

# University of Warsaw

# UW3

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#### 1 Contest

- 2 Mathematics
- 3 Data structures
- 4 Numerical
- 5 Number theory
- Combinatorial
- 7 Graph
- 8 Geometry
- 9 Strings
- 10 Various

#### Contest (1)

```
sol.cpp
                                                   27 lines
#include <bits/stdc++.h>
using namespace std;
#define rep(i, a, b) for (int i = (a); i < (b); i++)
#define all(x) begin(x), end(x)
#define sz(x) int((x).size())
using ll = long long;
using pii = pair<int, int>;
using vi = vector<int>;
auto operator<<(auto& o, auto x) -> decltype(x.first, o
auto operator<<(auto 6 o, auto x) -> decltype(x.end(), o | The extremum is given by x=-b/2a.
  for (int i = 0; auto y : x) o << ", " + !i++ * 2 << y
auto operator << (auto& o, auto x) -> decltype (x.first, o
  return o << "(" << x.first << ", " << x.second << ")"
void __print(auto... x) { ((cerr << x << " "), ...) <<</pre>
#define debug(x...) __print("[" #x "]:", x)
#define debug(...) 2137
#endif
int main()
 cin.tie(0)->sync_with_stdio(0);
```

#### .bashrc g++ -std=c++20 -fsanitize=address,undefined -g \ -DLOCAL -Wall -Wextra -Wshadow \$1.cpp -o \$1; nc() { g++ -std=c++20 -02 \$1.cpp -0 \$1; } alias rm='trash' alias mv='mv -i alias cp='cp -i' .vimrc

set nu et ts=2 sw=2 filetype indent on hi MatchParen ctermfg=66 ctermbg=234 cterm=underline inoremap {<cr> {<cr>}<esc>0 <bs>

#### 1 | hash.sh

3 lines # Hashes a file, ignoring all whitespace and comments. # verifying that code was correctly typed. cpp -dD -P -fpreprocessed | tr -d '[:space:]' | md5sum |

#### test.sh

for((i=1;i>0;i++)) do echo "\$i" echo "\$i" | ./gen > int diff -w <(./sol < int) <(./slow < int) || break

#### troubleshoot.txt

Czy na wejsciu pojawic sie moga long longi, np. Czy tresc jest w 100% jasna? Czy bardzo dokladnie przeczytana jest sekcja input? Daj komus innemu przeczytac tresc niezaleznie,

zadaj pytanie. Czy dziala dla brzegowych, np. n/m = 0/1, wszystkie a\_i = 0? Czy tablice za male?

Czy wszedzie modulujesz? Czy dobre modulo? Czy na wejsciu moga byc liczby poza [0, mod)? Czy zle parsujesz wejscie, np. zamiast wczytac double? WA z double nie oznacza bledu precyzji, tym bardziej z long double.

Czy napisales cos, co mogles przepisac z biblioteczki? Czy interakcja jest poprawna, np. brak znaku zapytania? Czy format wyjscia jest poprawny, np. brak YES? Czy uzywasz double gdzies, gdzie mozna tego uniknac?

# Mathematics (2)

#### 2.1 Equations

$$ax^2 + bx + c = 0 \Rightarrow x = \frac{-b \pm \sqrt{b^2 - 4ac}}{2a}$$

$$ax + by = e$$

$$cx + dy = f$$

$$x = \frac{ed - bf}{ad - bc}$$

$$y = \frac{af - ec}{ad - bc}$$

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

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

where  $A'_{i}$  is A with the i'th column replaced by

#### 2.2 Recurrences

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

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

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

#### 2.3 Trigonometry

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

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

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

where V, W are lengths of sides opposite angles

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

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

#### 2.4 Geometry

## 2.4.1 Triangles

Side lengths: a, b, c

Semiperimeter: 
$$p = \frac{a+b+c}{2}$$

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

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

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

Length of median (divides triangle into two equal-area triangles):

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

Length of bisector (divides angles in two):

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

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

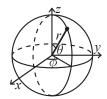
#### 2.4.2 Quadrilaterals

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

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

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

#### 2.4.3 Spherical coordinates



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

#### 2.5 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:

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

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

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

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

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

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

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

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

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

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

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

#### OrderStatisticTree HashMap SegmentTree LazySegmentTree UnionFind UnionFindRollback

#### 2.8 Probability theory

Let X be a discrete random variable with probability  $p_X(x)$  of assuming the value x. It will then have an expected value (mean)  $\mu = \mathbb{E}(X) = \sum_{x} x p_X(x)$  and variance  $\sigma^2 = V(X) = \mathbb{E}(X^2) - (\mathbb{E}(X))^2 =$  $\sum_{x} (x - \mathbb{E}(X))^2 p_X(x)$  where  $\sigma$  is the standard  $\overline{\text{deviation}}$ . If X is instead continuous it will have a probability density function  $f_X(x)$  and the sums above will instead be integrals with  $p_X(x)$  replaced by  $f_X(x)$ .

Expectation is linear:

$$\mathbb{E}(aX + bY) = a\mathbb{E}(X) + b\mathbb{E}(Y)$$

For independent X and Y.

$$V(aX + bY) = a^2V(X) + b^2V(Y).$$

#### 2.8.1 Discrete distributions Binomial distribution

The number of successes in n independent ves/no experiments, each which yields success with probability p is

Bin(n, p), n = 1, 2, ..., 0

$$p(k) = \binom{n}{k} p^k (1-p)^{n-k}$$

$$\mu = np, \, \sigma^2 = np(1-p)$$

Bin(n, p) is approximately Po(np) for small p.

#### First success distribution

The number of trials needed to get the first success in independent yes/no experiments, each which yields success with probability p is Fs(p),  $0 \le p \le 1$ .

$$p(k) = p(1-p)^{k-1}, k = 1, 2, \dots$$
  
$$\mu = \frac{1}{p}, \sigma^2 = \frac{1-p}{p^2}$$

#### Poisson distribution

The number of events occurring in a fixed period of time t if these events occur with a known average rate  $\kappa$  and independently of the time since the last event is  $Po(\lambda)$ ,  $\lambda = t\kappa$ .

$$p(k) = e^{-\lambda} \frac{\lambda^k}{k!}, k = 0, 1, 2, \dots$$
$$\mu = \lambda, \sigma^2 = \lambda$$

#### 2.8.2 Continuous distributions Uniform distribution

If the probability density function is constant between a and b and 0 elsewhere it is U(a, b), a < b.

$$f(x) = \begin{cases} \frac{1}{b-a} & a < x < b \\ 0 & \text{otherwise} \end{cases}$$
$$\mu = \frac{a+b}{2}, \ \sigma^2 = \frac{(b-a)^2}{12}$$

$$f(x) = \begin{cases} \frac{1}{b-a} & a < x \\ 0 & \text{otherw} \end{cases}$$

#### Exponential distribution

The time between events in a Poisson process is  $\text{Exp}(\lambda), \lambda > 0.$ 

$$f(x) = \begin{cases} \lambda e^{-\lambda x} & x \ge 0\\ 0 & x < 0 \end{cases}$$
$$\mu = \frac{1}{\lambda}, \sigma^2 = \frac{1}{\lambda^2}$$

#### Normal distribution

Most real random values with mean  $\mu$  and variance  $\sigma^2$  are well described by  $\mathcal{N}(\mu, \sigma^2)$ ,  $\sigma > 0$ .

$$f(x) = \frac{1}{\sqrt{2\pi\sigma^2}} e^{-\frac{(x-\mu)^2}{2\sigma^2}}$$

If  $X_1 \sim \mathcal{N}(\mu_1, \sigma_1^2)$  and  $X_2 \sim \mathcal{N}(\mu_2, \sigma_2^2)$  then

$$aX_1 + bX_2 + c \sim \mathcal{N}(\mu_1 + \mu_2 + c, a^2\sigma_1^2 + b^2\sigma_2^2)$$

#### 2.9 Markov chains

A Markov chain is a discrete random process with the property that the next state depends only on the current state. Let  $X_1, X_2, \ldots$  be a sequence of random variables generated by the Markov process. Then there is a transition matrix  $\mathbf{P} = (p_{ij})$ , with  $p_{ij} = \Pr(X_n = i | X_{n-1} = j), \text{ and } \mathbf{p}^{(n)} = \mathbf{P}^n \mathbf{p}^{(0)} \text{ is }$ the probability distribution for  $X_n$  (i.e.,  $p_i^{(n)} = \Pr(X_n = i)$ , where  $\mathbf{p}^{(0)}$  is the initial

 $\pi$  is a stationary distribution if  $\pi = \pi \mathbf{P}$ . If the Markov chain is irreducible (it is possible to get to any state from any state), then  $\pi_i = \frac{1}{\mathbb{E}(T_i)}$  where  $\mathbb{E}(T_i)$  is the expected time between two visits in state i.  $\pi_i/\pi_i$  is the expected number of visits in state j between two visits in state i.

For a connected, undirected and non-bipartite graph. where the transition probability is uniform among all neighbors,  $\pi_i$  is proportional to node i's degree.

A Markov chain is *ergodic* if the asymptotic distribution is independent of the initial distribution. A finite Markov chain is ergodic iff it is irreducible and aperiodic (i.e., the gcd of cycle lengths is 1).  $\lim_{k\to\infty} \mathbf{P}^k = \mathbf{1}\pi.$ 

A Markov chain is an A-chain if the states can be partitioned into two sets A and G, such that all states in **A** are absorbing  $(p_{ii} = 1)$ , and all states in **G** leads to an absorbing state in **A**. The probability for absorption in state  $i \in \mathbf{A}$ , when the initial state is j, is  $a_{ij} = p_{ij} + \sum_{k \in \mathbf{G}} a_{ik} p_{kj}$ . The expected time until absorption, when the initial state is i, is  $t_i = 1 + \sum_{k \in \mathbf{G}} p_{ki} t_k.$ 

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

```
b9b97b, 17 lines
#include <ext/pb_ds/assoc_container.hpp>
#include <ext/pb_ds/tree_policy.hpp>
using namespace __gnu_pbds;
template < class T>
using Tree = tree<T, null_type, less<T>, rb_tree_tag,
    tree_order_statistics_node_update>;
void example() {
   Tree<int> t, t2; t.insert(8);
 auto it = t.insert(10).first;
  assert(it == t.lower bound(9));
 assert (t.order_of_key(10) == 1);
  assert(t.order_of_key(11) == 2);
  assert(*t.find_by_order(0) == 8);
 t.join(t2); // assuming T < T2 or T > T2, merge t2
```

#### 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>
// To use most bits rather than just the lowest ones:
struct chash { // large odd number for C
 const uint64_t C = 11(4e18 * acos(0)) | 71;
 11 operator()(11 x) const { return __builtin_bswap64(
__gnu_pbds::gp_hash_table<11,11,chash> h({},{},{},{},{
     1<<16});
```

#### Segment Tree.h

Description: Zero-indexed max-tree. Bounds are inclusive to the left and exclusive to the right. Can be changed by modifving T, f and unit.

```
Time: \mathcal{O}(\log N)
                                                 0f4bdb, 19 lines
struct Tree {
  typedef int T:
  static constexpr T unit = INT_MIN;
  T f(T a, T b) { return max(a, b); } // (any
        associative fn)
  vector<T> s; int n;
 Tree(int n = 0, T def = unit) : s(2*n, def), n(n) {} void update(int pos, T val) {
    for (s[pos += n] = val; pos /= 2;)
       s[pos] = f(s[pos * 2], s[pos * 2 + 1]);
    query(int b, int e) { // query [b, e)
    T ra = unit, rb = unit;
    for (b += n, e += n; b < e; b /= 2, e /= 2) {
   if (b % 2) ra = f(ra, s[b++]);</pre>
      if (e % 2) rb = f(s[--e], rb);
    return f(ra, rb);
```

#### LazySegmentTree.h

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

```
Usage: Node* tr = new Node(v, 0, sz(v));
Time: \mathcal{O}(\log N).
```

```
34ecf5, 50 lines
"../various/BumpAllocator.h"
const int inf = 1e9;
struct Node {
 Node *1 = 0, *r = 0;
 int lo, hi, mset = inf, madd = 0, val = -inf;
 Node (int lo, int hi):lo(lo), hi(hi) {} // Large interval
 Node (vi& v, int lo, int hi) : lo(lo), hi(hi) {
    if (lo + 1 < hi) {
      int mid = lo + (hi - lo)/2;
      l = new Node(v, lo, mid); r = new Node(v, mid, hi
      val = max(1->val, r->val);
    else val = v[lo];
```

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

#### UnionFind.h

Description: Disjoint-set data structure.

```
Time: \mathcal{O}(\alpha(N))
                                            7aa27c, 14 lines
struct UF {
 UF(int n) : e(n, -1) {}
 bool sameSet(int a, int b) { return find(a) == find(b
  int size(int x) { return -e[find(x)]; }
  int find(int x) { return e[x] < 0 ? x : e[x] = find(e
        [x]); ]
  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);
    e[a] += e[b]; e[b] = a;
    return true;
```

#### UnionFindRollback.h

Description: Disjoint-set data structure with undo. If undo is not needed, skip st, time() and rollback().

```
Usage: int t = uf.time(); ...; uf.rollback(t);
Time: \mathcal{O}(\log(N))
                                             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;
};
```

#### DequeRollback.h

Description: Deque-like undoing on data structures with amortized O(log n) overhead for operations. Maintains a deque of objects alongside a data structure that contains all of them. The data structure only needs to support insertions and undoing of last insertion using the following interface: - insert(...) - insert an object to DS - time() - returns current version number - rollback(t) - undo all operations after t Assumes time() == 0 for empty DS. 57bab5, 38 lines

```
struct DequeUndo {
  // Argument for insert(...) method of DS.
  using T = tuple<int, int>;
  DataStructure ds; // Configure DS type here.
  vector<T> elems[2];
 vector<pii> his = {{0,0}};
// Push object to front or back of deque, depending
  on side arg.
void push(T val, bool side) {
    elems[side].push_back(val);
    doPush(0, side);
  ,
// Pop object from front or back of deque, depending
        on side arg.
  void pop(int side)
   auto &A = elems[side], &B = elems[!side];
int cnt[2] = {};
    if (A.emptv()) {
     assert(!B.empty());
      auto it = B.begin() + sz(B)/2 + 1;
     A.assign(B.begin(), it);
     B.erase(B.begin(), it);
     reverse(all(A)); his.resize(1);
      cnt[0] = sz(A); cnt[1] = sz(B);
    } else{
      do {
        cnt[his.back().y ^ side]++;
        his.pop_back();
      } while (cnt[0]*2 < cnt[1] && cnt[0] < sz(A));</pre>
    cnt[0]--; A.pop_back();
    ds.rollback(his.back().x);
    for (int i : {1, 0})
      while (cnt[i]) doPush(--cnt[i], i^side);
  void doPush (int i, bool s) {
    apply([&](auto... x) { ds.insert(x...); },elems[s].
         rbegin()[i]);
    his.push_back({ds.time(), s});
```

#### SubMatrix.h

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

Usage: SubMatrix<int> m(matrix);

m.sum(0, 0, 2, 2); // top left 4 elements

Time:  $\mathcal{O}\left(N^2+Q\right)$ 

```
c59ada, 13 lines
```

#### Matrix.h

Description: Basic operations on square matrices. Usage: Matrix<int, 3> A;

```
A.d = {{{{1,2,3}}, {{4,5,6}}, {{7,8,9}}}};
array<int, 3> vec = {1,2,3};
vec = (A^N) * vec; 6ab5db, 26 lines
template<class T. int N> struct Matrix {
```

```
template class T, int N> struct Matrix {
    typedef Matrix M;
    array < array < 7, N>, N> d{};
    M operator*(const M& m) const {
        M a;
        rep(i,0,N) rep(j,0,N)
        rep(k,0,N) a.d[i][j] += d[i][k]*m.d[k][j];
```

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

#### LineContainer.h

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

```
Time: \mathcal{O}\left(\log N\right)
                                                        8ec1c7, 30 lines
struct Line {
  mutable ll k, m, p;
  bool operator<(const Line& o) const { return k < o.k;
  bool operator<(11 x) const { return p < x; }</pre>
struct LineContainer : multiset<Line, less<>>> {
   // (for doubles, use inf = 1/.0, div(a,b) = a/b)
 // (for doubles, use inr = 1/.0, aiv(a, b) -
static const ll inf = LLONG_MAX;
ll div(ll a, ll b) { // floored division
return a / b - ((a ^ b) < 0 && a & b); }
bool isect(iterator x, iterator y) {
     if (y == end()) return x->p = inf, 0;
    if (x-)k == y-)k  x-)p = x-)m > y-)m ? inf : -inf; else x-)p = div(y-)m - x-)m, x-)k - y-)k;
     return x->p >= y->p;
  void add(ll k, ll m) {
     auto z = insert(\{k, m, 0\}), y = z++, x = y;
     while (isect(y, z)) z = erase(z);
     if (x != begin() && isect(--x, y)) isect(x, y =
     erase(y));

while ((y = x) != begin() && (--x)->p >= y->p)
       isect(x, erase(y));
  ll query(ll x) {
     assert(!empty());
     auto 1 = *lower_bound(x);
     return l.k * x + l.m;
```

#### Treap.h

**Description:** A short self-balancing tree. It acts as a sequential container with log-time splits/joins, and is easy to augment with additional data.

