

University of Warsaw

UW1

Adam Sołtan, Ivan Gechu, Franciszek Witt

AMPPZ 2024 2024-11-11 nnoremap : ;

2	Mathematics	1
3	Data structures	2
4	Numerical	4
5	Number theory	8
6	Combinatorial	11
7	Graph	12
8	Geometry	18
9	Strings	23
10 Various		

Contest (1)

set nu et ts=2 sw=2

colorscheme habamax

nnoremap ; :

filetype indent on syntax on

1 Contest

```
sol.cpp
#include <bits/stdc++.h>
using namespace std;
#define rep(i, a, b) for (int i = (a); i < (b); i++)
#define all(x) begin(x), end(x)
#define sz(x) int((x).size())
using 11 = long long;
using pii = pair<int, int>;
using vi = vector<int>;
#ifdef LOCAL
auto operator<<(auto& o, auto x) -> decltype(x.first, o);
auto operator<<(auto& o, auto x) -> decltype(x.end(), o) {
 for (int i = 0; auto y : x) o << ", " + !i++ * 2 << y;</pre>
 return o << "}"; }
auto operator<<(auto& o, auto x) -> decltype(x.first, o) {
 return o << "(" << x.first << ", " << x.second << ")"; }</pre>
void print(auto... x) { ((cerr << x << " "), ...) << endl; }</pre>
#define debug(x...) __print("[" #x "]:", x)
#define debug(...) 2137
#endif
int main() {
 cin.tie(0)->sync_with_stdio(0);
.vimrc
                                                             8 lines
```

hi MatchParen ctermfg=66 ctermbg=234 cterm=underline

```
inoremap {<cr> {<cr>}<esc>0 <bs>
.bashrc
c() {
  g++ -std=c++20 -fsanitize=address,undefined -g \
  -DLOCAL -Wall -Wextra -Wshadow $1.cpp -o $1;
nc() { q++ -std=c++20 -02 $1.cpp -0 $1; }
alias rm='trash'
alias mv='mv -i'
alias cp='cp -i'
hash.sh
# Hashes a file, ignoring all whitespace and comments. Use for
# verifying that code was correctly typed.
cpp -dD -P -fpreprocessed | tr -d '[:space:]' | md5sum |cut -c-6
for((i=1;i>0;i++)) do
  echo "$i"
  echo "$i" | ./gen > int
 diff -w <(./sol < int) <(./slow < int) || break
```

Mathematics (2)

2.1Trigonometry

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

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

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

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

Geometry

2.2.1 Triangles

Side lengths: a, b, cSemiperimeter: $p = \frac{a+b+c}{2}$

```
Area: A = \sqrt{p(p-a)(p-b)(p-c)}
```

Circumradius:
$$R = \frac{abc}{4A}$$

Inradius:
$$r = \frac{A}{p}$$

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

Length of bisector (divides angles in two):

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

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

Law of tangents: $\frac{a+b}{a-b} = \frac{\tan \frac{\alpha+\beta}{2}}{\tan \frac{\alpha-\beta}{2}}$

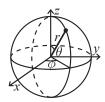
Quadrilaterals

With side lengths a, b, c, d, diagonals e, f, diagonals angle θ , 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 $A = \sqrt{(p-a)(p-b)(p-c)(p-d)}$.

Spherical coordinates



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

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

$$z = r \cos \theta \qquad \phi = a\tan(2(y, x))$$

2.3 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:

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

2.4 Sums

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

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

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

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

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

2.5 Series

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

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

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

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

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

2.6 Optimization

2.6.1 Lagrange multipliers

To optimize $f(x_1,...,x_n)$ subject to the constraints $g_k(x_1,...,x_n)=0$, a necessary condition for $(x_1,...,x_n)$ to be a local extremum is that the gradient $\nabla f(x_1,...,x_n)$ must be a linear combination of the gradients $\nabla g_k(x_1,...,x_n)$.

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

HashMap.h

Description: Hash map with mostly the same API as unordered_map, but ~3x faster. Uses 1.5x memory. Initial capacity must be a power of 2 (if provided).

```
#include <ext/pb_ds/assoc_container.hpp>
using namespace __gnu_pbds;
// 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); }
};
gp_hash_table<ll, l1, chash> h({},{},{},{},{1<<16});</pre>
```

LazySegtree.h

Description: Basic segment tree template with lazy propagation. Can be easily extended with advanced functionality.

```
if (mid < r) rec(u + (mid-lo) * 2, mid, hi, l, r, mod, f);
  if (mod) t[u] = join(t[u + 1], t[u + (mid - lo) * 2]);
}
Node get(int l, int r) {
  bool in = 0;
Node res;
  rec(0, 0, n, l, r, 0, [&](int u, int, int) {
    res = in ? join(res, t[u]) : t[u], in = 1;
  });
  return res;
}
void modify(int l, int r) {
  rec(0, 0, n, l, r, 1, [&](int u, int lo, int hi) {});
};</pre>
```

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: O(log(N))

de4ad0, 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]); }</pre>
 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;
```

LineContainer.h

Description: Container where you can add lines of the form kx+m, and query maximum values at points x. Useful for dynamic programming ("convex hull trick").

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

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

```
mutable 11 k, m, p;
bool operator<(const Line& o) const { return k < o.k; }
bool operator<(11 x) const { return p < x; }
};

struct LineContainer : multiset<Line, less<>> {
    // (for doubles, use inf = 1/.0, div(a,b) = a/b)
    static const 11 inf = LLONG_MAX;
    11 div(11 a, 11 b) { // floored division
        return a / b - ((a ^ b) < 0 && a % b); }
bool isect(iterator x, iterator y) {
    if (y == end()) return x->p = inf, 0;
    if (x->k == y->k) x->p = x->m > y->m ? inf : -inf;
    else x->p = div(y->m - x->m, x->k - y->k);
    return x->p > y->p;
}
void add(11 k, 11 m) {
    auto z = insert({k, m, 0}), y = z++, x = y;
```

Treap FenwickTree FenwickTree2d RMQ MoQueries

```
while (isect(y, z)) z = erase(z);
    if (x != begin() \&\& isect(--x, y)) isect(x, y = erase(y));
    while ((y = x) != begin() && (--x)->p >= y->p)
     isect(x, erase(v));
  11 query(ll x) {
    assert(!emptv());
    auto 1 = *lower bound(x);
    return l.k * x + l.m;
};
```

Treap.h

Description: Treap with lazy propagation and parent information. It acts as a sequential container with log-time splits/joins, and is easy to augment with additional data.

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

```
mt19937 rng(2137);
struct Node {
 Node *1 = 0, *r = 0, *p = 0;
 int val, pr, c = 1;
 Node(int x) : val(x), pr(rng()) {}
 void pull();
 void push();
int cnt(Node* n) { return n ? n->c : 0; }
void Node::pull() { c = cnt(1) + cnt(r) + 1; }
void Node::push() {}
pair<Node*, Node*> split(Node* n, int k) {
 if (!n) return {};
 n->push();
  if (cnt(n->1) >= k) { // "n->val >= k" for lower_bound(k)}
   auto pa = split(n->1, k);
   n->1 = pa.second;
   n->p = 0;
   if (n->1) n->1->p = n;
   n->pull();
   return {pa.first, n};
   auto pa = split(n->r, k - cnt(n->1) - 1); // and just "k"
   n->r = pa.first;
   n->p = 0;
   if (n->r) n->r->p = n;
   n->pull();
   return {n, pa.second};
Node* merge(Node* 1, Node* r) {
 if (!1 || !r) return 1 ?: r;
 if (1->pr > r->pr) {
   1->push();
   1->r = merge(1->r, r);
   1->r->p = 1;
   1->pull();
   return 1;
  } else {
    r->push();
   r->1 = merge(1, r->1);
   r->1->p = r;
   r->pull();
   return r;
```

```
FenwickTree.h
Description: Computes partial sums a[0] + a[1] + ... + a[pos - 1], and
```

updates single elements a[i], taking the difference between the old and new

```
Time: Both operations are \mathcal{O}(\log N).
struct FT {
 vector<11> s:
 FT(int n) : s(n) {}
 void update(int pos, ll dif) { // a[pos] += dif
```

for (; pos < sz(s); pos |= pos + 1) s[pos] += dif; 11 query (int pos) { // sum of values in [0, pos) 11 res = 0;for (; pos > 0; pos &= pos - 1) res += s[pos-1]; return res; int lower_bound(11 sum) {// min pos st sum of [0, pos] >= sum // Returns n if no sum is >= sum, or -1 if empty sum is. if $(sum \le 0)$ return -1; int pos = 0;**for** (int pw = 1 << 25; pw; pw >>= 1) { **if** (pos + pw \le sz(s) && s[pos + pw-1] \le sum) pos += pw, sum -= s[pos-1];

FenwickTree2d.h

};

return pos;

Description: Computes sums a[i,j] for all i<I, j<J, and increases single elements a[i,j]. Requires that the elements to be updated are known in advance (call fakeUpdate() before init()).

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

```
struct FT2 {
 vector<vi> ys; vector<FT> ft;
 FT2(int limx) : ys(limx) {}
 void fakeUpdate(int x, int y) {
   for (; x < sz(ys); x \mid = 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 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.

510c32, 16 lines

return res;

Usage: RMQ rmq(values); rmg.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 < g.first) del(L++, 0);
    while (R > q.second) del(--R, 1);
    res[qi] = 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/sgrt(0)
  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();
```

LinkCutTree.h

Description: Link-cut tree with path and subtree queries. Path operations can be arbitrary, but subtree operations need to be reversible. Current implementation supports subtree addition and sum.

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

6758b4, 105 lines

```
struct Node {
 Node *p, *c[2];
 Node() { p = c[0] = c[1] = 0; }
  // Vertex data (flip is required)
 bool rev = 0;
  11 v, s, vs = 0;
  int sz, vsz = 0;
  11 d = 0, vd = 0, cc = 0;
  void add(ll x) {
   d += x, vd += x;
   v += x, s += sz * x, vs += vsz * x;
  void flip() {
    swap(c[0], c[1]), rev ^= 1;
  // BST operations
  void push() {
   if (rev) {
     rep(i, 0, 2) if (c[i]) c[i]->flip();
     rev = 0;
     rep(i, 0, 2) if (c[i]) c[i]->add(d);
     d = 0:
  void pull() {
   s = v + vs;
   sz = 1 + vsz;
   if (c[0]) s += c[0]->s, sz += c[0]->sz;
   if (c[1]) s += c[1]->s, sz += c[1]->sz;
  // Virtual operations (vd = virtual delta, cc = cancel)
  void vadd(Node* x) { // Add x, x.cc = vd
   vs += x->s;
   vsz += x->sz;
   x->cc = vd;
  void vdel(Node* x) { // Push vd - x.cc, delete x, x.cc = 0
   x->add(vd - x->cc);
   vs -= x->s;
   vsz -= x->sz;
   x->cc = 0;
  void vswap (Node* x, Node* y) { // Swap x.cc, y.cc
    swap (x->cc, y->cc);
  // Splay operations
  int up() {
   if (!p) return -2;
    rep(i, 0, 2) if (p->c[i] == this) return i;
   return -1;
  bool isRoot() { return up() < 0; }</pre>
  friend void setLink(Node* x, Node* y, int d) {
   if (y) y->p = x;
   if (d >= 0) x -> c[d] = y;
  void rot() {
    int x = up(); Node* pp = p;
    setLink(pp->p, this, pp->up());
    setLink(pp, c[x ^ 1], x); setLink(this, pp, x ^ 1);
   if (pp->p) pp->p->vswap(pp, this);
```

```
pp->pull();
 void fix() { if (!isRoot()) p->fix(); push(); }
 void splay() {
   for (fix(); !isRoot();) {
     if (p->isRoot()) rot();
     else if (up() == p->up()) p->rot(), rot();
     else rot(), rot();
   pull();
};
struct LinkCut {
 vector<Node> t;
 LinkCut(int n) : t(n) {}
 void link(int u, int v) { // link u -> v
   makeRoot(&t[v]); access(&t[u]);
    setLink(&t[v], &t[u], 0); t[v].pull();
 void cut(int u, int v) { // cut u -> v
   makeRoot(&t[u]); access(&t[v]);
   t[v].c[0] = t[u].p = 0; t[v].pull();
 bool connected(int u, int v) {
   return lca(&t[u], &t[v]);
 Node* lca(Node* u, Node* v) {
   if (u == v) return u;
   access(u); access(v); if (!u->p) return 0;
   u->splay(); return u->p ?: u;
 void access(Node* u) {
   for (Node* x = u, *y = 0; x; x = x->p) {
     x->splay();
     if (y) x->vdel(y);
     if (x->c[1]) x->vadd(x->c[1]);
     x->c[1] = y; x->pull(); y = x;
   u->splav();
 void makeRoot(Node* u) { access(u), u->flip(), u->push(); }
```

Numerical (4)

4.1 Polynomials and recurrences

```
Polynomial.h
                                                     c9b7b0, 17 lines
struct Poly {
  vector<double> a:
  double operator()(double x) const {
    double val = 0;
    for (int i = sz(a); i--;) (val *= x) += a[i];
    return val;
  void diff() {
    rep(i, 1, sz(a)) a[i-1] = i*a[i];
    a.pop_back();
  void divroot(double x0) {
    double b = a.back(), c; a.back() = 0;
    for(int i=sz(a)-1; i--;) c = a[i], a[i] = a[i+1]*x0+b, b=c;
    a.pop_back();
};
```

```
PolyRoots
```

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

```
"Polynomial.h"
vector<double> polyRoots(Poly p, double xmin, double xmax) {
 if (sz(p.a) == 2) { return {-p.a[0]/p.a[1]}; }
 vector<double> ret;
 Poly der = p_i
 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 - 1 > 1e-8)
       double m = (1 + h) / 2, f = p(m);
       if ((f \le 0) ^ sign) 1 = m;
        else h = m;
     ret.push_back((l + h) / 2);
 return ret;
```

PolyInterpolate.h

Description: Given n points (x[i], y[i]), computes an n-1-degree polynomial p that passes through them: $p(x) = a[0] * x^0 + ... + a[n-1] * x^{n-1}$. For numerical precision, pick $x[k] = c * \cos(k/(n-1) * \pi), k = 0 \dots n-1$. **Time:** $\mathcal{O}(n^2)$

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

BerlekampMassev.h

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

```
Usage: berlekampMassey({0, 1, 1, 3, 5, 11}) // {1, 2}
```

```
Time: O(N²)
".../number-theory/ModInt.h"

vector<mint> berlekampMassey(vector<mint> s) {
   int n = sz(s), L = 0, m = 0;
   vector<mint> C(n), B(n), T;
   C[0] = B[0] = 1;

mint b = 1;
   rep(i,0,n) { ++m;
   mint d = s[i];
   rep(j,1,L+1) d += C[j] * s[i - j];
   if (d == 0) continue;
   T = C; mint coef = d / b;
```

```
rep(j,m,n) C[j] -= coef * B[j - m];
if (2 * L > i) continue;
L = i + 1 - L; B = T; b = d; m = 0;
}
C.resize(L + 1); C.erase(C.begin());
for (mint& x : C) x *= -1;
return C;
```

LinearRecurrence.h

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

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

4.2 Optimization

GoldenSectionSearch.h

Description: Finds the argument minimizing the function f in the interval [a,b] assuming f is unimodal on the interval, i.e. has only one local minimum 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 TernarySearch.h in the Various chapter for a discrete version.

