

Changxv's Competitive Programming Template Library

Mofish Automaton

 ${\rm Changxv}$

1 Contest	1 Makefile
2 Mathematics	CXXFLAGS = -02 -Wall -Wextra -Wconversion -Wno-unused-resultation -std=c++17 -fsanitize=undefined, address -fstack-protest
3 Data structures	vimre .vimre
4 Numerical	set hls ic is scs ru nu ls=2 lbr sc cin noswf aw ar et so=5 =2 ts=2 sts=2 sw=2 tm=50 mouse=a sy on
5 Number theory	no; : no:;
6 Combinatorial 1	
7 Graph	- 110 (12) .VS 111.CXC(C1)
8 Geometry 1	<pre>no <f9> :!make %:r<cr> no <f8> :!time ./%:r < in.txt<cr> com -range=% -nargs=1 P exe "<line1>, <line2>!".<q-args> y u echom @"</q-args></line2></line1></cr></f8></cr></f9></pre>
9 Strings 2	1 1 55
10 Various 2	troubleshoot.txt
$\underline{\text{Contest}}$ (1)	Pre-submit: Write a few simple test cases if sample is not enough. Are time limits close? If so, generate max cases.
template.cpp 12 line	Is the memory usage fine?
<pre>#include "bits/stdc++.h" #define sz(x) int((x).size()) #define all(x) begin(x), end(x)</pre>	Make sure to submit the right file. Wrong answer:
<pre>using namespace std; template<class t=""> using vc = vector<t>; using l1 = int64_t;</t></class></pre>	Print your solution! Print debug output, as well. Are you clearing all data structures between test cases? Can your algorithm handle the whole range of input? Read the full problem statement again.
<pre>int main() { cin.tie(0)->sync_with_stdio(0);</pre>	Do you handle all corner cases correctly? Have you understood the problem correctly? Any uninitialized variables? Any overflows?
}	Confusing N and M, i and j, etc.? Are you sure your algorithm works?
<pre>debug.cpp template<class class="" enable_if<!is_same<u,<="" pre="" t="typename" u,=""></class></pre>	What special cases have you not thought of? Are you sure the STL functions you use work as you think? Add some assertions, maybe resubmit.
<pre>string>::value, typename U::value_type>::type> ostream &operator << (ostream &os, U v) { os << "{ ";</pre>	Create some testcases to run your algorithm on. Go through the algorithm for a simple case. Go through this list again.
<pre>for (T x: v) os << x << ' '; return os << '}';</pre>	Explain your algorithm to a teammate. Ask the teammate to look at your code.
<pre>void dbg() { cerr << "\033[0m\n"; } template<class classu="" t,=""></class></pre>	Go for a small walk, e.g. to the toilet. Is your output format correct? (including whitespace) Rewrite your solution from the start or let a teammate do:
<pre>void dbg(T e, Ua) { cerr << e << ' '; dbg(a); } #define err(A) cerr << "\033[31;1m"#A ": ", dbg(A)</pre>	Runtime error: Have you tested all corner cases locally?
check.sh 9 line	Any uninitialized variables?
#!/bin/bash while true; do	Any assertions that might fail? Any possible division by 0? (mod 0 for example)
./gen	Any possible infinite recursion?
<pre>./a > myout.txt ./b > brute.txt</pre>	Invalidated pointers or iterators? Are you using too much memory?
diff myout.txt brute.txt	Debug with resubmits (e.g. remapped signals, see Various).
<pre>if [\$? -ne 0]; then break; fi echo AC</pre>	Time limit exceeded:
dans.	Do you have any neggible infinite leens?

done

```
Iakefile
```

XXFLAGS = -02 -Wall -Wextra -Wconversion -Wno-unused-result -q -std=c++17 -fsanitize=undefined,address -fstack-protector

zimrc

et hls ic is scs ru nu ls=2 lbr sc cin noswf aw ar et so=5 bs =2 ts=2 sts=2 sw=2 tm=50 mouse=a on on o ; : 0:; o , 、 m kj <esc> m ik <esc> o <c-a> ggVG"+ o <c-n> :noh<cr> o <f2> :vs in.txt<cr> o <f9> :!make %:r<cr> o <f8> :!time ./%:r < in.txt<cr> om -range=% -nargs=1 P exe "<line1>,<line2>!".<q-args> |y|sil om! -buffer -range=% Hash <line1>, <line2>P cpp -dD -P -

coubleshoot.txt

intime error:

Do you have any possible infinite loops?