```
Time: O(\log N)
                                              468a0f, 78 lines
mt19937 rng(2137);
struct Node (
 Node *1 = 0, *r = 0, *p = 0;
 int val, y, c = 1;
Node(int val) : val(val), y(rng()) {}
friend int cnt(Node* n) { return n ? n->c : 0; }
 void recalc() { c = cnt(l) + cnt(r) + 1; }
 void push() {}
void each (Node* n, auto f) {
 if (n) { each(n->1, f); f(n->val); each(n->r, f); }
pair<Node*, Node*> split(Node* n, int k) {
 if (!n) return {};
 n->push(); n->p = 0;
 if (cnt(n->1) >= k) { // "n->val >= k" for
       lower_bound(k)
    auto [L,R] = split(n->1, k);
    if (n->1) n->1->p = n;
    return n->recalc(), pair(L, n);
    auto [L,R] = split(n->r,k - cnt(n->l) - 1); // and
         just "k"
    if (n->r) n->r->p = n;
```

```
return n->recalc(), pair(n, R);
Node* merge(Node* 1, Node* r) {
 if (!l || !r) return l ?: r;
 if (1->y > r->y) {
    1->push():
    1->r = merge(1->r, r);
   if (1->r) 1->r->p = 1;
    return 1->recalc(), 1;
 } else {
    r->push();
    r->\bar{1} = merge(1, r->1);
    if (r->1) r->1->p = r;
    return r->recalc(), r;
Node* ins(Node* t, Node* n, int pos) {
 auto [l,r] = split(t, pos);
 return merge (merge (1, n), r);
 // Union of two sorted treaps, O(m log(n/m)) where m<=n
 / Makes small-to-large O(n log n) instead of log^2.
  Requires lower_bound split (not the default one).
Node* unite(Node* a, Node* b) {
 if (!a || !b) return a ?: b;
 if (a->y < b->y) swap(a, b);
 auto [1, r] = split(b, a->val); // lower_bound split
 return merge (unite(1, a->1), merge(a, unite(r, a->r)));
// Number of elements before n. If there are range
// reverse queries, recursively push the path to n.
int idx(Node* n) {
 int c = cnt(n->1);
 while (n->p) {
   if (n->p->l != n) c += cnt(n->p->l) + 1;
   n = n - > p:
 return c:
// Example application: move the range [1, r) to index
void move (Node * & t, int 1, int r, int k) {
 Node *a, *b, *c;
 tie(a,b) = split(t, 1); tie(b,c) = split(b, r - 1);
 if (k <= 1) t = merge(ins(a, b, k), c);
 else t = merge(a, ins(c, b, k - r));
```

#### IntSet.h

Description: bitset with fast predecessor and successor queries. Assumes x86 shift overflows. Extremely fast (50-200mln operations in 1 second). 85cd6f, 32 lines

```
template<int N>
struct IntSet {
 static constexpr int B = 64:
 uint64_t V[N / B + 1] = {};
IntSet<(N < B + 1 ? 0 : N / B + 1) > up;
  bool has(int i) { return (V[i / B] >> i) & 1; }
  void add(int i) {
    if (!V[i / B]) up.add(i / B);
V[i / B] |= 1ull << i;</pre>
  void del(int i) {
    if (!(V[i / B] &= ~(1ull << i))) up.del(i / B);
  int next(int i) { // j > i such that j inside or -1
    auto x = V[i / B] >> i;
    if (x &= ~1) return i + _builtin_ctzll(x);
return (i = up.next(i / B)) < 0 ? i :</pre>
       i * B + __builtin_ctzll(V[i]);
  int prev(int i) { // j < i such that j inside or -1
  auto x = V[i / B] << (B - i - 1);</pre>
    if (x &= INT64 MAX)
      return i-__builtin_clzll(x);
    return (i = up.prev(i / B)) < 0 ? i :
       i * B + B - 1 - __builtin_clzll(V[i]);
template<>
struct IntSet<0> {
  void add(int) {} void del(int) {}
  int next(int) { return -1; }
 int prev(int) { return -1; } };
```

#### FenwickTree.h

**Description:** Computes partial sums a[0] + a[1] + ... + a[pos - 1], and updates single elements a[i], taking the difference between the old and new value.

```
Time: Both operations are O(\log N).
                                           e62fac, 22 lines
struct FT {
 vector<ll> s:
 FT(int n) : s(n) {}
 void update(int pos, ll dif) { // a[pos] += dif
    for (; pos < sz(s); pos |= pos + 1) s[pos] += dif;
 ll query(int pos) { // sum of values in [0, pos)
    ll res = 0;
    for (; pos > 0; pos &= pos - 1) res += s[pos-1];
    return res;
  int lower_bound(ll sum) {// min pos st sum of [0, pos
    // Returns n if no sum is >= sum, or -1 if empty
         sum is.
    if (sum <= 0) return -1;</pre>
    int pos = 0;
    for (int pw = 1 << 25; pw; pw >>= 1) {
      if (pos + pw <= sz(s) && s[pos + pw-1] < sum)
        pos += pw, sum -= s[pos-1];
    return pos;
```

#### FenwickTree2d.h

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

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

"FenwickTree.h"

157f07, 22 lines

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

#### RMQ.h

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

rma.auerv(inclusive, exclusive);

**Time:**  $O(|V| \log |V| + Q)$  510c32, 16 lines

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

## 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();

#### PolvRoots.h

```
Description: Finds the real roots to a polynomial.
Usage:
                polyRoots(\{\{2,-3,1\}\},-1e9,1e9) // solve
x^2-3x+2 =
Time: \mathcal{O}\left(n^2\log(1/\epsilon)\right)
"Polynomial.h"
                                             b00bfe, 23 lines
vector<double> polyRoots(Poly p, double xmin, double
      xmax) {
  if (sz(p.a) == 2) { return {-p.a[0]/p.a[1]}; }
  vector<double> ret;
  Poly der = p_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 1 = 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 <= 0) ^ sign) l = m;
        else h = m;
      ret.push back((1 + h) / 2);
 return ret:
```

#### | PolvInterpolate.h

**Description:** 1. Interpolate set of points (i, vec[i]) and return it evaluated at x; 2. Given n points (x, f(x)) compute n-1-degree polynomial f that passes through them;

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

174038, 33 lines

```
template<class T>
 polyExtend(vector<T>& vec, T x) {
 int n = sz(vec);
 vector<T> fac(n, 1), suf(n, 1);
  rep(i, 1, n) fac[i] = fac[i-1] * i;
 for (int i=n; --i;) suf[i-1] = suf[i]*(x-i);
 T pref = 1, ret = 0;
   T d = fac[i] * fac[n-i-1] * ((n-i)%2*2-1);
   ret += vec[i] * suf[i] * pref / d;
   pref *= x-i;
 return ret:
template<class T>
vector<T> polyInterp(vector<pair<T, T>> P) {
 int n = sz(P);
 vector<T> ret(n), tmp(n);
 T last = 0;
 tmp[0] = 1;
 rep(k, 0, n-1) rep(i, k+1, n)
   P[i].second = (P[i].second-P[k].second) / (P[i].
         first-P[k].first);
 rep(k, 0, n) rep(i, 0, n) {
   ret[i] += P[k].second * tmp[i];
   swap(last, tmp[i]);
   tmp[i] -= last * P[k].first;
 return ret;
```

#### BerlekampMassev.h

ll d = s[i] % mod;

if (!d) continue;

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

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

Time:  $\mathcal{O}\left(N^2\right)$ ".../number-theory/ModPow.h" 96548b, 20 lines vector<1l> berlekampMassey(vector<1l> s) {
 int n = sz(s), L = 0, m = 0;
 vector<1l> c(n), B(n), T;
 C[0] = B[0] = 1;
 ll b = 1;
 rep(i,0,n) { ++m;

rep(j,1,L+1) d = (d + C[j] \* s[i - j]) % mod;

#### 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) f4e444, 26 lines
```

```
typedef vector<ll> Poly;
ll linearRec(Poly S, Poly tr, ll k) {
 int n = sz(tr):
 auto combine = [&] (Poly a, Poly b) {
Poly res(n * 2 + 1);
    rep(i,0,n+1) rep(j,0,n+1)
    res[i + j] = (res[i + j] + a[i] * b[j]) % mod;

for (int i = 2 * n; i > n; --i) rep(j,0,n)

res[i - 1 - j] = (res[i - 1 - j] + res[i] * tr[j]
             ]) % mod;
    res.resize(n + 1);
    return res:
  Polv 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 res = 0:
 rep(i, 0, n) res = (res + pol[i + 1] * S[i]) % mod;
 return res:
```

#### FastMulDet.h

**Description:** Given a matrix M, s.t. we can quickly compute f(v) = Mv for any vector v, computes  $\det(M)$ . Single iteration fails on identity matrix with probability around  $n^2/mod$ . For small mod you can modify this to use a field extension. **Time:** 4n calls to f

```
"BerlekampMassey.h"
                                          84262e, 30 lines
mt19937 64 rnd{2137};
vector<ll> rndVec(int n) {
 vector<ll> r(n);
 rep(i, 0, n) r[i] = rnd() % mod;
 return r;
ll dot(vector<ll> &a, vector<ll> &b) {
 11 r = 0;
 rep(i, 0, sz(a)) r += a[i] * b[i] % mod;
 return r % mod:
void pointwise (vector<ll> &a, vector<ll> &b) {
 rep(i, 0, sz(a)) a[i] = a[i] * b[i] % mod;
ll detOnce(int n, auto f) {
 auto v = rndVec(n), r = rndVec(n), a = rndVec(n);
 vector<ll> vals:
 rep(, 0, n*2) {
   pointwise(a, r);
   vals.push_back(dot(v, a = f(a)));
 auto ber = berlekampMassey(vals);
 if (sz(ber) != n) return 0;
 ll prod = 1:
 for (11 x : r) prod = prod * x % mod;
 int sq = n % 2 ? 1 : -1;
 return (mod + ber[n-1] * sg) * modpow(prod, mod-2) %
      mod:
il det(int n, auto f) {
 return detOnce(n, f) ?: detOnce(n, f); }
```

## 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 Ternary-
Search.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}\left(\log((b-a)/\epsilon)\right)
                                                 31d45b, 14 lines
double qss(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);
```

#### HillClimbing.h

return a;

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

8eeeaf, 14 lines

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

#### Integrate.h

**Description:** Simple integration of a function over an interval using Simpson's rule. The error should be proportional to  $h^4$ , although in practice you will want to verify that the result is stable to desired precision when epsilon changes 47566c. 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.

d quad(d a, d b, F f, d eps = 1e-8) {

return rec(f, a, b, eps, S(a, b));

#### Simplex.h

mod<P>...

typedef vector<T> vd;

typedef vector<vd> vvd;

const T eps = 1e-8, inf = 1/.0;

UW

Description: Solves a general linear maximization problem: maximize  $c^T x$  subject to  $Ax \leq b, x \geq 0$ . Returns -inf if there is no solution, inf if there are arbitrarily good solutions, or the maximum value of  $c^T x$  otherwise. The input vector is set to an optimal x (or in the unbounded case, an arbitrary solution fulfilling the constraints). Numerical stability is not guaranteed. For better performance, define variables such that x = 0is 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. aa8530, 68 lines typedef double T; // long double, Rational, double +

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

#### 4.3 Matrices

#### Determinant.h

Description: Calculates determinant of a matrix. Destroys the matrix, Time:  $\mathcal{O}\left(N^3\right)$ 4583fb, 16 lines

```
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;
```

#### IntDeterminant.h

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

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

3313dc, 18 lines

```
const 11 mod = 12345;
ll det(vector<vector<ll>>& a) {
 int n = sz(a); ll ans = 1;
 rep(i,0,n) {
    rep(j,i+1,n)
      while (a[j][i] != 0) { // gcd step
ll t = a[i][i] / a[j][i];
        if (t) rep(k,i,n)
         a[i][k] = (a[i][k] - a[j][k] * t) % mod;
        swap(a[i], a[j]);
        ans \star = -1;
    ans = ans * a[i][i] % mod;
   if (!ans) return 0;
 return (ans + mod) % mod;
```

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

4f0aa8, 41 lines

```
const double eps = 1e-12;
template<class T>
int solveLinear(vector<vector<T>>& 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: O(n^2m)
```

fa2d7a, 34 lines

```
typedef bitset<1000> bs;
int solveLinear(vector<bs>& A, vi& b, bs& x, int m) {
 int n = sz(A), rank = 0, br;
 assert (m \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;
    int bc = (int)A[br]._Find_next(i-1);
   swap(A[i], A[br]);
   swap(b[i], b[br]);
   swap(col[i], col[bc]);
rep(j,0,n) if (A[j][i] != A[j][bc]) {
     A[j].flip(i); A[j].flip(bc);
    rep(j,i+1,n) if (A[j][i]) {
     b[j] ^= b[i];
A[j] ^= A[i];
    rank++:
 x = bs():
 for (int i = rank; i--;) {
   if (!b[i]) continue;
x[col[i]] = 1;
   rep(j,0,i) b[j] ^= A[j][i];
 return rank; // (multiple solutions if rank < m)
```

#### MatrixInverse.h

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

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

d43579, 36 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;
    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;
      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];
```

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

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

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 ob-

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

Fails if the solution is not unique.

If  $|d_i| > |p_i| + |q_{i-1}|$  for all i, or  $|d_i| > |p_{i-1}| + |q_i|$ , or the matrix is positive definite, the algorithm is numerically stable and neither tr nor the check for diag[i] == 0 is needed.

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

```
typedef double T;
vector<T> tridiagonal (vector<T> diag, const vector<T>&
    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[
        il == 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;
```

#### 4.4 Fourier transforms

#### FastFourierTransform.h

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

Time:  $\mathcal{O}(N \log N)$  with  $N = |A| + |B| (\sim 1s \text{ for } 25)$ 

```
typedef complex<double> C:
typedef vector<double> vd;
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
  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-
```

```
a[i + j + k] = a[i + j] - z;
     a[i + j] += z;
vd conv(const vd& a, const vd& b) {
 if (a.empty() || b.empty()) return {};
 vd res(sz(a) + sz(b) - 1);
 int L = 32 - __builtin_clz(sz(res)), n = 1 << L;
vector<C> in(n), out(n);
  copy(all(a), begin(in));
  rep(i,0,sz(b)) in[i].imag(b[i]);
 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);
```

#### 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:  $O(N \log N)$ , 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</pre>
      (M));
  vector<C> L(n), R(n), outs(n), outl(n);
 rep(i,0,sz(a)) L[i] = C((int)a[i] / cut, (int)a[i] %
 rep(i,0,sz(b)) R[i] = C((int)b[i] / cut, (int)b[i] %
      cut);
 fft(L), fft(R);
 rep(i,0,n) {
   fft(outl), fft(outs);
 rep(i,0,sz(res)) {
   ll av = ll(real(outl[i])+.5), cv = ll(imag(outs[i])
       +.5);
   11 bv = 11(imag(outl[i])+.5) + 11(real(outs[i])+.5)
   res[i] = ((av % M * cut + bv) % M * cut + cv) % M;
 return res;
```

#### NumberTheoreticTransform.h

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

```
"../number-theory/ModPow.h"
const 11 mod = (119 << 23) + 1, root = 62; // =</pre>
     998244353
// For p < 2^30 there is also e.g. 5 << 25, 7 << 26,
     479 << 21
// and 483 << 21 (same root). The last two are > 10^9.
typedef vector<ll> vl;
void ntt(v1 &a) {
  int n = sz(a), L = 31 - __builtin_clz(n);
  static v1 rt(2, 1);
for (static int k = 2, s = 2; k < n; k *= 2, s++) {
   rt.resize(n);
    11 z[] = {1, modpow(root, mod >> s)};
    rep(i,k,2*k) rt[i] = rt[i / 2] * z[i & 1] % mod;
  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) {
        1 z = rt[j + k] * a[i + j + k] * mod, &ai = a[i</pre>
      + j];
a[i + j + k] = ai - z + (z > ai ? mod : 0);
ai += (ai + z >= mod ? z - mod : z);
vl conv(const vl &a, const vl &b) {
  if (a.empty() || b.empty()) return {};
  int inv = modpow(n, mod - 2);
  vl L(a), R(b), out(n);
  L.resize(n), R.resize(n);
  ntt(L), ntt(R);
    out [-i \& (n-1)] = (l1)L[i] * R[i] % mod * inv %
           mod;
  ntt(out);
  return {out.begin(), out.begin() + s};
```

#### FastSubsetTransform.h

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

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

## Number theory (5)

#### 5.1 Modular arithmetic

#### ModInt.h

Description: Operators for modular arithmetig<sub>f0e63</sub>, 29 lines

```
template<int M, int R>
struct Mod {
 static const int mod = M, rt = R;
 Mod(ll 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 *= a) if (n % 2) b *= a;
   return b; }
 Mod inv() const { assert(x); return pow(M - 2); }
 friend ostream& operator<<(ostream& os, Mod x) {</pre>
   return os << x.x; }
using mint = Mod<998244353, 3>;
```

#### ModInverse.h

Description: Pre-computation of modular inverses. Assumes LIM < mod and that mod is a prime. 6f684f, 3 lines