```
Usage: double func (double x) { return 4+x+.3*x*x; } double xmin = gss(-1000,1000,func); 
Time: \mathcal{O}(\log((b-a)/\epsilon))
```

31d45b, 14 lines

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

```
return a;
}

HillClimbing.h
Description: Poor man's optimization for unimodal functions
geeeaf, 14 lines

typedef array<double, 2> P;

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

Integrate.h
Description: Simple integration of a function over an interval using Simp-
```

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

4756fc, 7 lines

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

IntegrateAdaptive.h

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

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

template <class F>
d rec(F& f, d a, d b, d eps, d S) {
    d c = (a + b) / 2;
    d S1 = S(a, c), S2 = S(c, b), T = S1 + S2;
    if (abs(T - S) <= 15 * eps || b - a < 1e-10)
        return T + (T - S) / 15;
    return rec(f, a, c, eps / 2, S1) + rec(f, c, b, eps / 2, S2);
}
template < class F>
d quad(d a, d b, F f, d eps = 1e-8) {
    return rec(f, a, b, eps, S(a, b));
}
```

Simplex.h

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

```
Usage: vvd A = {{1,-1}, {-1,1}, {-1,-2}};
vd b = {1,1,-4}, c = {-1,-1}, x;
T val = LPSolver(A, b, c).solve(x);
```

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

```
typedef double T; // long double, Rational, double + mod<P>...
typedef vector<T> vd;
typedef vector<vd> vvd;
const T eps = 1e-8, inf = 1/.0;
#define MP make_pair
#define ltj(X) if (s == -1 \mid | MP(X[j], N[j]) < MP(X[s], N[s])) s=j
struct LPSolver {
  int m, n;
  vi N, B;
  vvd D;
  LPSolver (const vvd& A, const vd& b, const vd& c) :
    m(sz(b)), n(sz(c)), N(n+1), B(m), D(m+2), vd(n+2)) {
      rep(i, 0, m) rep(j, 0, n) D[i][j] = A[i][j];
      rep(i,0,m) { B[i] = n+i; D[i][n] = -1; D[i][n+1] = b[i]; }
      rep(j, 0, n) \{ N[j] = j; D[m][j] = -c[j]; \}
      N[n] = -1; D[m+1][n] = 1;
  void pivot(int r, int s) {
    T *a = D[r].data(), inv = 1 / a[s];
    rep(i, 0, m+2) if (i != r && abs(D[i][s]) > eps) {
      T *b = D[i].data(), inv2 = b[s] * inv;
      rep(j, 0, n+2) b[j] -= a[j] * inv2;
      b[s] = a[s] * inv2;
    rep(j,0,n+2) if (j != s) D[r][j] *= inv;
    rep(i,0,m+2) if (i != r) D[i][s] *= -inv;
    D[r][s] = inv;
    swap(B[r], N[s]);
  bool simplex(int phase) {
    int x = m + phase - 1;
    for (;;) {
      int s = -1;
      rep(j,0,n+1) if (N[j] != -phase) ltj(D[x]);
      if (D[x][s] >= -eps) return true;
      int r = -1:
      rep(i,0,m) {
        if (D[i][s] <= eps) continue;</pre>
        if (r == -1 || MP(D[i][n+1] / D[i][s], B[i])
                     < MP(D[r][n+1] / D[r][s], B[r])) r = i;
      if (r == -1) return false;
      pivot(r, s);
  T solve(vd &x) {
    int r = 0;
    rep(i,1,m) if (D[i][n+1] < D[r][n+1]) r = i;
    if (D[r][n+1] < -eps) {
      if (!simplex(2) || D[m+1][n+1] < -eps) return -inf;</pre>
      rep(i, 0, m) if (B[i] == -1) {
        int s = 0;
        rep(j,1,n+1) ltj(D[i]);
        pivot(i, s);
    bool ok = simplex(1); x = vd(n);
    rep(i,0,m) if (B[i] < n) x[B[i]] = D[i][n+1];
    return ok ? D[m][n+1] : inf;
```

```
UW
```

```
};
```

4.3 Matrices

Determinant.h

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

4583fb, 16 line

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

SolveLinear.h

Description: Solves Ax = b. If no solutions exist, returns -1. Otherwise, returns the rank of A and transforms it s.t. $\{A'_1, A'_2, \dots\}$ is a basis of the kernel of A.

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

8a34d1, 40 lines

```
const double eps = 1e-12;
template<class T>
int solveLinear(auto& A, vector<T>& b, vector<T>& 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) {
   T v, bv = 0;
   rep(r,i,n) rep(c,i,m)
     if ((v = abs(A[r][c])) > bv)
       br = r, bc = c, bv = v;
    if (bv <= eps) {
     rep(j,i,n) if (abs(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, 0, n) if (j != i) {
     T 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 - 1; i >= 0; i--) {
   b[i] /= A[i][i];
   x[col[i]] = b[i];
  vector<vector<T>> ker(m - rank, vector<T>(m));
  rep(i, rank, m) {
   ker[i - rank][col[i]] = 1;
```

```
rep(j, 0, rank) ker[i - rank][col[j]] -= A[j][i] / A[j][j];
}
return A = ker, rank;
```

SolveLinearBinary.h

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

```
typedef bitset<1000> bs;
int solveLinear(vector<bs>& A, vi& b, bs& x, int m) {
 int n = sz(A), rank = 0, br;
 assert(m \le sz(x));
 vi col(m); iota(all(col), 0);
 rep(i,0,n) {
   for (br=i; br<n; ++br) if (A[br].any()) break;</pre>
   if (br == n) {
     rep(j,i,n) if(b[j]) return -1;
     break;
    int bc = (int)A[br]._Find_next(i-1);
    swap(A[i], A[br]);
   swap(b[i], b[br]);
   swap(col[i], col[bc]);
   rep(j,0,n) if (A[j][i] != A[j][bc]) {
     A[j].flip(i); A[j].flip(bc);
   rep(j,i+1,n) if (A[j][i]) {
     b[j] ^= b[i];
     A[j] ^= A[i];
   rank++;
 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). For prime powers, repeatedly set $A^{-1} = A^{-1}(2I - AA^{-1}) \pmod{p^k}$ where A^{-1} starts as the inverse of A mod p, and k is doubled in each step.

Time: $\mathcal{O}\left(n^3\right)$ d43579, 33 lines

```
template<class T>
int matInv(vector<vector<T>>& A) {
 int n = sz(A); vi col(n);
 vector<vector<T>> tmp(n, vector<T>(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 (abs(A[j][k]) > abs(A[r][c]))
        r = j, c = k;
    if (abs(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]);
   T v = A[i][i];
    rep(j,i+1,n) {
```

```
T 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) {
  T v = A[j][i];
  rep(k,0,n) tmp[j][k] -= v*tmp[i][k];
}
rep(i,0,n) rep(j,0,n) A[col[i]][col[j]] = tmp[i][j];
return n;
```

Tridiagonal.h

Description: x = 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

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

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: $\mathcal{O}\left(N\right)$ 8f9fa8, 26 lines

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

BlackBoxDet.h

BlackBoxDet FFT FFTPoly FastFourierTransform

```
be a function s.t. f(v) = Mv. Can add multiple iterations if order of recur-
rence is less than n.
Time: \mathcal{O}\left(n^2 + 2n \text{ calls to } f\right)
"BerlekampMassey.h"
                                                          b316d2, 20 lines
mt19937_64 rng(2137);
mint det(int n, auto f) {
  auto rnd = [&]() {
    vector<mint> v(n);
    rep(i, 0, n) v[i] = rng() % mint::MOD;
    return v:
  auto a = rnd(), b = rnd(), c = rnd();
  vector<mint> s(2 * n);
  rep(i, 0, 2 * n) {
    rep(j, 0, n) s[i] += a[j] * b[j];
    rep(j, 0, n) b[j] *= c[j];
    b = f(move(b));
  auto v = berlekampMassey(s);
  if (sz(v) != n) return 0;
  mint p = 1;
  rep(i, 0, n) p *= c[i];
  return v[n - 1] / p * (n % 2 ? 1 : -1);
```

Description: Black-box algorithm for the determinant of a matrix. f should

4.4 Fourier transforms

```
FFT.h
```

```
Description: Multiply polynomials for any modulus. Works for n+m \leq 2^{24}
and c_k < 5 \cdot 10^{25}.
```

Time: $\overline{\mathcal{O}}((n+m)\log(n+m))$ "../number-theory/ModInt.h"

e993f5, 44 lines

```
template<class T>
void ntt(vector<T>& a, bool inv) {
  int n = sz(a); vector<T> b(n);
  for (int i = n / 2; i; i /= 2, swap(a, b)) {
   T w = T(T::ROOT).pow((T::MOD - 1) / n * i), m = 1;
    for (int j = 0; j < n; j += 2 * i, m *= w) rep(k, 0, i) {
     T u = a[j + k], v = a[j + k + i] * m;
     b[j / 2 + k] = u + v, b[j / 2 + k + n / 2] = u - v;
  if (inv) {
    reverse(1 + all(a));
    T z = T(n).inv(); rep(i, 0, n) a[i] *= z;
template < class T>
vector<T> conv(vector<T> a, vector<T> b) {
  int s = sz(a) + sz(b) - 1, n = 1 << __1g(2 * s - 1);
  a.resize(n); ntt(a, 0); b.resize(n); ntt(b, 0);
  rep(i, 0, n) a[i] *= b[i];
  ntt(a, 1); a.resize(s);
  return a;
template < class T>
vector<T> mconv(const auto& x, const auto& y) {
  auto con = [&] (const auto& v) {
   vector<T> w(sz(v)); rep(i, 0, sz(v)) w[i] = v[i].x;
    return w; };
  return conv(con(x), con(y));
template < class T>
vector<T> conv3(const vector<T>& a, const vector<T>& b) {
  using m0 = Mod<754974721, 11>; auto c0 = mconv<m0>(a, b);
```

```
using m1 = Mod<167772161, 3>; auto c1 = mconv<m1>(a, b);
  using m2 = Mod<469762049, 3>; auto c2 = mconv<m2>(a, b);
  int n = sz(c0); vector<T> d(n); m1 r01 = m1(m0::MOD).inv();
  m2 r02 = m2 (m0::MOD).inv(), r12 = m2 (m1::MOD).inv();
  rep(i, 0, n) {
    int x = c0[i].x, y = ((c1[i] - x) * r01).x,
        z = (((c2[i] - x) * r02 - y) * r12).x;
    d[i] = (T(z) * m1::MOD + y) * m0::MOD + x;
  return d:
FFTPolv.h
Description: Fast operations on polynomials.
Time: \mathcal{O}(n \log n), eval and interp are \mathcal{O}(n \log^2 n)
                                                     ed5af1, 117 lines
using Poly = vector<mint>;
Poly& operator+=(Poly& a, const Poly& b) {
 a.resize(max(sz(a), sz(b)));
  rep(i, 0, sz(b)) a[i] += b[i];
  return a;
Poly& operator = (Poly& a, const Poly& b) {
  a.resize (max(sz(a), sz(b)));
  rep(i, 0, sz(b)) a[i] -= b[i];
  return a;
Poly& operator *= (Poly& a, const Poly& b) {
  if (min(sz(a), sz(b)) < 50) {
    Polv c(sz(a) + sz(b) - 1);
    rep(i, 0, sz(a)) rep(j, 0, sz(b)) c[i + j] += a[i] * b[j];
    return a = c;
  return a = conv(move(a), b);
Poly operator+(Poly a, const Poly& b) { return a += b; }
Poly operator-(Poly a, const Poly& b) { return a -= b; }
Poly operator* (Poly a, const Poly& b) { return a *= b; }
Poly modK(Poly a, int k) { return a.resize(min(sz(a), k)), a; }
Poly inv(const Poly& a) { // a[0] != 0
  Poly b = \{1 / a[0]\};
  while (sz(b) < sz(a))
    b = modK(b * (Poly{2} - modK(a, 2 * sz(b)) * b), 2*sz(b));
  return modK(b, sz(a));
Poly deriv(Poly a) {
  if (!sz(a)) return {};
  rep(i, 1, sz(a)) a[i - 1] = a[i] * i;
  return a.pop_back(), a;
Poly integr(const Poly& a) {
  if (!sz(a)) return {};
  Polv b(sz(a) + 1);
  rep(i, 1, sz(b)) b[i] = a[i - 1] / i;
  return b:
Poly log(const Poly& a) { // a[0] = 1
  Poly b = integr(deriv(a) * inv(a));
  return b.resize(sz(a)), b;
Poly exp(const Poly& a) { // a[0] = 0
  Poly b = \{1\};
  if (!sz(a)) return b;
  while (sz(b) < sz(a)) {
    b.resize(sz(b) * 2);
    b \star = Poly\{1\} + modK(a, sz(b)) - log(b);
    b.resize(sz(b) / 2 + 1);
```

return modK(b, sz(a));

