clz(n);
 static v1 rt(2, 1);
 for (static int k = 2, s = 2; k < n; k \neq 2, s++) {
   rt.resize(n);
   11 z[] = {1, modpow(root, mod >> s)};
   rep(i,k,2*k) rt[i] = rt[i / 2] * z[i & 1] % mod;
 vi rev(n):
 rep(i, 0, n) rev[i] = (rev[i / 2] | (i & 1) << L) / 2;
 rep(i,0,n) if (i < rev[i]) swap(a[i], a[rev[i]]);
 for (int k = 1; k < n; k *= 2)
    for (int i = 0; i < n; i += 2 * k) rep(j,0,k) {
     11 z = rt[j + k] * a[i + j + k] % mod, &ai = a[i + j];
     a[i + j + k] = ai - z + (z > ai ? mod : 0);
     ai += (ai + z >= mod ? z - mod : z);
vl conv(const vl &a, const vl &b) {
 if (a.empty() || b.empty()) return {};
 int s = sz(a) + sz(b) - 1, B = 32 - _builtin_clz(s),
     n = 1 << B;
 int inv = modpow(n, mod - 2);
 vl L(a), R(b), out(n);
 L.resize(n), R.resize(n);
 ntt(L), ntt(R);
 rep(i,0,n)
   out[-i \& (n - 1)] = (ll)L[i] * R[i] % mod * inv % mod;
 return {out.begin(), out.begin() + s};
```

#### FastSubsetTransform.h

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

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

```
d41d8c 16 lines
void FST(vi& a, bool inv) {
 for (int n = sz(a), step = 1; step < n; step *= 2) {
    for (int i = 0; i < n; i += 2 * step) rep(i, i, i+step) {
      int &u = a[j], &v = a[j + step]; tie(u, v) =
        inv ? pii(v - u, u) : pii(v, u + v); // AND
        inv ? pii(v, u - v) : pii(u + v, u); // OR
        pii(u + v, u - v);
 if (inv) for (int& x : a) x /= sz(a); // XOR only
vi conv(vi a, vi b) {
 FST(a, 0); FST(b, 0);
 rep(i, 0, sz(a)) a[i] *= b[i];
 FST(a, 1); return a;
```

# Number theory (5)

#### 5.1 Modular arithmetic

#### ModInverse.h

**Description:** Pre-computation of modular inverses. Assumes LIM < mod and that mod is a prime. d41d8c, 3 lines

```
const 11 mod = 1000000007, LIM = 200000;
11* inv = new l1[LIM] - 1; inv[1] = 1;
rep(i,2,LIM) inv[i] = mod - (mod / i) * inv[mod % i] % mod;
```

```
ModPow.h
```

```
d41d8c, 8 lines
```

```
const 11 mod = 1000000007; // faster if const

11 modpow(11 b, 11 e) {
    11 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.  $\operatorname{modLog}(a,1,m)$  can be used to calculate the order of a. Time:  $\mathcal{O}(\sqrt{m})$ 

11 modLog(l1 a, l1 b, l1 m) {
 ll n = (l1) sqrt(m) + 1, e = 1, f = 1, j = 1;
 unordered\_map<l1, l1> A;

while (j <= n && (e = f = e \* a % m) != b % m)
 A[e \* b % m] = j++;

if (e == b % m) return j;

if (\_gcd(m, e) == \_gcd(m, b))
 rep(i,2,n+2) if (A.count(e = e \* f % m))
 return n \* i - A[e];

return -1;

#### ModSum.h

**Description:** Sums of mod'ed arithmetic progressions.

modsum(to, c, k, m) =  $\sum_{i=0}^{\rm to-1} (ki+c)\%m$ . divsum is similar but for floored division.

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

d41d8c, 16 lines

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

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

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

#### ModMulLL.h

**Description:** Calculate  $a \cdot b \mod c$  (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 d41d8c, 11 lines

typedef unsigned long long ull;
ull modmul(ull a, ull b, ull M) {
 ll ret = a \* b - M \* ull(1.L / M \* a \* b);
 return ret + M \* (ret < 0) - M \* (ret >= (11)M);
}
ull modpow(ull b, ull e, ull mod) {
 ull ans = 1;
 for (; e; b = modmul(b, b, mod), e /= 2)
 if (e & 1) ans = modmul(ans, b, mod);
 return ans;

#### ModSgrt.h

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

Time:  $\mathcal{O}\left(\log^2 p\right)$  worst case,  $\mathcal{O}\left(\log p\right)$  for most p"ModPow.h" d41d8c, 24 lines

```
ll sgrt(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} \text{ or } 2^{(n+3)/8} * 2^{(n-1)/4} \text{ works if } p \% 8 == 5
 11 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;
 11 x = modpow(a, (s + 1) / 2, p);
  ll b = modpow(a, s, p), g = modpow(n, s, p);
  for (;; r = m) {
    11 t = b;
    for (m = 0; m < r && t != 1; ++m)
     t = t * t % p;
    if (m == 0) return x;
    11 \text{ gs} = \text{modpow}(g, 1LL \ll (r - m - 1), p);
    q = qs * qs % p;
    x = x * qs % p;
    b = b * q % p;
```

## 5.2 Primality

FastEratosthenes.h

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

```
d41d8c, 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 <= S; j += 2 * i) sieve[j] = 1;</pre>
  for (int L = 1; L <= R; L += S) {</pre>
    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 extend A randomly.

**Time:** 7 times the complexity of  $a^b \mod c$ .

```
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

 $\begin{array}{ll} \textbf{Description:} & \text{Pollard-rho randomized factorization algorithm.} & \text{Returns} \\ \text{prime factors of a number, in arbitrary order (e.g. 2299 -> \{11, 19, 11\}).} \\ \end{array}$ 

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

```
"ModMulLL.h", "MillerRabin.h"
                                                     d41d8c, 18 lines
ull pollard(ull n) {
 ull x = 0, y = 0, t = 30, prd = 2, i = 1, q;
 auto f = [\&](ull x) \{ return modmul(x, x, n) + i; \};
 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 1 = factor(x), r = factor(n / x);
 l.insert(l.end(), all(r));
 return 1;
```

## 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 (mod b).

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

#### CRT.h

**Description:** Returns  $(x, \operatorname{lcm}(m, n))$  such that  $x \equiv a \pmod{m}$ ,  $x \equiv b \pmod{n}$ . If |a| < m and |b| < n, x will obey  $0 \le x < \operatorname{lcm}(m, n)$ . Returns (0,0) if there is no solution. Assumes  $mn < 2^{62}$ . Time:  $\log(n)$ 

x = (b - a) % n \* x % n / g \* m + a;

# return {x < 0 ? x + m\*n/g : x, m\*n/g}; 5.3.1 Bézout's identity</pre>

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

```
ax + by = d
```

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

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

#### phiFunction.h

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

**Fermat's little thm**:  $p \text{ prime } \Rightarrow a^{p-1} \equiv 1 \pmod{p} \ \forall a.$ 

d41d8c, 8 lines

```
const int LIM = 50000000;
int phi[LIM];

void calculatePhi() {
  rep(i,0,LIM) phi[i] = i&1 ? i : i/2;
  for (int i = 3; i < LIM; i += 2) if(phi[i] == i)
    for (int j = i; j < LIM; j += i) phi[j] -= phi[j] / i;
}</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 < N. It will obey |p/q - x| < 1/qN.

For consecutive convergents,  $p_{k+1}q_k - q_{k+1}p_k = (-1)^k$ .  $(p_k/q_k$  alternates between > x and < x.) If x is rational, y eventually becomes  $\infty$ ; if x is the root of a degree 2 polynomial the a's eventually become cyclic.

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

```
typedef double d; // for N \sim 1e7; long double for N \sim 1e9
pair<11, 11> approximate(d x, 11 N) {
  11 LP = 0, LO = 1, P = 1, O = 0, inf = LLONG MAX; dv = x;
  for (;;) {
    ll lim = min(P ? (N-LP) / P : inf, Q ? (N-LQ) / Q : 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)NO) < abs(x - (d)P / (d)O)) ?
       make_pair(NP, NQ) : make_pair(P, Q);
    if (abs(y = 1/(y - (d)a)) > 3*N) {
      return {NP, NQ};
    LP = P; P = NP;
    LQ = Q; Q = NQ;
```

#### FracBinarySearch.h

**Description:** Given f and N, finds the smallest fraction  $p/q \in [0,1]$  such that f(p/q) is true, and  $p,q \leq N$ . You may want to throw an exception from f if it finds an exact solution, in which case N can be removed.

Usage: fracBS([](Frac f) { return f.p>=3\*f.q; }, 10); // {1,3} Time:  $\mathcal{O}(\log(N))$  d41d8c, 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 (30-bit number), 31443539979727 (45-bit), 3006703054056749 (52-bit),  $2^{61}-1$  (61-bit). There are 78498 primes less than 1000000.

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  $\mathbb{Z}_{2a}^{\times}$  is instead isomorphic to  $\mathbb{Z}_2 \times \mathbb{Z}_{2a-2}$ .

#### 5.7 Estimates

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

The number of divisors of n is at most around 250 for n < 1e6, 1500 for n < 1e9, 100 000 for n < 1e18.

#### 5.8 Mobius Function

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

Mobius Inversion:

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

Other useful formulas/forms:

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

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

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

# Combinatorial (6)

#### 6.1 Permutations

#### 6.1.1 Factorial

#### IntPerm.h

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

#### 6.1.2 Derangements

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

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

#### 6.1.3 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

#### 6.2.1 Partition function

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

$$p(0) = 1, \ p(n) = \sum_{k \in \mathbb{Z} \setminus \{0\}} (-1)^{k+1} p(n - k(3k - 1)/2)$$
$$p(n) \sim 0.145/n \cdot \exp(2.56\sqrt{n})$$

#### multinomial BellmanFord FloydWarshall PushRelabel

#### 6.2.2 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  $\binom{n}{m} \equiv \prod_{i=0}^{k} \binom{n_i}{m_i} \pmod{p}$ .

#### 6.2.3 Binomials

multinomial.h

#### General purpose numbers

#### 6.3.1 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 $c(n,2) = 0, 0, 1, 3, 11, 50, 274, 1764, 13068, 109584, \dots$ 

#### 6.3.2 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.  $\pi(j) > j$ .

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

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

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

#### 6.3.3 Stirling numbers of the second kind

Partitions of n distinct elements into exactly k groups.

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

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

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

#### 6.3.4 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,

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

#### 6.3.5 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)!)
```

#### 6.3.6 Catalan numbers

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

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

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

- 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.