```
const 11 mod = 1000000007, LIM = 200000;
                  l1* inv = new l1[LIM] - 1; inv[1] = 1;
rep(i,2,LIM) inv[i] = mod - (mod / i) * inv[mod % i] %
                  ModPow.h
                                                                   b83e45, 8 lines
                  const 11 mod = 1000000007; // faster if const
                  ll modpow(ll b, ll e) {
                    ll ans = 1;
                    for (; e; b = b * b % mod, e /= 2)
                      if (e & 1) ans = ans * b % mod;
                    return ans;
                  ModLog.h
                  Description: Returns the smallest x > 0 s.t. a^x = b
                  (\text{mod } m), or -1 if no such x exists. \text{modLog}(a,1,m) can be
                  used to calculate the order of a.
                  Time: \mathcal{O}\left(\sqrt{m}\right)
                  ll modLog(ll a, ll b, ll m) {
                   ll n = (ll) sqrt(m) + 1, e = 1, f = 1, j = 1;
                    unordered map<ll, ll> A:
                    while (j <= n && (e = f = e * a % m) != b % m)
A[e * b % m] = j++;</pre>
                    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))
464cf3, 16 lines
                        return n * i - A[e];
                    return -1;
                  ModSum.h
                  Description: Sums of mod'ed arithmetic progressions.
                  modeum (to, c, k, m) = \sum_{i=0}^{t_0-1} (ki+c)\%m. divsum is similar but for floor d
                  but for floored division.
                  Time: \log(m), with a large constant.
                                                                  5c5bc5, 16 lines
                  typedef unsigned long long ull;
                  ull sumsq(ull to) { return to / 2 * ((to-1) | 1); }
                  ull divsum(ull to, ull c, ull k, ull m) {
                   ull res = k / m * sumsq(to) + c / m * to;
                   k %= m; c %= m;
                   if (!k) return res;
                   ull to2 = (to * k + c) / m;

return res + (to - 1) * to2 - divsum(to2, m-1 - c, m,
                  11 modsum(ull 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,
                  ModLinear.h
                  Description: Statistics on a mod'ed arithmetic sequence.
                  Time: O(\log m)
                  "euclid.h"
                                                                  45f202, 34 lines
                  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 \text{ y2} = (n * a + b) / m;
```

```
11 x2 = cdiv(m2 * y2 - b, a);
 if (x2 * a - m2 * y2 + b >= m2) --x2;
 return minRemainder(a, b, m2, x2);
// min x >= 0 s.t. 1 <= (ax + b) % m <= r
ll minBetween(ll a, ll b, ll m, ll l, ll r) {
 ll x, y, g = euclid(a, m, x, y);
 if (g > 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);
Description: Calculate a \cdot b \mod c (or a^b \mod c) for 0 <
a, b \le c \le 7.2 \cdot 10^{18}.
Time: \mathcal{O}(1) for modmul, \mathcal{O}(\log b) for modpow bbbd8f, 11 lines
typedef unsigned long long ull;
ull modmul(ull a, ull b, ull M) {
 11 ret = a * b - M * ull(1.L / M * a * b);
 return ret + M * (ret < 0) - M * (ret >= (11)M);
ull modpow(ull b, ull e, ull mod) {
  for (; e; b = modmul(b, b, mod), e /= 2)
   if (e & 1) ans = modmul(ans, b, mod);
```

#### ModSart.h

Description: Tonelli-Shanks algorithm for modular square roots. Finds x s.t.  $x^2 = a \pmod{p} (-x \text{ gives the other solu-}$ tion) or -1 if no such x exists.

```
Time: \mathcal{O}\left(\log^2 p\right) worst case, \mathcal{O}\left(\log p\right) for most p
"ModMulLL.h"
ll sgrt(ll a, ll p) {
 a %= p; if (a < 0) a += p;
  if (a == 0) return 0;
  if (modpow(a, (p-1)/2, p) != 1) return -1;
if (p % 4 == 3) return modpow(a, (p+1)/4, p);
// a^(n+3)/8 or 2^(n+3)/8 * 2^(n-1)/4 works if p % 8
        == 5
  11 s = p - 1, n = 2;
  int r = 0, m;
  while (s % 2 == 0)
     ++r, s /= 2;
   while (modpow(n, (p - 1) / 2, p) != p - 1) ++n;
  11 x = modpow(a, (s + 1) / 2, p);

11 b = modpow(a, s, p), g = modpow(n, s, p);
  for (;; r = m) {
     11 t = b;
     for (m = 0; m < r && t != 1; ++m)
      t = t * t % p;
     if (m == 0) return x;
     11 \text{ gs} = \text{modpow}(g, 1LL << (r - m - 1), p);
     q = qs * qs % p;
     x = x * gs % p;
     b = b * g % p;
```

#### PrimitiveRoot.h

**Description:** Finds a primitive root modulo p.

```
"Factor.h", "ModMulLL.h"
                                                      d63925, 9 lines
mt19937 64 rng(2137);
11 primitiveRoot(11 p) {
  auto f = factor(p - 1); sort(all(f));
  f.resize(unique(all(f)) - f.begin()); rep:
  11 g = rng() % (p - 1) + 1;

for (auto x : f) if (modpow(q, (p - 1) / x, p) == 1)
    goto rep;
  return a:
```

#### 5.2 Primality

#### FastEratosthenes.h

Description: Prime sieve for generating all primes smaller than LIM.

```
Time: LIM=1e9 \approx 1.5s
                                               6b2912, 20 lines
const int LIM = 1e6;
bitset<LIM> isPrime;
vi eratosthenes() {
  const int S = (int) round(sqrt(LIM)), R = LIM / 2;
 vi pr = {2}, sieve(S+1); pr.reserve(int(LIM/log(LIM)
        *1.1));
  vector<pii> cp;
  for (int i = 3; i <= S; i += 2) if (!sieve[i]) {</pre>
    cp.push_back({i, i * i / 2});
for (int j = i * i; j <= S; j += 2 * i) sieve[j] =</pre>
  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]</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 7.10<sup>18</sup>; for larger numbers, use Python and extend A randomly.

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

```
60dcd1, 12 lines
bool isPrime(ull n) {
   if (n < 2 | | n % 6 % 4 != 1) return (n | 1) == 3;
ull A[] = {2, 325, 9375, 28178, 450775, 9780504,
              1795265022},
   179320322;
s = _builtin_ctzll(n-1), d = n >> s;
for (ull a : A) { // ^ count trailing zeroes
ull p = modpow(a%n, d, n), i = s;
while (p != 1 && p != n - 1 && a % n && i--)
       p = modmul(p, p, n);
if (p != n-1 && i != s) return 0;
   return 1;
```

#### Factor.h

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

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

```
"ModMulLL.h", "MillerRabin.h"
  ull x = 0, y = 0, t = 30, prd = 2, i = 1, q;
  auto f = [&](ull x) { return modmul(x, x, n) + i; };
  while (t++ % 40 || __gcd(prd, n) == 1) {
   if (x == y) x = ++i, y = f(x);
   if ((q = modmul(prd, max(x,y) - min(x,y), n))) prd
    = q;
x = f(x), y = f(f(y));
  return __gcd(prd, n);
vector<ull> factor(ull n) {
  if (n == 1) return {};
  if (isPrime(n)) return {n};
  ull x = pollard(n);
  auto l = factor(x), r = factor(n / x);
  l.insert(l.end(), all(r));
  return 1:
```

#### 5.3 Divisibility

#### euclid.h

**Description:** Finds two integers x and y, such that ax + by =gcd(a, b). If you just need gcd, use the built in  $\_gcd$  instead. If a and b are coprime, then x is the inverse of  $a_3$  (mage b) lines

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

#### CRT.h

Description: Chinese Remainder Theorem.

crt (a, m, b, n) computes x such that  $x \equiv a \pmod{m}$ ,  $x \equiv b$ (mod n). If |a| < m and |b| < n, x will obey 0 < x < nlcm(m, n). Assumes  $mn < 2^{62}$ .

Time:  $\log(n)$ 04d93a, 7 lines

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

#### 5.3.1 Bézout's identity

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

$$ax + by = d$$

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

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

#### phiFunction.h

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

Fermat's little thm:  $p \text{ prime} \Rightarrow a^{p-1} \equiv 1 \pmod{8}$  (most weakly selines

```
const int LIM = 5000000;
int phi[LIM];
void calculatePhi() {
 rep(i,0,LIM) phi[i] = i&1 ? i : i/2;
  for (int i = 3; i < LIM; i += 2) if(phi[i] == i)</pre>
    for (int j = i; j < LIM; j += i) phi[j] -= phi[j] /</pre>
```

#### Min25.h

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

Time: 
$$\mathcal{O}\left(\frac{n^{3/4}}{\log n}\right)$$

"FastEratosthenes.h" 4c2ea6, 47 lines

```
template<class T>
struct Min25 {
 ll n, sq, s, hls; vi p;
Min25(ll N) : n(N) {
    sq = sqrt(n) + 5;
    while (sq * sq > n) sq--;
    hls = quo(n, sq);
    if (hls != 1 && quo(n, hls - 1) == sq) hls--;
   s = hls + sq;
rep(i, 2, sq + 1) if (isPrime[i]) p.push_back(i);
  vector<T> sieve(auto f) {
   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 (ll x : p) {
      T xt = f(x) - f(x - 1), pi = h[s - x + 1];
      11 x2=x*x, imax=min(hls, quo(n, x2) + 1), ix=x;
      for (int i = 1; i < imax; i++, ix += x)</pre>
        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(vector<T>& fp, auto f) {
    vector<ll> ns(s);
    rep(i, 1, hls) ns[i] = quo(n, i);
    rep(i, 1, sq + 1) ns[s - i] = i;
    auto F = fp, G = fp;
    for (ll P : p | views::reverse) {
      for (ll pk=P, k=1; quo(n,P)>=pk; k++, pk*=P) {
        T = fp[idx(P)], y=f(P,k,pk), z=f(P,k+1,pk*P);
          if (P * pk > m) break;
          G[i] += y * (F[idx(quo(m, pk))] - x) + z;
      copy_n(G.begin(), min(s,idx(P*P)+1),F.begin());
    rep(i, 1, sz(ns)) F[i] += 1;
```

```
il quo(ll x, ll y) { return (double)x / y; }
ll idx(ll x) { return x <= sq ? s - x : quo(n, x); }
```

#### 5.4 Fractions

#### ContinuedFractions.h

**Description:** Given N and a real number x > 0, finds the closest rational approximation p/q with  $p, q \leq \overline{N}$ . It will obey

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

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

#### FracBinarySearch.h

LQ = Q; Q = NQ;

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

Usage: fracBS([](Frac f) { return f.p>=3\*f.q; }, 10); // {1,3}

```
Time: \mathcal{O}(\log(N))
                                              27ab3e, 25 lines
struct Frac { 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;
```

#### 5.5 Pythagorean Triples

The Pythagorean triples are uniquely generated

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

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

#### 5.6 Primes

p = 962592769 is such that  $2^{21} | p - 1$ , which may be useful. For hashing use 970592641 (31-bit number), 31443539979727 (45-bit), 3006703054056749 (52-bit). There are 78498 primes less than 1000000.

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

#### 5.7 Estimates

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

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

#### 5.8 Mobius Function

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

Mobius Inversion:

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

Other useful formulas/forms:

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

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

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

## Combinatorial (6)

#### 6.1 Permutations

#### 6.1.1 Factorial

n	1 2 3	4	5 6	7	8	9	10	
$\overline{n!}$	1 2 6	24 1	20 720	5040	40320	362880	3628800	
n	11	12	13	14	15	16	17	
n!	4.0e7	4.8e	8 6.2e	9 8.7e	10 1.3e	12 2.1e	13 3.6e14	
n	20	25	30	40	50 10	00   15	0 171	
n!	2e18	2e25	3e32	$8e47 \ 3$	e64 9e1	157  6e2	$62 > DBL_M$	ΑZ

#### IntPerm.h

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

```
int permToInt(vi& v) {
 int use = 0, i = 0, r = 0;
 for(int x:v) r = r * ++i + __builtin_popcount(use &
       -(1<<x)),
    use |= 1 << x;
                                       // (note: minus
          not ∼!)
```

#### **6.1.2** Cycles

UW

Let  $q_S(n)$  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.

$$D(n) = (n-1)(D(n-1) + D(n-2)) = nD(n-1) + (-1)^n$$

#### 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 q

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.

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

#### 6.2.2 Lucas' Theorem

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

#### 6.2.3 Binomials

multinomial.h

return c;

Description: Computes 
$$\binom{\kappa_1 + \cdots + \kappa_n}{k_1, k_2, \dots, k_n} = \frac{(\sum \kappa_i)!}{k_1 \log 12 / \kappa_n \text{ lines}}$$
.

11 multinomial (vi& v) {
 ll c = 1, m = v.empty() ? 1 : v[0];
 rep(i,1,sz(v)) rep(j,0,v[i]) c = c \* ++m / (j+1);

6.3 General purpose numbers

#### 6.3.1 Bernoulli numbers

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

Sums of powers:

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

Euler-Maclaurin formula for infinite sums:

$$D(n) = (n-1)(D(n-1) + D(n-2)) = nD(n-1) + (-1)^n = \left\lfloor \frac{n!}{e} \right\rfloor_{k=m}^{\infty} f(i) = \int_m^{\infty} f(x) dx - \sum_{k=1}^{\infty} \frac{B_k}{k!} f^{(k-1)}(m)$$

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

# 6.3.2 Stirling numbers of the first

Number of permutations on n items with kcycles.

$$c(n,k) = c(n-1,k-1) + (n-1)c(n-1,k), \ c(0,0) = 1$$
 
$$\sum_{k=0}^{n} c(n,k)x^{k} = x(x+1)\dots(x+n-1)$$
 
$$c(8,k) = 8,0,5040,13068,13132,6769,1960,322,28,1$$
 
$$c(n,2) = 0,0,1,3,11,50,274,1764,13068,109584,\dots$$

#### 6.3.3 Eulerian numbers

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

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

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

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

# 6.3.4 Stirling numbers of the second

Partitions of n distinct elements into exactly k

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

#### 6.3.5 Bell numbers

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

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

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

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

- sub-diagonal monotone paths in an  $n \times n$  grid.
- strings with n pairs of parenthesis, correctly
- 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

#### 6.4 Other

#### DeBruiin h

**Description:** Recursive FKM, given alphabet [0, k) constructs cyclic string of length kn that contains every length n string as substr. a7faa5, 13 lines

```
vi dseq(int k, int n) {
if (k == 1) return {0};
 vi res, aux(n+1);
 function<void(int,int)> gen = [&](int t, int p) {
   if (t > n) { // consider lyndon word of len p
  if (n%p == 0) FOR(i,1,p+1) res.pb(aux[i]);
      aux[t] = aux[t-p]; gen(t+1,p);
FOR(i,aux[t-p]+1,k) aux[t] = i, gen(t+1,t);
gen(1,1); return res;
```

#### PermGroup.h

Description: Schreier-Sims lets you add a permutation to a group, count number of permutations in a group, test whether a permutation is a member of a group. Works well for n < 15, maybe for larger too. Construct PermGroup() and run order() to get order of the group.

```
Time: \mathcal{O}\left(n^6\right)
                                             d6edf4, 54 lines
vi inv(vi v) { vi V(sz(v)); rep(i,sz(v)) V[v[i]]=i;
     return V: }
vi id(int n) { vi v(n); iota(all(v),0); return v; }
vi operator* (const vi& a, const vi& b)
  vi c(sz(a)); rep(i,sz(a)) c[i] = a[b[i]];
 return c:
struct PermGroup {
  struct Group {
    vector<vi> gen, sigma;
    Group(int n, int p) : flag(n), sigma(n) {
        flag[p] = 1; sigma[p] = id(n);
  int n = 0; vector<Group> q;
 bool check (const vi& cur, int k) {
```

```
if (!k) return 1;
  int t = cur[k]:
  return g[k].flag[t] ? check(inv(g[k].sigma[t])*cur,
void updateX(const vi& cur, int k) {
  int t = cur[k]; // if flag, fixes k -> k
  if (g[k].flag[t]) ins(inv(g[k].sigma[t])*cur,k-1);
       g[k].flag[t] = 1, g[k].sigma[t] = cur;
      for (auto x: q[k].gen)
        undateX(x*cur,k);
void ins(const vi& cur, int k) {
  if (check(cur,k)) return;
  a[k].gen.pb(cur);
  rep(i,n) if (g[k].flag[i]) updateX(cur*g[k].sigma[i
        1.k);
ll order(vector<vi> gen) {
  if(sz(gen) == 0) return 1;
  n = sz(gen[0]);
  rep(i,n) g.pb(Group(n,i));
  for (auto a: gen)
      ins(a, n-1); // insert perms into group one by
  11 tot = 1; // watch out for overflows, can be up
        to n!
      int cnt = 0;
      rep(j,i+1) cnt += g[i].flag[j];
  return tot;
```

#### GravCode.h

**Description:** Gray code:  $gray(0), \ldots, gray(2^n-1)$  - permutation in which each two consecutive (cyclically) numbers. differ in exactly one bit.

```
using ull = unsigned long long;
ull grav(ull i) { return i^i>>1;
ull invg(ull i) { // i=invg(gray(i))=gray(invg(i))
  i^=i>>1; i^=i>>2; i^=i>>4;
 i^=i>>8; i^=i>>16; i^=i>>32; return i;
```

## Graph (7)

#### 7.1 Fundamentals

#### BellmanFord.h

**Description:** Calculates shortest paths from s in a graph that might have negative edge weights. Unreachable nodes get dist = inf; nodes reachable through negative-weight cycles get dist = -inf. Assumes  $V^2 \max |w_i| < \sim 2^{63}$ .