```
Poly pow (Poly a, 11 m) {
  int p = 0, n = sz(a);
  if (!m) { a.assign(n, 0); a[0] = 1; return a; }
  while (p < n \&\& !a[p].x) p++;
  if (p >= (n + m - 1) / m) return Poly(n);
  mint j = a[p];
  a = Poly(p + all(a)) * Poly{1 / j};
  a.resize(n);
  Poly res = exp(log(a) * Poly{m}) * Poly{j.pow(m)};
  res.insert(res.begin(), p * m, 0);
  return modK(res, n);
Poly& operator/=(Poly& a, Poly b) {
  if (sz(a) < sz(b)) return a = {};</pre>
  int s = sz(a) - sz(b) + 1;
  reverse(all(a)), reverse(all(b));
  a.resize(s), b.resize(s);
  a \star = inv(b);
  a.resize(s), reverse(all(a));
  return a:
Poly operator/(Poly a, Poly b) { return a /= b; }
Poly& operator%=(Poly& a, const Poly& b) {
  if (sz(a) < sz(b)) return a;</pre>
  return a = modK(a - (a / b) * b, sz(b) - 1);
Poly operator% (Poly a, const Poly& b) { return a %= b; }
vector<mint> eval(const Poly& a, const vector<mint>& x) {
  int n = sz(x);
  if (!n) return {};
  vector<Poly> up(2 * n);
  rep(i, 0, n) up[i + n] = Poly\{0 - x[i], 1\};
  for (int i = n - 1; i > 0; i--)
    up[i] = up[2 * i] * up[2 * i + 1];
  vector<Poly> down(2 * n);
  down[1] = a % up[1];
  rep(i, 2, 2 * n) down[i] = down[i / 2] % up[i];
  vector<mint> v(n);
  rep(i, 0, n) y[i] = down[i + n][0];
  return y;
Poly interp(vector<mint> x, vector<mint> y) {
  int n = sz(x);
  vector<Poly> up(2 * n);
  rep(i, 0, n) up[i + n] = Poly\{0 - x[i], 1\};
  for (int i = n - 1; i > 0; i--)
    up[i] = up[2 * i] * up[2 * i + 1];
  vector<mint> a = eval(deriv(up[1]), x);
  vector<Poly> down(2 * n);
  rep(i, 0, n) down[i + n] = Poly\{y[i] / a[i]\};
  for(int i = n - 1; i > 0; i--)
    down[i] = down[2*i] * up[2*i+1] + down[2*i+1] * up[2*i];
  return down[1];
Poly subsetSum(Poly a) { // a[0] = 0
  int n = sz(a);
  Poly b(n);
  rep(i, 1, n) b[i] = mint(i).inv() * (i % 2 ? 1 : -1);
  for (int i = n - 2; i > 0; i--)
    for (int j = 2; i * j < n; j++)
     a[i * j] += b[j] * a[i];
  return exp(a);
```

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: $\mathcal{O}(N \log N)$ with N = |A| + |B| (~1s for $N = 2^{22}$)

```
typedef complex<double> C;
typedef vector<double> vd;
void fft(vector<C>& a) {
  int n = sz(a), L = 31 - __builtin_clz(n);
  static vector<complex<long double>> R(2, 1);
  static vector<C> rt(2, 1); // (^ 10% faster if double)
  for (static int k = 2; k < n; k \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];
  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.emptv() || b.emptv()) return {};
  vd res(sz(a) + sz(b) - 1);
  int L = 32 - \underline{\text{builtin\_clz}(\text{sz(res)})}, n = 1 << L;
  vector<C> in(n), out(n);
  copy(all(a), begin(in));
  rep(i,0,sz(b)) in[i].imag(b[i]);
  fft(in);
  for (C& x : in) x *= x;
  rep(i, 0, n) out[i] = in[-i & (n - 1)] - conj(in[i]);
  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: $\mathcal{O}\left(N\log N\right)$, where N=|A|+|B| (twice as slow as NTT or FFT) "FastFourierTransform.h" b82773, 22 lines

```
typedef vector<ll> v1;
template<int M> vl convMod(const vl &a, const vl &b) {
  if (a.empty() || b.empty()) return {};
  vl res(sz(a) + sz(b) - 1);
  int B=32-__builtin_clz(sz(res)), n=1<<B, cut=int(sqrt(M));</pre>
  vector<C> L(n), R(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}(\text{outl}[i]) + .5) + 11(\text{real}(\text{outs}[i]) + .5);
    res[i] = ((av % M * cut + bv) % M * cut + cv) % M;
  return res;
```

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

8fd8c5, 18 lines

Number theory (5)

5.1 Modular arithmetic

ModInt.h

Description: Operators for modular arithmetic.

a902ca, 28 lines

```
template<int M, int R>
struct Mod {
 static const int MOD = M, ROOT = R;
  Mod(11 y = 0) : x(y % M) { x += (x < 0) * M; }
 Mod& operator+= (Mod o) {
    if ((x += 0.x) >= M) x -= M;
    return *this; }
  Mod& operator = (Mod o) {
    if ((x -= 0.x) < 0) x += M;
    return *this; }
  Mod& operator *= (Mod o) {
   x = 111 * x * o.x % M;
   return *this: }
  Mod& operator/=(Mod o) { return *this *= o.inv(); }
  friend Mod operator+ (Mod a, Mod b) { return a += b; }
  friend Mod operator-(Mod a, Mod b) { return a -= b; }
  friend Mod operator* (Mod a, Mod b) { return a *= b; }
  friend Mod operator/(Mod a, Mod b) { return a /= b; }
  auto operator<=>(const Mod&) const = default;
  Mod pow(ll n) const {
    Mod a = x, b = 1;
    for (; n; n /= 2, a \star= a) if (n & 1) b \star= a;
    return b:
 Mod inv() const { assert(x != 0); return pow(M - 2); }
using mint = Mod<998244353, 3>;
```

ModInverse.h

Description: Pre-computation of modular inverses. Assumes LIM \leq mod and that mod is a prime.

```
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;
```

ModLog.h

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

```
11 modLog(11 a, 11 b, 11 m) {
    11 n = (11) sqrt(m) + 1, e = 1, f = 1, j = 1;
    unordered_map<11, 11> 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;
}</pre>
```

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. 9c796e, 15 lines

```
11 sumsq(11 to) { return to / 2 * ((to-1) | 1); }
11 divsum(11 to, 11 c, 11 k, 11 m) {
    11 res = k / m * sumsq(to) + c / m * to;
    k %= m; c %= m;
    if (!k) return res;
    11 to2 = (to * k + c) / m;
    return res + (to - 1) * to2 - divsum(to2, m-1 - c, m, k);
}
11 modsum(11 to, 11 c, 11 k, 11 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 $a^b \mod c$). **Time:** $\mathcal{O}(1)$ for modmul, $\mathcal{O}(\log b)$ for modpow

02ea06, 9 lines

```
11 modmul(11 a, 11 b, 11 M) {
   return (_int128)a * b % M;
}
11 modpow(11 b, 11 e, 11 mod) {
   11 ans = 1;
   for (; e; b = modmul(b, b, mod), e /= 2)
      if (e & 1) ans = modmul(ans, b, mod);
   return ans;
```

ModSqrt.h

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

```
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}(q, 1LL \ll (r - m - 1), p);
 q = qs * qs % p;
 x = x * gs % p;
 b = b * g % p;
```

ModGen.h

Description: Finds a primitive root modulo p.

```
"Factor.h", "ModMulLL.h"
                                                         ff3110, 8 lines
mt19937 64 rng(2137);
11 modGen(11 n) {
  map<11, int> f; factor(n - 1, f); rep:
 11 g = rng() % (n - 1) + 1;
  for (auto [p, _] : f)
   if (modpow(g, (n - 1) / p, n) == 1) goto rep;
  return q;
```

ModArith.h

Description: Statistics on a mod'ed arithmetic sequence.

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

```
ll cdiv(ll x, ll y) { return x / y + ((x ^ y) > 0 && x % y); }
// \min (ax + b) % m for 0 <= x <= n
11 minRemainder(ll a, ll b, ll m, ll n) {
  assert (a >= 0 && m > 0 && b >= 0 && n >= 0);
  a \% = m, b \% = m; n = min(n, m - 1);
  if (a == 0) return b;
  if (b >= a) {
   11 \text{ ad} = \text{cdiv}(m - b, a);
   n -= ad; if (n < 0) return b;
   b += ad * a - m;
  11 q = m / a, m2 = m % a;
  if (m2 == 0) return b;
  if (b / m2 > n / q) return b - n / q * m2;
  n -= b / m2 * q; b %= m2;
  11 \ v2 = (n * a + b) / m;
  11 \times 2 = cdiv(m2 * y2 - b, a);
  if (x2 * a - m2 * y2 + b >= m2) --x2;
  return minRemainder(a, b, m2, x2);
// \min x >= 0 \text{ s.t. } 1 <= (ax + b) % m <= r
11 minBetween(ll a, ll b, ll m, ll l, ll r) {
 ll x, y, g = euclid(a, m, x, y);
  if (a > 1)
   return minBetween(a/g,b/g,m/g,1/g+(1%g>b%g),r/g-(r%g<b%g));
  if (1 > r) return -1; // no solution
  if ((x \% = m) < 0) x += m;
 11 b2 = (1 - b) * x % m;
  return minRemainder(x, b2 < 0 ? b2 + m : b2, m, r - 1);
```

5.2 Primality

FastEratosthenes.h

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

```
6<u>b2912</u>, 20 lines
const int LIM = 1e6;
bitset<LIM> isPrime:
vi eratosthenes() {
 const int S = (int)round(sqrt(LIM)), R = LIM / 2;
 vi pr = \{2\}, sieve(S+1); pr.reserve(int(LIM/log(LIM) *1.1));
  vector<pii> cp;
  for (int i = 3; i <= S; i += 2) if (!sieve[i]) {</pre>
    cp.push_back(\{i, i * i / 2\});
    for (int j = i * i; j \le S; j += 2 * i) sieve[j] = 1;
 for (int L = 1; L <= R; L += S) {
    array<bool, S> block{};
    for (auto &[p, idx] : cp)
      for (int i=idx; i < S+L; idx = (i+=p)) block[i-L] = 1;</pre>
    rep(i, 0, min(S, R - L))
      if (!block[i]) pr.push_back((L + i) * 2 + 1);
  for (int i : pr) isPrime[i] = 1;
 return pr;
```

MillerRabin.h

Description: Deterministic Miller-Rabin primality test. Guaranteed to work for numbers up to $7 \cdot 10^{18}$; for larger numbers, use Python and extend A randomly.

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

```
"ModMulLL.h"
                                                       418a8d, 12 lines
bool isPrime(ll n) {
 if (n < 2 || n % 6 % 4 != 1) return (n | 1) == 3;</pre>
 11 A[] = \{2, 325, 9375, 28178, 450775, 9780504, 1795265022\},
      s = \underline{\quad} builtin_ctzll(n-1), d = n >> s;
 for (ll a : A) { // ^ count trailing zeroes
   ll p = modpow(a%n, d, n), i = s;
    while (p != 1 && p != n - 1 && a % n && i--)
     p = modmul(p, p, n);
    if (p != n-1 && i != s) return 0;
 return 1;
```

Factor.h

Description: Pollard-rho randomized factorization algorithm.

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

```
"ModMulLL.h", "MillerRabin.h"
                                                     0750d1, 16 lines
11 pollard(ll n) {
 11 \times = 0, y = 0, t = 30, prd = 2, i = 1, q;
  auto f = [&](11 k) { return modmul(k, k, 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);
void factor(ll n, map<ll, int>& cnt) {
 if (n == 1) return;
 if (isPrime(n)) { cnt[n]++; return; }
 11 x = pollard(n);
 factor(x, cnt); factor(n / x, cnt);
```

Min25.h

Description: Fast multiplicative function prefix sums. Requires isPrime calculated up to \sqrt{n} .