# Graph (7)

#### 7.1 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  $V^2 \max |w_i| < \sim 2^{63}$ . Time:  $\mathcal{O}(VE)$ 

```
const ll inf = LLONG_MAX;
struct Ed { int a, b, w, s() { return a < b ? a : -a; }};
struct Node { ll dist = inf; int prev = -1; };
void bellmanFord(vector<Node>& nodes, vector<Ed>& eds, int s) {
 nodes[s].dist = 0;
 sort(all(eds), [](Ed a, Ed b) { return a.s() < b.s(); });</pre>
 int lim = sz(nodes) / 2 + 2; // /3+100 with shuffled vertices
 rep(i,0,lim) for (Ed ed : eds) {
   Node cur = nodes[ed.a], &dest = nodes[ed.b];
    if (abs(cur.dist) == inf) continue;
   11 d = cur.dist + ed.w;
    if (d < dest.dist) {</pre>
     dest.prev = ed.a;
     dest.dist = (i < lim-1 ? d : -inf);
 rep(i,0,lim) for (Ed e : eds) {
   if (nodes[e.a].dist == -inf)
     nodes[e.b].dist = -inf;
```

#### 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)
                                                        d41d8c, 12 lines
const 11 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;
```

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#### 7.2 Network flow

#### PushRelabel.h

**Description:** Push-relabel using the highest label selection rule and the gap heuristic. Quite fast in practice. To obtain the actual flow, look at positive values only.

```
Time: \mathcal{O}\left(V^2\sqrt{E}\right)
                                                      d41d8c, 48 lines
struct PushRelabel {
  struct Edge {
    int dest, back;
    11 f, c;
  vector<vector<Edge>> q;
  vector<11> ec;
  vector<Edge*> cur;
  vector<vi> hs; vi H;
  PushRelabel(int n): q(n), ec(n), cur(n), hs(2*n), H(n) {}
  void addEdge(int s, int t, ll cap, ll rcap=0) {
    if (s == t) return;
    g[s].push_back({t, sz(g[t]), 0, cap});
    g[t].push_back({s, sz(g[s])-1, 0, rcap});
  void addFlow(Edge& e, ll f) {
    Edge &back = g[e.dest][e.back];
    if (!ec[e.dest] && f) hs[H[e.dest]].push_back(e.dest);
    e.f += f; e.c -= f; ec[e.dest] += f;
    back.f -= f; back.c += f; ec[back.dest] -= f;
 11 calc(int s, int t) {
    int v = sz(g); H[s] = v; ec[t] = 1;
    vi co(2*v); co[0] = v-1;
    rep(i,0,v) cur[i] = g[i].data();
    for (Edge& e : g[s]) addFlow(e, e.c);
    for (int hi = 0;;) {
      while (hs[hi].empty()) if (!hi--) return -ec[s];
      int u = hs[hi].back(); hs[hi].pop_back();
      while (ec[u] > 0) // discharge u
        if (cur[u] == g[u].data() + sz(g[u])) {
          H[u] = 1e9;
          for (Edge& e : g[u]) if (e.c && H[u] > H[e.dest]+1)
            H[u] = H[e.dest]+1, cur[u] = &e;
          if (++co[H[u]], !--co[hi] && hi < v)</pre>
            rep(i, 0, v) if (hi < H[i] && H[i] < v)
              --co[H[i]], H[i] = v + 1;
          hi = H[u];
        } else if (cur[u]->c && H[u] == H[cur[u]->dest]+1)
          addFlow(*cur[u], min(ec[u], cur[u]->c));
        else ++cur[u];
```

bool leftOfMinCut(int a) { return H[a] >= sz(g); }

d41d8c, 21 lines

```
UCI
```

#### MinCostMaxFlow.h

**Description:** Min-cost max-flow. 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:**  $\mathcal{O}(FE\log(V))$  where F is max flow.  $\mathcal{O}(VE)$  for setpi.<sub>d41d8c, 80 lines</sub>

```
#include <bits/extc++.h>
const 11 INF = numeric_limits<11>::max() / 4;
struct MCMF {
  struct edge {
   int from, to, rev;
   11 cap, cost, flow;
  };
  int N;
  vector<vector<edge>> ed;
  vector<ll> dist, pi;
  vector<edge*> par;
  MCMF(int N) : N(N), ed(N), seen(N), dist(N), pi(N), par(N) {}
  void addEdge(int from, int to, 11 cap, 11 cost) {
   if (from == to) return;
   ed[from].push_back(edge{ from, to, sz(ed[to]), cap, cost, 0 });
   ed[to].push_back(edge{ to,from,sz(ed[from])-1,0,-cost,0 });
  void path(int s) {
    fill(all(seen), 0);
   fill(all(dist), INF);
   rep(i, 0, sz(par)) par[i] = NULL;
   dist[s] = 0; ll di;
    __gnu_pbds::priority_queue<pair<11, int>> q;
    vector<decltype(g)::point iterator> its(N);
   q.push({ 0, s });
    while (!q.empty()) {
     s = q.top().second; q.pop();
     seen[s] = 1; di = dist[s] + pi[s];
     for (edge& e : ed[s]) if (!seen[e.to]) {
       11 val = di - pi[e.to] + e.cost;
       if (e.cap - e.flow > 0 && val < dist[e.to]) {</pre>
         dist[e.to] = val;
         par[e.to] = &e;
         if (its[e.to] == q.end())
           its[e.to] = q.push({ -dist[e.to], e.to });
          else
            q.modify(its[e.to], { -dist[e.to], e.to });
    rep(i, 0, N) pi[i] = min(pi[i] + dist[i], INF);
  pair<11, 11> maxflow(int s, int t) {
    11 totflow = 0, totcost = 0;
    while (path(s), seen[t]) {
     11 fl = INF;
     for (edge* x = par[t]; x; x = par[x->from])
       fl = min(fl, x->cap - x->flow);
     totflow += fl;
     for (edge* x = par[t]; x; x = par[x->from]) {
       x->flow += fl;
```

```
ed[x->to][x->rev].flow -= fl;
   rep(i,0,N) for(edge& e : ed[i]) totcost += e.cost * e.flow;
   return {totflow, totcost/2};
  // 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 (edge& e : ed[i]) if (e.cap)
          if ((v = pi[i] + e.cost) < pi[e.to])</pre>
            pi[e.to] = v, ch = 1;
    assert(it >= 0); // negative cost cycle
};
Description: Flow algorithm with complexity O(VE \log U) where U =
\max |\operatorname{cap}|. O(\min(E^{1/2}, V^{2/3})E) if U = 1; O(\sqrt{V}E) for bipartite match-
                                                     d41d8c, 42 lines
struct Dinic {
 struct Edge {
   int to, rev;
   11 flow() { return max(oc - c, OLL); } // if you need flows
 };
 vi lvl, ptr, q;
 vector<vector<Edge>> adj;
 Dinic(int n) : lvl(n), ptr(n), q(n), adj(n) {}
 void addEdge(int a, int b, ll c, ll rcap = 0) {
    adj[a].push_back({b, sz(adj[b]), c, c});
    adj[b].push_back({a, sz(adj[a]) - 1, rcap, rcap});
 11 dfs(int v, int t, 11 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 (ll p = dfs(e.to, t, min(f, e.c))) {
          e.c -= p, adj[e.to][e.rev].c += p;
          return p;
    return 0;
 11 calc(int s, int t) {
   11 flow = 0; q[0] = s;
    rep(L,0,31) do { // 'int L=30' maybe faster for random data
     lvl = ptr = vi(sz(q));
     int qi = 0, qe = lvl[s] = 1;
     while (qi < qe && !lvl[t]) {
       int v = q[qi++];
       for (Edge e : adj[v])
          if (!lvl[e.to] && e.c >> (30 - L))
            q[qe++] = e.to, lvl[e.to] = lvl[v] + 1;
     while (ll p = dfs(s, t, LLONG_MAX)) flow += p;
    } while (lvl[t]);
    return flow;
 bool leftOfMinCut(int a) { return lvl[a] != 0; }
};
```

#### | MinCu

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

#### GlobalMinCut.h

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

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

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

#### GomoryHu.h

return best;

**Description:** Given a list of edges representing an undirected flow graph, returns edges of the Gomory-Hu tree. The max flow between any pair of vertices is given by minimum edge weight along the Gomory-Hu tree path.

**Time:**  $\mathcal{O}(V)$  Flow Computations

 $mat[0][t] = INT_MIN;$ 

rep(i, 0, n) mat[i][s] = mat[s][i];

## 7.3 Matching

B[b] = 0;

#### hopcroftKarp.h

**Description:** Fast maximum bipartite matching algorithm. Let the left and right partitions have size n, m respectively. Then graph g  $(g[0] \dots g[n-1])$  contains the adjacent vertices in the right partition  $(0 \dots m-1)$ . Returns the size of the matching. btoa[i] will be the match for vertex i of the right partition, or -1 if it's not matched.

```
Usage: vi btoa(m, -1); hopcroftKarp(g, btoa); 

\frac{\textbf{Time: }\mathcal{O}\left(\sqrt{V}E\right)}{\textbf{bool } dfs(\textbf{int a, int }L, \ \text{vvi\& g, vi\& btoa, vi\& A, vi\& B) } \{ \\ \textbf{if } (A[a] \ != \ L) \ \textbf{return } 0; \\ A[a] \ = \ -1;
```

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

#### 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. n and m are the sizes of the left and right partitions respectively. Then graph  $g(g[0] \dots g[n-1])$  contains the adjacent vertices in the right partition  $(0 \dots m-1)$ .

```
"hopcroftKarp.h"
                                                     d41d8c, 20 lines
vi cover(vector<vi>& g, int n, int m) {
  vi match (m, -1);
  int res = hopcroftKarp(q, 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;
```

Weighted Matching.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. Requires  $N \leq M$ . **Time:**  $\mathcal{O}(N^2M)$ 

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

#### GeneralMatching.h

**Description:** Matching for general graphs. Fails with probability N/mod. Time:  $\mathcal{O}\left(N^3\right)$ 

```
d41d8c, 40 lines
"../numerical/MatrixInverse-mod.h"
vector<pii> generalMatching(int N, vector<pii>& ed) {
  vector<vector<ll>> mat(N, vector<ll>(N)), A;
  for (pii pa : ed) {
    int a = pa.first, b = pa.second, r = rand() % mod;
    mat[a][b] = r, mat[b][a] = (mod - r) % mod;
  int r = matInv(A = mat), M = 2*N - r, fi, fj;
  assert (r % 2 == 0);
  if (M != N) do {
    mat.resize(M, vector<ll>(M));
    rep(i,0,N) {
     mat[i].resize(M);
      rep(j,N,M) {
        int r = rand() % mod;
        mat[i][j] = r, mat[j][i] = (mod - r) % mod;
 } while (matInv(A = mat) != M);
 vi has (M, 1); vector<pii> ret;
 rep(it,0,M/2) {
    rep(i,0,M) if (has[i])
      rep(j,i+1,M) if (A[i][j] && mat[i][j]) {
       fi = i; fj = j; goto done;
```

```
} assert(0); done:
if (fj < N) ret.emplace_back(fi, fj);
has[fi] = has[fj] = 0;
rep(sw,0,2) {
    ll a = modpow(A[fi][fj], mod-2);
    rep(i,0,M) if (has[i] && A[i][fj]) {
        ll b = A[i][fj] * a % mod;
        rep(j,0,M) A[i][j] = (A[i][j] - A[fi][j] * b) % mod;
    }
    swap(fi,fj);
}
return ret;</pre>
```

#### 7.4 DFS algorithms

#### SCC.h

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

**Usage:**  $scc(graph, [\&](vi\& v) \{ ... \})$  visits all components in reverse topological order. comp(i] holds the component index of a node (a component only has edges to components with lower index). ncomps will contain the number of components. **Time:**  $\mathcal{O}(E+V)$ 

```
vi val, comp, z, cont;
int Time, ncomps;
template < class G, class F> int dfs (int j, G& g, F& f) {
  int low = val[j] = ++Time, x; z.push_back(j);
 for (auto e : g[j]) if (comp[e] < 0)</pre>
    low = min(low, val[e] ?: dfs(e,q,f));
 if (low == val[i]) {
     x = z.back(); z.pop_back();
     comp[x] = ncomps;
     cont.push_back(x);
    } while (x != j);
    f(cont); cont.clear();
    ncomps++;
 return val[j] = low;
template < class G, class F > void scc(G& g, F f) {
 int n = sz(q);
 val.assign(n, 0); comp.assign(n, -1);
 Time = ncomps = 0;
 rep(i,0,n) if (comp[i] < 0) dfs(i, g, f);
```

#### BiconnectedComponents.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.