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

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

#### FlovdWarshall.h

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

```
Time: \mathcal{O}\left(N^3\right)
                                             531245, 12 lines
const 11 inf = 1LL << 62;</pre>
void floydWarshall(vector<vector<ll>>& m) {
  int n = sz(m);
  rep(i, 0, n) m[i][i] = min(m[i][i], OLL);
  rep(k,0,n) rep(i,0,n) rep(j,0,n)
    if (m[i][k] != inf && m[k][j] != inf) {
      auto newDist = max(m[i][k] + m[k][j], -inf);
     m[i][j] = min(m[i][j], newDist);
  rep(k,0,n) if (m[k][k] < 0) rep(i,0,n) rep(j,0,n)
    if (m[i][k] != inf && m[k][j] != inf) m[i][j] = -
```

#### TopoSort.h

**Description:** Topological sorting. Given is an oriented graph. Output is an ordering of vertices, such that there are edges only from left to right. If there are cycles, the returned list will have size smaller than n – nodes reachable from cycles will not be returned

```
Time: \mathcal{O}(|V| + |E|)
vi topoSort(const vector<vi>& gr) {
  vi indeg(sz(gr)), q;
 for (auto& li : gr) for (int x : li) indeg[x]++;
rep(i,0,sz(gr)) if (indeg[i] == 0) q.push_back(i);
  rep(j, 0, sz(q)) for (int x : gr[q[j]])
    if (-indeg[x] == 0) q.push_back(x);
  return q;
```

Description: SPFA with subtree erasure heuristic. Returns array of distances or empty array if negative cycle is reachable from source. par[v] = parent in shortest path tree

```
Time: \mathcal{O}(VE) but fast on random
                                              bf4a3f, 33 lines
using Edge = pair<int, 11>;
vector<11> spfa(vector<vector<Edge>>& G,
                vi& par, int src) {
  int n = sz(G); vi que, prv(n+1);
  iota(all(prv), 0); vi nxt = prv;
  vector<ll> dist(n, INT64_MAX);
 par.assign(n, -1);
auto add = [&](int v, int p, ll d) {
   par[v] = p; dist[v] = d;
prv[n] = nxt[prv[v] = prv[nxt[v] = n]] = v;
  auto del = [&](int v) {
    nxt[prv[nxt[v]] = prv[v]] = nxt[v];
    prv[v] = nxt[v] = v;
  for (add(src, -2, 0); nxt[n] != n;) {
    int v = nxt[n]; del(v);
    for (auto e : G[v]) {
      11 alt = dist[v] + e.y;
      if (alt < dist[e.x]) {</pre>
        que = \{e.x\};
        rep(i, sz(que)) {
          int w = que[i]; par[w] = -1;
          del(w);
          for (auto f : G[w])
            if (par[f.x] == w) que.pb(f.x);
        if (par[v] == -1) return {};
        add(e.x, v, alt);
  return dist; }
```

#### 7.2 Network flow

#### PushRelabel.h

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

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

0ae1d4, 48 lines

```
struct PushRelabel {
 struct Edge {
   int dest, back;
   11 f. c:
 vector<vector<Edge>> g:
 vector<11> ec:
 vector<Edge*> cur:
 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.
    e.f += f; e.c -= f; ec[e.dest] += f;
   back.f -= f; back.c += f; ec[back.dest] -= f;
 ill calc(int s, int t) {
   int v = sz(g); H[s] = v; ec[t] = 1;
   vi co(2*v); co[0] = v-1;
    rep(i,0,v) cur[i] = g[i].data();
    for (Edge& e : g[s]) addFlow(e, e.c);
    for (int hi = 0;;) {
      while (hs[hi].empty()) if (!hi--) return -ec[s];
      int u = hs[hi].back(); hs[hi].pop_back();
      while (ec[u] > 0) // discharge u
       if (cur[u] == g[u].data() + sz(g[u])) {
         H[u] = 1e9;
         for (Edge& e : g[u]) if (e.c && H[u] > H[e.
               dest]+1)
           H[u] = H[e.dest]+1, cur[u] = &e;
         if (++co[H[u]], !--co[hi] && hi < v)</pre>
           rep(i, 0, v) if (hi < H[i] && H[i] < v)
               -co[H[i]], H[i] = v + 1;
         hi = H[u];
        } else if (cur[u]->c && H[u] == H[cur[u]->dest
         addFlow(*cur[u], min(ec[u], cur[u]->c));
       else ++cur[u];
 bool leftOfMinCut(int a) { return H[a] >= sz(g); }
```

#### MinCostMaxFlow.h

Description: Min-cost max-flow. If costs can be negative, call setpi before maxflow, but note that negative cost cycles are not supported. To obtain the actual flow, look at positive values only.

```
Time: \mathcal{O}(FE \log(V)) where F is max flow. \mathcal{O}(FE \log(V))
#include <ext/pb_ds/priority_queue.hpp>
const 11 INF = numeric_limits<11>::max() / 4;
struct MCMF {
  struct edge {
    int from, to, rev;
    ll 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, ll cap, ll 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;
    __qnu_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]) {
         ll val = di - pi[e.to] + e.cost;
```

```
if (e.cap - e.flow > 0 && val < dist[e.to]) {</pre>
        dist[e.to] = val;
par[e.to] = &e;
        if (its[e.to] == q.end())
          its[e.to] = q.push({ -dist[e.to], e.to });
        else
          q.modify(its[e.to], { -dist[e.to], e.to });
  rep(i,0,N) pi[i] = min(pi[i] + dist[i], INF);
pair<ll, ll> maxflow(int s, int t) {
  11 totflow = 0, totcost = 0;
  while (path(s), seen[t]) {
    11 fl = INF;
    for (edge* x = par[t]; x; x = par[x->from])
      fl = min(fl, x->cap - x->flow);
    totflow += fl:
    for (edge* x = par[t]; x; x = par[x->from]) {
      x \rightarrow flow += fl:
      ed[x->to][x->rev].flow -= fl;
  rep(i,0,N) for(edge& e : ed[i]) totcost += e.cost *
        e.flow;
  return {totflow, totcost/2};
// If some costs can be negative, call this before
      maxflow:
void setpi(int s) { // (otherwise, leave this out)
  fill(all(pi), INF); pi[s] = 0;
  int it = N, ch = 1; ll v;
  while (ch-- && it--)
    rep(i,0,N) if (pi[i] != INF)
      for (edge& e : ed[i]) if (e.cap)
        if ((v = pi[i] + e.cost) < pi[e.to])
  pi[e.to] = v, ch = 1;
assert(it >= 0); // negative cost cycle
```

#### EdmondsKarp.h

Description: Flow algorithm with guaranteed complexity  $O(VE^2)$ . To get edge flow values, compare capacities before and after, and take the positive values only. \_482fe0, 36 lines

```
template < class T > T edmonds Karp (vector < unordered_map <
     int, T>>&
    graph, int source, int sink) {
  assert (source != sink);
 T flow = 0:
  vi par(sz(graph)), q = par;
 for (;;) {
    fill(all(par), -1);
    par[source] = 0;
    int ptr = 1:
    q[0] = source;
    rep(i,0,ptr) {
      int x = q[i];
      for (auto e : graph[x]) {
        if (par[e.first] == -1 && e.second > 0) {
          par[e.first] = x;
          g[ptr++] = e.first;
          if (e.first == sink) goto out;
    return flow;
0111:
    T inc = numeric limits<T>::max();
    for (int y = sink; y != source; y = par[y])
      inc = min(inc, graph[par[y]][y]);
    flow += inc;
    for (int y = sink; y != source; y = par[y]) {
  int p = par[y];
      if ((graph[p][y] -= inc) <= 0) graph[p].erase(y);</pre>
      graph[y][p] += inc;
```

#### 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. d7f0f1, 42 lines

```
struct Dinic {
 struct Edge {
   int to, rev;
```

```
ll flow() { return max(oc - c, OLL); } // if you
        need flows
vi lvl, ptr, q;
vector<vector<Edge>> adj;
Dinic(int n) : lvl(n), ptr(n), q(n), adj(n) {}
void addEdge(int a, int b, ll c, ll rcap = 0) {
  adj[a].push_back({b, sz(adj[b]), c, c});
adj[b].push_back({a, sz(adj[a]) - 1, rcap, rcap});
11 dfs(int v, int t, 11 f) {
  if (v == t || !f) return f;
  for (int& i = ptr[v]; i < sz(adj[v]); i++) {</pre>
    Edge& e = adj[v][i];
    if (lvl[e.to] == lvl[v] + 1)
      if (ll p = dfs(e.to, t, min(f, e.c))) {
        e.c -= p, adj[e.to][e.rev].c += p;
         return p;
  return 0;
il 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; }
```

#### MinCut.h

Description: After running max-flow, the left side of a mincut 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

8b0e19, 21 lines

```
graph, as represented by an adjacency matrix.
Time: \mathcal{O}\left(V^3\right)
pair<int, vi> globalMinCut(vector<vi> mat) {
 pair<int, vi> best = {INT_MAX, {}};
  int n = sz(mat);
  vector<vi> co(n);
  rep(i, 0, n) co[i] = {i};
  rep(ph,1,n) {
    vi w = mat[0];
    size_t s = 0, t = 0;
    rep(it, 0, n-ph) { // O(V^2) -> O(E log V) with prio.
       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;
```

#### GomoryHu.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
```

vector<Edge> tree;

vi par(N);

```
0418b3, 1<u>3 lines</u>
"PushRelabel.h"
typedef array<11, 3> Edge;
vector<Edge> gomoryHu(int N, vector<Edge> ed) {
```

#### 7.3 Matching

#### Matching.h

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

```
Time: \mathcal{O}\left(E\sqrt{V}\right)
                                            daf32c, 21 lines
int matching(vector<vi>& g, vi& r) {
 int n = sz(g), res = 0;
  vi l(n, -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 \mid | (d[r[v]] == t && f(f, r[v])))
       return 1[u] = v, r[v] = u, 1;
  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 : q[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 res;
   rep(i, 0, n) if ([i] == -1) res += 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.

#### 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 < M.

```
Time: \mathcal{O}\left(N^2M\right)
```

```
pair<11, vi> hungarian(const vector<vector<11>>> &a) {
    if (a.empty()) return {0, {}};
    int n = sz(a) + 1, m = sz(a[0]) + 1;
    vi p(m), ans(n - 1); vector<11> u(n), v(m);
    rep(i,1,n) {
```

#### Blossom.h

**Description:** Matching for general graphs using Blossom algorithm. Time:  $\mathcal{O}(NM)$ , fast in practice

```
int blossom(vector<vi>& G, vi& match) {
  int n = sz(G), cnt = -1, ans = 0; match.assign(n, -1);
 vi lab(n), par(n), orig(n), aux(n, -1), q;

auto blos = [&] (int v, int w, int a) {
    while (orig[v] != a) {
      par[v] = w; w = match[v];
if (lab[w] == 1) lab[w] = 0, q.push_back(w);
      orig[v] = orig[w] = a; v = par[w];
 };
 rep(i, 0, n) if (match[i] == -1)
    for (auto e : G[i]) if (match[e] == -1) {
      match[match[e] = i] = e; ans++; break;
  rep(root, 0, n) if (match[root] == -1) {
    fill(all(lab), -1);
    iota(all(orig), 0);
    lab[root] = 0;
    q = \{root\};
    rep(i, 0, sz(q)) {
      int v = q[i];
      for (auto x : G[v]) if (lab[x] == -1) {
         lab[x] = 1; par[x] = v;
        if (match[x] == -1) {
  for (int y = x; y+1;) {
             int p = par[y], w = match[p];
match[match[p] = y] = p; y = w;
           ans++;
           goto nxt;
         lab[match[x]] = 0; q.push_back(match[x]);
       } else if (lab[x] == 0 && orig[v]!=orig[x]) {
         int a = orig[v], b = orig[x];
         for (cnt++;; swap(a, b)) if (a+1) {
           if (aux[a] == cnt) break;
           a = (match[a]+1)?
             orig[par[match[a]]]: -1);
         blos(x, v, a); blos(v, x, a);
    nxt:;
  return ans; }
```

#### WeightedBlossom.h

**Description:** Edmond's Blossom algorithm for weighted maximum matching in general graphs. Weights must be positive. **Time:**  $\mathcal{O}\left(N^3\right)$  998ff6, 228 lines

```
struct WeightedBlossom {
   struct edge { int u, v, w; };
   int n, s, nx;
   vector<vector<edge>> g;
   vi lab, match, slack, st, pa, S, vis;
   vector<vi>> flo, floFrom;
```

```
queue<int> q;
// Initialize for k vertices
WeightedBlossom(int k)
    : n(k), s(n*2+1).
      g(s, vector<edge>(s)),
      lab(s), match(s), slack(s), st(s),
      pa(s), S(s), vis(s), flo(s),
  floFrom(s, vi(n+1)) {
rep(u, 1, n+1) rep(v, 1, n+1)
g[u][v] = {u, v, 0};
// Add edge between u and v with weight w
void addEdge(int u, int v, int w) {
  u++; v++;
  q[u][v].w = q[v][u].w = max(q[u][v].w, w);
// Compute max weight matching.
   'count' is set to matching size,
   'weight' is set to matching weight.
// Returns vector 'match' such that:
// match[v] = vert matched to v or -1
vi solve(int& count, ll& weight) {
  fill(all(match), 0);
  nx = n;
  weight = count = 0;
  rep(u, 0, n+1) flo[st[u] = u].clear();
  int tmp = 0;
rep(u, 1, n+1) rep(v, 1, n+1) {
    floFrom[u][v] = (u-v ? 0 : v);
    tmp = max(tmp, g[u][v].w);
  rep(u, 1, n+1) lab[u] = tmp;
  while (matching()) count++;
  rep(u, 1, n+1)
    if (match[u] && match[u] < u)</pre>
      weight += g[u][match[u]].w;
  vi ans(n);
  rep(i, 0, n) ans[i] = match[i+1]-1;
  return ans;
int delta(edge& e) {
  return lab[e.u]+lab[e.v]-g[e.u][e.v].w*2;
void updateSlack(int u, int x)
  void setSlack(int x) {
  slack[x] = 0;
  rep(u, 1, n+1) if (g[u][x].w > 0 &&
    st[u] != x && !S[st[u]])
      updateSlack(u, x);
void push(int x) {
  if (x <= n) q.push(x);
else rep(i, 0, sz(flo[x])) push(flo[x][i]);</pre>
void setSt(int x, int b) {
  st[x] = b;
if (x > n) rep(i, 0, sz(flo[x]))
  setSt(flo[x][i],b);
int getPr(int b, int xr) {
  int pr = int(find(all(flo[b]), xr) -
    flo[b].begin());
  if (pr % 2) {
    reverse(flo[b].begin()+1, flo[b].end());
    return sz(flo[b]) - pr;
  } else return pr:
void setMatch(int u, int v) {
  match[u] = q[u][v].v;
  if (n <= n) return:
  edge e = g[u][v];
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(flo[u].begin(), flo[u].begin()+pr,
  flo[u].end());
void augment(int u, int v) {
  while (1) {
    int xnv = st[match[u]];
    setMatch(u, v);
    if (!xnv) return;
    setMatch(xnv, st[pa[xnv]]);
    u = st[pa[xnv]], v = xnv;
int getLca(int u, int v) {
  static int t = 0;
```

```
for (++t; u||v; swap(u, v)) {
    if (!u) continue;
    if (vis[u] == t) return u;
    vis[u] = t;
    u = st[match[u]];
    if (u) u = st[pa[u]];
  return 0:
void blossom(int u, int lca, int v) {
  int h = n+1:
  while (b <= nx && st[b]) ++b;
  if (b > nx) ++nx;
  lab[b] = S[b] = 0;
  match[b] = match[lca];
  flo[b].clear();
  flo[b].push_back(lca);
  for (int x=u, y; x != lca; x = st[pa[y]]) {
  flo[b].push_back(x);
    flo[b].push_back(y = st[match[x]]);
    push(y);
  reverse(flo[b].begin()+1, flo[b].end());
  for (int x=v, y; x != lca; x = st[pa[y]]) {
    flo[b].push_back(x);
    flo[b].push_back(y = st[match[x]]);
    push (y);
  setSt(b, b);
  rep(x, 1, nx+1) g[b][x].w = g[x][b].w = 0;
  rep(x, 1, n+1) floFrom[b][x] = 0;
  rep(i, 0, sz(flo[b])) {
    int xs = flo[b][i];
    rep(x, 1, nx+1) if (!q[b][x].w ||
      delta(g[xs][x]) < delta(g[b][x]))
        q[b][x]=q[xs][x], q[x][b]=q[x][xs];
    rep(x, 1, n+1) if (floFrom[xs][x])
       floFrom[b][x] = xs;
  setSlack(b):
void blossom(int b) {
  for (auto &e : flo[b]) setSt(e, e);
  int xr = floFrom[b][g[b][pa[b]].u];
  int pr = getPr(b, xr);
  int pr getric, xr,
for (int i = 0; i < pr; i += 2) {
  int xs = flo[b][i], xns = flo[b][i+1];</pre>
    pa[xs] = q[xns][xs].u;
    S[xs] = 1; S[xns] = slack[xs] = 0;
    setSlack(xns); push(xns);
  S[xr] = 1; pa[xr] = pa[b];
  rep(i, pr+1, sz(flo[b])) {
    int xs = flo[b][i];
    S[xs] = -1; setSlack(xs);
  st[b] = 0;
bool found (const edge& e) {
  int u = st[e.u], v = st[e.v];
  if (S[v] == -1) {
  pa[v] = e.u; S[v] = 1;
  int nu = st[match[v]];
    slack[v] = slack[nu] = S[nu] = 0;
    push (nu):
  } else if (!S[v]) {
    int lca = getLca(u, v);
    if (!lca) return augment(u, v),
      augment (v, u), 1;
    else blossom(u, lca, v);
  return 0:
bool matching() {
  fill(S.begin(), S.begin()+nx+1, -1);
  fill(slack.begin(), slack.begin()+nx+1, 0);
  q = {};
  rep(x, 1, nx+1)
 if (st[x] == x && !match[x])
pa[x] = S[x] = 0, push(x);
if (q.empty()) return 0;
  while (1) {
    while (q.size()) {
      int u = q.front(); q.pop();
      if (S[st[u]] == 1) continue;
      rep(v, 1, n+1)
        if (g[u][v].w > 0 && st[u] != st[v]){
           if (!delta(g[u][v])) {
             if (found(g[u][v])) return 1;
           } else updateSlack(u, st[v]);
```

```
int d = INT_MAX;
  rep(b, n+1, nx+1)
    if (st[b] == b && S[b] == 1)
      d = min(d, lab[b]/2);
  rep(x, 1, nx+1)
    if (st[x] == x && slack[x]) {
   if (S[x] == -1)
        d = min(d, delta(g[slack[x]][x]));
      else if (!S[x])
d = min(d,delta(g[slack[x]][x])/2);
  rep(u, 1, n+1) {
    if (!S[st[u]]) {
   if (lab[u] <= d) return 0;</pre>
      lab[u] -= d;
    } else if (S[st[u]] == 1) lab[u] += d;
  rep(b, n+1, nx+1) if (st[b] == b) {
    if (!S[st[b]]) lab[b] += d*2;
    else if (S[st[b]] == 1) lab[b] -= d*2;
  rep(x, 1, nx+1)
    if (st[x] == x && slack[x] &&
      st[slack[x]] != x &&
      !delta(g[slack[x]][x]) &&
      found(g[slack[x]][x])) return 1;
  rep(b, n+1, nx+1)
    if (st[b] == b && S[b] == 1 && !lab[b])
      blossom(b):
return 0;
```