```
Time: \mathcal{O}\left(\frac{n^{3/4}}{\log n}\right)
                                                                                             c0b399, 47 lines
template < class T > struct Min25 {
```

```
ll n, sq, s; vi p;
Min25(11 _n) : n(_n) {
  sq = sqrtl(n) + 5;
  while (sq * sq > n) sq--;
  11 \text{ hls} = quo(n, sq);
  while (hls != 1 && quo(n, hls - 1) == sq) hls--;
  s = hls + sa;
  rep(i, 2, sq + 1) if (isPrime[i]) p.push_back(i);
vector<T> sieve(auto f) {
  11 \text{ hls} = s - sq;
  vector<T> h(s);
  rep(i, 1, hls) h[i] = f(quo(n, i)) - 1;
  rep(i, 1, sq + 1) h[s - i] = f(i) - 1;
  for (int x : p) {
   T xt = f(x) - f(x - 1), pi = h[s - x + 1];
    11 \times 2 = 111 \times \times \times \times, imax = min(hls, quo(n, x2) + 1);
    for (11 i = 1, ix = x; i < imax; i++, ix += x)
     h[i] = ((ix < hls ? h[ix] : h[s-quo(n, ix)])-pi) * xt;
    for (int i = sq; i >= x2; i--)
     h[s - i] = (h[s - quo(i, x)] - pi) * xt;
  return h;
vector<T> unsieve(const vector<T>& fp, auto f) {
  vector<11> ns = {0};
  rep(i, 1, s - sq) ns.push_back(quo(n, i));
  rep(i, 0, sq) ns.push back(sq - i);
  auto F = fp, G = fp;
  for (ll P : p | views::reverse) {
    for (11 pk = P, k = 1; quo(n, P) >= pk; k++, pk *= P) {
      T x = fp[idx(P)], y = f(P, k, pk), z = f(P, k+1, pk*P);
      rep(i, 1, s) {
        11 m = ns[i];
        if (P * pk > m) break;
        G[i] += y * (F[idx(quo(m, pk))] - x) + z;
    copy_n(G.begin(), min((int)s, idx(P*P) + 1), F.begin());
  rep(i, 1, sz(ns)) F[i] += 1;
  return F;
11 quo(11 a, 11 b) { return (double) a / b; }
int idx(11 a) { return a <= sq ? s - a : quo(n, a); }</pre>
```

5.3 Divisibility

Euclid.h

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

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

CRT.h

```
Description: Chinese Remainder Theorem.
```

crt (a, m, b, n) computes x such that $x \equiv a \pmod{m}$, $x \equiv b \pmod{n}$. If |a| < m and |b| < n, x will obey $0 \le x < \operatorname{lcm}(m, n)$. Assumes $mn < 2^{62}$ Time: $\log(n)$

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

SameDiv.h

Description: Divides the interval $[1, \infty)$ into constant division intervals. For a significant speedup, get rid of v and do the calculations directly in the

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

24617c, 13 lines

```
vector<ll> sameFloor(ll n) {
  vector<11> v:
  for (ll i = 1; i <= n; i = n / (n / i) + 1) v.push_back(i);</pre>
  return v.push_back(n + 1), v;
vector<ll> sameCeil(ll n) {
 vector<ll> v:
  for (ll i = 1, j; i < n; i = (n + j - 2) / (j - 1)) {
   j = (n + i - 1) / i;
   v.push_back(i);
  return v.push back(n), v;
```

Fractions

ContinuedFractions.h

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

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

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

FracBinarySearch.h

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

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}

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

Fraction.h

Description: Arithmetic on rational numbers. Consider using python or doubles.

```
template<class T>
struct Frac {
 typedef Frac F;
 T p, q;
 Frac(T x = 0) : Frac(x, 1) \{ \}
 Frac(T x, T y) : p(x), q(y) {
   T g = \underline{gcd}(p, q); p /= g, q /= g;
    if (q < 0) p *= -1, q *= -1;
 friend F abs(F f) { return F(abs(f.p), f.q); }
 F operator-() const { return F(-p, q); }
 friend F operator+(F a, F b) {
   return F(a.p * b.q + b.p * a.q, a.q * b.q); }
 friend F operator-(F a, F b) {
   return F(a.p * b.q - b.p * a.q, a.q * b.q); }
  friend F operator*(F a, F b) { return F(a.p*b.p, a.q*b.q); }
 friend F operator/(F a, F b) { return F(a.p*b.q, a.q*b.p); }
 friend F& operator+=(F& a, F b) { return a = a + b; }
 friend F& operator-=(F& a, F b) { return a = a - b; }
 friend F& operator*=(F& a, F b) { return a = a * b; }
 friend F& operator/=(F& a, F b) { return a = a / b; }
 auto operator <=> (const F& f) const {
   return p * f.q - f.p * q <=> 0; }
 bool operator == (const F& f) const {
   return p == f.p && q == f.q; }
```

Pythagorean Triples

The Pythagorean triples are uniquely generated by

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

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

5.6 Primes

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

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

Estimates

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

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

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:

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

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

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

Combinatorial (6)

6.1 Permutations

6.1.1 Factorial

n	1 2 3	4	5 6	7	8	9	10
							3628800
n	11	12	13	14	15	16	17
n!	4.0e7	4.8e	8 6.2e	9 8.7e	10 1.3e	12 2.1e1	3 3.6e14
n) 171
n!	2e18	2e25	3e32	$8e47 \ 3$	8e64 9e	157 6e26	32 >dbl_max

IntPerm.h

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

Integer -> permutation can use a lookup table **Time:** $\mathcal{O}(n)$

6.1.2 Cycles

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

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

6.1.3 Derangements

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

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

6.1.4 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.

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

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

$$\left(\sum_{n=0}^{\infty} p(n)x^n\right)^{-1} = 1 + \sum_{n=1}^{\infty} (-1)^n \left(x^{n(3n+1)/2} + x^{n(3n-1)/2}\right)$$

6.2.2 Lucas' Theorem

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

6.3 General purpose numbers

6.3.1 Bernoulli numbers

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

Sums of powers:

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

Euler-Maclaurin formula for infinite sums:

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

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

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

6.3.3 Eulerian numbers

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

11

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

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

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

6.3.4 Stirling numbers of the second kind

Partitions of n distinct elements into exactly k groups.

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

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

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

6.3.5 Bell numbers

Total number of partitions of n distinct elements. B(n) = 1, 1, 2, 5, 15, 52, 203, 877, 4140, 21147, For <math>p prime,

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

6.3.6 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 : $(n-2)!/((d_1-1)!\cdots(d_n-1)!)$

6.3.7 Catalan numbers

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

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

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

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

MatroidIsect (M1 m1, M2 m2, int n) : n(n), iset(n + 1), m1(m1)

rep(i,0,n) if (m1.check(i) && m2.check(i))

iset[i] = true, m1.add(i), m2.add(i);

rep(i,0,n) if (iset[i]) ans.push_back(i);

queue<int> q({n}); // starts at dummy node

ans.push_back(b), frm[b] = a;

rep(v, 0, n) if (iset[v] && v != a) m1.add(v);

while ((c = backE(b)) >= 0) if (c == n) {

while (b != n) iset[b] ^= 1, b = frm[b];

rep(b, 0, n) if (!iset[b] && frm[b] == -1 && m1.check(b))

if ((v == b || iset[v]) && (frm[v] == -1) == cas) {

return cas ? q.push(v), frm[v] = b, v : -1;

, m2(m2) {}

while (augment());

vector<int> frm(n, -1);

auto fwdE = [&](int a) {

auto backE = [&](int b) {

m2.add(v);

while (!q.empty()) {

for (int b : fwdE(a))

return true;

rep(cas, 0, 2) rep(v, 0, n)

int a = q.front(), c; q.pop();

if (!m2.check(v))

vi solve() {

return ans;

bool augment() {

vi ans;

};

m1.clear();

return ans;

m2.clear();

return n;

return false:

};

6.3.8 LGV lemma

Let $A=(a_1,\ldots,a_n),\ B=(b_1,\ldots,b_n)$ be subsets of vertices of a DAG. By $\omega(P)$ denote a path weight, the product of edge weights in that path. Let $M_{i,j}$ be the sum of path weights over all possible paths from a_i to b_j (when unit weights, note this is the number of paths).

$$\det(M) = \sum_{(P_1, \dots, P_n) \in S_{\pi}(A, B)} \operatorname{sgn}(\pi) \prod_{i=1}^n \omega(P_i)$$

Where $S_{\pi}(A, B)$ is the set of *n*-tuples of vertex disjoint paths (including endpoints) where the *k*-th path is from a_k to $b_{\pi(k)}$. Particularly useful when only identity permutation is possible.

6.4 Other

NimProduct.h

Description: Nim product.

Time: 64² xors per multiplication.

9bba25, 16 lines

MatroidIntersection.h

Description: Given two matroids, finds the largest common independent set. Pass the matroid with more expensive add/clear operations to M1. **Time:** R^2N (M2.add + M1.check + M2.check) + R^3 M1.add + R^2 M1.clear + RN M2.clear, where R is the size of the largest independent set.

```
"../data-structures/UnionFind.h"
                                                     9812a7, 60 lines
struct ColorMat {
  vi cnt, clr;
  ColorMat(int n, vector<int> clr) : cnt(n), clr(clr) {}
 bool check(int x) { return !cnt[clr[x]]; }
  void add(int x) { cnt[clr[x]]++; }
  void clear() { fill(all(cnt), 0); }
struct GraphMat {
  UF uf;
  vector<array<int, 2>> e;
  GraphMat(int n, vector<array<int, 2>> e) : uf(n), e(e) {}
  bool check(int x) { return !uf.sameSet(e[x][0], e[x][1]); }
  void add(int x) { uf.join(e[x][0], e[x][1]); }
  void clear() { uf = UF(sz(uf.e)); }
template <class M1, class M2> struct MatroidIsect {
  vector<char> iset;
  M1 m1; M2 m2;
```

```
DeBruijnSeg.h
```

Description: Constructs a cyclic string from the alphabet [0, k) of length k^n that contains every length n string as a substring.

```
vi deBruijnSeq(int k, int n) {
   if (k == 1) return {0};
   vi seq, aux(n + 1);
   auto rec = [&] (auto f, int t, int p) -> void {
      if (t > n) {
        if (n % p == 0) rep(i, 1, p + 1) seq.push_back(aux[i]);
      } else {
        aux[t] = aux[t - p]; f(f, t + 1, p);
      while (++aux[t] < k) f(f, t + 1, t);
    }
   ;
   return rec(rec, 1, 1), seq;
}</pre>
```

Graph (7)

7.1 Network flow

Dinic.h

Description: Flow algorithm with complexity $O(VE \log U)$ where $U = \max |\text{cap}|$. $O(\min(E^{1/2}, V^{2/3})E)$ if U = 1; $O(\sqrt{V}E)$ for bipartite matching.

```
struct Dinic {
 struct Edge {
    int to, rev;
    11 c, oc;
    11 flow() { return max(oc - c, OLL); } // if you need flows
  vi lvl, ptr, q;
  vector<vector<Edge>> adi:
  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 (11 p = dfs(e.to, t, min(f, e.c))) {
          e.c -= p, adj[e.to][e.rev].c += p;
          return p;
    return 0;
 ll 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; }
```

PushRelabel.h

Time: $\mathcal{O}\left(V^2\sqrt{E}\right)$

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.

0ae1d4, 48 lines

```
struct PushRelabel {
    struct Edge {
        int dest, back;
        ll f, c;
    };
    vector<vector<Edge>> g;
    vector<ll> ec;
    vector<vector< ur;
    vector<vi> hs; vi H;
```

```
PushRelabel(int n) : g(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;
ll calc(int s, int t) {
 int v = sz(q); H[s] = v; ec[t] = 1;
 vi co(2*v); co[0] = v-1;
  rep(i,0,v) cur[i] = q[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 : q[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(q); }
```

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: $\mathcal{O}\left(FE\log(V)\right)$ where F is max flow. $\mathcal{O}\left(VE\right)$ for setpi. _{f8c012, 79 lines}

```
#include <ext/pb_ds/priority_queue.hpp>
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;
  vi seen;
  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);
    dist[s] = 0; ll di;
    __gnu_pbds::priority_queue<pair<ll, int>> q;
    vector<decltype(q)::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 });
            g.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 \text{ 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; 11 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
};
```

MinCut.h

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

GlobalMinCut.h

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

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

```
pair<int, vi> globalMinCut(vector<vi> mat) {
```

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

GomorvHu.h

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

7.2 Matching

Matching.h

Description: Fast biparite matching algorithm. Graph g should be a list of neighbors of the left partition. Returns the match for every left vertex.

```
Time: \mathcal{O}\left(E\sqrt{V}\right)
```

8b0e19, 21 lines

797692, 19 line

```
vi match(int n, int m, vector<vi>& g) {
  vi l(n, -1), r(m, -1), q(n), d(n);
  auto dfs = [&] (auto f, int u) -> bool {
    int t = exchange(d[u], 0) + 1;
    for (int v : g[u])
      if (r[v] == -1 || (d[r[v]] == t && f(f, r[v])))
        return l[u] = v, r[v] = u, 1;
    return 0;
};
for (int t = 0, f = 0;; t = f = 0, d.assign(n, 0)) {
    rep(i, 0, n) if (l[i] == -1) q[t++] = i, d[i] = 1;
    rep(i, 0, t) for (int v : g[q[i]]) {
      if (r[v] == -1) f = 1;
        else if (!d[r[v]]) d[r[v]] = d[q[i]] + 1, q[t++] = r[v];
    }
    if (!f) return l;
    rep(i, 0, n) if (l[i] == -1) dfs(dfs, i);
}
```

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.

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

WeightedMatching.h

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

```
Time: \mathcal{O}(N^2M)
                                                     b35132, 31 lines
pair<11, vi> hungarian(const vector<vi> &a) {
  if (a.emptv()) return {0, {}};
  int n = sz(a) + 1, m = sz(a[0]) + 1;
  vi p(m), ans(n - 1); vector<ll> u(n), v(m);
  rep(i,1,n) {
   p[0] = i;
    int j0 = 0; // add "dummy" worker 0
    vi pre(m, -1); vector<ll> dist(m, LLONG_MAX);
    vector<bool> done(m + 1);
    do { // dijkstra
     done[i0] = true;
      int i0 = p[j0], j1; l1 delta = LLONG_MAX;
      rep(j,1,m) if (!done[j]) {
       11 cur = a[i0 - 1][j - 1] - u[i0] - v[j];
        if (cur < dist[j]) dist[j] = cur, pre[j] = j0;
       if (dist[j] < delta) delta = dist[j], j1 = j;</pre>
      rep(j,0,m) {
       if (done[j]) u[p[j]] += delta, v[j] -= delta;
        else dist[j] -= delta;
      i0 = 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
```

Blossom.h

Description: Matching for general graphs.

Time: $\mathcal{O}(nm)$, fast in practice

4e943d, 46 lines