```
Usage: int eid = 0; ed.resize(N);
for each edge (a,b) {
  ed[a].emplace.back(b, eid);
  ed[b].emplace.back(a, eid++); }
  bicomps([&](const vi& edgelist) {...});
  Time: O(E+V)
```

```
vi num, st;

vector<vector<pii>> ed;

int Time;

template<class F>

int dfs(int at, int par, F& f) {
```

#### BlockCutTree 2sat EulerWalk EdgeColoring

```
int me = num[at] = ++Time, top = me;
  for (auto [y, e] : ed[at]) if (e != par) {
    if (num[v]) {
      top = min(top, num[y]);
      if (num[y] < me)
       st.push back(e);
    } else {
      int si = sz(st);
      int up = dfs(y, e, f);
      top = min(top, up);
      if (up == me) {
       st.push_back(e);
        f(vi(st.begin() + si, st.end()));
       st.resize(si);
     else if (up < me) st.push_back(e);</pre>
      else { /* e is a bridge */ }
  return top;
template<class F>
void bicomps(F f) {
 num.assign(sz(ed), 0);
  rep(i,0,sz(ed)) if (!num[i]) dfs(i, -1, f);
```

#### BlockCutTree.h

Description: Returns block-cut tree of graph. The block-cut tree has a node for each articulation point and each biconnected component. There is an edge between each biconnected component and articulation point it contains. Let i be a node in the original graph. artp[i] == true means i is an articulation point. id[i] is the node i is a part of in the block-cut tree. Let nartp = accumulate(all(artp), 0), then nodes 0...nartp-1 are articulation points. Can pass artp, id with any contents. Time:  $\mathcal{O}(n)$ .

```
vvi blockCutTree(vvi& g, vector<bool>& artp, vi& id) {
 int n = sz(q), timer = 0;
 vvi comps; vi stk, num(n), low(n);
 artp.assign(n, 0); id.assign(n, 0);
  function<void(int,int,int&)> dfs = [&](int no,int p,int &t) {
   num[no] = low[no] = ++t; stk.push back(no);
   for (int son : q[no]) {
     if (son == p) continue;
     if (num[son]) low[no] = min(low[no], num[son]);
       dfs(son, no, t); low[no] = min(low[no], low[son]);
       if (low[son] >= num[no]){
         artp[no] = (num[no]>1 || num[son]>2);
         comps.push back({no});
         while (comps.back().back() != son) {
           comps.back().push_back(stk.back());
           stk.pop_back();
 } } } };
 dfs(0, -1, timer);
 rep(i, 0, n) if (artp[i]) t.push_back({}), id[i] = sz(t)-1;
 for (auto &comp : comps) {
   t.push back({});
   for (int u : comp)
     if (!artp[u]) id[u] = sz(t)-1;
     else t[sz(t)-1].push_back(id[u]), t[id[u]].push_back(sz(t)
 return t;
```

#### 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
ts.atMostOne(\{0, \sim 1, 2\}); // <= 1 of vars 0, \sim 1 and 2 are true
ts.solve(); // Returns true iff it is solvable
ts.values[0..N-1] holds the assigned values to the vars
Time: \mathcal{O}(N+E), where N is the number of boolean variables, and E is the
number of clauses.
```

d41d8c, 56 lines

```
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 = ~li[0];
   rep(i,2,sz(li)) {
     int next = addVar();
     either(cur, ~li[i]);
     either(cur, next);
     either(~li[i], next);
     cur = ~next;
   either(cur, ~li[1]);
 vi val, comp, z; int time = 0;
 int dfs(int i) {
   int low = val[i] = ++time, x; z.push_back(i);
   for(int e : gr[i]) if (!comp[e])
     low = min(low, val[e] ?: dfs(e));
    if (low == val[i]) do {
     x = z.back(); z.pop_back();
     comp[x] = low;
     if (values[x>>1] == -1)
       values[x>>1] = x&1;
    } while (x != i);
    return val[i] = low;
 bool solve() {
   values.assign(N, -1);
   val.assign(2*N, 0); comp = val;
   rep(i,0,2*N) if (!comp[i]) dfs(i);
   rep(i,0,N) if (comp[2*i] == comp[2*i+1]) return 0;
   return 1;
```

#### EulerWalk.h

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

```
vi eulerWalk (vector<vector<pii>> & gr, int nedges, int src=0) {
  int n = sz(qr);
  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(v, 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()};
```

#### 7.5 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:  $\mathcal{O}(NM)$ 

d41d8c, 31 lines

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

#### 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

d41d8c, 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 & ~eds[q];
  rep(i,0,sz(eds)) if (cands[i]) {
   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.

d41d8c, 49 lines

```
typedef vector<br/>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>
     g.push_back(R.back().i);
     for(auto v:R) if (e[R.back().i][v.i]) T.push_back({v.i});
       if (S[lev]++ / ++pk < limit) init(T);</pre>
       int j = 0, mxk = 1, mnk = max(sz(qmax) - sz(q) + 1, 1);
       C[1].clear(), C[2].clear();
        for (auto v : T) {
         int k = 1;
          auto f = [&](int i) { return e[v.i][i]; };
          while (any_of(all(C[k]), f)) k++;
         if (k > mxk) mxk = k, C[mxk + 1].clear();
          if (k < mnk) T[j++].i = v.i;
          C[k].push_back(v.i);
        if (j > 0) T[j - 1].d = 0;
        rep(k, mnk, mxk + 1) for (int i : C[k])
         T[j].i = i, T[j++].d = k;
        expand(T, lev + 1);
      } else if (sz(q) > sz(qmax)) qmax = q;
      q.pop_back(), R.pop_back();
```

```
vi maxClique() { init(V), expand(V); return qmax; }
 Maxclique(vb conn) : e(conn), C(sz(e)+1), S(sz(C)), old(S) {
   rep(i,0,sz(e)) V.push_back({i});
};
```

#### MaximumIndependentSet.h

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

#### 7.7Trees

#### LCA.h

**Description:** Data structure for computing lowest common ancestors in a tree (with 0 as root). C should be an adjacency list of the tree, either directed or undirected.

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

```
"../data-structures/RMQ.h"
                                                      d41d8c, 21 lines
struct LCA {
  int T = 0;
  vi time, path, ret;
  RMQ<int> rmq;
  LCA(vector < vi > \& C) : time(sz(C)), rmq((dfs(C, 0, -1), ret)) {}
 void dfs(vector<vi>& C, int v, int par) {
   time[v] = T++;
    for (int y : C[v]) if (y != par) {
      path.push_back(v), ret.push_back(time[v]);
      dfs(C, y, v);
  int lca(int a, int b) {
    if (a == b) return a;
    tie(a, b) = minmax(time[a], time[b]);
    return path[rmg.query(a, b)];
 //dist(a,b){return depth[a] + depth[b] - 2*depth[lca(a,b)];}
```

#### CompressTree.h

Description: Given a rooted tree and a subset S of nodes, compute the minimal subtree that contains all the nodes by adding all (at most |S|-1) pairwise LCA's and compressing edges. Returns a list of (par, orig\_index) representing a tree rooted at 0. The root points to itself.

Time:  $\mathcal{O}(|S| \log |S|)$ 

```
d41d8c, 21 lines
typedef vector<pair<int, int>> vpi;
vpi compressTree(LCA& lca, const vi& subset) {
 static vi rev; rev.resize(sz(lca.time));
 vi li = subset, &T = lca.time;
 auto cmp = [&](int a, int b) { return T[a] < T[b]; };</pre>
 sort(all(li), cmp);
 int m = sz(li)-1;
 rep(i,0,m) {
   int a = li[i], b = li[i+1];
   li.push_back(lca.lca(a, b));
 sort(all(li), cmp);
 li.erase(unique(all(li)), li.end());
 rep(i, 0, sz(li)) rev[li[i]] = i;
 vpi ret = {pii(0, li[0])};
 rep(i, 0, sz(li)-1) {
   int a = li[i], b = li[i+1];
```

```
ret.emplace_back(rev[lca.lca(a, b)], b);
return ret;
```

#### HLD.h

**Description:** Decomposes a tree into vertex disjoint heavy paths and light edges such that the path from any leaf to the root contains at most log(n) light edges. Code does additive modifications and max queries, but can support commutative segtree modifications/queries on paths and subtrees. Takes as input the full adjacency list. VALS\_EDGES being true means that values are stored in the edges, as opposed to the nodes. All values initialized to the segtree default. Root must be 0.