#### 7.4 DFS algorithms

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

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

```
Time: \mathcal{O}\left(E+V\right)
                                               76b5c9, 24 lines
vi val, comp, z, cont;
int Time, ncomps:
template<class G, class F> int dfs(int j, G& g, F& f) {
 int low = val[j] = ++Time, x; z.push_back(j);
  for (auto e : g[j]) if (comp[e] < 0)</pre>
 low = min(low, val[e] ?: dfs(e,g,f));
if (low == val[j]) {
    do {
      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(g);
  val.assign(n, 0); comp.assign(n, -1);
  Time = ncomps = 0;
```

#### BiconnectedComponents.h

rep(i,0,n) if (comp[i] < 0) dfs(i, g, f);

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

```
Usage: int eid = 0; ed.resize(N);
for each edge (a,b) {
ed[a].emplace_back(b, eid);
ed[b].emplace_back(a, eid++); }
bicomps([&](const vi& edgelist) {...});
Time: \mathcal{O}\left(E+V\right)
                                              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 up = dfs(y, e, f);
      top = min(top, up);
      if (up == me) {
        st.push back(e);
        f(vi(st.begin() + si, st.end()));
        st.resize(si);
      else if (up < me) st.push_back(e);
      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);
```

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 ( $\sim x$ ).

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

Time:  $\mathcal{O}(N+E)$ , where N is the number of boolean variables, and E is the number of clauses. 5f9706, 56 lines

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

```
if (low == val[i]) do {
       x = z.back(); z.pop_back();
       comp[x] = low;
       if (values[x>>1] == -1)
         values[x>>1] = x&1;
    } while (x != i);
    return val[i] = low;
  bool solve() {
    values.assign(N, -1);
values.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, incoming edge) in the Eulerian path/cycle with src at both start and end, or empty list if no cycle/path exists.

```
Time: \mathcal{O}(V+E)
                                             c62d93, 16 lines
vector<pii> eulerWalk(vector<vector<pii>>& gr, int
     nedges, int src=0) {
 int n = sz(qr);
 vi D(n), its(n), eu(nedges); 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(gr[x]);
    if (it == end) {
     ret.push back(s.back()); s.pop back(); continue;
   tie(y, e) = gr[x][it++];
if (!eu[e]) {
     D[x]--, D[y]++;
eu[e] = 1; s.push_back({y, e});
 for (int x : D) if (x < 0 \mid \mid sz(ret) != nedges+1)
       return {};
 return {ret.rbegin(), ret.rend()};
```

#### Dominators.h

Description: Tarjan's dominators in directed graph. Returns tree (as array of parents) of immediate dominators idom. idom[root] = root, idom[v] = -1 if v is unreachable from root. Time:  $\mathcal{O}\left(|E|log|V|\right)$ 

```
vi dominators(vector<vi>& G, int root) {
 int n = sz(G); vector<vi> in(n), bucket(n);
 vi pre(n, -1), anc(n, -1), par(n), best(n);
 vi ord, idom(n, -1), sdom(n, n), rdom(n);
 auto dfs = [&] (auto f, int v, int p) -> void {
  if (pre[v] == -1) {
      par[v] = p; pre[v] = sz(ord);
      ord.push back(v);
      for (auto e : G[v])
        in[e].push_back(v), f(f, e, v);
 };
 auto find = [&] (auto f, int v) -> pii {
   if (anc[v] == -1) return {best[v], v};
int b; tie(b, anc[v]) = f(f, anc[v]);
    if (sdom[b] < sdom[best[v]]) best[v] = b;</pre>
    return {best[v], anc[v]};
  rdom[root] = idom[root] = root;
 iota(all(best), 0); dfs(dfs, root, -1);
 rep(i, 0, sz(ord)) {
    int v = ord[sz(ord)-i-1], b = pre[v];
    for (auto e : in[v])
     b = min(b, pre[e] < pre[v] ? pre[e] :
      sdom[find(find, e).first]);
   for (auto u : bucket[v])rdom[u]=find(find,u).first;
    sdom[v] = b; anc[v] = par[v];
    bucket[ord[sdom[v]]].push_back(v);
  for (auto v : ord) idom[v] = (rdom[v] == v ?
   ord[sdom[v]] : idom[rdom[v]]);
  return idom; }
```

#### KthShortest.h

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

constexpr ll INF = 1e18;

Description: Given directed weighted graph with nonnegative edge weights gets K-th shortest walk (not necessarily simple) or -1 if no next path (can only happen in DAG). **Memory:**  $\mathcal{O}(m \log m + k \log m)$  (uses persistent heaps)

11

e29f5a, 57 lines

```
struct Eppstein {
 using T = 11; using Edge = pair<int, T>;
struct Node { int E[2] = {}, s = 0; Edge x; };
T shortest; // Shortest path length
  priority_queue<pair<T, int>> Q;
  vector<Node> P{1}; vi h;
  Eppstein(vector<vector<Edge>>& G, int s, int t) {
    int n = sz(G); vector<vector<Edge>> H(n);
    rep(i,0,n) for(auto &[j, w] : G[i])
     H[j].push_back({i,w});
    vi ord, par(n, -1); vector<T> d(n, -INF);
Q.push({d[t] = 0, t});
    while (!Q.empty()) {
      auto [dd, v] = Q.top(); Q.pop();
      if (d[v] == dd)
        ord.push back(v);
         for (auto &[u, w] : H[v])
        if (dd-w > d[u]) {
           Q.push({d[u] = dd-w, u});
           par[u] = v;
    if ((shortest = -d[s]) >= INF) return;
    h.resize(n);
    for (auto &v : ord)
      int p = par[v]; if (p+1) h[v] = h[p];
      for(auto &[u, w] : G[v]) if (d[u] > -INF) {
        T k = w - d[u] + d[v];
        if (k || u != p)
        h[v] = push(h[v], \{u, k\});
else p = -1;
    P[0].x.first = s; Q.push({0, 0});
  int push (int t, Edge x) {
    P.push back (P[t]);
    if (!P[t = sz(P)-1].s \mid\mid P[t].x.second >= x.second)
      swap(x, P[t].x);
    if (P[t].s) {
      int i = P[t].E[0], j = P[t].E[1];
      int d = P[i].s > P[j].s;
int k = push(d ? j : i, x);
P[t].E[d] = k; // Don't inline k!
    P[t].s++; return t;
  il nextPath() { // next length, -1 if no next path
    if (Q.empty()) return -1;
    auto [d, v] = Q.top(); Q.pop();
    for (int i : P[v].E) if (i)
      Q.push({ d-P[i].x.second+P[v].x.second, i });
    int t = h[P[v].x.first];
    if (t) Q.push({d - P[t].x.second, t });
    return shortest - d; } };
```

#### DenseDFS.h

Description: DFS over dense graph. Suddenly DFS over N <= 1000 graph many times becomes feasible 6e9645, 68 lines

```
// DFS over bit-packed adjacency matrix
// G = NxN adjacency matrix of graph
       G(i,j) \iff (i,j) \text{ is edge}
// V = 1xN matrix containing unvisited vertices
       V(0,i) <=> i-th vertex is not visited
// Total DFS time: O(n^2/64)
using ull = uint64_t;
// Matrix over Z_2 (bits and xor)
// TODO: arithmetic operations //!HIDE
struct BitMatrix {
 vector<ull> M;
 int rows, cols, stride;
  // Create matrix with n rows and m columns
 BitMatrix(int n = 0, int m = 0) {
   rows = n; cols = m; stride = (m+63)/64;
    M.resize(n*stride);
  // Get pointer to bit-packed data of i-th row
  ull* row(int i) { return &M[i*stride]; }
  // Get value in i-th row and j-th column
  bool operator()(int i, int j) {
```

```
return (row(i)[j/64] >> (j%64)) & 1;
 // Set value in i-th row and j-th column void set(int i, int j, bool val) {
    ull &w = row(i)[j/64], m = 1ull << (j\%64);
    if (val) w |= m;
    else w &= \sim m;
struct DenseDFS {
  BitMatrix G, V; // space: O(n^2/64)
  // Initialize structure for n vertices
  DenseDFS(int n = 0) : G(n, n), V(1, n) {
    reset():
  // Mark all vertices as unvisited
  void reset() { for(auto &x : V.M) x = -1; }
  // Get/set visited flag for i-th vertex
 void setVisited(int i) { V.set(0, i, 0); }
bool isVisited(int i) { return !V(0, i);
  // DFS step: func is called on each unvisited
  // neighbour of i. You need to manually call
  // setVisited(child) to mark it visited
  // or this function will call the callback
  // with the same vertex again.
  void step(int i, auto func) {
    ull* E = G.row(i);
    for (int w = 0; w < G.stride;) {</pre>
      ull x = E[w] & V.row(0)[w];
      if (x) func((w<<6) | __builtin_ctzll(x));</pre>
      else w++;
```

#### 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, 24 lines

```
template<class P>
vector<vi>planarFaces(vector<vi>& q, 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;
```

#### 7.5 Coloring

#### EdgeColoring.h

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

```
while (d = free[v], !loc[d] && (v = adj[u][d]) !=
    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
       1 [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:** Calculates chromatic number of a graph represented by a vector of bitmasks. Self loops are not allowed. **Usage:** chromaticNumber({6, 5, 3}) // 3-clique

Time:  $\mathcal{O}\left(2^n n\right)$  07ea3d, 20 lines

```
const int MOD = 1000500103; // big prime
int chromaticNumber(vi g) {
 int n = sz(q);
 if (!n) return 0;
 vi ind(1 << n, 1), s(1 << n);
 rep(i, 0, 1 << n) s[i] = __popcount(i) & 1 ? -1 : 1;
 rep(i, 1, 1 << n) {
   int ctz = __builtin_ctz(i);
   ind[i] = ind[i - (1 << ctz)] + ind[(i - (1 << ctz))]
          & ~g[ctz]];
   if (ind[i] >= MOD) ind[i] -= MOD;
  rep(k, 1, n) {
   11 sum = 0;
   rep(i, 0, 1 << n) {
      s[i] = int((ll)s[i] * ind[i] % MOD);
      sum += s[i];
   if (sum % MOD) return k;
 return n: }
```

#### 7.6 Heuristics

#### MaximalCliques.h

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

Time:  $\mathcal{O}\left(3^{n/3}\right)$ , much faster for sparse graphs b0d5b1, 12 lines

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

#### MaximumClique.h

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

Time: Runs in about 1s for n=155 and worst case random graphs (p=.90). Runs faster for sparse graphs. f7c0bc, 49 lines

```
typedef vector<bitset<200>> vb;
struct Maxclique {
   double limit=0.025, pk=0;
   struct Vertex { int i, d=0; };
   typedef vector<Vertex> vv;
```

```
vb e;
vv V:
vector<vi> C;
vi qmax, q, S, old;
void init (vv& r) {
  for (auto& v : r) v.d = 0;
  for (auto& v : r) for (auto j : r) v.d \neq e[v.i][j.
  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);
    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;</pre>
        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)),
  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.

#### 7.7 Trees

#### BinaryLifting.h

**Description:** Calculate power of two jumps in a tree, to support fast upward jumps and LCAs. Assumes the root node points to itself. **Time:** construction  $\mathcal{O}(N \log N)$ , queries  $\mathcal{O}(\log_b N)$  in Eq. (2) lines

```
vector<vi> treeJump(vi& P) {
 int on = 1, d = 1;
 while (on < sz(P)) on *= 2, d++;
 vector<vi> jmp(d, P);
 rep(i,1,d) rep(j,0,sz(P))
    jmp[i][j] = jmp[i-1][jmp[i-1][j]];
 return jmp;
int jmp(vector<vi>& tbl, int nod, int steps){
 rep(i,0,sz(tbl))
    if(steps&(1<<i)) nod = tbl[i][nod];
int lca(vector<vi>& tbl, vi& depth, int a, int b) {
  if (depth[a] < depth[b]) swap(a, b);</pre>
  a = jmp(tbl, a, depth[a] - depth[b]);
 if (a == b) return a;
 for (int i = sz(tbl); i--;) {
    int c = tbl[i][a], d = tbl[i][b];
    if (c != d) a = c, b = d;
 return tbl[0][a];
```

#### LCA.

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

Time:  $O(N \log N + Q)$ 

```
"../data-structures/RMO.h"
                                            0f62fb, 21 lines
struct LCA {
 int T = 0;
  vi time, path, ret;
  RMQ<int> rmq;
 LCA(vector < vi > \& C) : time(sz(C)), rmq((dfs(C, 0, -1)),
  void dfs(vector<vi>& C, int v, int par) {
    time[v] = T++;
    for (int y : C[v]) if (y != par) {
     path.push_back(v), ret.push_back(time[v]);
      dfs(C, y, v);
  int lca(int a, int b) {
   if (a == b) return a;
    tie(a, b) = minmax(time[a], time[b]);
    return path[rmq.query(a, b)];
  //dist(a,b) {return depth[a] + depth[b] - 2*depth[lca(
       a,b)];}
```

#### CompressTree.h

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

"LCA.h" 9775a0, 21 lines

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

#### HLD.h

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

Time:  $\mathcal{O}\left((\log N)^2\right)$ 

```
"../data-structures/LazySegmentTree.h"
                                           9547af, 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)
    for (int& u : adj[v]) {
      adj[u].erase(find(all(adj[u]), v));
      par[u] = v;
      dfsSz(u);
      siz[v] += siz[u];
      if (siz[u] > siz[adj[v][0]]) swap(u, adj[v][0]);
```

vi seen(n, -1), path(n), par(n);

```
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 (;; v = par[rt[v]]) {
   if (pos[u] > pos[v]) swap(u, v);
   if (rt[u] == rt[v]) break;
   op(pos[rt[v]], pos[v] + 1);
 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] +
```

#### LCT.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: O(\log n)
                                          6758b4, 104 lines
struct Node {
 Node *p, *c[2];
 Node() \{ p = c[0] = c[1] = 0; \}
 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; }
 void push() {
   if (rev) {
     rep(i, 0, 2) if (c[i]) c[i]->flip();
     rev = 0:
   if (d) {
     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;
  // Add a virtual edge, cancel current virtual delta
  void vadd (Node* x) {
   vs += v->s:
   vsz += x->sz:
   x->cc = vd;
  // Delete a virutal edge, push the virtual delta
  void vdel(Node* x) {
   x \rightarrow add(vd - x \rightarrow cc):
    vs -= x->s;
   vsz -= x->sz;
   x->cc = 0;
  // Swap the cancels because of a splay rotation
  void vswap (Node* x, Node* y) {
    swap (x->cc, y->cc);
  int up() {
   if (!p) return -2;
    rep(i, 0, 2) if (p->c[i] == this) return i;
```

```
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 splav() {
   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) {
   makeRoot(&t[v]); access(&t[u]);
   setLink(&t[v], &t[u], 0); t[v].pull();
 void cut(int u, int 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 (v) x->vdel(v);
      if (x->c[1]) x->vadd(x->c[1]);
     x - c[1] = y; x - c[1]; y = x;
   u->splay();
 void makeRoot(Node* u) {
   access(u), u->flip(), u->push();
DirectedMST.h
Description: Finds a minimum spanning tree/arborescence
of a directed graph, given a root node. If no MST exists, re-
turns -1
Time: \mathcal{O}\left(E\log V\right)
 ../data-structures/UnionFindRollback.h"
                                            39e620, 60 lines
struct Edge { int a, b; ll w; };
struct Node
 Edge key;
 Node *1, *r:
 11 delta:
 void prop() {
   key.w += delta;
   if (1) 1->delta += delta;
   if (r) r->delta += delta;
   delta = 0:
 Edge top() { prop(); return key; }
Node *merge(Node *a, Node *b) {
 if (!a || !b) return a ?: b;
 a->prop(), b->prop();
 if (a->key.w > b->key.w) swap(a, b);
 swap(a->1, (a->r = merge(b, a->r)));
void pop(Node*& a) { a->prop(); a = merge(a->1, a->r);
pair<ll, vi> dmst(int n, int r, vector<Edge>& g) {
 RollbackUF uf(n);
 vector<Node*> heap(n);
```

for (Edge e : g) heap[e.b] = merge(heap[e.b], new

Node(e));