```
vi blossom(vector<vi>& q) {
 int n = sz(q), t = -1;
 vi m(n, -1), l(n), p(n), o(n), b(n, -1), q;
 auto lca = [&](int x, int y) {
   for (t++; ; swap(x, y)) {
     if (x == -1) continue;
     if (b[x] == t) return x;
     b[x] = t;
     x = (m[x] == -1 ? -1 : o[p[m[x]]]);
 };
 auto blossom = [&](int v, int w, int a) {
   while (o[v] != a) {
     p[v] = w; w = m[v];
     if (1[w] == 1) 1[w] = 0, q.push_back(w);
     o[v] = o[w] = a; v = p[w];
 };
 auto augment = [&](int v) {
   while (v != -1) {
     int pv = p[v], nv = m[pv];
     m[v] = pv; m[pv] = v; v = nv;
 };
 auto bfs = [&](int r) {
   fill(all(1), -1); iota(all(0), 0); q.clear();
   l[r] = 0; q.push_back(r);
   rep(i, 0, sz(q)) {
     int v = q[i];
     for (auto x : g[v]) {
       if (1[x] == -1) {
         1[x] = 1; p[x] = v;
         if (m[x] == -1) return augment (x), 1;
         l[m[x]] = 0; q.push_back(m[x]);
       } else if (1[x] == 0 && o[v] != o[x]) {
         int a = lca(o[v], o[x]);
         blossom(x, v, a); blossom(v, x, a);
   return 0;
 // Time halves if you start with (any) maximal matching.
 rep(i, 0, n) if (m[i] == -1) bfs(i);
 return m;
```

WeightedBlossom.h

Description: General max weight matching. Edge weights must be positive. **Time:** $\mathcal{O}(n^3)$, faster in practice

```
812025, 147 lines
template<int N> struct WeightedBlossom {
 struct Edge { int u, v, w; } g[N*2][N*2];
 int n,m,lab[N*2],match[N*2],slack[N*2],st[N*2];
 int par[N*2],floFrom[N*2][N],s[N*2],aux[N*2];
 vi flo[N*2]; queue<int> q;
 void init(int _n) { n = _n;
    rep(u,1,n+1) rep(v,1,n+1) g[u][v] = \{u,v,0\}; }
 void ae(int u, int v, int w) { u++, v++;
    q[u][v].w = q[v][u].w = max(q[u][v].w, w);}
 int eDelta(Edge e) {
   return lab[e.u]+lab[e.v]-g[e.u][e.v].w*2; }
 void updSlack(int u, int x) {
   if (!slack[x] || eDelta(g[u][x]) < eDelta(g[slack[x]][x]))</pre>
     slack[x] = u; }
 void setSlack(int x) {
```

```
slack[x] = 0; rep(u,1,n+1) if (q[u][x].w > 0
   && st[u] != x && s[st[u]] == 0) updSlack(u,x); }
void gPush(int x) {
  if (x <= n) q.push(x);
  else for (int t : flo[x]) qPush(t); }
void setSt(int x, int b) {
  st[x] = b; if (x > n) for (int t : flo[x]) setSt(t,b); }
int getPr(int b, int xr) {
  int pr = find(all(flo[b]),xr)-begin(flo[b]);
  if (pr&1) { reverse(1+all(flo[b])); return sz(flo[b])-pr; }
  return pr; }
void setMatch(int u, int v) {
  Edge e = g[u][v]; match[u] = e.v; if (u <= n) return;
  int xr = floFrom[u][e.u], pr = getPr(u,xr);
  rep(i, 0, pr) setMatch(flo[u][i],flo[u][i^1]);
  setMatch(xr,v); rotate(begin(flo[u]),pr+all(flo[u])); }
void augment(int u, int v) {
  while (1) {
    int xnv = st[match[u]]; setMatch(u,v);
    if (!xnv) return;
    setMatch(xnv,st[par[xnv]]);
    u = st[par[xnv]], v = xnv;
int lca(int u, int v) {
  static int t = 0;
  for (++t;u||v;swap(u,v)) {
   if (!u) continue;
    if (aux[u] == t) return u;
    aux[u] = t; u = st[match[u]];
    if (u) u = st[par[u]];
  return 0;
void addBlossom(int u, int anc, int v) {
  int b = n+1; while (b <= m && st[b]) ++b;</pre>
  if (b > m) ++m;
  lab[b] = s[b] = 0; match[b] = match[anc]; flo[b] = {anc};
  auto blossom = [&](int x) {
    for (int y; x != anc; x = st[par[y]]) {
      flo[b].push_back(x), flo[b].push_back(y=st[match[x]]);
  };
  blossom(u); reverse(1+all(flo[b])); blossom(v); setSt(b,b);
  rep(x,1,m+1) q[b][x].w = q[x][b].w = 0;
  rep(x,1,n+1) floFrom[b][x] = 0;
  for(int xs : flo[b]) {
    rep(x,1,m+1) if (q[b][x].w == 0 || eDelta(q[xs][x]) <
      eDelta(g[b][x])) g[b][x]=g[xs][x], g[x][b]=g[x][xs];
    rep(x,1,n+1) if (floFrom[xs][x]) floFrom[b][x] = xs;
  setSlack(b);
void expandBlossom(int b) {
  for (int t : flo[b]) setSt(t,t);
  int xr = floFrom[b][q[b][par[b]].u], pr = getPr(b,xr);
  for(int i = 0; i < pr; i += 2) {</pre>
    int xs = flo[b][i], xns = flo[b][i+1];
    par[xs] = g[xns][xs].u; s[xs] = 1;
    s[xns] = slack[xs] = slack[xns] = 0; qPush(xns);
  s[xr] = 1, par[xr] = par[b];
  rep(i,pr+1,sz(flo[b])) {
    int xs = flo[b][i]; s[xs] = -1, setSlack(xs); }
  st[b] = 0;
bool onFoundEdge (Edge e) {
  int u = st[e.u], v = st[e.v];
```

SCC BiconnectedComponents 2sat

```
if (s[v] == -1) {
      par[v] = e.u, s[v] = 1; slack[v] = 0;
      int nu = st[match[v]]; s[nu] = slack[nu] = 0; qPush(nu);
    } else if (s[v] == 0) {
      int anc = lca(u, v);
      if (!anc) return augment(u,v), augment(v,u),1;
     addBlossom(u,anc,v);
   return 0;
  bool matching() {
   q = queue<int>();
    rep(x, 1, m+1) {
     s[x] = -1, slack[x] = 0;
     if (st[x] == x \&\& !match[x]) par[x] = s[x] = 0, qPush(x);
    if (!sz(q)) return 0;
    while (1) {
     while (sz(q)) {
        int u=q.front(); q.pop(); if (s[st[u]] == 1) continue;
        rep(v, 1, n+1) if (q[u][v].w > 0 && st[u] != st[v]) {
          if (eDelta(g[u][v]) == 0) {
            if (onFoundEdge(g[u][v])) return 1;
          } else updSlack(u,st[v]);
      int d = INT_MAX;
      rep(b, n+1, m+1) if (st[b] == b \&\& s[b] == 1)
       d = \min(d, lab[b]/2);
      rep(x, 1, m+1) if (st[x] == x \&\& slack[x]) {
       if (s[x] == -1) d=min(d,eDelta(g[slack[x]][x]));
        else if (s[x] == 0) d=min(d,eDelta(g[slack[x]][x])/2);
      rep(u, 1, n+1) {
        if (s[st[u]] == 0) {
          if (lab[u] <= d) return 0;</pre>
          lab[u] -= d;
       } else if (s[st[u]] == 1) lab[u] += d;
      rep(b, n+1, m+1) if (st[b] == b \&\& s[b] != -1)
        lab[b] += (s[b] == 0 ? 1 : -1) *d*2;
      q = queue<int>();
      rep(x, 1, m+1) if (st[x] == x \&\& slack[x]
        && st[slack[x]] != x && eDelta(g[slack[x]][x]) == 0)
          if (onFoundEdge(q[slack[x]][x])) return 1;
      rep(b, n+1, m+1) if (st[b]==b && s[b]==1 && lab[b]==0)
        expandBlossom(b);
    return 0;
  pair<11, vi> calc() {
   m = n; st[0] = 0; rep(i, 1, 2*n+1) aux[i] = 0;
    rep(i,1,n+1) match[i] = 0, st[i] = i, flo[i].clear();
    int wMax = 0;
    rep(u, 1, n+1) rep(v, 1, n+1)
     floFrom[u][v] = (u==v ? u : 0), wMax=max(wMax,q[u][v].w);
    rep(u,1,n+1) lab[u] = wMax;
    ll w = 0; vi mt(n, -1); while (matching());
    rep(i,1,n+1) if (match[i])
     w += g[i][match[i]].w, mt[i-1] = match[i]-1;
    return {w/2,mt};
};
```

7.3 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: $sc(graph, [\&](vi\& v) \{ ... \})$ 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:** $\mathcal{O}(E+V)$

```
76b5c9, 24 lines
vi val, comp, z, cont;
int Time, ncomps;
template < class G, class F> int dfs (int j, G& q, F& f) {
 int low = val[j] = ++Time, x; z.push_back(j);
  for (auto e : q[i]) 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

Usage: int eid = 0; ed.resize(N);

st.resize(si);

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.

```
for each edge (a,b) {
ed[a].emplace_back(b, eid);
ed[b].emplace_back(a, eid++); }
bicomps([&](const vi& edgelist) {...});
Time: \mathcal{O}(E+V)
                                                      c6b7c7, 32 lines
vi num, st:
vector<vector<pii>> ed;
int Time;
template<class F>
int dfs(int at, int par, F& f) {
  int me = num[at] = ++Time, top = me;
  for (auto [y, e] : ed[at]) if (e != par) {
    if (num[y]) {
      top = min(top, num[y]);
      if (num[y] < me)</pre>
        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()));
```

```
}
else if (up < me) st.push_back(e);
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);
}
```

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 unsatisfiable. Negated variables are represented by bit-inversions $(\sim \times)$.

```
Usage: TwoSat ts(number of boolean variables); ts.either(0, \sim3); // Var 0 is true or var 3 is false ts.setValue(2); // Var 2 is true ts.atMostOne(\{0, \sim 1, 2\}); // <= 1 of vars 0, \sim1 and 2 are true ts.solve(); // Returns true iff it is solvable ts.values[0..N-1] holds the assigned values to the vars Time (2(N+E)) where N is the number of boolean variables and E is the property of the
```

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

```
struct TwoSat {
 int N:
 vector<vi> gr;
 vi values; // 0 = false, 1 = true
 TwoSat(int n = 0) : N(n), gr(2*n) {}
 int addVar() { // (optional)
   gr.emplace back();
   gr.emplace_back();
    return N++;
 void either(int f, int j) {
   f = \max(2 * f, -1 - 2 * f);
   j = \max(2*j, -1-2*j);
   gr[f].push_back(j^1);
   gr[j].push_back(f^1);
 void setValue(int x) { either(x, x); }
 void atMostOne(const vi& li) { // (optional)
   if (sz(li) <= 1) return;</pre>
   int cur = \simli[0];
   rep(i,2,sz(li)) {
     int next = addVar();
     either(cur, ~li[i]);
     either(cur, next);
     either(~li[i], next);
     cur = ~next;
   either(cur, ~li[1]);
 vi val, comp, z; int time = 0;
 int dfs(int i) {
   int low = val[i] = ++time, x; z.push_back(i);
   for(int e : gr[i]) if (!comp[e])
     low = min(low, val[e] ?: dfs(e));
    if (low == val[i]) do {
     x = z.back(); z.pop_back();
```

```
comp[x] = low;
    if (values[x>>1] == -1)
        values[x>>1] = x&1;
} while (x != i);
return val[i] = low;
}

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

EulerWalk.h

Description: Eulerian undirected/directed path/cycle algorithm. Input should be a vector of (dest, global edge index), where for undirected graphs, forward/backward edges have the same index. Returns a list of pairs (node, edge) in the Eulerian path/cycle with src at the start, or empty list if no cycle/path exists.

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

428fec, 16 lines

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

PlanarFaces.h

Description: Finds the faces of a simple planar graph and returns the vertex indices for each face in either clockwise (inner) or counterclockwise (outer) order. Disconnected graphs may have multiple outer faces and require careful handling.

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

```
"../geometry/Point.h", "../geometry/AngleCmp.h" 2c9685, 23 lines
```

```
template<class P>
vector<vi> planarFaces(vector<vi>& g, vector<P>& p) {
 int n = sz(q); P o;
 auto cmp = [&](int x,int y){return angleCmp(p[x]-o,p[y]-o);};
 vector<vi> vis(n);
  rep(i, 0, n) {
   o = p[i], sort(all(g[i]), cmp);
   vis[i].resize(sz(g[i]));
 vector<vi> f;
  rep(i, 0, n) rep(j, 0, sz(adj[i])) {
   if (vis[i][j]) continue;
   vi s; int u = i, k = j;
   while (!vis[u][k]) {
     vis[u][k] = 1; s.push_back(u);
     int v = adj[u][k]; o = p[v];
     int kk = lower_bound(all(g[v]), u, cmp) - g[v].begin();
     u = v, k = (kk + 1) % sz(adj[u]);
    f.push_back(s);
```

```
return f;
```

7.4 Coloring

EdgeColoring.h

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

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

ChromaticNumber.h

Description: Fast computation of chromatic number. Time: $\mathcal{O}(n2^n)$

const int MOD = 1034865179; // random large prime
int chromaticNumber(vi& g) {
 int n = sz(g);
 vi dp(1 << n), f(n); dp[0] = 1;
 rep(i, 1, 1 << n) {
 int ctz = __builtin_ctz(i), j = i - (1 << ctz);
 dp[i] = dp[j] + dp[j & ~g[ctz]];
 if (dp[i] >= MOD) dp[i] -= MOD;
}

rep(i, 0, 1 << n) {
 ll x = (n - __builtin_parity(i)) & 1 ? MOD - 1 : 1;
 rep(j, 0, n) {
 if ((f[j] += x) >= MOD) f[j] -= MOD;
 x = x * dp[i] % MOD;
 }
}
rep(i, 0, n) if (f[i]) return i;
return n;

7.5 Heuristics

MaximalCliques.h

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

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

b0d5b1, 12 lines

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

MaximumClique.h

b2180<u>6</u>, 19 lines

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

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

f7c0bc, 49 lines

```
typedef vector<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>
      q.push_back(R.back().i);
      vv T;
      for(auto v:R) if (e[R.back().i][v.i]) T.push_back({v.i});
      if (sz(T)) {
        if (S[lev]++ / ++pk < limit) init(T);</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 MinimumVertexCover.