Time:  $\mathcal{O}\left((\log N)^2\right)$ "../data-structures/LazySegmentTree.h"

```
template <bool VALS EDGES> struct HLD {
 int N, tim = 0;
 vector<vi> adj;
 vi par, siz, rt, pos;
 Node *tree;
 HLD(vector<vi> adj )
   : N(sz(adj_)), adj(adj_), par(N, -1), siz(N, 1),
     rt(N),pos(N),tree(new Node(0, N)) { dfsSz(0); dfsHld(0); }
 void dfsSz(int v) {
   if (par[v] != -1) adj[v].erase(find(all(adj[v]), par[v]));
   for (int& u : adj[v]) {
     par[u] = v;
     dfsSz(u);
     siz[v] += siz[u];
     if (siz[u] > siz[adj[v][0]]) swap(u, adj[v][0]);
 void dfsHld(int v) {
   pos[v] = tim++;
   for (int u : adj[v]) {
     rt[u] = (u == adj[v][0] ? rt[v] : u);
     dfsHld(u);
 template <class B> void process(int u, int v, B op) {
   for (; rt[u] != rt[v]; v = par[rt[v]]) {
     if (pos[rt[u]] > pos[rt[v]]) swap(u, v);
     op(pos[rt[v]], pos[v] + 1);
   if (pos[u] > pos[v]) swap(u, v);
   op(pos[u] + VALS_EDGES, pos[v] + 1);
 void modifyPath(int u, int v, int val) {
   process(u, v, [&](int 1, int r) { tree->add(1, r, val); });
 int queryPath(int u, int v) { // Modify depending on problem
   int res = -1e9;
   process(u, v, [&](int 1, int r) {
       res = max(res, tree->query(1, r));
   });
   return res;
 int querySubtree(int v) { // modifySubtree is similar
    return tree->query(pos[v] + VALS_EDGES, pos[v] + siz[v]);
};
```

#### CentroidDecomposition.h

**Description:**  $O(\log(n))$  depth. Builds into cent. path $(u, v) = \operatorname{path}(u, w) +$ path(w, v), w = lca(u, v) in centroid decomposition. Usage: {initialize adj/cent; build(cent, root);}

d41d8c, 25 lines

Time:  $\mathcal{O}(n \log n)$ vector<vi> adj;

#### LinkCutTree DirectedMST DominatorTree

```
vector<bool> rem(size(adj)); vi sub(size(adj));
int dfs_s(int c, int p = -1) {
    sub[c] = 1;
    for (int ch: adj[c]) {
       if (ch == p || rem[ch]) continue;
        sub[c] += dfs_s(ch, c);
    return sub[c];
int get_c(int c, int n, int p = -1) {
    for (int ch: adj[c]) {
       if (ch == p || rem[ch]) continue;
        if (sub[ch] * 2 > n) return get_c(ch, n, c);
    return c;
int build(vector<vi> &cent, int c = 0) {
    int cen = get_c(c, dfs_s(c)); // do something
    rem[cen] = true;
    for (int ch: adj[centroid]) {
        if (not rem[ch]) cent[cen].push_back(build(cent, ch));
    return cen;
```

#### 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.

**else** p->p->rot(c2, c1 != c2);

```
Time: All operations take amortized \mathcal{O}(\log N).
                                                     d41d8c, 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];
     y - > c[h ^ 1] = x;
    z->c[i ^ 1] = this;
    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);
```

```
Node* first() {
    pushFlip();
    return c[0] ? c[0]->first() : (splay(), this);
};
struct LinkCut {
  vector<Node> node;
  LinkCut(int N) : node(N) {}
  void link(int u, int v) { // add an edge (u, v)
    assert(!connected(u, v));
    makeRoot (&node[u]);
    node[u].pp = &node[v];
  void cut (int u, int v) { // remove an edge (u, v)
    Node *x = &node[u], *top = &node[v];
    makeRoot(top); x->splay();
    assert(top == (x->pp ?: x->c[0]));
    if (x->pp) x->pp = 0;
    else {
      x->c[0] = top->p = 0;
      x->fix();
  bool connected (int u, int v) { // are u, v in the same tree?
    Node* nu = access(&node[u])->first();
    return nu == access(&node[v])->first();
  void makeRoot(Node* u) {
    access(u);
    u->splay();
    if(u->c[0]) {
      u - c[0] - p = 0;
      u - c[0] - flip ^= 1;
      u - c[0] - pp = u;
      u - > c[0] = 0;
      u \rightarrow fix();
  Node* access(Node* u) {
    u->splay();
    while (Node* pp = u->pp) {
      pp \rightarrow splay(); u \rightarrow 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 11:
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"
                                                        d41d8c, 60 lines
struct Edge { int a, b; ll w; };
struct Node {
 Edge key;
 Node *1, *r;
 ll delta;
  void prop() {
```

```
key.w += delta;
if (1) 1->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->1, (a->r = merge(b, a->r)));
  return a:
void pop(Node*\& a) { a->prop(); a = merge(a->1, a->r); }
pair<11, vi> dmst(int n, int r, vector<Edge>& g) {
  RollbackUF uf(n);
  vector<Node*> heap(n);
  for (Edge e : g) heap[e.b] = merge(heap[e.b], new Node{e});
  11 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) {</pre>
      if (!heap[u]) return {-1,{}};
      Edge e = heap[u] \rightarrow 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};
```

#### DominatorTree.h

**Description:** Given a directed graph, we say b is a dominator of a if all paths  $1 \to a$  go through b. Dominator tree satisfies: (b is an ancestor of a)  $\iff$  (b is a dominator of a)

```
Usage: {initialize adj 1 indexed; build();}
Time: \mathcal{O}((n+m) \cdot \log(n))
```

void Union(int u, int v) { dsu[v] = u; }

```
d41d8c, 37 lines
const int N=0; vi adj[N];
vi dom tree[N], rg[N], buc[N];
int sdom[N], par[N], dom[N], dsu[N], lab[N], ar[N], rev[N], T;
int Find(int u, int x=0) {
    if (u == dsu[u]) return x ? -1 : u;
    int v = Find(dsu[u], x+1);
    if (v < 0) return u;</pre>
    if (sdom[lab[dsu[u]]] < sdom[lab[u]]) lab[u] = lab[dsu[u]];</pre>
    dsu[u] = v;
    return x ? v : lab[u];
```

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```
void dfs0(int u) {
    ar[u]=++T; rev[T]=u; lab[T]=T; sdom[T]=T; dsu[T]=T;
    for (int w: adj[u]) {
        if (!ar[w]) dfs0(w), par[ar[w]]=ar[u];
        rg[ar[w]].push_back(ar[u]);
void build() {
   dfs0(1);
    for (int i = T; i >= 1; i--) {
        for (int j = 0; j < sz(rq[i]); j++)</pre>
            sdom[i] = min(sdom[i], sdom[Find(rg[i][j])]);
        if (i > 1) buc[sdom[i]].push_back(i);
        for (int w : buc[i]) {
            int v = Find(w);
            if (sdom[v] == sdom[w]) dom[w] = sdom[w];
            else dom[w] = v;
        if (i > 1) Union(par[i], i);
    for (int i = 2; i <= T; i++) {</pre>
       if (dom[i] != sdom[i]) dom[i] = dom[dom[i]];
        dom_tree[rev[dom[i]]].push_back(rev[i]);
```

#### 7.8 Math

#### 7.8.1 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).

#### 7.8.2 Erdős–Gallai theorem

A simple graph with node degrees  $d_1 \ge \cdots \ge d_n$  exists iff  $d_1 + \cdots + d_n$  is even and for every  $k = 1 \dots n$ ,  $\sum_{i=1}^k d_i \le k(k-1) + \sum_{i=k+1}^n \min(d_i, k).$ 

# Geometry (8)

#### 8.1 Geometric primitives

#### Point.h

**Description:** Class to handle points in the plane. T can be e.g. double or long long. (Avoid int.)  $$_{\rm d41d8c,\ 28\ lines}$$ 

```
template <class T> int sgn(T x) { return (x > 0) - (x < 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 a, P b) const { return (a-*this): cross(b-*this); }
    T dist2() const { return x*x + y*y; }</pre>
```

```
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(); }
// returns point rotated 'a' radians ccw around the origin
P rotate(double a) const {
  return P(x*cos(a)-y*sin(a),x*sin(a)+y*cos(a)); }
friend ostream& operator<<(ostream& os, P p) {
  return os << "(" << p.x << "," << p.y << ")"; }
};</pre>
```

#### lineDistance.h

#### Description:

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



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

#### SegmentDistance.h

#### Description:

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

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

d41d8c, 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<II> 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.

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 \$1,e1\$ and \$2,e2 exists \$\{1, point\}\$ is returned. If no intersection point exists \$\{0, (0,0)\}\$ is returned and if infinitely many exists \$\{-1, (0,0)\}\$ is returned. The wrong position will be returned if P is Point<||> and the intersection point does not have integer coordinates. Products of three coordinates are used in intermediate steps so watch out for overflow if using int or ll.



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

#### sideOf.h

**Description:** Returns where p is as seen from s towards e.  $1/0/-1 \Leftrightarrow left/on line/right$ . 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.

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

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

#### OnSegment.h

"Point.h"

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

"Point.h" d41d8c, 3 lines

```
template < class P > bool on Segment (P s, P e, P p) {
  return p.cross(s, e) == 0 && (s - p).dot(e - p) <= 0;
}</pre>
```

# linearTransformation.h Description:

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



```
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();
}
```

#### LineProjectionReflection.h

**Description:** Projects point p onto line ab. Set refl=true to get reflection of point p across line ab instead. The wrong point will be returned if P is an integer point and the desired point doesn't have integer coordinates. Products of three coordinates are used in intermediate steps so watch out for overflow.

"Point.h" d41d8c, 5 lines

```
template<class P>
P lineProj(P a, P b, P p, bool refl=false) {
 P v = b - a;
 return p - v.perp()*(1+refl)*v.cross(p-a)/v.dist2();
```

#### Angle.h

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

Usage: vector<Angle>  $v = \{w[0], w[0].t360() ...\}; // sorted$ int j = 0; rep(i,0,n) { while (v[j] < v[i].t180()) ++j; } // sweeps j such that (j-i) represents the number of positively oriented triangles with vertices at 0 and i d41d8c, 35 lines

```
struct Angle {
  int x, y;
  int t:
  Angle(int x, int y, int t=0) : x(x), y(y), t(t) {}
  Angle operator-(Angle b) const { return {x-b.x, y-b.y, t}; }
  int half() const {
   assert(x || v);
    return y < 0 || (y == 0 && x < 0);
  Angle t90() const { return \{-y, x, t + (half() \&\& x >= 0)\}; \}
  Angle t180() const { return {-x, -y, t + half()}; }
  Angle t360() const { return {x, y, t + 1}; }
bool operator < (Angle a, Angle b) {
  // add a.dist2() and b.dist2() to also compare distances
  return make_tuple(a.t, a.half(), a.y * (ll)b.x) <</pre>
         make_tuple(b.t, b.half(), a.x * (11)b.y);
// Given two points, this calculates the smallest angle between
// them. i.e., the angle that covers the defined line segment.
pair<Angle, Angle> segmentAngles(Angle a, Angle b) {
  if (b < a) swap(a, b);
  return (b < a.t180() ?
          make_pair(a, b) : make_pair(b, a.t360()));
Angle operator+(Angle a, Angle b) { // point a + vector b
  Angle r(a.x + b.x, a.y + b.y, a.t);
  if (a.t180() < r) r.t--;
  return r.t180() < a ? r.t360() : r;
Angle angleDiff(Angle a, Angle b) { // angle b - angle a
  int tu = b.t - a.t; a.t = b.t;
 return {a.x*b.x + a.y*b.y, a.x*b.y - a.y*b.x, tu - (b < a)};
```

#### 8.2 Circles

#### CircleIntersection.h

P vec = b - a;

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

```
d41d8c, 11 lines
typedef Point < double > P;
bool circleInter(P a,P b,double r1,double r2,pair<P, P>* out) {
  if (a == b) { assert(r1 != r2); return false; }
```

double d2 = vec.dist2(), sum = r1+r2, dif = r1-r2,