11 res = 0;

```
seen[r] = r:
  vector<Edge> Q(n), in(n, {-1,-1}), comp;
deque<tuple<int, int, vector<Edge>>> cycs;
  rep(s,0,n) {
    int u = s, qi = 0, w;
    while (seen[u] < 0) {
      if (!heap[u]) return {-1,{}};
      Edge e = heap[u]->top();
heap[u]->delta -= e.w, pop(heap[u]);
      Q[qi] = e, path[qi++] = u, seen[u] = s;
res += e.w, u = uf.find(e.a);
      if (seen[u] == s) {
        Node* cyc = 0;
        int end = qi, time = uf.time();
        do cyc = merge(cyc, heap[w = path[--qi]]);
        while (uf.join(u, w));
        u = uf.find(u), heap[u] = cyc, seen[u] = -1;
        cycs.push_front({u, time, {&Q[qi], &Q[end]}});
    rep(i,0,qi) in[uf.find(Q[i].b)] = Q[i];
 for (auto& [u,t,comp] : cycs) { // restore sol (
        optional)
    uf.rollback(t);
    Edge inEdge = in[u];
    for (auto& e : comp) in[uf.find(e.b)] = e;
    in[uf.find(inEdge.b)] = inEdge;
  rep(i,0,n) par[i] = in[i].a;
 return {res, par};
CentroidTree.h
Description: Centroid decomposition tree. Example usage
can be extended for weighted trees with a BST.
Time: O\left(n\log n + q\log^2 n\right)
"../data-structures/FenwickTree.h"
                                              fd433b, 60 lines
struct CT {
 vi sub, cp, d; // centroid subtree, parent, depth
 vector<vi> g, dst; // dst[depth][descendant]
 CT(vector<vi>& G): sub(sz(G)), cp(sz(G), -2),
d(sz(G)), q(G), dst(_lq(sz(G)) + 1, vi(sz(G))) {
    rec(0, 0):
 void dfs(int u, int p) {
   sub[u] = 1;
    for (int v : q[u]) if (v != p && cp[v] == -2)
      dfs(v, u), sub[u] += sub[v];
  void gen(int u, int p, int lev) {
    dst[lev][u] = dst[lev][p] + 1;

for (int v : g[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 : g[u]) if (cp[v] == -2)
  gen(v, u, d[u]), cp[rec(v, dd + 1)] = u;
    return ii:
 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(q)) d[i] = c[i] = FT(sub[i] + 1);
  // Add x to verts whose distance from p is in [1, r)
  void add(int p, int l, 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(int p) {
    11 ans = 0:
    path(p, [&](int u, int v, int dd) {
     ans += d[u].query(dd + 1);
     if (v != -1) ans -= c[v].query(dd + 1);
    return ans:
};
```

#### RerootDP.h

Description: Calculates a DP from every root in a tree. Use dp, rdp and p for edge dp. Time:  $\mathcal{O}(\sum d \log d)$ 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(q), t = 1;
  vi q(n), p(n);
  for (int u : q) for (int v : q[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 = a[i], k = -1;
     dp[u].init(u);
     rep(j, 0, sz(q[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 11 : a)
    int d = sz(q[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 >= 0; i--) e[i] = e[i / 2];
       Tor (int 1 - d - 1; 1 > -0; 1 - 1 e[1] - e[1 / 2] rep(i, 0, d - (d & !b)) {
    S& s = g[u][i] != p[u] ? dp[g[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:
```

#### 7.8 Math

#### 7.8.1 Number of Spanning Trees

Create an  $N \times N$  matrix mat, and for each edge  $a \to b \in G$ , do mat[a][b]--, mat[b][b]++ (and mat[b][a] --, mat[a][a] ++ if G is undirected). Remove the ith row and column and take the determinant; this yields the number of directed spanning trees rooted at i (if G is undirected, remove

#### 7ng.20w Erdősm Gallai theorem

A simple graph with node degrees  $d_1 > \cdots > d_n$ exists iff  $d_1 + \cdots + d_n$  is even and for every  $k=1\ldots n$ ,

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

Point lineDistance SegmentDistance SegmentIntersection lineIntersection sideOf UnSegment linearTransformation LineProjectionReflection Angle AngleOmp directedSegment CircleInterse UW

#### Geometry (8)

#### 8.1 Geometric primitives

#### Point.h

```
Description: Class to handle points in the plane. T can be
e.g. double or long long. (Avoid int.)
                                            3e64f3, 26 lines
template <class T> int sqn(T x) \{ return (x > 0) - (x < 0) \}
      0); }
template<class T>
struct Point {
 typedef Point P;
  T x, v:
```

```
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 sqrt((double)dist2()); }
// angle to x-axis in interval (-pi, pi)
double angle() const { return atan2(y, x); }
P unit() const { return *this/dist(); } // makes dist
P perp() const { return P(-y, x); } // rotates +90
      dearees
P normal() const { return perp().unit(); }
// returns point rotated 'a' radians ccw around the
      origin
P rotate (double a) const {
  return P(x*cos(a)-y*sin(a),x*sin(a)+y*cos(a)); }
friend ostream& operator<<(ostream& os, P p) {
  return os << "(" << p.x << "," << p.y << ")"; }</pre>
```

#### lineDistance.h

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

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

#### SegmentDistance.h

Description: Returns the shortest distance between point p and the line segment from point s to e. Usage: Point < double > a, b(2,2), p(1,1); bool onSegment = segDist(a,b,p) < 1e-10; 5c88f4, 6 lines

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

#### SegmentIntersection.h

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

```
so watch out for overflow if using int or long long.
Usage: vector <P > inter = segInter(s1,e1,s2,e2);
if (sz(inter) == 1)
cout << "segments intersect at " << inter[0] <<
endl;
"Point.h", "OnSegment.h"
                                             9d57f2, 13 lines
```

```
template<class P> vector<P> segInter(P a, P b, P c, P d | LineProjectionReflection.h
 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) };</pre>
 if (onSegment(c, d, a)) s.insert(a);
if (onSegment(c, d, b)) s.insert(b);
 if (onSegment(a, b, c)) s.insert(c);
 if (onSegment(a, b, d)) s.insert(d);
 return {all(s)};
```

#### lineIntersection.h

Description: If a unique intersection point of the lines going through s1,e1 and s2,e2 exists {1, point} is returned. If no intersection point exists  $\{0, (0,0)\}$  is returned and if infinitely many exists  $\{-1, (0,0)\}$  is returned. The wrong position will be returned if P is Point < ll> and the intersection point does not have integer coordinates. Products of three coordinates are used in intermediate steps so watch out for overflow if using int or 11 Usage: auto res = lineInter(s1,e1,s2,e2);

```
if (res.first == 1)
cout << "intersection point at " << res.second <<
endl;
"Point.h"
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;

```
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 on Segment (P s, P e, P p) {
 return p.cross(s, e) == 0 && (s - p).dot(e - p) <= 0;
```

#### linearTransformation.h Description:

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

```
03a306, 6 lines
typedef Point < double > P;
P linearTransformation(const P& p0, const P& p1,
   const P& q0, const P& q1, const P& r) {
 P dp = p1-p0, dq = q1-q0, num(dp.cross(dq), dp.dot(dq)
 return q0 + P((r-p0).cross(num), (r-p0).dot(num))/dp.
```

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 over-

```
b55<u>62d</u>, 5 lines
template<class P>
P lineProj(P a, P b, P p, bool refl=false) {
 P v = b - a;
 return p - v.perp()*(1+refl)*v.cross(p-a)/v.dist2();
```

#### Angle.h

sorted

Description: A class for ordering angles (as represented by int points and a number of rotations around the origin). Useful for rotational sweeping. Sometimes also represents points or vectors. **Usage:** vector < Angle  $> v = \{w[0], w[0], t360() ...\}; //$ 

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

```
++j; }
// sweeps j such that (j-i) represents the number of
positively oriented triangles with vertice 90f0602, 350dines
struct Angle {
 int x, y;
 int t;
 Angle(int x, int y, int t=0) : x(x), y(y), t(t) {}
 Angle operator-(Angle b) const { return {x-b.x, y-b.y
 int half() const {
    assert(x || y);
    return y < 0 | | (y == 0 && x < 0);
 Angle t90() const { return {-y, x, t + (half() && x
       >= 0) }; }
 Angle t180() const { return {-x, -y, t + half()}; }
Angle t360() const { return {x, y, t + 1}; }
bool operator < (Angle a, Angle b) {
 // add a.dist2() and b.dist2() to also compare
        distances
 return make_tuple(a.t, a.half(), a.y * (11)b.x) <
    make_tuple(b.t, b.half(), a.x * (11)b.y);</pre>
// Given two points, this calculates the smallest angle
      between
// them, i.e., the angle that covers the defined line
     seament.
pair<Angle, Angle> segmentAngles(Angle a, Angle b) {
 if (b < a) swap(a, b);
 return (b < a.t180() ?
          make_pair(a, b) : make_pair(b, a.t360()));
Angle operator+(Angle a, Angle b) { // point a + vector
 Angle r(a.x + b.x, a.y + b.y, a.t);
 if (a.t180() < r) r.t--;
 return r.t180() < a ? r.t360() : r;
Angle angleDiff(Angle a, Angle b) { // angle b - angle
 int tu = b.t - a.t; a.t = b.t;
 return {a.x*b.x + a.y*b.y, a.x*b.y - a.y*b.x, tu - (b
         < a) }:
```

## AngleCmp.h

Description: Sorts points in ascending order by angle within the interval  $(-\pi, \pi]$ . The point (0,0) has an angle of 0. Equivalent to sorting by atan2(y, x).

```
496e9b, 6 lines
template<class P>
bool angleCmp(P a, P b) {
 auto half = [](P p){ return sgn(p.y) ?: -sgn(p.x); };
int A = half(a), B = half(b);
```

return A == B ? a.cross(b) > 0 : A < B;

## directedSegment.h

Description: Segment representation usefull for sweeping. Compares two disjoint (can touch) segments on the sweep line (OY projection). Transitivity breaks if three segments touch in one point and are on different side of the sweep line. Can be easily fixed by comparing mean X coordinates.

```
template<class P>
struct dirSeg {
 P s, e; int rev;
 dirSeg(P _s, P _e) : s(_s), e(_e), rev(0) {
   if(e < s) swap(s, e), rev = 1;
 P getY(P X) { // takes x * 2, returns y * 2 as a
       fraction
    P d = (e - s);
    return ! sgn(d.x) ? P(s.y+e.y, 1) : P(d.cross(s*2-X)
         . d.x);
  int cmp(dirSeg b) { // needs ~64 * M^3 !
   P \times (\max(s.x, b.s.x) + \min(e.x, b.e.x), 0);
    return sgn(getY(X).cross(b.getY(X)));
```

#### 8.2 Circles

#### CircleIntersection.h

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

```
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 = \text{vec.dist2}(), sum = r1+r2, dif = r1-r2,
        p = (d2 + r1*r1 - r2*r2)/(d2*2), h2 = r1*r1 -
              p*p*d2;
  if (sum*sum < d2 || dif*dif > d2) return false;
  P mid = a + vec*p, per = vec.perp() * sqrt(fmax(0, h2
      ) / d2);
  *out = {mid + per, mid - per};
 return true;
```

#### CircleTangents.h

Description: Finds the external tangents of two circles, or internal if r2 is negated. Can return 0, 1, or 2 tangents - 0 if one circle contains the other (or overlaps it, in the internal case, or if the circles are the same); 1 if the circles are tangent to each other (in which case .first = .second and the tangent line is perpendicular to the line between the centers). .first and .second give the tangency points at circle 1 and 2 respectively. To find the tangents of a circle with a point set r2 to 0. b0153d, 13 lines

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

#### CircleLine.h

"Point.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>.

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\};
```

## CirclePolygonIntersection.h

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

```
Time: \mathcal{O}(n)
"../../content/geometry/Point.h"
                                                19add1, 19 lines
typedef Point<double> P;
#define arg(p, q) atan2(p.cross(q), p.dot(q))
double circlePoly(P c, double r, vector<P> ps) {
 auto tri = [&](P p, P q) {
  auto r2 = r * r / 2;
    P d = q - p;
    auto a = d.dot(p)/d.dist2(), b = (p.dist2()-r*r)/d.
          dist2();
    auto det = a * a - b;
    if (det <= 0) return arg(p, q) * r2;</pre>
    auto s = max(0., -a-sqrt(det)), t = min(1., -a+sqrt
          (det));
    if (t < 0 || 1 <= s) return arg(p, g) * r2;</pre>
    P u = p + d * s, v = q + d * (t-1);

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. 1caa3a, 9 lines

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

#### MinimumEnclosingCircle.h

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

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

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

#### CirclesUnionArea.h

Description: Returns the area of the sum of circles.

```
Time: \mathcal{O}\left(n^2 \log n\right)
```

"CircleIntersection.h" 2be987, 34 lines

```
template<typename T> // double or long double
T circlesArea(vector<pair<P, T>> c) {
 const T PI = acos((T)-1);
  sort(all(c)); c.erase(unique(all(c)), c.end());
   res = 0;
  for(auto &[p, r]: c) {
   int cnt = 0, cover = 0;
```

```
vector<pair<T, int>> eve = {{-PI, 0}};
  for(auto &[q, s]: c) if(make_pair(p, r) !=
       make_pair(q, s)) {
    T dst = (p - q).dist();
if(r + dst <= s) { cover = 1; break; }</pre>
    pair<P, P> inters;
    if(!circleInter(p, q, r, s, &inters)) continue;
    T le = (inters.first - p).angle();
    T re = (inters.second - p).angle();
    cnt += le > re:
    eve.pb({le, 1}), eve.pb({re, -1});
  if(cover) continue:
  sort(eve.begin() + 1, eve.end());
  eve.pb({PI, 0});
  T loc = 0;
  rep(i, 1, SZ(eve)) {
    if(!cnt) {
      T = eve[i-1].first, b = eve[i].first;
      loc += r * (b - a) +
        p.cross(P(cos(b)-cos(a), sin(b)-sin(a)));
    cnt += eve[i].second;
  res += r * loc;
return res / 2;
```

#### 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

```
Usage: vector < P > v = {P{4,4}, P{1,2}, P{2,1}};
bool in = inPolygon(v, P\{3, 3\}, false);
Time: \mathcal{O}(n)
"Point.h", "OnSegment.h", "SegmentDistance.h"
                                            2bf504, 11 lines
template<class P>
bool inPolygon(vector<P> &p, P a, bool strict = true)
 int cnt = 0, n = sz(p);
 rep(i,0,n) {
   P q = p[(i + 1) % n];
   if (onSegment(p[i], q, a)) return !strict;
    //or: if (segDist(p[i], q, a) <= eps) return !
   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:  $\mathcal{O}(n)$ 

```
9706dc, 9 lines
typedef Point<double> P;
 polygonCenter(const vector<P>& v) {
  porygoneence: (coins)
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]);</pre>
     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));
                                           d07181, 13 lines
typedef Point < double > P;
vector<P> polygonCut (const vector<P>& poly, P s, P e) {
 vector<P> res;
 rep(i,0,sz(polv)) {
   P cur = poly[i], prev = i ? poly[i-1] : poly.back()
    auto a = s.cross(e, cur), b = s.cross(e, prev);
    if ((a < 0) != (b < 0))
      res.push_back(cur + (prev - cur) * (a / (a - b)))
    if (a < 0)
      res.push_back(cur);
 return res:
```

#### 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
```

```
typedef Point < double > P;
double rat(P a, P b) { return sgn(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])) {
   \bar{P} A = poly[i][v], \bar{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 && sgn((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 = segs[0].second;
    rep(j,1,sz(segs)) {
      if (!cnt) sum += segs[j].first - segs[j - 1].
           first;
      cnt += segs[j].second;
    ret += A.cross(B) * sum;
 return ret / 2;
```

## PolygonTangents.h

Description: Polygon tangents from a given point. The polygon must be ccw and have no collinear points. Returns a pair of indices of the given polygon. Should work for a point on border (for a point being polygon vertex returns previous and next one).

```
Time: O(\log n)
                                           096fab, 21 lines
#define pdir(i) (ph ? p - poly[(i)%n] : poly[(i)%n] - p
#define cmp(i, j) sqn(pdir(i).cross(poly[(i)%n]-poly[(j)
#define extr(i) cmp(i + 1, i) >= 0 && cmp(i, i - 1 + n)
template <class P>
array<int, 2> polygonTangents(vector<P>& poly, P p) {
 auto bs = [&](int ph) {
    int n = sz(poly), lo = 0, hi = n;
```

```
if(extr(0)) return 0;
 while (lo + 1 < hi) {
    int m = (lo + hi) / 2:
   if(extr(m)) return m;
    int ls = cmp(lo + 1, lo), ms = cmp(m + 1, m);
    (ls < ms \mid | (ls == ms && ls == cmp(lo, m)) ? hi:
         lo) = m;
 return lo;
array<int, 2> res = {bs(0), bs(1)};
if(res[0] == res[1]) res[0] = (res[0] + 1) % sz(poly)
if(poly[res[0]] == p) res[0] = (res[0] + 1) % sz(poly
return res;
```

#### ConvexHull.h

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

```
310954, 13 lines
```

```
typedef Point<11> P;
vector<P> convexHull(vector<P> pts) {
 if (sz(pts) <= 1) return pts;
  sort (all (pts));
  vector<P> h(sz(pts)+1);
  int s = 0, t = 0;
  for (int it = 2; it--; s = --t, reverse(all(pts)))
    for (P p : pts) {
     while (t >= s + 2 \&\& h[t-2].cross(h[t-1], p) <=
           0) t--;
     h[t++] = p;
 return {h.begin(), h.begin() + t - (t == 2 && h[0] ==
        h[1])};
```

#### ConvexHullOnline.h

Description: Allows online point insertion. If exists, left vertical segment is included; right one is excluded. To get a lower hull add (-x, -y) instead of (x, y).