7.6 Trees

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"
                                                     03139d, 46 lines
template <bool VALS_EDGES> struct HLD {
  int N, tim = 0;
  vector<vi> adi;
 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]);
};
RerootDP.h.
Description: Calculates a DP from every root in a tree.
Time: \mathcal{O}\left(\sum d \log d\right)
                                                      c0a8b6, 40 lines
struct S {
 void init(int u) {}
  void join(int u, int i, const S& c) {}
  void push(int u, int i) \{\} // i = -1 if root
vector<S> reroot(vector<vi>& g) {
 int n = sz(g), t = 1;
 vi q(n), p(n);
  for (int u : q) for (int v : g[u]) if (p[u] != v) {
   p[v] = u, q[t++] = v;
  vector<S> dp(n), rdp(n), ans(n);
  for (int i = n - 1; i >= 0; i--) {
    int u = q[i], k = -1;
    dp[u].init(u);
    rep(j, 0, sz(g[u])) {
      if (g[u][j] != p[u]) dp[u].join(u, j, dp[g[u][j]]);
    ans[u] = dp[u], dp[u].push(u, k);
  if (n == 1) return dp;
  for (int u : q) {
    int d = sz(g[u]); vector<S> e(d);
    rep(i, 0, d) e[i].init(u);
    for (int b = __lg(d); b >= 0; b--) {
      for (int i = d - 1; i \ge 0; i--) e[i] = e[i / 2];
      rep(i, 0, d - (d \& !b)) {
        S\& s = q[u][i] != p[u] ? dp[q[u][i]] : rdp[u];
        e[(i >> b) ^ 1].join(u, i, s);
    rep(i, 0, sz(g[u])) {
      if (p[u] != g[u][i]) (rdp[g[u][i]] = e[i]).push(u, i);
      else ans[u].join(u, i, rdp[u]);
    ans[u].push(u, -1);
 return ans;
```

CentroidTree.h

Description: Centroid decomposition tree. Example usage can be extended for weighted trees with a BST. **Time:** $\mathcal{O}(n \log n + q \log^2 n)$

```
dst[lev][u] = dst[lev][p] + 1;
    for (int v : q[u]) if (v!=p && cp[v]==-2) gen(v, u, lev);
 int rec(int u, int dd) {
    dfs(u, -1);
    int p = -1, s = sub[u]; rep:
    for (int v : g[u]) if (v!=p && cp[v]==-2 && sub[v]>s/2) {
     p = u, u = v; goto rep;
    sub[u] = s, d[u] = dd, cp[u] = -1;
    for (int v : q[u]) if (cp[v] == -2)
     gen(v, u, d[u]), cp[rec(v, dd + 1)] = u;
    return u:
 void path(int u, auto f) { // f(centroid, son, dist)
    for (int x = u, y = -1; x != -1; y = x, x = cp[x])
      f(x, y, dst[d[x]][u]);
};
struct ContourAdd : CT {
 vector<FT> d, c;
 ContourAdd(vector<vi>G): CT(G), d(sz(g), FT(0)), c(d) {
   rep(i, 0, sz(g)) d[i] = c[i] = FT(sub[i] + 1);
 // Add x to all vertices whose distance from p is in [1, r)
 void add(int p, int 1, int r, int x) {
   path(p, [&](int u, int v, int dd) {
     d[u].update(max(0, 1 - dd), x);
     if (r - dd < sub[u]) d[u].update(max(0, r - dd), -x);
     if (v != −1) {
       c[v].update(max(0, 1 - dd), x);
       if (r - dd < sub[u]) c[v].update(max(0, r - dd), -x);
   });
 ll get(ll p) {
   11 \text{ ans} = 0;
   path(p, [&](int u, int v, int dd) {
     ans += d[u].querv(dd + 1);
     if (v != -1) ans -= c[v].query(dd + 1);
   }):
    return ans;
};
```

7.7 Advanced

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)
```

KthWalk DominatorTree Point

```
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 \text{ 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]->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};
```

KthWalk.h

Description: Eppstein's algorithm for the k-th shortest walk in a directed graph with non-negative edge weights.

Memory: $\mathcal{O}((n+m)\log n + k)$ Time: $\mathcal{O}((n+m)\log n + k\log k)$

f7b9b0, 53 lines

```
struct KthWalk {
  using Edge = pair<int, 11>;
  struct Node { // persistent leftist heap node
   Node *1 = 0, *r = 0;
   int s; Edge e;
   Node (Edge _e) : e(_e) {}
  11 d0;
  priority_queue<pair<11, Node*>> q;
  vector<Node*> h;
  KthWalk(vector<vector<Edge>>& g, int s, int t) {
    int n = sz(q); vector<vector<Edge>> r(n);
    rep(i, 0, n) for (auto [j,w] : g[i]) r[j].push_back({i,w});
    vector<ll> d(n, LLONG_MAX);
   vi ord, p(n, -1);
   priority_queue<pair<ll, int>> pq;
   pq.push({d[t] = 0, t});
```

```
while (sz(pq)) {
     auto [dd, u] = pq.top(); pq.pop();
     if (d[u] != -dd) continue;
     ord.push_back(u);
     for (auto [v, w] : r[u]) if (d[u] + w < d[v])
       pq.push({-(d[v] = d[u] + w), v}), p[v] = u;
   if ((d0 = d[s]) == LLONG_MAX) return;
   h.resize(n);
   for (int u : ord) {
     int pp = p[u]; if (pp != -1) h[u] = h[pp];
     for (auto [v, w] : g[u]) if (d[v] != LLONG_MAX) {
       11 x = w + d[v] - d[u];
       if (x || v != pp) h[u] = merge(h[u], new Node({v, x}));
       else pp = -1;
   q.push({0, new Node({s, 0})});
 Node* merge(Node* a, Node* b) {
   if (!a || !b) return a ?: b;
   if(a->e.second > b->e.second) swap(a, b);
   Node* c = new Node(*a); c->r = merge(c->r, b);
   if (!c->1 || c->1->s < c->r->s) swap(c->1, c->r);
   c->s = (c->r ? c->r->s : 0) + 1; return c;
 11 next() { // -1 if no path
   if (!sz(q)) return -1;
   auto [d, a] = q.top(); q.pop();
   if (a->1) q.push({d - a->1->e.second + a->e.second, a->1});
   if (a->r) q.push({d - a->r->e.second + a->e.second, a->r});
   Node* t = h[a->e.first];
   if (t) q.push({d - t->e.second, t});
   return d0 - d;
};
```

DominatorTree.h

Description: Finds the parent of each vertex in the dominator tree of g. Vertex a dominates vertex b iff every path from src to b passes through a. **Time:** $\mathcal{O}(m \log n)$

```
vi dominatorTree(vector<vi>& q, int src) {
 int n = sz(q), tt = 0;
 vi ans(n, -1), t(n, -1), rt(n), s(n), p(n), d(n), b(n);
 vector < vi > c(n), r(n), sc(n);
 auto get = [&] (auto f, int u) -> int {
    if (p[u] != u) {
     int v = f(f, p[u]); p[u] = p[p[u]];
     if (s[v] < s[b[u]]) b[u] = v;
   return b[u];
 auto dfs = [&] (auto f, int u) -> void {
   t[u] = tt, rt[tt] = u;
    s[tt] = p[tt] = b[tt] = tt; tt++;
   for (int v : q[u]) {
     if (t[v] == -1) f(f, v), c[t[u]].push_back(t[v]);
     r[t[v]].push_back(t[u]);
 };
 dfs(dfs, src);
 for (int i = tt - 1; i >= 0; i--) {
    for (int j : r[i]) s[i] = min(s[i], s[get(get, j)]);
    if (i) sc[s[i]].push_back(i);
    for (int j : sc[i]) {
     int k = get(get, j);
     d[j] = s[j] == s[k] ? s[j] : k;
```

```
for (int j : c[i]) p[j] = i;
}
rep(i, 1, tt) {
  if (d[i] != s[i]) d[i] = d[d[i]];
  ans[rt[i]] = rt[d[i]];
}
return ans;
```

7.8 Math

7.8.1 Matrix tree theorem

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 $d_1 \geq \cdots \geq d_n$ exists iff $d_1 + \cdots + d_n$ is even and for every $k = 1 \dots n$,

$$\sum_{i=1}^{k} d_i \le k(k-1) + \sum_{i=k+1}^{n} \min(d_i, k).$$

7.8.3 Gale-Ryser theorem

A simple bipartite graph with degree sequences $a_1 \ge \cdots \ge a_n$ and b_1, \ldots, b_m exists iff $\sum a_i = \sum b_i$ and for every $1 \le k \le n$

$$\sum_{i=1}^{k} a_i \le \sum_{i=1}^{m} \min(b_i, k).$$

7.8.4 BEST theorem

The number of Eulerian circuits on an Eulerian graph equals

$$t(v)\prod_{u}(\deg(u)-1)!$$

where t(v) is the number of spanning trees directed towards an arbitrary root v, and deg(u) is the outdegree of vertex u.

Geometry (8)

8.1 Geometric primitives

19

Point.h

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

```
template \langle class T \rangle int sqn(T x) \{ return (x > 0) - (x < 0); \}
template<class T>
struct Point {
  typedef Point P;
 Тх, у;
  auto operator<=>(const P&) const = default;
  P operator+(P p) const { return P(x+p.x, y+p.y); }
  P operator-(P p) const { return P(x-p.x, y-p.y); }
  P operator*(T d) const { return P(x*d, y*d); }
  P operator/(T d) const { return P(x/d, y/d); }
 T dot(P p) const { return x*p.x + y*p.y; }
  T cross(P p) const { return x*p.y - y*p.x; }
 T cross(P a, P b) const { return (a-*this).cross(b-*this); }
  T dist2() const { return x*x + y*y; }
  double dist() const { return sgrt((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(-v, 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) {</pre>
    return os << "(" << p.x << "," << p.y << ")"; }
};
```

LineDist.h

Description:

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



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

SegDist.h

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

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

5c88f4, 6<u>lines</u> 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;
```

SegInter.h Description:

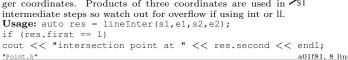
If a unique intersection point between the line segments going from s1 to e1 and from s2 to e2 exists then it is returned. If no intersection point exists an empty vector is returned. If infinitely many exist a vector with 2 elements is returned. containing the endpoints of the common line segment. The tersection 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.

```
wrong position will be returned if P is Point<|l> and the in-
Usage: vector<P> inter = segInter(s1,e1,s2,e2);
if (sz(inter) == 1)
cout << "segments intersect at " << inter[0] << endl;</pre>
"Point.h", "OnSegment.h"
template<class P> vector<P> segInter(P a, P b, P c, P d) {
  auto oa = c.cross(d, a), ob = c.cross(d, b),
       oc = a.cross(b, c), od = a.cross(b, d);
  // Checks if intersection is single non-endpoint point.
 if (sgn(oa) * sgn(ob) < 0 && sgn(oc) * sgn(od) < 0)
    return { (a * ob - b * oa) / (ob - oa) };
  set<P> s;
  if (onSegment(c, d, a)) s.insert(a);
 if (onSegment(c, d, b)) s.insert(b);
  if (onSegment(a, b, c)) s.insert(c);
  if (onSegment(a, b, d)) s.insert(d);
  return {all(s)};
```

LineInter.h

Description:

If a unique intersection point of the lines going through s1,e1 and s2.e2 exists {1, point} is returned. If no intersection point exists $\{0, (0,0)\}$ is returned and if infinitely many exists $\{-1,$ (0,0)} is returned. The wrong position will be returned if P is Point<|l> and the intersection point does not have integer coordinates. Products of three coordinates are used in \(\sigma)\) 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. $1/0/-1 \Leftrightarrow \text{left/on}$ line/right. If the optional argument eps is given 0 is returned if p is within distance eps from the line. P is supposed to be Point<T> where T is e.g. double or long long. It uses products in intermediate steps so watch out for overflow if using int or long long.

```
Usage: bool left = sideOf(p1,p2,q)==1;
"Point.h"
                                                      3af81c, 9 lines
template<class P>
int sideOf(P s, P e, P p) { return sgn(s.cross(e, p)); }
template<class P>
int sideOf(const P& s, const P& e, const P& p, double eps) {
 auto a = (e-s).cross(p-s);
 double 1 = (e-s).dist()*eps;
 return (a > 1) - (a < -1);
```

OnSegment.h

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

```
"Point.h"
                                                        c597e8, 3 lines
template<class P> bool onSegment(P s, P e, P p) {
 return p.cross(s, e) == 0 && (s - p).dot(e - p) <= 0;
```

LinearTransformation.h Description:

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



```
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();
```

LineProj.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"
```

```
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();
```

AngleCmp.h

Description: Sorts points in ascending order by angle within the interval $(-\pi, \pi]$. The point (0, 0) has an angle of 0.

```
"Point.h"
                                                                           496e9b, 6 lines
```

```
template<class P>
bool angleCmp(P a, P b) {
 auto half = [](P p) { return sqn(p.y) ?: -sqn(p.x); };
 int A = half(a), B = half(b);
 return A == B ? a.cross(b) > 0 : A < B;
```

HalfPlane.h

Description: Computes the intersection of a set of half-planes. Input is given as a set of planes, facing left. Output is the convex polygon representing the intersection. The points may have duplicates and be collinear. Will not fail catastrophically if 'eps > sqrt(2)(line intersection error)'. Likely to work for more ranges if 3 half planes are never guaranteed to intersect at the same point.