```
if (sum*sum < d2 || dif*dif > d2) return false;
 P mid = a + vec*p, per = vec.perp() * sqrt(fmax(0, h2) / d2);
 *out = {mid + per, mid - per};
 return true;
CircleTangents.h
```

p = (d2 + r1\*r1 - r2\*r2)/(d2\*2), h2 = r1\*r1 - p\*p\*d2;

**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"
                                                     d41d8c, 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;
```

#### CircleLine.h

Description: Finds the intersection between a circle and a line. Returns a vector of either 0, 1, or 2 intersection points. P is intended to be Point<double>.

```
"Point.h"
                                                       d41d8c, 9 lines
template<class P>
vector<P> circleLine(P c, double r, P a, P b) {
 P \ ab = b - a, \ p = a + ab * (c-a).dot(ab) / ab.dist2();
 double s = a.cross(b, c), h2 = r*r - s*s / ab.dist2();
  if (h2 < 0) return {};
 if (h2 == 0) return {p};
 P h = ab.unit() * sqrt(h2);
 return {p - h, p + h};
```

#### CirclePolygonIntersection.h

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

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

```
"../../content/geometry/Point.h"
                                                       d41d8c, 19 lines
typedef Point < double > P;
#define arg(p, q) atan2(p.cross(q), p.dot(q))
double circlePoly(P c, double r, vector<P> ps) {
  auto tri = [&] (P p, P q) {
    auto r2 = r * r / 2;
    P d = q - p;
    auto a = d.dot(p)/d.dist2(), b = (p.dist2()-r*r)/d.dist2();
    auto det = a * a - b;
    if (det <= 0) return arg(p, q) * r2;</pre>
    auto s = max(0., -a-sqrt(det)), t = min(1., -a+sqrt(det));
    if (t < 0 || 1 <= s) return arg(p, g) * r2;</pre>
    P u = 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:

"Point.h"

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.



```
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"
                                                     d41d8c, 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};
```

### 8.3 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\langle P \rangle 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"
                                                         d41d8c, 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) \le eps) return !strict;
    cnt ^= ((a.y<p[i].y) - (a.y<q.y)) * a.cross(p[i], q) > 0;
  return cnt;
```

#### PolygonArea.h

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

"Point.h" d41d8c, 6 lines

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

#### PolygonCenter.h

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

Time:  $\mathcal{O}\left(n\right)$ 

"Point.h" d41d8c, 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;
}</pre>
```

# 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"



d41d8c, 13 lines

```
typedef Point<double> P;
vector<P> polygonCut(const vector<P>& poly, P s, P e) {
  vector<P> res;
  rep(i,0,sz(poly)) {
    P cur = poly[i], prev = i ? poly[i-1] : poly.back();
    bool side = s.cross(e, cur) < 0;
    if (side != (s.cross(e, prev) < 0))
      res.push_back(lineInter(s, e, cur, prev).second);
    if (side)
      res.push_back(cur);
  }
  return res;
}</pre>
```

#### HalfPlane.h

**Description:** Computes the area of the intersection of a set of half-planes. Each half-plane Line(s, e) is represented as the area to the right of ray s->e. To return the actual intersection, make small edits. It will be a ccw convex polygon, but may have duplicate/collinear points.

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

```
"Point.h", "lineIntersection.h"
                                                       d41d8c, 43 lines
#define eps 1e-8 // \ check > \sim sqrt(2)*error(lineInter) ?
typedef Point<double> P;
struct Line {
  P s, e; // Right of ray s \rightarrow e
  explicit Line(P = P(), P = P()) : s(s), e(e) {};
  P intpo(Line y) { return lineInter(s, e, y.s, y.e).S; }
  P dir() { return e - s; }
  bool contains(P x) { return (e - s).cross(x - s) < eps; }</pre>
  bool out(P x) { return !contains(x); }
bool mycmp(P a, P b) {
  if (a.x * b.x < 0) return a.x < 0;
  if (abs(a.x) < eps) {
    if (abs(b.x) < eps) return a.y > 0 && b.y < 0;
    return b.x < 0 ? a.y > 0 : 1;
```

```
if (abs(b.x) < eps) return a.x < 0 ? b.y < 0 : 0;</pre>
 return a.cross(b) > 0;
double halfPlaneIntersectionArea(ve<Line> b) {
 sort(all(b), [&](Line &a, Line &b)
    { return mycmp(a.dir(), b.dir()); });
  int n = sz(b), q = 1, h = 0;
  ve < Line > c(all(b) + 10);
  rep(i, 0, n) {
    while (q < h && b[i].out(c[h].intpo(c[h - 1]))) h--;</pre>
    while (q < h && b[i].out(c[q].intpo(c[q + 1]))) q++;</pre>
    c[++h] = b[i];
    if (q < h && abs(c[h].dir().cross(c[h - 1].dir())) < eps)</pre>
      if (c[h].dir().dot(c[h - 1].dir()) > 0) { h--;
        if (b[i].out(c[h].s)) c[h] = b[i];
      } else return 0; // 0 or infinity, bounding box <math>\Rightarrow 0
  while (q < h - 1 \&\& c[q].out(c[h].intpo(c[h - 1]))) h--;
  while (q < h - 1 \&\& c[h].out(c[q].intpo(c[q + 1]))) q++;
  if (h - q <= 1) return 0; // Intersection is empty</pre>
 c[h + 1] = c[q];
  ve<P> s; // s is the intersection as a ccw convex polygon
  rep(i, q, h+1) s.push_back(c[i].intpo(c[i + 1]));
  double area = s.back().cross(s[0]);
  rep(i, 0, sz(s)-1) area += s[i].cross(s[i + 1]);
  return area / 2;
```

#### PolygonUnion.h

**Description:** Calculates the area of the union of n polygons (not necessarily convex). The points within each polygon must be given in CCW order. (Epsilon checks may optionally be added to sideOf/sgn, but shouldn't be needed.)

**Time:**  $\mathcal{O}(N^2)$ , where N is the total number of points "Point.h", "sideOf.h"

```
typedef Point < double > P;
double rat(P a, P b) { return sqn(b.x) ? a.x/b.x : a.y/b.y; }
double polyUnion(vector<vector<P>>& poly) {
 double ret = 0:
 rep(i, 0, sz(poly)) rep(v, 0, sz(poly[i])) {
   P A = poly[i][v], B = poly[i][(v + 1) % sz(poly[i])];
   vector<pair<double, int>> seqs = {{0, 0}, {1, 0}};
   rep(j,0,sz(poly)) if (i != j) {
      rep(u,0,sz(poly[j])) {
       P C = poly[j][u], D = poly[j][(u + 1) % sz(poly[j])];
       int sc = sideOf(A, B, C), sd = sideOf(A, B, D);
       if (sc != sd) {
         double sa = C.cross(D, A), sb = C.cross(D, B);
         if (\min(sc, sd) < 0)
            segs.emplace_back(sa / (sa - sb), sgn(sc - sd));
        } else if (!sc && !sd && j<i && sqn((B-A).dot(D-C))>0) {
          segs.emplace_back(rat(C - A, B - A), 1);
          segs.emplace_back(rat(D - A, B - A), -1);
    sort (all (segs));
    for (auto& s : seqs) s.first = min(max(s.first, 0.0), 1.0);
   double sum = 0;
   int cnt = segs[0].second;
   rep(j,1,sz(segs)) {
     if (!cnt) sum += seqs[j].first - seqs[j - 1].first;
     cnt += segs[j].second;
    ret += A.cross(B) * sum;
 return ret / 2;
```

#### 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.



```
Time: \mathcal{O}(n \log n)
```

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

#### HullDiameter.h

**Description:** Returns the two points with max distance on a convex hull (ccw, no duplicate/collinear points).

Time:  $\mathcal{O}\left(n\right)$ 

d41d8c, 33 lines

#### 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"

typedef Point<11> P;

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

```
"Point.h"
                                                     d41d8c, 39 lines
#define cmp(i,j) sgn(dir.perp().cross(poly[(i)%n]-poly[(j)%n]))
#define extr(i) cmp(i + 1, i) >= 0 && cmp(i, i - 1 + n) < 0
template <class P> int extrVertex(vector<P>& poly, P dir) {
  int n = sz(poly), lo = 0, hi = n;
  if (extr(0)) return 0;
  while (lo + 1 < hi) {
   int m = (lo + hi) / 2;
   if (extr(m)) return m;
   int 1s = cmp(1o + 1, 1o), ms = cmp(m + 1, m);
    (1s < ms \mid | (1s == ms \&\& 1s == cmp(1o, m)) ? hi : 1o) = m;
  return lo;
#define cmpL(i) sgn(a.cross(poly[i], b))
template <class P>
array<int, 2> lineHull(P a, P b, vector<P>& poly) {
  int endA = extrVertex(poly, (a - b).perp());
  int endB = extrVertex(poly, (b - a).perp());
  if (cmpL(endA) < 0 \mid \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;
```

#### LinePolyEdgeIntersection.h

**Description:** For use in computing line-polygon intersection between some line s-e and polygon side p1-p2. To use, apply this function to all polygon edges, sort outputs by first key and maintain running sum of second key. Segments for which the running sum was nonzero are in the polygon, 8 lines

```
vector<pair<P, 11>> linePoly(P s, P e, P p1, P p2) {
    11 s1 = sideOf(s,e,p1), s2 = sideOf(s,e,p2);
    if(s1==s2) return {};
    11 d = 1;
    if(s1*s2)d=2;
    if(s1<s2)d*=-1;
    return {{lineInter(s,e,p1,p2).second,d}};
}</pre>
```

#### MinkowskiSum.h

**Description:** Computes Minkowski sum of two strictly convex polygons P, Q (no collinear). This is the ccw convex hull of  $\{p+q:p\in P,q\in Q\}$ . If P, Q have  $\geq 3$  sides, remove first if and ConvexHull.h.