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

```
10c55b, 16 lines
using P = Point<11>;
struct UpperHull : set<P> {
  bool rm(auto it) {
    if (it==begin() || it==end() || next(it)==end() ||
        it->cross(*prev(it), *next(it)) > 0)
      return false;
    erase(it); return true;
 bool add(P p) { // true iff added
auto [it, ok] = emplace(p);
    if (!ok || rm(it)) return false;
    while (rm(next(it)));
    while (it != begin() && rm(prev(it)));
    return true:
```

#### ConvexHullOnline2.h

1), i});

sort(all(pts));

};

Description: Fully dynamic upper / lower convex hull, can be used for computing onion layers. All points should be known in advance. Points on the edges are included in the hull. Return indices are the same as in the input.

```
Time: \mathcal{O}(\log^2 n), as fast as other \mathcal{O}(\log n) hulls
```

```
78db53, 70 lines
template<class T>
struct DynHull {
  using P = Point<T>;
  struct Node { int 1, r; }; vector<P> ps;
  int n; vi in, id; int s; vector<Node> t; vector<T> m;
  DynHull(vector<P> _ps, bool lower = 0, int start =
    : ps(_ps), n(sz(ps)), in(n), id(n) {
    if(start == -1) start = n;
    s = 1; while(s < n) s *= 2;
    t.resize(s * 2, \{-1, -1\}); m.resize(s);
    vector<pair<P, int>> pts;
    rep(i, 0, n) pts.push_back({ps[i] * (lower ? -1 :
```

```
rep(i, 0, n) {
      Lep(1, 0, n) {
    tie(ps[i], id[i]) = pts[i]; in[id[i]] = i;
    int p = i + s; while((p & 1) ^ 1) p >>= 1;
    m[p >> 1] = ps[i].x;
    rep(i, 0, start) t[s + in[i]] = {in[i], in[i]};
for (int i = s - 1; i >= 1; i --) pull(i);
    while (t[v].1 < 0) v = v * 2 + t[v].1 + 3;
    return v:
  void pull(int v) {
    auto crossNegX = [](Pa, Pb, Pc, Pd, Tx) {
   // change __int128 if using doubles!
         _{int128} p = a.cross(b, c), q = b.cross(a, d);
       return p + q == 0 || (d.x - x) * p + (c.x - x) *
             q <= 0;
     int p = v * 2, q = p + 1;
    else if(t[q].1 == -1) t[v] = {-1, -1};
else if(t[q].1 == -1) t[v] = {-2, -2};
else if(t[q].1 == -1) t[v] = {-3, -3};
     else {
      p = go(p), q = go(q);
       while(p < s || q < s) {
         auto [a, b] = t[p]; auto [c, d] = t[q];
         if (a != b && ps[a].cross(ps[b], ps[c]) > 0) {
           p = qo(p * 2); }
         else if(c != d && ps[b].cross(ps[c], ps[d]) >
               0) {
           q = qo(q * 2 + 1); }
         else if (a == b) q = go(q * 2);
         else if(c == d ||
           crossNegX(ps[a], ps[b], ps[c], ps[d], m[v]))
           p = go(p * 2 + 1); }
         else q = go(q * 2);
      t[v] = \{p - s, q - s\};
  void add(int i) {
    i = in[i]; int v = i + s; t[v] = {i, i};
     while(v >>= 1) pull(v);
  void del(int i) {
    i = in[i]; int v = i + s; t[v] = \{-1, -1\};
     while(v >>= 1) {
      if(t[v].1 < 0 || t[v].1 == i || t[v].r == i) pull</pre>
              (v); }
  void dfs(int v, int l, int r, vi &h) {
    if(v >= s) return h.push_back(id[t[v].1]);
    if(1 \le t[v].1) dfs(go(v * 2), 1, min(t[v].1, r), h
     if(t[v].r \le r) dfs(go(v * 2 + 1), max(t[v].r, 1),
          r, h);
  vi hull() {
    vi h; if(~t[1].1) dfs(go(1), 0, n - 1, h); return h
};
```

#### HullDiameter.h

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

c571b8, 12 lines

```
typedef Point<ll> P:
arrav<P, 2> hullDiameter(vector<P> S) {
 int n = sz(S), j = n < 2 ? 0 : 1;
pair<11, 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 con-

```
if point lies within the hull. If strict is true, points on the
boundary aren't included.
Time: O(\log N)
"Point.h", "sideOf.h", "OnSegment.h"
                                                 71446b, 14 lines
typedef Point<ll> P;
bool inHull(const vector<P>& 1, P p, bool strict = true
 int a = 1, b = sz(l) - 1, r = !strict;
if (sz(l) < 3) return r && onSegment(l[0], l.back(),</pre>
        ; (a
  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) \ll -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(l[a].cross(l[b], p)) < r;</pre>
```

vex hull (CCW order, with no collinear points). Returns true

#### LineHullIntersection.h

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

```
Time: \mathcal{O}(\log n)
"Point.h"
                                              7cf45b, 39 lines
#define cmp(i,j) sqn(dir.perp().cross(poly[(i)%n]-poly
#define extr(i) cmp(i + 1, i) >= 0 && cmp(i, i - 1 + n)
template <class P> int extrVertex(vector<P>& polv, P
    dir) {
  int n = sz(polv), lo = 0, hi = n;
  if (extr(0)) return 0;
  while (lo + 1 < hi) {
    int m = (lo + hi) / 2;
    if (extr(m)) return m;
    int ls = cmp(lo + 1, lo), ms = cmp(m + 1, m);
    (ls < ms | | (ls == ms && ls == cmp(lo, m)) ? hi :
          lo) = m;
  return lo:
#define cmpL(i) sqn(a.cross(polv[i], b))
template <class P>
array<int, 2> lineHull(P a, P b, vector<P>& polv) {
 int endA = extrVertex(poly, (a - b).perp());
int endB = extrVertex(poly, (b - a).perp());
  if (cmpL(endA) < 0 || cmpL(endB) > 0)
   return {-1, -1};
  array<int, 2> res;
  rep(i,0,2) {
    int lo = endB, hi = endA, n = sz(poly);
    while ((lo + 1) % n != hi) {
      int m = ((lo + hi + (lo < hi ? 0 : n)) / 2) % n;</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; Minkowski.h

Description: Computes Minkowski sum of two convex polygons in ccw order. Vertices are required to be in ccw order. Time:  $\mathcal{O}(n+m)$ 7d8a28, 18 lines "Point.h", "Angle.h"

```
P edgeSeg(vector<P> p, vector<P>& edges) {
 int i = 0, n = sz(p);
 rep(j, 0, n) if (tie(p[i].y, p[i].x) > tie(p[j].y, p[
        j].x)) i = j;
  rep(j, 0, n) edges.push_back(p[(i+j+1)%n] - p[(i+j)%n
 return p[i];
vector<P> hullSum(vector<P> A, vector<P> B) {
  vector<P> sum, e1, e2, es(sz(A) + sz(B));
  P pivot = edgeSeq(A, e1) + edgeSeq(B, e2);
 merge(all(e1), all(e2), es.begin(), [&](Pa, Pb){
    return Angle(a.x, a.y) < Angle(b.x,b.y);
  sum.push back(pivot);
 for (auto e: es) sum.push_back(sum.back() + e);
  sum.pop_back();
 return sum; //can have collinear vertices!
HalfplaneIntersection.h
Description: Online half plane intersection. Works both for
ll and long double. Bounding box is optional, but needed for
distinguishing bounded vs unbounded. Halfplanes are sorted
ccw in HPI.s. Time: O(log n) per add.
 Ta,b,c;
       //ax + by + c >= 0 (coords <= 10^9)
```

```
5e6600, 98 lines
using T = 11; // has to fit 2*|pts|**2
using P = Point<T>; // only cross needed
using SuperT = __int128_t; // has to fit 6*/pts/**3
const SuperT EPS = 1e-12; // |pts/ <= 10^6 (for T=dbl)</pre>
  Line (T a_=0, T b_=0, T c_=0): a(a_), b(b_), c(c_) {}
  Line (P p, P q): a(p.y-q.y), b(q.x-p.x), c(p.cross(q))
 {} //p->q cw (coords <= 10^6)
Line operator- () const {return Line(-a, -b, -c); }
bool up() const { return a?(a<0):(b>0);}
  P v() const {return P(a,b);}
  P vx() {return P(b,c);} P vy() {return P(a,c);}
  T wek(Line p) const {return v().cross(p.v());}
  bool operator<(Line b) const {
    if (up() != b.up()) return up() > b.up();
    return wek(b) > 0;
bool parallel (Line a, Line b) {return !a.wek(b);}
bool same (Line a, Line b) {
 return parallel(a,b) && !a.vy().cross(b.vy()) && !a.
       vx().cross(b.vx()):
T weaker (Line a. Line b) {
 if (abs(a.a) > abs(a.b)) return a.c*abs(b.a) - b.c*
       abs(a.a):
  return a.c*abs(b.b) - b.c*abs(a.b);
array<SuperT, 3> intersect(Line a, Line b) {
 SuperT det = a.wek(b):
  SuperT x = a.vx().cross(b.vx());
  SuperT y = a.vy().cross(b.vy());
  // if (T=dbl) return {x / det, -y / det, 1.0};
  if (det > 0) return {x, -y, det};
  return {-x, y, -det};
struct HPI {
 bool empty=0, pek=0;
  set < Line > s:
  typedef set<Line>::iterator iter;
  iter next(iter it) {return ++it == s.end() ? s.begin()
         : it;}
  iter prev(iter it) {return it == s.begin() ? --s.end()
         : --it;}
 bool hide (Line a, Line b, Line c) { // do a, b hide c?
    if (parallel(a,b)) {
      if (weaker(a, -b) < 0) empty = 1;
      return 0:
    if (a.wek(b) < 0) swap(a,b);
    auto [rx, ry, rdet] = intersect(a,b);
    auto v = rx*c.a + ry*c.b + rdet*c.c;
    if (a.wek(c) >=0 && c.wek(b) >=0 && v >= -EPS)
         return 1;
    if (a.wek(c) < 0 && c.wek(b) < 0) {
      if (v < -EPS) empty = 1;
      else if (v <= EPS) pek = 1;
    return 0;
  void delAndMove(iter& i, int nxt) {
    iter j = i;
```

```
if(nxt==1) i = next(i);
    else i = prev(i):
    s.erase(j);
  void add(Line 1) {
   if (empty) return;
    if (1.a == 0 && 1.b == 0) {
     if (1.c < 0) empty = 1;
      return;
    iter it = s.lower_bound(1); //parallel
    if(it != s.end() && parallel(*it, 1) && it->up() ==
          1.up()) {
      if (weaker(1, *it)>=0) return;
      delAndMove(it,1);
   if(it == s.end()) it = s.begin(); //*it>p
while(sz(s) >= 2 && hide(l, *next(it), *it))
      delAndMove(it,1);
    if(sz(s)) it = prev(it); //*it < p
    while (sz(s) \ge 2 && hide(1, *prev(it), *it))
     delAndMove(it,0);
    if(sz(s) < 2 \mid \mid !hide(*it, *next(it), 1)) s.insert(
  int type() { // 0=empty, 1=point, 2=segment,
    if(empty) return 0; // 3=halfline, 4=line,
    if(sz(s) <= 4) { // 5=polygon or unbounded
      vector<Line> r(all(s));
      if(sz(r) == 2 \&\& parallel(r[0], r[1]) \&\& weaker(r
           [0],-r[1])<0)
        return 0;
      rep(i, 0, sz(r)) rep(j, 0, i) if(same(r[i], r[j])
        if(sz(r) == 2) return 4;
        if(sz(r) == 3) return 3;
        if(sz(r) == 4 \&\& same(r[0], r[2]) \&\& same(r[1],
              r[3])) return 1;
        return 2:
      if(sz(r) == 3 && pek) return 1;
    return 5;
};
```

#### PointLocation.h

**Description:** Computes (not necessarily convex) polygon tree structure. Also for each query point computes its location (including boundaries).

```
Time: \mathcal{O}(n \log n)
"directedSegment.h"
template<class P>
pair<vi, vi> pointLoc(vector<vector<P>> polys, vector<P>
      pts) {
  vector<tuple<P, int, int>> eve; // {point, event_type
       , id}
  vector<pair<dirSeg<P>, int>> segs; // {s, e, poly_id
  rep(i, 0, sz(polys)) rep(j, 0, sz(polys[i])) {
    dirSeg<P> seg(polys[i][j], polys[i][(j+1)%sz(polys[
         il)]);
    eve.push_back({seg.s,0,sz(segs)}), eve.push_back({
         seg.e,2,sz(segs)});
    segs.push_back({seg, i});
  rep(i, 0, sz(pts)) eve.push_back({pts[i], 1, i});
  sort(all(eve));
  vi par(sz(polys), -2), ans(sz(pts), -1);
  auto cmp = [](auto a, auto b) {
    return make_pair(a.first.cmp(b.first), a.second) <</pre>
         make_pair(0, b.second);
  set<pair<dirSeq<P>, int>, decltype(cmp)> s(cmp);
  for (auto &[_, eve_tp, id]: eve) {
   if (eve_tp == 1) { // point query}
      P p = pts[id];
      auto it = s.lower_bound({dirSeg(p, p), 0});
      if(it != s.begin()) { // on vertical segment?
        auto prv = prev(it);
        if(!sgn(p.cross(prv->first.s, prv->first.e)))
```

if(it == s.end()) ans[id] = -1;

strictness there!

int poly\_id = segs[seg\_id].second; //

ans[id] = !seg.rev && sgn(p.cross(seg.s, seg.e)

auto [seg, seg\_id] = \*it;

else {

#### ClosestPair ManhattanMST kdTree FastDelaunay PolyhedronVolume Point3D 3dHull

```
? par[poly_id] : poly_id;
  if(eve_tp == 0) { // add segment
    auto it = next(s.insert({segs[id].first, id}).
          first):
    int poly_id = segs[id].second;
if(par[poly_id] == -2) {
   if(it == s.end()) par[poly_id] = -1;
        int up_rev = it->first.rev, up_id = segs[it->
               secondl.second:
        par[poly_id] = !up_rev ? par[up_id] : up_id;
  if(eve_tp == 2) s.erase({seqs[id].first, id}); //
        del seament
return {par, ans};
```

#### 8.4 Misc. Point Set Problems

#### ClosestPair.h

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

```
ac41a6, 17 lines
typedef Point<11> P;
pair<P, P> closest(vector<P> v) {
 assert(sz(v) > 1);
 sort(all(v), [](P a, P b) { return a.v < b.v; });
 pair<ll, pair<P, P>> ret{LLONG_MAX, {P(), P()}};
 int j = 0;
  for (P p : v) {
   P d{1 + (ll)sqrt(ret.first), 0};
    while (v[j].y <= p.y - d.x) S.erase(v[j++]);</pre>
   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);
```

#### ManhattanMST.h

return ret.second;

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.yq.y-. Edges are in the form (distance, src, dst). Use a standard MST algorithm on the result to find the final MST. Time:  $\mathcal{O}(N \log N)$ 

```
df6f59, 23 lines
"Point.h"
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;});
    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) bac5b0, 63 lines

```
typedef long long T;
typedef Point<T> P;
const T INF = numeric limits<T>::max();
bool on_x(const P& a, const P& b) { return a.x < b.x; }
bool on v(const P& a, const P& b) { return a.v < b.v; }
```

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

#### Fast Delaunav.h

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

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