What is the complexity of your algorithm? Are you copying a lot of unnecessary data? (References) How big is the input and output? (consider scanf) Avoid vector, map. (use arrays/unordered_map) What do your teammates think about your algorithm?

Memory limit exceeded:

What is the max amount of memory your algorithm should need? Are you clearing all data structures between test cases?

Mathematics (2)

2.1 Recurrences

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

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

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

2.2 Trigonometry

$$\sin v + \sin w = 2\sin\frac{v+w}{2}\cos\frac{v-w}{2}$$

$$\sin v - \sin w = 2\cos\frac{v+w}{2}\sin\frac{v-w}{2}$$

$$\cos v + \cos w = 2\cos\frac{v+w}{2}\cos\frac{v-w}{2}$$

$$\cos v - \cos w = -2\sin\frac{v+w}{2}\sin\frac{v-w}{2}$$

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

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

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

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

2.3 Geometry

2.3.1 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):

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

template debug check Makefile .vimrc troubleshoot

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

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

2.3.2 Quadrilaterals

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

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

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

2.4 Derivatives/Integrals

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

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

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

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

2.5 Sums

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

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

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

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

2.6 Series

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

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

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

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

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

2.7 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 σ is the standard 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,

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

2.7.1 Discrete distributions Binomial distribution

The number of successes in n independent yes/no experiments, each which yields success with probability p is $Bin(n, p), n = 1, 2, ..., 0 \le p \le 1.$

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

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

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 wich 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 κ and independently of the time since the last event is $Po(\lambda)$, $\lambda = t\kappa$.

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

$$\mu = \lambda, \, \sigma^2 = \lambda$$

2.7.2 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.

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

$$\mu = \frac{a+b}{2}, \, \sigma^2 = \frac{(b-a)^2}{12}$$

Exponential distribution

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

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

Normal distribution

Most real random values with mean μ and variance σ^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

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

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, \dots be a sequence of random variables generated by the Markov process. Then there is a transition matrix $\mathbf{P} = (p_{ij})$, with $p_{ij} = \Pr(X_n = i | X_{n-1} = j)$, and $\mathbf{p}^{(n)} = \mathbf{P}^n \mathbf{p}^{(0)}$ is the probability distribution for X_n (i.e., $p_i^{(n)} = \Pr(X_n = i)$), where $\mathbf{p}^{(0)}$ is the initial distribution.

 π 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. π_j/π_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, π_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:** $\mathcal{O}(\log N)$

```
#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>;
/* basic */
begin(), end(), size(), empty(), clear()
insert(pair<key, T>), erase(iter), erase(key), operator[]
find(key), lower_bound(key), upper_bound(key)
/* advanced */
order_of_key(key); // number of element less than key
// return the iter of order+1, end() if too large
find by order(size type order);
// join the other if their ranges don't intersect
join(tree &other);
// split tr and overide other with element larger than key
tr.split(key, tree &other);
```

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).

26e9d1f2, 14 lines

```
#include <ext/pb_ds/assoc_container.hpp>
#include <ext/pb_ds/hash_policy.hpp>
// const int RANDOM = chrono::high_resolution_clock::now().\
time_since_epoch().count();
struct chash {
  const uint64_t C = ll(4e18 * acos(0)) | 71;
   ll operator()(ll x) const { return __builtin_bswap64(x*C); }
   /*
   ll operator()(ll x) const {
    return __builtin_bswap64((x*RANDOM)*C);
   }
   */
};
__gnu_pbds::gp_hash_table<1l,int,chash> h({},{},{},{},{},{1<<16});</pre>
```