```
Time: \mathcal{O}(n+m) without if
"Point.h", "ConvexHull.h", "sideOf.h"
                                                       d41d8c, 22 lines
typedef Point<11> P;
void reorder(vector<P> &p) {
  if (sideOf(p[0], p[1], p[2]) < 0) reverse(all(p));</pre>
  rotate(p.begin(), min_element(all(p)), p.end());
vector<P> minkowskiSum(vector<P> p, vector<P> q) {
  if (min(sz(p), sz(q)) < 3) {
    vector<P> v;
    for (P pp : p) for (P qq : q) v.push_back(pp + qq);
    return convexHull(v);
  reorder(p), reorder(q);
  rep(i, 0, 2) p.push_back(p[i]), q.push_back(q[i]);
  vector<P> r;
  11 i = 0, j = 0;
  while (i + 2 < sz(p) | | j + 2 < sz(q)) {
    r.push_back(p[i] + q[j]);
    ll cross = (p[i + 1] - p[i]).cross(q[j + 1] - q[j]);
    i += cross >= 0, j += cross <= 0;
  return r;
```

#### 8.4 Misc. Point Set Problems

#### ClosestPair.h

**Description:** Finds the closest pair of points.

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

```
"Point.h"
                                                      d41d8c, 17 lines
typedef Point<11> 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<11, pair<P, P>> ret{LLONG MAX, {P(), P()}};
 int j = 0;
 for (P p : v) {
   P d{1 + (ll)sqrt(ret.first), 0};
   while (v[j].y \le 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)

return (P(x,y) - p).dist2();

```
typedef long long T;
typedef Point<T> P;
const T INF = numeric_limits<T>::max();

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

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

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

```
Node(vector<P>&& vp) : pt(vp[0]) {
    for (P p : vp) {
      x0 = min(x0, p.x); x1 = max(x1, p.x);
      y0 = min(y0, p.y); y1 = max(y1, p.y);
    if (vp.size() > 1) {
      // split on x if width >= height (not ideal...)
      sort(all(vp), x1 - x0 >= y1 - y0 ? on_x : on_y);
      // divide by taking half the array for each child (not
      // best performance with many duplicates in the middle)
      int half = sz(vp)/2;
      first = new Node({vp.begin(), vp.begin() + half});
      second = new Node({vp.begin() + half, vp.end()});
};
struct KDTree {
 Node* root;
 KDTree(const vector<P>& vp) : root(new Node({all(vp)})) {}
 pair<T, P> search (Node *node, const P& p) {
    if (!node->first) {
      // uncomment if we should not find the point itself:
      // if (p = node \rightarrow 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)
  pair<T, P> nearest (const P& p) {
    return search(root, p);
};
```

#### FastDelaunav.h

} \*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  $\{t[0][0], t[0][1], t[0][2], t[1][0], \ldots\}$ , all counter-clockwise.

```
Time: O(nlog n)

"Point.h"

typedef Point<11> P;

typedef struct Quad* Q;

typedef __int128_t 111; // (can be ll if coords are < 2e4)

P arb(LLONG_MAX, LLONG_MAX); // not equal to any other point

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

#### VoronoiDiagram PolyhedronVolume Point3D 3dHull

```
B = b.dist2()-p2, C = c.dist2()-p2;
  return p.cross(a,b)*C + p.cross(b,c)*A + p.cross(c,a)*B > 0;
Q makeEdge(P orig, P dest) {
  O r = H ? H : new Ouad{new Ouad{new Ouad{new Ouad{0}}}};
  H = r -> 0; r -> r() -> r() = r;
  rep(i,0,4) r = r->rot, r->p = arb, r->o = i & 1 ? <math>r : r->r();
  r->p = orig; r->F() = dest;
  return r;
void splice(Q a, Q b) {
  swap(a->o->rot->o, b->o->rot->o); swap(a->o, b->o);
Q connect(Q a, Q b) {
  Q = makeEdge(a->F(), b->p);
  splice(q, a->next());
  splice(q->r(), b);
  return q;
pair<0,0> rec(const vector<P>& s) {
  if (sz(s) <= 3) {
    Q = makeEdge(s[0], s[1]), b = makeEdge(s[1], s.back());
    if (sz(s) == 2) return { a, a->r() };
    splice(a->r(), b);
    auto side = s[0].cross(s[1], s[2]);
    Q c = side ? connect(b, a) : 0;
    return {side < 0 ? c->r() : a, side < 0 ? c : b->r() };
#define H(e) e->F(), e->p
#define valid(e) (e->F().cross(H(base)) > 0)
  O 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 \rightarrow 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 \rightarrow dir; \
      splice(e, e->prev()); \
      splice(e->r(), e->r()->prev()); \
      e->o = 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());
    else
      base = connect(base->r(), LC->r());
  return { ra, rb };
vector<P> triangulate(vector<P> pts) {
  sort(all(pts)); assert(unique(all(pts)) == pts.end());
  if (sz(pts) < 2) return {};
  Q e = rec(pts).first;
  vector<Q> q = \{e\};
  int qi = 0;
  while (e->o->F().cross(e->F(), e->p) < 0) e = e->o;
```

bool circ(P p, P a, P b, P c) { // is p in the circumcircle?

111 p2 = p.dist2(), A = a.dist2()-p2,

```
#define ADD { Q c = e; do { c->mark = 1; pts.push_back(c->p); \
  q.push\_back(c->r()); c = c->next(); } while (c != e); }
  ADD; pts.clear();
  while (qi < sz(q)) if (!(e = q[qi++])->mark) ADD;
  return pts;
VoronoiDiagram.h
Description: Given points, returns map[p] = \{is\_inf, ccw\_voronoi\_region(p)\}
for each p in points. is inf=true if the region is infinite, in which case the
two rays will have length INF. There must be no duplicate points and they
should not be all collinear.
Time: \mathcal{O}(n \log n)
"FastDelaunay.h"
                                                      d41d8c, 28 lines
typedef Point<double> Pd;
map<P, pair<bool, ve<Pd>>>> voronoi(ve<P>& pts, double INF=1e9) {
 ve<P> d = triangulate(pts); map<P, ve<P>> adj;
 for (int i = 0; i < sz(d); i += 3) rep(j, 0, 3) {
   P = d[i+j], b = d[i+(j+1)%3];
    adj[a].push_back(b-a); adj[b].push_back(a-b);
 map<P, pair<bool, ve<Pd>>> v;
  auto c = [](P a) { return Pd(a.x, a.y); };
 for (auto &[p, a]: adj) {
    auto &[inf, b] = v[p];
    sort(all(a), [](P& a, P& b) {
      return pii(a.y < 0 || (a.y == 0 && a.x < 0), a.y * b.x)
        < pii(b.y < 0 | | (b.y == 0 && b.x < 0), a.x * b.y); });
    a.erase(unique(all(a)), end(a)); int m = sz(a);
    rep(i, 0, m) if (a[(i+m-1)%m].cross(a[i]) <= 0)
      \inf = 1, rotate(begin(a), begin(a)+i, end(a));
    rep(i, 0, m-inf) {
      Pd d = c(a[(i+1)%m]), e = c(a[i]);
      b.push_back(
      c(p) + (d*e.dist2()-e*d.dist2()).perp()/d.cross(e)/2);
    if (inf)
      b.push\_back(b[m-2] + c(a[m-1]).perp().unit() * INF),
      b.insert(begin(b), b[0] - c(a[0]).perp().unit() * INF);
 return v;
8.5 3D
PolyhedronVolume.h
Description: Magic formula for the volume of a polyhedron. Faces should
point outwards.
                                                       d41d8c, 6 lines
template<class V, class L>
double signedPolyVolume(const V& p, const L& trilist) {
  for (auto i : trilist) v += p[i.a].cross(p[i.b]).dot(p[i.c]);
  return v / 6;
Point3D.h
Description: Class to handle points in 3D space. T can be e.g. double or
                                                      d41d8c, 32 lines
template<class T> struct Point3D {
 typedef Point3D P;
  typedef const P& R;
```

explicit Point3D(T x=0, T y=0, T z=0) : x(x), y(y), z(z) {}

**return** tie(x, y, z) < tie(p.x, p.y, p.z); }

return tie(x, y, z) == tie(p.x, p.y, p.z); }

bool operator<(R p) const {</pre>

bool operator==(R p) const {

```
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 sgrt((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"
                                                     d41d8c, 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[1]) > q.dot(A[i]))
     q = q * -1;
    F f{q, i, j, k};
    E(a,b).ins(k); E(a,c).ins(j); E(b,c).ins(i);
   FS.push_back(f);
 rep(i,0,4) rep(j,i+1,4) rep(k,j+1,4)
   mf(i, j, k, 6 - i - j - k);
  rep(i,4,sz(A)) {
    rep(j,0,sz(FS)) {
     F f = FS[j];
      if(f.q.dot(A[i]) > f.q.dot(A[f.a])) {
       E(a,b).rem(f.c);
        E(a,c).rem(f.b);
        E(b,c).rem(f.a);
        swap(FS[j--], FS.back());
        FS.pop_back();
    int nw = sz(FS);
```

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

#### sphericalDistance.h

**Description:** Returns the shortest distance on the sphere with radius radius between the points with azimuthal angles (longitude) f1  $(\phi_1)$  and f2  $(\phi_2)$  from x axis and zenith angles (latitude) f1  $(\theta_1)$  and t2  $(\theta_2)$  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);
```

# Strings (9)

#### 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:  $\mathcal{O}\left(n\right)$ 

```
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;
}
```

#### Zfunc.h

**Description:** z[i] computes the length of the longest common prefix of s[i:] and s, except z[0] = 0. (abacaba -> 0010301)

```
Time: \mathcal{O}\left(n\right) d41d8c, 12 lines
```

```
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 - 1]);
  while (i + z[i] < sz(S) && S[i + z[i]] == S[z[i]])
    z[i]++;
  if (i + z[i] > r)
```

```
1 = i, r = i + z[i];
}
return z;
```

#### Manacher.h

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

```
Time: \mathcal{O}\left(N\right) d41d8c, 13 lines
```

```
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=0,1=0,r=0; i < n; i++) {
    int t = r-i+!z;
    if (i<r) p[z][i] = min(t, p[z][1+t]);
    int L = i-p[z][i], R = i+p[z][i]-!z;
    while (L>=1 && R+1<n && s[L-1] == s[R+1])
      p[z][i]++, L--, R++;
    if (R>r) l=L, r=R;
  }
  return p;
}
```

#### MinRotation.h

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

: O(N) d41d8c, 8 lines

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

#### SuffixArray.h

**Description:** Builds suffix array for a string. lcp(i, j) returns lcp(s[i..], s[j..]) in O(1). 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 alcp array contains longest common prefixes for neighbouring strings in the suffix array: alcp[i] = lcp(sa[i], sa[i-1]), alcp[0] = 0. The input string must not contain any zero bytes. Can remove RMQ.h if you don't need lcp().

**Time:**  $\mathcal{O}(n \log n)$  to construct,  $\mathcal{O}(1)$  for lcp.

```
"../data-structures/RMQ.h"
                                                     d41d8c, 32 lines
struct SuffixArray {
 vi sa, alcp, rsa;
 RMO<int> st;
 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 = alcp = y, iota(all(sa), 0);
   for (int j = 0, p = 0; p < n; j = max(1, j * 2), lim = p) {
     p = j, iota(all(y), n - j);
     rep(i,0,n) if (sa[i] >= j) y[p++] = sa[i] - j;
     fill(all(ws), 0);
     rep(i, 0, n) ws[x[i]] ++;
     rep(i,1,lim) ws[i] += ws[i-1];
     for (int i = n; i--;) sa[--ws[x[y[i]]]] = y[i];
      swap(x, y), p = 1, x[sa[0]] = 0;
     rep(i,1,n) a = sa[i - 1], b = sa[i], x[b] =
        (y[a] == y[b] && y[a + j] == y[b + j]) ? p - 1 : p++;
    rep(i,1,n) rank[sa[i]] = i;
    for (int i = 0, j; i < n - 1; alcp[rank[i++]] = k)
```

#### SuffixTree.h

**Description:** Ukkonen's algorithm for online suffix tree construction. Each node contains indices [l, r) into the string, and a list of child nodes. Suffixes are given by traversals of this tree, joining [l, r) substrings. The root is 0 (has l=-1, r=0), non-existent children are -1. To get a complete tree, append a dummy symbol – otherwise it may contain an incomplete path (still useful for substring matching, though).