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

"Point.h", "SideOf.h", "LineInter.h"

sort(all(vs), cmp);

```
typedef Point<double> P;
typedef array<P, 2> Line;
#define sp(a) a[0], a[1]
#define ang(a) (a[1] - a[0]).angle()
int angDiff(Line a, Line b) { return sgn(ang(a) - ang(b)); }
bool cmp (Line a, Line b) {
 int s = angDiff(a, b);
 return (s ? s : sideOf(sp(a), b[0])) < 0;
vector<P> halfPlaneIntersection(vector<Line> vs) {
 const double EPS = sqrt(2) * 1e-8;
```

2bf504, 11 lines

```
vector<Line> deg(sz(vs) + 5);
vector < P > ans(sz(vs) + 5);
deq[0] = vs[0];
int ah = 0, at = 0, n = sz(vs);
rep(i,1,n+1) {
 if (i == n) vs.push_back(deq[ah]);
 if (angDiff(vs[i], vs[i - 1]) == 0) continue;
 while (ah<at && sideOf(sp(vs[i]), ans[at-1], EPS) < 0)</pre>
  while (i!=n && ah<at && sideOf(sp(vs[i]),ans[ah],EPS)<0)</pre>
  auto res = lineInter(sp(vs[i]), sp(deq[at]));
 if (res.first != 1) continue;
 ans[at++] = res.second, deg[at] = vs[i];
if (at - ah <= 2) return {};</pre>
return {ans.begin() + ah, ans.begin() + at};
```

Circles

CircleInter.h

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

```
84d6d3, 11 lines
typedef Point<double> P;
bool circleInter(P a, P b, double r1, double r2, pair < P, P > * out) {
  if (a == b) { assert(r1 != r2); return false; }
  P \text{ vec} = b - a;
  double d2 = vec.dist2(), sum = r1+r2, dif = r1-r2,
         p = (d2 + r1*r1 - r2*r2)/(d2*2), h2 = r1*r1 - p*p*d2;
  if (sum*sum < d2 || dif*dif > d2) return false;
  P mid = a + \text{vec*p}, per = \text{vec.perp}() * \text{sqrt}(\text{fmax}(0, h2) / d2);
  *out = {mid + per, mid - per};
  return true;
```

CircleTangents.h

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

```
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"
                                                                           e0cfba, 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};
CirclePolv.h
Description: Returns the area of the intersection of a circle with a ccw
```

polygon.

```
Time: \mathcal{O}(n)
"Point.h"
```

```
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, q) * r2;</pre>
    P u = p + d * s, v = p + d * t;
    return arg(p,u) * r2 + u.cross(v)/2 + arg(v,q) * r2;
  auto sum = 0.0;
  rep(i, 0, sz(ps))
    sum += tri(ps[i] - c, ps[(i + 1) % sz(ps)] - c);
  return sum;
```

Circumcircle.h

Description:

The 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.



a1ee63, 19 lines

```
"Point.h"
typedef Point<double> P;
double ccRadius (const P& A, const P& B, const P& C) {
  return (B-A).dist()*(C-B).dist()*(A-C).dist()/
      abs((B-A).cross(C-A))/2;
P ccCenter (const P& A, const P& B, const P& C) {
  P b = C-A, c = B-A;
  return A + (b*c.dist2()-c*b.dist2()).perp()/b.cross(c)/2;
```

MinEnclosingCircle.h

Description: Computes the minimum circle that encloses a set of points. **Time:** expected $\mathcal{O}(n)$

```
"Circumcircle.h"
                                                      09dd0a, 17 lines
pair<P, double> mec(vector<P> ps) {
 shuffle(all(ps), mt19937(time(0)));
 P \circ = ps[0];
 double r = 0, EPS = 1 + 1e-8;
 rep(i, 0, sz(ps)) if ((o - ps[i]).dist() > r * EPS) {
   o = ps[i], r = 0;
    rep(j,0,i) if ((o - ps[j]).dist() > r * EPS) {
     o = (ps[i] + ps[j]) / 2;
      r = (o - ps[i]).dist();
      rep(k, 0, j) if ((o - ps[k]).dist() > r * EPS) {
        o = ccCenter(ps[i], ps[j], ps[k]);
        r = (o - ps[i]).dist();
```

```
return {o, r};
```

8.3 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}\left(n\right)
```

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

PolygonArea.h

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

"Point.h" f12300, 6 lines

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

PolygonCenter.h

Description: Returns the center of mass for a polygon.

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

"Point.h"

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

PolygonCut.h

Description:

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

```
Usage: vector<P> p = ...;
p = polygonCut(p, P(0,0), P(1,0));
"Point.h", "LineInter.h"
```



9706dc, 9 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();
```

249823, 27 lines

```
UW
```

```
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>
```

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

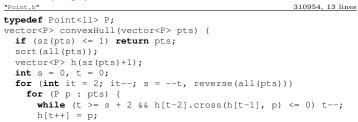
```
"Point.h", "SideOf.h"
                                                     3931c6, 33 lines
typedef Point < double > P;
double rat(P a, P b) { return sqn(b.x) ? a.x/b.x : a.y/b.y; }
double polyUnion(vector<vector<P>>& poly) {
  double ret = 0;
  rep(i, 0, sz(poly)) rep(v, 0, sz(poly[i])) {
   P A = poly[i][v], B = poly[i][(v + 1) % sz(poly[i])];
    vector<pair<double, int>> segs = {{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 : segs) s.first = min(max(s.first, 0.0), 1.0);
   double sum = 0;
   int cnt = seqs[0].second;
   rep(j,1,sz(segs)) {
     if (!cnt) sum += seqs[j].first - seqs[j - 1].first;
     cnt += seqs[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.





```
return {h.begin(), h.begin() + t - (t == 2 && h[0] == h[1])};
```

HullDiameter.h

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

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

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

PointInsideHull.h

Description: Determine whether a point t lies inside a convex hull (CCW order, with no collinear points). Returns true if point lies within the hull. If strict is true, points on the boundary aren't included.

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

"Point.h", "SideOf.h", "OnSegment.h"

```
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}(\log n)
```

```
#define cmpL(i) sqn(a.cross(polv[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};
  arrav<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;</pre>
      (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;
```

HullTangents.h

Description: Finds the left and right tangent vertices of a convex polygon relative to point a. The polygon must have at least 3 vertices, be CCW-ordered, and no collinear points. Returns the vertex indices.

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

71446b, 14 lines

```
template<class P>
pii hullTangents (const vector < P > & p, P a) {
 int n = sz(p), t[2];
 rep(it, 0, 2) {
    auto dir = [&](int i) {
      P u = p[i] - a, v = p[(i + 1) % n] - a;
      auto c = u.cross(v);
      if (c != 0) return c < 0;
      if (u.dot(v) <= 0) return true;</pre>
      return u.dist2() > v.dist2();
    auto idir = [&](int i) { return dir(i) ^ it; };
    if (idir(0) && !idir(n - 1)) { t[it] = 0; continue; }
    int s[2] = \{0, n - 1\};
    while (s[1] - s[0] > 2) {
      int mid = (s[0] + s[1]) / 2, x = idir(mid);
      if (idir(s[x ^ 1]) == (x ^ 1)) {
        s[x] = mid;
      } else {
        bool b = a.cross(p[mid], p[s[1]]) < 0;
        s[b ^ x ^ it ^ 1] = mid;
    t[it] = s[0] + 1 + (idir(s[0] + 1) == 0);
 return {t[0], t[1]};
```

8.4 Misc. Point Set Problems

ClosestPair.h

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

```
"Point.h"
```

ac41a6, 17 lines

```
typedef Point<ll> P;
```

ManhattanMST kdTree FastDelaunay

```
pair<P, P> closest(vector<P> v) {
    assert(sz(v) > 1);
    set<P> S;
    sort(all(v), [](P a, P b) { return a.y < b.y; });
    pair<11, pair<P, P>> ret{LLONG_MAX, {P(), P()}};
    int j = 0;
    for (P p : v) {
        P d{1 + (l1)sqrt(ret.first), 0};
        while (v[j].y <= p.y - d.x) S.erase(v[j++]);
        auto lo = S.lower_bound(p - d), hi = S.upper_bound(p + d);
        for (; lo != hi; ++lo)
            ret = min(ret, {(*lo - p).dist2(), {*lo, p}});
        S.insert(p);
    }
    return ret.second;
}</pre>
```

ManhattanMST.h

Description: Given N points, returns up to 4*N edges, which are guaranteed to contain a minimum spanning tree for the graph with edge weights w(p,q) = |p.x-q.x| + |p.y-q.y|. Edges are in the form (distance, src, dst). Use a standard MST algorithm on the result to find the final MST.

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

```
"Point.h"
                                                      df6f59, 23 lines
typedef Point<int> P;
vector<array<int, 3>> manhattanMST(vector<P> ps) {
  vi id(sz(ps));
  iota(all(id), 0);
  vector<array<int, 3>> edges;
  rep(k,0,4) {
    sort(all(id), [&](int i, int j) {
         return (ps[i]-ps[j]).x < (ps[j]-ps[i]).y;});</pre>
    map<int, int> sweep;
    for (int i : id) {
      for (auto it = sweep.lower_bound(-ps[i].y);
                it != sweep.end(); sweep.erase(it++)) {
        int j = it->second;
       P d = ps[i] - ps[j];
       if (d.y > d.x) break;
        edges.push_back(\{d.y + d.x, i, j\});
     sweep[-ps[i].y] = i;
    for (P& p : ps) if (k & 1) p.x = -p.x; else swap(p.x, p.y);
  return edges;
```

kdTree.h

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

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

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

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

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

```
Node (vector<P>&& vp) : pt(vp[0]) {
   for (P p : vp) {
     x0 = min(x0, p.x); x1 = max(x1, p.x);
     y0 = min(y0, p.y); y1 = max(y1, p.y);
   if (vp.size() > 1) {
     // split on x if width >= height (not ideal...)
     sort(all(vp), x1 - x0 >= y1 - y0 ? on_x : on_y);
      // divide by taking half the array for each child (not
      // best performance with many duplicates in the middle)
     int half = sz(vp)/2;
     first = new Node({vp.begin(), vp.begin() + half});
     second = new Node({vp.begin() + half, vp.end()});
 }
};
struct KDTree {
 Node* root;
 KDTree(const vector<P>& vp) : root(new Node({all(vp)})) {}
 pair<T, P> search(Node *node, const P& p) {
   if (!node->first) {
     // uncomment if we should not find the point itself:
     // if (p == node->pt) return {INF, P()};
     return make_pair((p - node->pt).dist2(), node->pt);
   Node *f = node->first, *s = node->second;
   T bfirst = f->distance(p), bsec = s->distance(p);
   if (bfirst > bsec) swap(bsec, bfirst), swap(f, s);
   // search closest side first, other side if needed
   auto best = search(f, p);
   if (bsec < best.first)</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

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

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

```
111 p2 = p.dist2(), A = a.dist2()-p2,
      B = b.dist2()-p2, C = c.dist2()-p2;
  return p.cross(a,b) \starC + p.cross(b,c) \starA + p.cross(c,a) \starB > 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<Q,Q> 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)
  Q A, B, ra, rb;
  int half = sz(s) / 2;
  tie(ra, A) = rec({all(s) - half});
  tie(B, rb) = rec(\{sz(s) - half + all(s)\});
  while ((B->p.cross(H(A)) < 0 && (A = A->next())) | 
          (A->p.cross(H(B)) > 0 && (B = B->r()->o)));
  Q base = connect(B->r(), A);
  if (A->p == ra->p) ra = base->r();
  if (B->p == rb->p) rb = base;
#define DEL(e, init, dir) O e = init->dir; if (valid(e)) \
    while (circ(e->dir->F(), H(base), e->F())) { \
      0 t = e \rightarrow dir; \setminus
      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());
      base = connect(base->r(), LC->r());
  return { ra, rb };
vector<P> triangulate(vector<P> pts) {
  sort(all(pts)); assert(unique(all(pts)) == pts.end());
  if (sz(pts) < 2) return {};
  Q e = rec(pts).first;
  vector<Q> q = \{e\};
  int qi = 0;
  while (e->o->F().cross(e->F(), e->p) < 0) e = e->o;
#define ADD { Q c = e; do { c->mark = 1; pts.push_back(c->p); \
```

```
q.push_back(c->r()); c = c->next(); } while (c != e); }
ADD; pts.clear();
while (qi < sz(q)) if (!(e = q[qi++])->mark) ADD;
return pts;
}
```

$8.5 \quad 3D$

PolyhedronVolume.h

Description: Magic formula for the volume of a polyhedron. Faces should point outwards. $$_{3058c3,\ 6$\ lines}$$

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

Point3D.h

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

8058ae, 32 lines

```
template < class T > struct Point 3D {
  typedef Point3D P;
  typedef const P& R:
  T x, y, z;
  explicit Point3D(T x=0, T y=0, T z=0) : x(x), y(y), z(z) {}
  bool operator<(R p) const {</pre>
   return tie(x, y, z) < tie(p.x, p.y, p.z); }
  bool operator==(R p) const {
   return tie(x, y, z) == tie(p.x, p.y, p.z); }
  P operator+(R p) const { return P(x+p.x, y+p.y, z+p.z); }
  P operator-(R p) const { return P(x-p.x, y-p.y, z-p.z); }
  P operator*(T d) const { return P(x*d, y*d, z*d); }
  P operator/(T d) const { return P(x/d, y/d, z/d); }
  T dot(R p) const { return x*p.x + y*p.y + z*p.z; }
  P cross(R p) const {
    return P(y*p.z - z*p.y, z*p.x - x*p.z, x*p.y - y*p.x);
  T dist2() const { return x*x + y*y + z*z; }
  double dist() const { return sqrt((double)dist2()); }
  //Azimuthal angle (longitude) to x-axis in interval [-pi, pi]
  double phi() const { return atan2(y, x); }
  //Zenith angle (latitude) to the z-axis in interval [0, pi]
  double theta() const { return atan2(sqrt(x*x+y*y),z); }
  P unit() const { return *this/(T)dist(); } //makes dist()=1
  //returns unit vector normal to *this and p
  P normal(P p) const { return cross(p).unit(); }
  //returns point rotated 'angle' radians ccw around axis
  P rotate (double angle, P axis) const {
    double s = sin(angle), c = cos(angle); P u = axis.unit();
    return u*dot(u)*(1-c) + (*this)*c - cross(u)*s;
};
```