Rope.h

Description: STL BST Time: $O(\log N)$

3bbc1517, 16 lines

};

```
#include <ext/rope>
using namespace __gnu_cxx;
```

```
rope<T> rp;
crope rp; // same for rope<char>
/* basic */
size(), operator[]
push_back(key);
insert (pos, x);
erase(pos, x);
replace(pos, x);
substr(pos, x);
/* advanced */
rope<int> *his[N];
his[i] = new rope<int> (*his[i - 1]);
SegmentTree.h
Description: Zero-indexed segment tree. Bounds are inclusive to the left
and exclusive to the right.
Time: \mathcal{O}(\log N)
                                                    dea81b51, 49 lines
struct Seq {
 int n, h;
 vc<int> s, add;
 Seg(int n): n(n), h(32 - \underline{builtin_clz(n)}), s(n * 2),
              add(n) {}
 void update(int x, int v, int 1) {
   s[x] += v * 1;
   if (x < n) add[x] += v;
 void push(int 1, int r) {
   int b = h, k = 1 << (h - 1);
    for (1 += n, r += n - 1; b > 0; --b, k /= 2)
      for (int x = 1 >> b; x <= r >> b; ++x) if (add[x]) {
        update(x * 2, add[x], k);
        update(x * 2 + 1, add[x], k);
        add[x] = 0;
 void pull(int 1, int r) {
   int k = 2;
    for (1 += n, r += n - 1; 1 > 1; k *= 2) {
     1 /= 2, r /= 2;
      for (int x = r; x >= 1; --x) {
       s[x] = s[x * 2] + s[x * 2 + 1] + add[x] * k;
 void modify(int 1, int r, int v) {
   push(1, 1 + 1);
   push(r - 1, r);
    int 10 = 1, r0 = r, k = 1;
    for (1 += n, r += n; 1 < r; 1 /= 2, r /= 2, k *= 2) {
     if (1 & 1) update(1++, v, k);
     if (r & 1) update(--r, v, k);
   pull(10, 10 + 1);
   pull(r0 - 1, r0);
  } //3fc22174
 int query(int 1, int r) {
   push(1, 1 + 1);
   push(r - 1, r);
   int res = 0;
   for (1 += n, r += n; 1 < r; 1 /= 2, r /= 2) {
     if (1 & 1) res += s[1++];
     if (r & 1) res += s[--r];
    return res;
```

```
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).
"../various/BumpAllocator.h"
                                                        86334bd5, 52 lines
constexpr int INF = numeric_limits<int>::max() / 2;
struct Seq {
  Seq *ls = 0, *rs = 0;
  int 1, r, mset = INF, madd = 0, val = -INF;
  Seg(int 1, int r): 1(1), r(r){}
  Seg(vc<int> &v, int 1, int r): 1(1), r(r) {
    if (1 + 1 < r) {
      int mi = 1 + (r - 1)/2;
      ls = new Seg(v, l, mi), rs = new Seg(v, mi, r);
    } else val = v[1];
  } // 972dfe1b
  void pull() { val = max(ls->val, rs->val); }
  void push() {
    if (!ls) {
      int mi = 1 + (r - 1)/2;
      ls = new Seg(1, mi), rs = new Seg(mi, r);
    if (mset != INF)
      ls->set(l,r,mset), rs->set(l,r,mset), mset = INF;
    else if (madd)
      ls \rightarrow add(l, r, madd), rs \rightarrow add(l, r, madd), madd = 0;
  int query(int L, int R) {
    if (R <= 1 | | r <= L) return -INF;</pre>
    if (L <= 1 && r <= R) return val;</pre>
    return max(ls->query(L, R), rs->query(L, R));
  void set(int L, int R, int x) {
    if (R <= 1 || r <= L) return;
    if (L <= 1 && r <= R) mset = val = x, madd = 0;</pre>
      push();
      ls \rightarrow set(L, R, x), rs \rightarrow set(L, R, x);
      pull();
  void add(int L, int R, int x) {
    if (R <= 1 || r <= L) return;</pre>
    if (L <= 1 && r <= R) {
      if (mset != INF) mset += x;
      else madd += x;
      val += x;
    else {
      push();
      ls \rightarrow add(L, R, x), rs \rightarrow add(L, R, x);
      pull();
};
```