```
Time: \mathcal{O}\left(26N\right)
```

```
struct SuffixTree {
 enum { N = 200010, ALPHA = 26 }; //N \sim 2*maxlen+10
 int toi(char c) { return c - 'a'; }
 string a; //v = cur \ node, q = cur \ position
 int t[N][ALPHA],1[N],r[N],p[N],s[N],v=0,q=0,m=2;
 void ukkadd(int i, int c) { suff:
    if (r[v]<=a) {
     if (t[v][c]==-1) { t[v][c]=m; l[m]=i;
       p[m++]=v; v=s[v]; q=r[v]; goto suff; }
     v=t[v][c]; q=l[v];
    if (q==-1 || c==toi(a[q])) q++; else {
     l[m+1]=i; p[m+1]=m; l[m]=l[v]; r[m]=q;
     p[m]=p[v]; t[m][c]=m+1; t[m][toi(a[q])]=v;
     l[v]=q; p[v]=m; t[p[m]][toi(a[l[m]])]=m;
     v=s[p[m]]; q=l[m];
      while (q<r[m]) { v=t[v][toi(a[q])]; q+=r[v]-l[v]; }</pre>
     if (q==r[m]) s[m]=v; else s[m]=m+2;
      q=r[v]-(q-r[m]); m+=2; goto suff;
 SuffixTree(string a) : a(a) {
    fill(r,r+N,sz(a));
   memset(s, 0, sizeof s);
   memset(t, -1, sizeof t);
    fill(t[1],t[1]+ALPHA,0);
   s[0] = 1; 1[0] = 1[1] = -1; r[0] = r[1] = p[0] = p[1] = 0;
    rep(i,0,sz(a)) ukkadd(i, toi(a[i]));
  // example: find longest common substring (uses ALPHA = 28)
 pii best;
  int lcs(int node, int i1, int i2, int olen) {
   if (1[node] <= i1 && i1 < r[node]) return 1;</pre>
   if (1[node] <= i2 && i2 < r[node]) return 2;</pre>
   int mask = 0, len = node ? olen + (r[node] - l[node]) : 0;
   rep(c, 0, ALPHA) if (t[node][c] != -1)
     mask |= lcs(t[node][c], i1, i2, len);
    if (mask == 3)
     best = max(best, {len, r[node] - len});
    return mask;
  static pii LCS(string s, string t) {
   SuffixTree st(s + (char) ('z' + 1) + t + (char) ('z' + 2));
    st.lcs(0, sz(s), sz(s) + 1 + sz(t), 0);
```

};

Eertree.h

return st.best;

#### Eertree Depam Hashing AhoCorasick

```
drome which ends at i'th position in the string.
Time: \mathcal{O}(n \log \alpha)
                                                       d41d8c, 31 lines
struct enode{int len; enode *suf, *ne[26];
  enode(int 1, enode *s=nullptr):len(1), suf(s){
    for(int i{}; i<26; ++i) ne[i] = nullptr;</pre>
struct eertree{
  string s;
  enode *curr, *root, *iroot;
  vector<enode*> p;
  eertree(){
    iroot = new enode(-1),iroot->suf = iroot;
    root = new enode(0),root->suf = iroot;
    curr = root;
  // returns true if new palindrome added
  bool append(char c) {
    bool add{}; s += c; enode *t = curr; int ind = s.size()-1;
    while(ind-1-t->len<0 || s[ind-1-t->len] != c)t=t->suf;
    if(!t->ne[c-'a']){
      enode *ne = t->ne[c-'a'] = new enode(t->len + 2);
      if(ne->len == 1)ne->suf = root;
      else{
        do t = t -> suf;
        while(ind-1-t->len<0||s[ind-1-t->len]!=c);
        ne->suf = t->ne[c-'a'];
      curr = ne, add = true;
    }else curr = t->ne[c-'a'];
    p.push_back(curr); return add;
```

**Description:** Builds eartree for a string, p[i] is the node for longest palin-

#### Depam.h

**Description:** alpha is [OS, OS+SIZE), sentinel (OSET-1) no[0]: len 0(null in nxt), node[1]: len -1, node[2, nodLen] pali str ts[pos,pos+len] pre/suff when created, fail: longest proper pali quick[a]: longes proper pre/suff followed/preceded by a nos->nodes, noLen->nodeLength,his->history,OSET/OS->OFFSET

```
template <class T, int SIZE, T OSET> struct Depam {
  struct Node {int len, pos, fail, nxt[SIZE], quick[SIZE];};
  int noLen, pre, suf, nL, nR, 1, r, hisLen;
  vector<Node>nos;vector<T>tsBuffer;T*ts;vector<pii>his;
  // curr whole str: ts[l, r) (-nL <= l <= 0 <= r <= nR)
  //ts \rightarrow ((\sim pre)/suf before pushF/B, par of made node or -1)
  void setNode(Node &g, int 1, int p, int f){
   g.len=1, g.pos=p, g.fail=f; memset(g.nxt, 0, sizeof(g.nxt));}
  Depam(int nL_, int nR_) : nL(nL_), nR(nR_) {
   noLen = 2; hisLen = 1 = r = pre = suf = <math>0;
   nos.resize(2 + nL + nR); his.resize(nL + nR);
   setNode(nos[0], 0, 0, 1); setNode(nos[1], -1, 0, 1);
   rep(a, 0, SIZE) nos[0].quick[a]=nos[1].quick[a] = 1;
   tsBuffer.assign(1 + nL + nR + 1, OSET - 1);
   ts = tsBuffer.data() + (1 + nL);}
  const Node &operator[](int u) const {return nos[u];}
  void pushFront(T t) {
   const int a=t-OSET;his[hisLen++]={~pre, -1};ts[--1]=t;
   if (ts[l+1+nos[pre].len]!=t) pre=nos[pre].quick[a];
   Node &f = nos[pre]; if (!f.nxt[a]) {
     his[hisLen-1].S = pre; Node &g=nos[noLen];
```

```
setNode(g,f.len+2,l,nos[f.quick[a]].nxt[a]);
    memcpy(q.quick, nos[q.fail].quick, sizeof(q.quick));
    g.quick[ts[l + nos[g.fail].len] - OSET] = g.fail;
    f.nxt[a]=noLen++;/*needs to be after setting g.fail*/}
  if (nos[pre = f.nxt[a]].len == r - 1) suf = pre;}
void pushBack(T t) {
  const int a=t-OSET; his[hisLen++]={suf,-1}; ts[r++]=t;
  if (ts[r-2-nos[suf].len]!=t) suf=nos[suf].quick[a];
  Node &f = nos[suf]; if (!f.nxt[a]) {
   his[hisLen-1].S = suf; Node &g=nos[noLen];
    setNode(q,f.len+2,r-q.len,nos[f.quick[a]].nxt[a]);
    memcpy(g.quick, nos[g.fail].quick, sizeof(g.quick));
    g.quick[ts[r - 1 - nos[g.fail].len] - OSET] = g.fail;
    f.nxt[a]=noLen++;/*needs to be after setting g.fail*/}
  if (nos[suf = f.nxt[a]].len == r - 1) pre = suf;}
void undo() {
  const pii h = his[--hisLen]; if (h.F < 0) \{//pF\}
    if(nos[pre].len==r-l)suf=nos[suf].fail;
    pre=\sim h.F; if(\sim h.S) \{--noLen; nos[h.S].nxt[ts[l]-OSET]=0; \}
    ts[l++]=OSET-1; else{/*pB*/if(nos[suf].len==r-1)
    pre=nos[pre].fail; suf=h.F; if (~h.S) {--noLen;
    nos[h.S].nxt[ts[r-1]-OSET]=0;}ts[--r]=OSET-1;}};
```

#### Hashing.h

**Description:** Self-explanatory methods for string hashing. d41d8c, 44 line

```
// Arithmetic mod 2^64-1. 2x slower than mod 2^64 and more
// code, but works on evil test data (e.g. Thue-Morse, where
// ABBA... and BAAB... of length 2^10 hash the same mod 2^64).
// "typedef ull H;" instead if you think test data is random,
// or work mod 10^9+7 if the Birthday paradox is not a problem.
typedef uint64_t ull;
struct H {
 ull x; H(ull x=0) : x(x) {}
 H operator+(H o) { return x + o.x + (x + o.x < x); }
 H operator-(H o) { return *this + ~o.x; }
 H operator*(H o) { auto m = ( uint128 t)x * o.x;
   return H((ull)m) + (ull)(m >> 64); }
 ull get() const { return x + !~x; }
 bool operator==(H o) const { return get() == o.get(); }
 bool operator<(H o) const { return get() < o.get(); }</pre>
static const H C = (11)1e11+3; // (order \sim 3e9; random \ also \ ok)
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];
vector<H> getHashes(string& str, int length) {
 if (sz(str) < length) return {};</pre>
 H h = 0, pw = 1;
 rep(i,0,length)
   h = h * C + str[i], pw = pw * C;
  vector<H> ret = {h};
  rep(i,length,sz(str)) {
    ret.push_back(h = h * C + str[i] - pw * str[i-length]);
  return ret;
```

```
H hashString(string& s){H h{}; for(char c:s) h=h*C+c; return h;}
```

#### AhoCorasick.h

**Description:** Aho-Corasick automaton, used for multiple pattern matching. Initialize with AhoCorasick ac(patterns); the automaton start node will be at index 0. find(word) returns for each position the index of the longest word that ends there, or -1 if none. findAll(-, word) finds all words (up to  $N\sqrt{N}$  many if no duplicate patterns) that start at each position (shortest first). Duplicate patterns are allowed; empty patterns are not. To find the longest words that start at each position, reverse all input. For large alphabets, split each symbol into chunks, with sentinel bits for symbol boundaries.

**Time:** construction takes  $\mathcal{O}(26N)$ , where N = sum of length of patterns. find(x) is  $\mathcal{O}(N)$ , where N = length of x. findAll is  $\mathcal{O}(NM)$ .

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

#### SuffixAutomaton.h

d41d8c, 26 lines