3dHull k

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.

```
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);
 \label{eq:vector} $\operatorname{vector} \operatorname{PR}>> \ E\left(\operatorname{sz}\left(A\right), \ \operatorname{vector} \operatorname{PR}>\left(\operatorname{sz}\left(A\right), \ \left\{-1, \ -1\right\}\right)\right)$;
#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[i];
#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.g) <= 0) swap(it.c, it.b);
 return FS;
```

sphericalDistance.h

Description: Returns the shortest distance on the sphere with radius radius between the points with azimuthal angles (longitude) f1 (ϕ_1) and f2 (ϕ_2) from x axis and zenith angles (latitude) t1 (θ_1) and t2 (θ_2) from z axis (0 = north pole). All angles measured in radians. The algorithm starts by converting the spherical coordinates to cartesian coordinates so if that is what you have you can use only the two last rows. dx*radius is then the difference between the two points in the x direction and d*radius is the total distance between the points.

```
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.

Zfunc h

Description: z[i] computes the length of the longest common prefix of s[i:] and s.

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

```
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)
    l = i, r = i + z[i];
  }
  if (sz(S)) z[0] = sz(S);
  return z;
}
```

Manacher.h

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

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

e7ad79, 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][l+t]);
    int L = i-p[z][i], R = i+p[z][i]-!z;
    while (L>=1 && R+1<n && s[L-1] == s[R+1])
      p[z][i]++, L--, R++;
    if (R>r) l=L, r=R;
}
return p;
```

MinRotation.h

 $\begin{array}{lll} \textbf{Description:} \ \ \text{Finds the lexicographically smallest rotation of a string.} \\ \textbf{Usage:} \ \ \text{rotate(v.begin(), v.begin()+minRotation(v), v.end());} \\ \textbf{Time:} \ \ \mathcal{O}\left(N\right) & & & & & & & \\ \end{array}$

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

SuffixArrav.h

Description: Builds suffix array for a string. sa[i] is the starting index of the suffix which is i'th in the sorted suffix array. The returned vector is of size n + 1, and sa[0] = n. The lcp array contains longest common prefixes for neighbouring strings in the suffix array: lcp[i] = lcp(sa[i], sa[i-1]), lcp[0] = 0. The input string must not contain any zero bytes. Time: $\mathcal{O}(n \log n)$ 38db9f, 23 lines

```
struct SuffixArrav {
  vi sa, lcp;
  SuffixArray(string& s, int lim=256) { // or basic_string<int>
    int n = sz(s) + 1, k = 0, a, b;
    vi \times (all(s)+1), v(n), ws(max(n, lim)), rank(n);
    sa = lcp = y, iota(all(sa), 0);
    for (int j = 0, p = 0; p < n; j = max(1, j * 2), lim = p) {
     p = j, iota(all(y), n - j);
     rep(i,0,n) if (sa[i] >= j) y[p++] = sa[i] - j;
      fill(all(ws), 0);
     rep(i, 0, n) ws[x[i]] ++;
      rep(i, 1, lim) ws[i] += ws[i - 1];
      for (int i = n; i--;) sa[--ws[x[y[i]]]] = y[i];
     swap(x, y), p = 1, x[sa[0]] = 0;
     rep(i,1,n) a = sa[i - 1], b = sa[i], x[b] =
        (y[a] == y[b] \&\& y[a + j] == y[b + j]) ? p - 1 : p++;
    rep(i,1,n) rank[sa[i]] = i;
    for (int i = 0, j; i < n - 1; lcp[rank[i++]] = k)</pre>
     for (k \& \& k--, j = sa[rank[i] - 1];
          s[i + k] == s[j + k]; 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: $\mathcal{O}(26N)$

aae0b8, 50 lines

```
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]<=q) {
     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] - 1[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);
    return st.best;
};
```

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. $\operatorname{find}(x)$ 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.emptv());
   int n = 0;
   for (char c : s) {
     int& m = N[n].next[c - first];
     if (m == -1) { n = m = sz(N); N.emplace_back(-1); }
     else n = m;
   if (N[n].end == -1) N[n].start = j;
   backp.push_back(N[n].end);
   N[n].end = j;
   N[n].nmatches++;
 AhoCorasick(vector<string>& pat) : N(1, -1) {
   rep(i,0,sz(pat)) insert(pat[i], i);
   N[0].back = sz(N);
```

```
N.emplace_back(0);
  queue<int> q:
  for (q.push(0); !q.empty(); q.pop()) {
    int n = q.front(), prev = N[n].back;
    rep(i,0,alpha) {
      int &ed = N[n].next[i], y = N[prev].next[i];
      if (ed == -1) ed = y;
      else {
        N[ed].back = y;
        (N[ed].end == -1 ? N[ed].end : backp[N[ed].start])
          = N[y].end;
        N[ed].nmatches += N[y].nmatches;
        q.push(ed);
vi find(string word) {
  int n = 0;
  vi res; // 11 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;
```

Duval.h

Description: Constructs Lyndon factorization of s. A word is called simple iff it is strictly smaller than any of its nontrivial suffixes. The Lyndon factorization of a string is the unique division into non-increasing simple words.

Time: $\mathcal{O}(n)$ cab391, 12 lines

```
vi duval(const string& s) {
 int n = sz(s); vi f;
 for (int i = 0; i < n;) {</pre>
    int j = i + 1, k = i;
    for (; j < n && s[k] <= s[j]; j++) {</pre>
      if (s[k] < s[j]) k = i;
      else ++k:
    for (; i <= k; i += j - k) f.push_back(i);</pre>
 return f.push_back(n), f;
```

WildcardMatching.h

Description: Finds all occurrences of T in S over an alphabet with wildcards. Requires |T| < |S|.

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

```
"../numerical/FFT.h"
                                                        639ceb, 20 lines
mt19937 rng(2137);
vector<bool> match(string s, string t, char w = '*') {
 int n = sz(s), m = sz(t); mint d = rng();
```

```
vector<mint> f1(n), f2(n), f3(n), g1(m), g2(m), g3(m);
rep(i, 0, n) f1[i] = s[i] == w ? 0 : s[i] + d;
rep(i, 0, n) f2[i] = f1[i] * f1[i], f3[i] = f2[i] * f1[i];
rep(i, 0, m) gl[i] = t[i] == w ? 0 : t[i] + d;
rep(i, 0, m) g2[i] = g1[i] * g1[i], g3[i] = g2[i] * g1[i];
auto mul = [&](auto a, auto b) {
 a.resize(sz); ntt(a, 0); b.resize(sz); ntt(b, 0);
 rep(i, 0, sz) a[i] *= b[i];
 ntt(a, 1); a.erase(a.begin(), a.begin() + m - 1);
 return a:
};
auto a = mul(f1, g3), b = mul(f2, g2), c = mul(f3, g1);
vector \langle bool \rangle ans (n - m + 1);
rep(i, 0, n - m + 1) ans[i] = a[i] - b[i] - b[i] + c[i] == 0;
return ans;
```

Hash61.h

Description: Arithmetic for fast hashing modulo $2^{61} - 1$ (prime). Time: About 30% faster than naive modulo.

```
51cf65, 7 lines
const 11 MOD = (111 << 61) - 1;</pre>
11 add(11 a, 11 b) { return a+b >= MOD ? a+b - MOD : a+b; }
11 sub(11 a, 11 b) { return add(a, MOD - b); }
ll mul(ll a, ll b) {
 auto c = (\underline{\quad}int128)a * b;
  return add(c & MOD, c >> 61);
```

PalTree.h

Description: Palindrome tree. Can be used for counting number of occurrences, just add 1 to suffix link path. Replace array with map if ML is

```
Time: \mathcal{O}(26N), split is \mathcal{O}(n \log n)
```

2b16ac, 43 lines

```
const int A = 26;
struct PalTree {
  int last = 0;
  vi len = \{0, -1\}, link = \{1, 0\}, s = \{-1\};
  vector<array<int, A>> to = {{}}, {{}};
  int find(int u) {
    while (s.back() != s[sz(s) - len[u] - 2]) u = link[u];
   return u:
  int add(int x) { // x in [0, A)
    s.push_back(x); last = find(last);
    if (!to[last][x]) {
     to.push_back({});
     len.push_back(len[last] + 2);
     link.push_back(to[find(link[last])][x]);
     to[last][x] = sz(to) - 1;
    return last = to[last][x];
// dp[i] = min even/odd palindromic split of prefix of size i
const int INF = 1e9;
struct F { int e, o; };
F op (F x, F y) { return {min(x.e, y.e), min(x.o, y.o)}; }
vector<F> split(vi v) {
  PalTree t:
  vector\langle F \rangle s(2), ans(sz(v) + 1, {INF, INF});
  vi go(2), d(2); ans[0] = s[0] = s[1] = {0, INF};
  rep(i, 0, sz(v)) {
    int x = t.add(v[i]), y = t.link[x];
    if (x >= sz(qo)) {
      d.push_back(t.len[x] - t.len[y]);
      go.push_back(d[x] == d[y] ? go[y] : y);
```

```
s.push_back(ans[0]);
  for (int u = x; t.len[u] > 0; u = qo[u]) {
   s[u] = ans[i + 1 - t.len[go[u]] - d[u]];
    if (d[u] == d[t.link[u]]) s[u] = op(s[u], s[t.link[u]]);
    ans[i + 1] = op(ans[i + 1], {s[u].o + 1, s[u].e + 1});
return ans;
```

Squares.h

Description: Main-Lorentz algorithm for finding all squares in a string. Returns triples (l, r, t) signifying that for each $i \in [l, r)$ there is a square at position i of size 2t. Triples with the same t don't intersect or touch.

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

```
"Zfunc.h"
                                                       51e311, 23 lines
vector<array<int, 3>> squares(const string& s) {
 vector<array<int, 3>> ans; vi pos(sz(s) / 2 + 2, -1);
 rep(m, 1, sz(s)) {
    int p = m \& \sim (m - 1), o = m - p, e = min(m + p, sz(s));
    auto a = s.substr(o, p), b = s.substr(m, e - m);
    auto ra = a, rb = b; reverse(all(ra)); reverse(all(rb));
   rep(j, 0, 2) {
     vi z1 = Z(ra), z2 = Z(b + ' \setminus 0' + a);
     z1.push_back(0); z2.push_back(0);
     rep(c, 0, sz(a)) {
        int 1 = sz(a) - c;
        int x=c-min(1-1, z1[1]), y=c-max(1-z2[sz(b)+c+1], j);
       if (x > y) continue;
        int sb = \frac{1}{2} ? e-y-1*2 : o+x, se = \frac{1}{2} ? e-x-1*2+1 : o+y+1;
        int& k = pos[1];
        if (k != -1 \&\& ans[k][1] == sb) ans[k][1] = se;
        else k = sz(ans), ans.push_back({sb, se, 1});
      swap(a, rb); swap(b, ra);
 return ans;
```

Various (10)

10.1 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: $\mathcal{O}(\log(b-a))$ 9155b4, 11 lines

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

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

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;

while (L--) ans[L] = cur, cur = prev[cur];

FastKnapsack.h

return ans;

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

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

b20ccc, 16 lines

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

SubsetSum.h

Time: $\mathcal{O}\left(\frac{s\sqrt{s}}{32}\right)$

Description: Finds all subset sums of a multiset with sum s. cnt[k] should be the number of elements with value k.

Usage: subsetSum<MAXN>(s, cnt);

```
6aa4ae, 13 lines
template<int N>
vi subsetSum(int s, vi& cnt) {
  if (s < 3 * N / 4) return subsetSum<3 * N / 4>(s, cnt);
  bitset<N> b; b[0] = 1;
  rep(i, 1, sz(cnt)) {
    for (int j = 1; j <= cnt[i]; j *= 2)</pre>
     b \mid = b \ll (j * i), cnt[i] -= j;
    if (cnt[i]) b |= b << (cnt[i] * i);</pre>
  rep(i, 0, s + 1) if (b[i]) ans.push_back(i);
  return ans:
```

10.2 Dynamic programming

KnuthDP.h

Description: When doing DP on intervals: $a[i][j] = \min_{i < k < j} (a[i][k] + a[k][j]) + f(i,j)$, where the (minimal) optimal k increases with both i and j, one can solve intervals in increasing order of length, and search k = p[i][j] for a[i][j] only between p[i][j-1] and p[i+1][j]. This is known as Knuth DP. Sufficient criteria for this are if $f(b,c) \le f(a,d)$ and $f(a,c)+f(b,d) \le f(a,d)+f(b,c)$ for all $a \le b \le c \le d$. Consider also: LineContainer (ch. Data structures), monotone queues, ternary search. **Time:** $\mathcal{O}\left(N^2\right)$

DivideAndConquerDP.h

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

Time: $\mathcal{O}\left(\left(N+(hi-lo)\right)\log N\right)$ d38d2b, 18 lines

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

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

10.3 Optimization tricks

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;
  }
};
```

FastInput.h

 $\bf Description:$ Read an integer from stdin. Usage requires your program to pipe in input from file.

Usage: ./a.out < input.txt

Time: About 5x as fast as cin/scanf.

7b3c70, 17 lines

```
inline char gc() { // like getchar()
  static char buf[1 << 16];
  static size_t bc, be;
  if (bc >= be) {
    buf[0] = 0, bc = 0;
    be = fread(buf, 1, sizeof(buf), stdin);
  }
  return buf[bc++]; // returns 0 on EOF
}
```

```
int readInt() {
   int a, c;
   while ((a = gc()) < 40);
   if (a == '-') return -readInt();
   while ((c = gc()) >= 48) a = a * 10 + c - 480;
   return a - 48;
}

Pragmas.h
Description: GCC optimization pragmas. Should be added before bits a fines
#include <bits/allocator.h>
#pragma GCC optimize("03, unroll-loops")
#pragma GCC target("avx2, bmi, bmi2, lzcnt, popent")
```