```
typedef Point<ll> P;
typedef struct Quad* 0;
typedef __int128_t lll; // (can be ll if coords are < 2</pre>
P arb(LLONG_MAX, LLONG_MAX); // not equal to any other
     point
struct Ouad {
 Q rot, o; P p = arb; bool mark;
 P& F() { return r()->p; }
 Q& r() { return rot->rot; }
 Q prev() { return rot->o->rot;
 Q next() { return r()->prev(); }
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)*C + p.cross(b,c)*A + p.cross(c,a)
        *B > 0:
Q makeEdge(P orig, P dest) {
 Q r = H ? H : new Quad{new Quad{new Quad{new Quad{0}}}
```

```
H = r -> 0; r -> r() -> r() = r;
 rep(i,0,4) r = r->rot, r->p = arb, r->o = i & 1 ? r :
        r->r();
 r->p = orig; r->F() = dest;
 return r:
void splice(Q a, Q b) {
 swap(a->o->rot->o, b->o->rot->o); swap(a->o, b->o);
Q connect(Q a, Q b) {
 Q q = makeEdge(a->F(), b->p);
 splice(q, a->next());
 splice(q->r(), b);
 return a:
pair<Q,Q> rec(const vector<P>& s) {
 if (sz(s) <= 3) {
    Q a = makeEdge(s[0], s[1]), b = makeEdge(s[1], s.
    back());
if (sz(s) == 2) return { a, a->r() };
    splice(a->r(), b);
    auto side = s[0].cross(s[1], s[2]);
    Q c = side ? connect(b, a) : 0;
    return {side < 0 ? c->r() : a, side < 0 ? c : b->r
          () };
#define H(e) e->F(), e->p
#define valid(e) (e->F().cross(H(base)) > 0)
 Q A, B, ra, rb;
 int half = sz(s) / 2;
tie(ra, A) = rec({all(s) - half});
tie(B, rb) = rec({sz(s) - half + all(s)});
  while ((B->p.cross(H(A)) < 0 && (A = A->next())) | |
         (A->p.cross(H(B)) > 0 && (B = B->r()->o)));
 O base = connect(B->r(), A);
 if (A->p == ra->p) ra = base->r();
 if (B->p == rb->p) rb = base;
#define DEL(e, init, dir) Q e = init->dir; if (valid(e)
    while (circ(e->dir->F(), H(base), e->F())) { \
     Q t = e->dir; \
      splice(e, e->prev()); \
      splice(e->r(), e->r()->prev()); \
      e->0 = H; H = e; e = t; \setminus
    DEL(LC, base->r(), o); DEL(RC, base, prev());
if (!valid(LC) && !valid(RC)) break;
    if (!valid(LC) || (valid(RC) && circ(H(RC), H(LC)))
      base = connect(RC, base->r());
    else
     base = connect(base->r(), LC->r());
 return { ra, rb };
vector<P> triangulate(vector<P> pts) {
 sort(all(pts)); assert(unique(all(pts)) == pts.end()
 if (sz(pts) < 2) return {};
 Q e = rec(pts).first;
  vector < Q > q = \{e\};
 int gi = 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 3D

#### Polyhedron Volume.h

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

```
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

```
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.h
Description: Computes all faces of the 3-dimension hull of a
point set. *No four points must be coplanar*, or else random
results will be returned. All faces will point outwards.
Time: \mathcal{O}\left(n^2\right)
"Point3D.h"
                                              5b45fc, 49 lines
typedef Point3D<double> P3;
struct PR {
  void ins(int x) { (a == -1 ? a : b) = x; }
  void rem(int x) { (a == x ? a : b) = -1; }
int cnt() { return (a != -1) + (b != -1); }
  int a, b;
struct F { P3 q; int a, b, c; };
vector<F> hull3d(const vector<P3>& A) {
 assert(sz(A) >= 4);
  vector<vector<PR>> E(sz(A), vector<PR>(sz(A), {-1, -1
        }));
#define E(x,y) E[f.x][f.y]
  vector<F> FS;
auto mf = [&](int i, int j, int k, int l) {
    P3 q = (A[j] - A[i]).cross((A[k] - A[i]));
    if (q.dot(A[1]) > q.dot(A[i]))
    q = q * -1;
F f{q, i, j, k};
    E(a,b).ins(k); E(a,c).ins(j); E(b,c).ins(i);
    FS.push back(f):
  rep(i,0,4) rep(j,i+1,4) rep(k,j+1,4)
    mf(i, j, k, 6 - i - j - k);
  rep(i,4,sz(A)) {
    rep(j,0,sz(FS)) {
      F f = FS[j];

if(f.q.dot(A[i]) > f.q.dot(A[f.a])) {
        E(a,b).rem(f.c);
        E(a,c).rem(f.b);
        E(b,c).rem(f.a);
        swap(FS[j--], FS.back());
        FS.pop_back();
    int nw = sz(FS);
    rep(j,0,nw) {
      F f = FS[j];
```

#define C(a, b, c) if (E(a,b).cnt() != 2) mf(f.a, f.b,

template<class T> struct Point3D {

typedef Point3D P: typedef const P& R:

```
C(a, b, c); C(a, c, b); C(b, c, a);
}
for (F& it : FS) if ((A[it.b] - A[it.a]).cross(
    A[it.c] - A[it.a]).dot(it.q) <= 0) swap(it.c, it.b)
return FS;
};</pre>
```

#### sphericalDistance.h

Description: Returns the shortest distance on the sphere with radius radius between the points with azimuthal angles (longitude) f1  $(\phi_1)$  and f2  $(\phi_2)$  from x axis and zenith angles (latitude) t1  $(\theta_1)$  and t2  $(\theta_2)$  from z axis (0= north pole). All angles measured in radians. The algorithm starts by converting the spherical coordinates to cartesian coordinates so if that is what you have you can use only the two last rows. dx\*radius is then the difference between the two points in the x direction and d\*radius is the total distance between the points7, 8 lines

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

#### Strings (9)

#### KMP.h

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

```
Time: O(n)

vi pi(const string& s) {
  vi p(sz(s));
  rep(i,1,sz(s)) {
    int g = p[i-1];
    while (g && s[i] != s[g]) g = p[g-1];
    p[i] = g + (s[i] == s[g]);
  }
  return p;
}
vi match(const string& s, const string& pat) {
  vi p = pi(pat + '\0' + s), res;
  rep(i,sz(p)-sz(s),sz(p))
    if (p[i] == sz(pat)) res.push_back(i - 2 * sz(pat))
    return res;
```

#### Zfunc.h

 $\begin{array}{ll} \textbf{Description:} \ z[i] \ computes the length of the longest common \\ prefix of s[i:] \ and s \ (abacaba -> 7010301). \\ \textbf{Time:} \ \mathcal{O}\left(n\right) \\ & 584523, \ 13 \ lines \end{array}$ 

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

```
p[z][i]++, L--, R++;
if (R>r) l=L, r=R;
}
return p;
```

#### MainLorentz.h

"Zfunc.h"

**Description:** Main-Lorentz algorithm for finding all squares in given word; Results are in compressed form: (b, e, l) means that for each b <= i < e there is square at position i of size 2l. Each square is present in only one interval. **Time:**  $\mathcal{O}(nlqn)$ 

```
struct Sqr {
 int begin, end, len;
vector<Sqr> lorentz(const string &s) {
 vector<Sqr> ans;
 vi pos(sz(s) / 2 + 2, -1);
 rep(mid, 1, sz(s)) {
    int part = mid & \sim(mid - 1), off = mid - part;
    int end = min(mid + part, sz(s));
    auto a = s.substr(off, part);
    auto b = s.substr(mid, end - mid);
    string ra(a.rbegin(), a.rend());
    string rb(b.rbegin(), b.rend());
    rep(j, 0, 2) {
        Set # to some unused character!
      vi z1 = Z(ra);
      vi z2 = Z(b + "#" + a);
      z1.push_back(0);
      z2.push_back(0);
      rep(c, 0, sz(a)) {
        int l = sz(a) - c;
        int x = c - min(1 - 1, z1[1]);
        int y = c - max(1 - z2[sz(b) + c + 1], j);
        if (x > y)
         continue;
        int sb = (j ? end - y - 1 * 2 : off + x);
        int se = (j ? end - x - 1 * 2 + 1 : off + y +
             1);
        int &p = pos[1];
        if (p != -1 && ans[p].end == sb)
         ans[p].end = se;
         p = sz(ans), ans.push back({sb, se, 1});
      a.swap(rb);
     b.swap(ra);
 return ans;
```

#### Lvndon.h

**Description:** Compute Lyndon factorization for s; Word is simple iff it's stricly smaller than any of it's nontrivial suffixes. Lyndon factorization is division of string into non-increasing simple words. It is unique.

```
Time: \mathcal{O}(n) fa3adf, 12 lines

vector<string> duval(const string &s) {
  int n = sz(s), i = 0;
  vector<string> ret;
  while (i < n) {
    int j = i + 1, k = i;
    while (j < n && s[k] <= s[j])
    k = (s[k] < s[j]? i : k + 1), j++;
  while (i <= k)
    ret.push_back(s.substr(i, j - k)), i += j - k;
  }

return ret;
```

#### MinRotation.h

 $\bf Description:$  Finds the lexicographically smallest rotation of a string.

```
return a;
```

#### SuffixArray.h

**Description:** Builds suffix array for a string. sa[i] is the starting index of the suffix which is i'th in the sorted suffix array. The returned vector is of size n+1, and sa[0]=n. The lcp array contains longest common prefixes for neighbouring strings in the suffix array: lcp[i] = lcp(sa[i], sa[i-1]), lcp[0] = 0. The input string must not contain any nul chars. **Time:**  $\mathcal{O}(n \log n)$  635552, 22 lines

```
e01cd1, 46 lines | struct SuffixArray {
                vi sa. lcp:
                SuffixArray(string s, int lim=256) { // or vector<int
                  s.push_back(0); int n = sz(s), k = 0, a, b;
                  vi x(all(s)), y(n), ws(max(n, lim));
                  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];
                    : p++;
                  for (int i = 0, j; i < n - 1; lcp[x[i++]] = k)
                    for (k && k--, j = sa[x[i] - 1];
    s[i + k] == s[j + k]; k++);
              };
```

#### SuffixTree.h

pii best;

]) : 0;

**Description:** Ukkonen's algorithm for online suffix tree construction. Each node contains indices [1, r] into the string, and a list of child nodes. Suffixes are given by traversals of this tree, joining [1, r] substrings. The root is 0 (has 1 = -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:  $\mathcal{O}(26N)$ 

```
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]
           ]; }
      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)
```

int lcs(int node, int i1, int i2, int olen) {

mask |= lcs(t[node][c], i1, i2, len);

rep(c, 0, ALPHA) if (t[node][c] != -1)

if (1[node] <= i1 && i1 < r[node]) return 1;
if (1[node] <= i2 && i2 < r[node]) return 2;</pre>

int mask = 0, len = node ? olen + (r[node] - 1[node

```
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;
}
```

#### Hashing.h

```
Description: Self-explanatory methods for stringalershingnes
// Arithmetic mod 2^64-1. 2x slower than mod 2^64 and
// code, but works on evil test data (e.g. Thue-Morse,
     where
// ABBA... and BAAB... of length 2^10 hash the same mod
// "typedef ull H;" instead if you think test data is
     random,
// or work mod 10^9+7 if the Birthday paradox is not a
     problem.
typedef uint64_t ull;
struct H {
 ull x; H(ull x=0) : x(x) {}
 H operator+(H o) { return x + o.x + (x + o.x < x); }
   operator-(H o) { return *this + ~o.x; }
  H operator*(H o) { auto m = (_uint128_t)x * o.x;
    return H((ull)m) + (ull)(m >> 64); }
  ull get() const { return x + !~x; }
 bool operator==(H o) const { return get() == o.get();
 bool operator<(H o) const { return get() < o.get(); }</pre>
static const H C = (11)1e11+3; // (order ~ 3e9; random
     also ok)
struct HashInterval {
  vector<H> ha, pw:
  HashInterval(string& str) : ha(sz(str)+1), pw(ha) {
    pw[0] = 1:
    ren(i.0.sz(str))
     ha[i+1] = ha[i] * C + str[i],
     pw[i+1] = pw[i] * C;
  H hashInterval(int a, int b) { // hash [a, b)
    return ha[b] - ha[a] * pw[b - a];
vector<H> getHashes(string& str, int length) {
 if (sz(str) < length) return {};</pre>
 H h = 0, pw = 1;
 rep(i,0,length)
   h = h * C + str[i], pw = pw * C;
  vector<H> ret = {h};
 rep(i,length,sz(str)) {
    ret.push\_back(h = h * C + str[i] - pw * str[i-
         length]);
```

#### Hash61.h

return ret;

return h:}

Description: Arithmetic for fast hashing modulo  $2^{61} - 1$  (prime).

Time: About 30% faster than naive modulo.

51cf65, 8 lines

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

```
const 11 MOD = (111 << 61) - 1;
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); }
11 mul(11 a, 11 b) {
   auto c = (_int128) a * b;
   return add(c & MOD, c >> 61);
```

#### AhoCorasick.h

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

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

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

#### Various (10)

#### 10.1 Intervals

#### IntervalContainer.h

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

```
Time: O(\log N)
                                          edce47, 23 lines
set<pii>::iterator addInterval(set<pii>& is, int L, int
      R) {
 if (L == R) return is.end();
 auto it = is.lower_bound({L, R}), before = it;
  while (it != is.end() && it->first <= R) {
   R = max(R, it->second);
   before = it = is.erase(it);
 if (it != is.begin() && (--it)->second >= L) {
   L = min(L, it->first);
   R = max(R, it->second);
   is.erase(it):
 return is.insert(before, {L,R});
void removeInterval(set<pii>& is, int L, int R) {
 if (L == R) return;
 auto it = addInterval(is, L, R);
 auto r2 = it->second;
 if (it->first == L) is.erase(it);
 else (int&)it->second = L;
 if (R != r2) is.emplace(R, r2);
```

#### IntervalCover.h

```
template<class T>
vi cover(pair<T, T> G, vector<pair<T, T>> I) {
 vi S(sz(T)). R:
 iota(all(S), 0);
 sort(all(S), [&](int a, int b) { return I[a] < I[b];</pre>
      });
 T cur = G.first;
 int at = 0;
 while (cur < G.second) { // (A)
   pair<T, int> mx = make_pair(cur, -1);
   while (at < sz(I) && I[S[at]].first <= cur) {
     mx = max(mx, make_pair(I[S[at]].second, S[at]));
     at++;
   if (mx.second == -1) return {};
   cur = mx.first:
   R.push back(mx.second);
 return R:
```

#### ConstantIntervals.h

Time:  $\mathcal{O}\left(k\log\frac{n}{L}\right)$ 

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

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

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

```
rec(from, to-1, f, g, i, p, q); g(i, to, q);
```

# 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:  $O(N^2)$ 

```
DivideAndConquerDP.h
```

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

```
d38d2b, 18 lines
struct DP { // Modify at will:
 int lo(int ind) { return 0; }
 int hi(int ind) { return ind; }
 11 f(int ind, int k) { return dp[ind][k]; }
 void store(int ind, int k, ll v) { res[ind] = pii(k,
      v); }
 void rec(int L, int R, int LO, int HI) {
   if (L >= R) return;
   int mid = (L + R) >> 1;
   pair<11, int> best (LLONG_MAX, LO);
   rep(k, max(LO, lo(mid)), min(HI, hi(mid)))
     best = min(best, make pair(f(mid, k), k));
   store (mid, best.second, best.first);
   rec(L, mid, LO, best.second+1);
   rec(mid+1, R, best.second, HI);
 void solve (int L, int R) { rec(L, R, INT MIN, INT MAX
```

#### 10.3 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: O(log(b - a)) 9155b4, 11 lines
template < class F >
int ternSearch(int a, int b, F f) {
    assert(a <= b);
    while (b - a >= 5) {
        int mid = (a + b) / 2;
        if (f (mid) < f (mid) + 1) a = mid; // (A)
        else b = mid+1;
    }
    rep(i, a+1, b+1) if (f (a) < f (i)) a = i; // (B)</pre>
```

#### LIS.h

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

```
int L = sz(res), cur = res.back().second;
vi ans(L);
while (L--) ans[L] = cur, cur = prev[cur];
return ans;
```

#### FastKnapsack.h

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

```
Time: \mathcal{O}(N \max(w_i)) b20ccc, 16 lines
```

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

#### 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  $[0, 2b]_{a02, 8}$  lines

```
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

**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;
}
```

#### BumpAllocator.h

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

745db2, 8 lines

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

#### SmallPtr.h

"BumpAllocator.h

**Description:** A 32-bit pointer that points into BumpAllocator memory.

2dd6c9, 10 lines

```
template<class T> struct ptr {
```

```
unsigned ind;
ptr(T* p = 0) : ind(p ? unsigned((char*)p - buf) : 0)

    assert(ind < sizeof buf);
}
T& operator*() const { return *(T*) (buf + ind); }
T& operator->() const { return &**this; }
T& operator[] (int a) const { return (&**this) [a]; }
explicit operator bool() const { return ind; }
};

BumpAllocatorSTL.h
Description: BumpAllocator for STL containers.
Usage: vector<vector<int, small<int>>>biscb(N),114 lines
char buf[450 << 20] alignas(16);
size_t buf_ind = sizeof buf;
template<class T> struct small {
    typedef T value_type;
    small() {}
    template<class U> small(const U&) {}
    T* allocate(size_t n) {
        buf_ind -= n * sizeof(T);
        buf_ind &= 0 - alignof(T);
        return (T*) (buf + buf_ind);
    }
void deallocate(T*, size_t) {}
}
```