```
struct SuffixAutomaton {
 struct state {
   int len, link; map<char, int> next;
   state() { len = link = 0; next.clear(); } };
  ve<state> st; int sz, last;
 void sa init() {
   st.resize(1); st[0].len = 0;
   st[0].link = -1; sz++; last = 0; }
  void sa extend(char c) {
   int cur = sz++; st.emplace back();
   st[cur].len = st[last].len + 1; int p = last;
   while (p != -1 \&\& !st[p].next.count(c))
     st[p].next[c] = cur, p = st[p].link;
   if (p == -1) st[cur].link = 0;
   else {
     int q = st[p].next[c];
     if (st[p].len + 1 == st[q].len) st[cur].link = q;
     else { st.emplace_back();
       int clone = sz++; st[clone].len = st[p].len + 1;
       st[clone].next= st[q].next; st[clone].link= st[q].link;
       while (p != -1 && st[p].next[c] == q)
            st[p].next[c] = clone, p = st[p].link;
       st[q].link = st[cur].link = clone;
   } last = cur;
 } };
```

# Dynamic Programming (10)

#### KnuthOpt.h

**Description:** Solves  $dp[j] = \min_{0 < i < j} (dp[i] + c(i+1,j)) \ c(i,j)$  must satisfy Quadrangle Inequality:  $\operatorname{convex}(c(b,d) - c(b,c) \le c(a,d) - c(a,c))$  for all a < b < c < d. For concave flip < to >. To use, fill out c, f, T. Time:  $\mathcal{O}(N \log N)$ d41d8c, 47 lines

```
struct Transition { int i, l, r; };
T knuth_convex(int n) {
  ve<T> dp(n+1);
  auto c = [&] (int 1, int r) \rightarrow 11 {}; // 1 <= l <= r <= n
  auto f = [&](int i, int j) -> T {return dp[i] + c(i+1, j);};
  deque < Transition > dq = \{\{0, 1, n\}\};
  rep(i, 1, n+1) {
    dp[i] = f(dq.front().i, i);
    if (dq.front().r == i) dq.pop_front(); else dq.front().l++;
    while (sz(dg) && f(dg.back().i, dg.back().l)
                     >= f(i, dq.back().1)) dq.pop_back();
    if (!sz(dq)) dq.push_back({i, i+1, n});
    else {
      auto [pi, 1, r] = dq.back();
      while (1 < r) {
        int m = (1+r+1)/2;
        if (f(pi, m) >= f(i, m)) r = m-1;
        else 1 = m;
```

```
if ((dq.back().r=r) != n) dq.push_back({i, r+1, n});
 return dp[n];
T knuth_concave(int n) {
 ve<T> dp(n+1);
  auto c = [&] (int 1, int r) -> 11 {}; // 1 <= l <= r <= n
  auto f = [\&] (int i, int j) \rightarrow T \{return dp[i] + c(i+1, j);\};
  stack<Transition> s; s.push({0, 1, n});
 rep(i, 1, n+1) {
    dp[i] = f(s.top().i, i);
    if (s.top().r == i) s.pop(); else s.top().l++;
    while (sz(s) && f(s.top().i, s.top().r)
                    > f(i, s.top().r)) s.pop();
    if (!sz(s)) s.push({i, i+1, n});
      auto [pi, l, r] = s.top();
      while (1 < r) {
       int m = (1+r)/2;
       if (f(pi, m) > f(i, m)) 1 = m+1;
        else r = m;
      if ((s.top().l=1) != i+1) s.push({i, i+1, l-1});
  return dp[n];
```

#### Knuth2d1d.h

**Description:** (Array Merging) When doing DP on intervals: a[i][j] = $\min_{i < k < j} (a[i][k] + a[k+1][j]) + f(i,j)$ . (Array Partition) For min cost to split an array of length n into m partitions. Need Quadrangle Inequality,  $f(a,c) + f(b,d) \leq f(a,d) + f(b,c)$ , and for array partition, also need f(b,c) < f(a,d) for all a < b < c < d. (i, j) are 0 indexed. Time:  $\mathcal{O}(NM, N^2)$ 

```
int knuth_array_partition(int n, int m, F f){
 // dp[index][partition number]
 // f(i, j, k) = cost \ of \ [i, j] \ as \ the \ kth \ partition
 vvi dp(n+1, vi(m+1, 1e18)), opt(n+1, vi(m+1, n));
 dp[0][0] = opt[0][0] = 0;
 irep(i, n, 0) opt[i][1] = 0, dp[i][1] = f(0, i-1, 1);
 rep(j, 2, m+1) irep(i, n, 0)
    rep(k, opt[i][j-1], i==n?n+1:min(i+1, opt[i+1][j]+1))
      if(dp[i][j] >= dp[k][j-1] + f(k, i-1, j))
       dp[i][j]=dp[k][j-1]+f(k, i-1, j), opt[i][j] = k;
 return dp.back().back();
int knuth_array_merging(int n, F f) {
 // dp[left bound][right bound], inclusive on both sides
 // f(i, j) calls to determine cost of [i, j]
 vvi dp(n, vi(n, le18)), opt(n, vi(n));
 rep(i, 0, n) dp[i][i] = f(i, i), opt[i][i] = i;
 irep(i, n-1, -1) rep(j, i+1, n)
   rep(k, opt[i][j-1], min(j-1, opt[i+1][j])+1)
     if(dp[i][j] >= dp[i][k] + dp[k+1][j] + f(i, j))
       dp[i][j]=dp[i][k]+dp[k+1][j]+f(i, j), opt[i][j] = k;
 return dp[0].back();
```

#### DivideAndConquerDP.h

**Description:** Given  $a[i] = \min_{lo(i) \le k < hi(i)} (f(i, k))$  where the (minimal) optimal k non-decreasing with i, computes a[i] for i = L...R - 1. Time:  $\mathcal{O}((N + (hi - lo)) \log N)$ 

```
d41d8c, 18 lines
struct DP { // Modify at will:
 int lo(int ind) { return 0; }
 int hi(int ind) { return ind; }
```

```
11 f(int ind, int k) { return dp[ind][k]; }
void store(int ind, int k, ll v) { res[ind] = pii(k, v); }
void rec(int L, int R, int LO, int HI) {
 if (L >= R) return;
  int mid = (L + R) \gg 1;
  pair<11, int> best (LLONG MAX, LO);
  rep(k, max(LO,lo(mid)), min(HI,hi(mid)))
   best = min(best, make_pair(f(mid, k), k));
  store(mid, best.second, best.first);
  rec(L, mid, LO, best.second+1);
  rec(mid+1, R, best.second, HI);
void solve(int L, int R) { rec(L, R, INT_MIN, INT_MAX); }
```

#### AliensTrick.h

**Description:** For problems asking for min/max cost using k or  $\leq k$  'groups'. Replaces the groups state of the DP with a binary search. Let f(x) be the best cost using x groups. f(x) must be convex/concave. Ensure  $\max(\|f(x) - f(x)\|_{L^2(x)})$  $f(x-1)\|$   $< \max(\|l\|, \|r\|)$ : almost guaranteed for gen\_eq=0. gen\_eq=0 is:  $\operatorname{convex}(\operatorname{nonincreasing} = k \mid \operatorname{general} \leq k), \operatorname{concave}(\operatorname{nondecreasing} = k \mid \operatorname{general} \leq k)$ eral  $\leq k$ ), gen\_eq=1 is: convex/concave(general = k). Fill out best: add +lambda extra cost for each group and return min{cost, groups} for convex, max for concave. If gen\_eq=1 and  $n \cdot \max(\|f(x) - f(x-1)\|) > 4 \cdot 10^{18}$ , remove /n + use = int128 cost.

```
Time: O(\log(10^{18})/N)
                                                             d41d8c, 12 lines
int n, k; // global
```

```
pii best(int lambda) { }
11 aliens_trick(bool convex, bool gen_eq=0) {
    11 1 = gen eg ? -4e18/n : -4e18*!convex;
    11 r = gen_eq ? 4e18/n : 4e18*convex;
    int sgn = convex ? 1 : -1;
    while (1 < r) {
       11 m = 1 + (r-1)/2;
        sgn*best(m).S \le sgn*k ? r = m : 1 = m+1;
    return best(r).F - r*k;
```

# Various (11)

#### 11.1 Intervals

#### IntervalContainer.h

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

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

d41d8c, 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) {</pre>
   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);
}
```

#### 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  $\mid \mid$  R.empty(). Returns empty set on failure (or if G is empty).

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

d41d8c, 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]; });
    T cur = G.first;
    int at = 0;
    while (cur < G.second) { // (A)
        pair < T, int > mx = make_pair(cur, -1);
        while (at < sz(I) && I[S[at]].first <= cur) {
            mx = max(mx, make_pair(I[S[at]].second, S[at]));
            at++;
        }
        if (mx.second == -1) return {};
        cur = mx.first;
        R.push_back(mx.second);
    }
    return R;</pre>
```

#### 11.2 Misc. algorithms

#### 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;
}</pre>
```

#### LIS.h

**Description:** Compute indices for the longest increasing subsequence. **Time:**  $\mathcal{O}\left(N\log N\right)$ 

```
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  $S \le t$  such that S is the sum of some subset of the weights.

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

d41d8c, 16 lines

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

#### Random.h

 ${\bf Description:}\ {\rm For}\ {\rm rng}$ 

d41d8c, 2 lines

```
mt19937_64 rng();
int a_to_b = uniform_int_distribution<int>(a, b) (rng);
```

## 11.3 Optimization tricks

#### 11.3.1 Bit hacks

- for (int x = m; x; ) { --x &= 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.
- rep(b,0,K) rep(i,0,(1 << K))
   if (i & 1 << b) D[i] += D[i^(1 << b)];
  computes all sums of subsets.</pre>

#### 11.3.2 Pragmas

- #pragma GCC optimize ("Ofast") will make GCC auto-vectorize loops and optimizes floating points better.
- #pragma GCC optimize ("unroll-loops") reduces number of branches and optimizes parallel computation, but increased code size can lead to instruction cache misses.
- #pragma GCC target ("avx2") can double performance of vectorized code, but causes crashes on old machines.

FastMod.h

**Description:** Compute a%b about 5 times faster than usual, where b is constant but not known at compile time. Returns a value congruent to  $a \pmod{b}$  in the range [0, 2b).

```
typedef unsigned long long ull;
struct FastMod {
  ull b, m;
  FastMod(ull b) : b(b), m(-1ULL / b) {}
  ull reduce(ull a) { // a % b + (0 or b)
    return a - (ull) ((__uint128_t(m) * a) >> 64) * b;
  }
};
```

#### BumpAllocator.h

**Description:** When you need to dynamically allocate many objects and don't care about freeing them. "new X" otherwise has an overhead of something like 0.05us + 16 bytes per allocation.

```
// Either globally or in a single class:
static char buf[450 << 20];
void* operator new(size_t s) {
    static size_t i = sizeof buf;
    assert(s < i);
    return (void*) &buf[i -= s];
} void operator delete(void*) {}</pre>
```

#### BumpAllocatorSTL.h

**Description:** BumpAllocator for STL containers.

```
char buf[450 << 20] alignas(16);
size_t buf_ind = sizeof buf;
template<class T> struct small {
    typedef T value_type; small() {}
    template<class U> small(const U&) {}
    T* allocate(size_t n) {
        buf_ind &= 0 - alignof(T);
        return (T*)(buf + buf_ind);
    } void deallocate(T*, size_t) {};
```