PersistentSegmentTree.h

Description: Segment tree with persistent ability. 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 = null; tr->insert(-INF, INF, pos, val);

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

"../various/BumpAllocator.h" 7569d05b, 25 lines
```

```
CSU
struct Seq {
  static Seg *null;
  Seq *ls = this, *rs = this;
  int val = 0;
  Sea() {}
  Seg(int val): ls(null), rs(null), val(val) {}
  Seq(Seq *ls, Seq *rs): ls(ls), rs(rs) {
   val += ls->val;
   val += rs->val;
  Seg *insert(int 1, int r, int pos, int v) {
    if (pos < 1 || pos >= r) return this;
    if (1 + 1 == r) return new Seg(val + v);
   int mi = 1 + (r - 1) / 2;
    return new Seg(ls->insert(l, mi, pos, v),
                   rs->insert(mi, r, pos, v));
  int query(Seg *p, int 1, int r, int k) {
   if (1 + 1 == r) return 1;
   int mi = 1 + (r - 1) / 2, cnt = ls->val - p->ls->val;
   return cnt >= k ? ls->query(p->ls, l, mi, k) :
                      rs->query(p->rs, mi, r, k - cnt);
};
Seg *Seg::null = new Seg;
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)
                                                    005e7e70, 22 lines
  vc<int> e;
  vc<pair<int, int>> st;
```

```
struct RollbackUF {
  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 = sz(st); 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 0;
   if (e[a] > e[b]) swap(a, b);
   st.emplace_back(a, e[a]);
   st.emplace back(b, e[b]);
   e[a] += e[b]; e[b] = a;
    return 1;
};
```

Matrix.h

Description: Basic operations on square matrices.

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

ed7aad73, 31 lines

```
template<class T, int N>
struct Matrix {
 using M = Matrix;
  array<array<T, N>, N> d{};
 M operator * (const M &m) const {
   Ma;
   for (int i = 0; i < N; ++i)
```

```
for (int k = 0; k < N; ++k)
        for (int j = 0; j < N; ++j)
         a.d[i][j] += d[i][k] * m.d[k][j];
 vc<T> operator * (const vc<T> &vec) const {
   vc<T> ret(N);
    for (int i = 0; i < N; ++i)
     for (int j = 0; j < N; ++j)
        ret[i] += d[i][j] * vec[j];
    return ret;
 M operator ^ (int p) const {
   assert (p >= 0);
   M a, b(*this);
    for (int i = 0; i < N; ++i) a.d[i][i] = 1;</pre>
    while (p) {
     if (p\&1) a = a * b;
     b = b * b;
     p /= 2;
    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}(\log N)
```

```
7f54e44f, 30 lines
```

```
struct Line {
  mutable 11 k, m, p;
  bool operator < (const Line &o) const { return k < o.k; }</pre>
  bool operator < (11 x) const { return p < x; }</pre>
struct LineContainer: multiset<Line, less<>>> {
  // for doubles, use INF = 1 / .0, div(a, b) = a / b
  static const ll INF = numeric_limits<ll>::max();
  ll div(ll a, ll b) { // floored division
    return a / b - ((a ^ b) < 0 && a % b);
  bool isect(iterator x, iterator y) {
    if (y == end()) return x \rightarrow p = INF, 0;
    if (x->k == y->k) x->p = x->m > y->m ? INF : -INF;
    else x -> p = div(y -> m - x -> m, x -> k - y -> k);
    return x->p >= v->p;
  void add(ll k, ll m) {
    auto z = insert(\{k, m, 0\}), y = z++, x = y;
    while (isect(v, 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 1.k * x + 1.m;
};
```

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

0dd1013a, 54 lines

```
struct Node {
 Node *1 = 0, *r = 0;
```

```
if (r) c += r->c;
};
struct Treap {
 Node *tr = 0;
 int cnt(Node *n) { return n ? n->c : 0; }
 // split is [, )
 pair<Node*, Node*> split(Node* n, int k) {
   if (!n) return {};
    if (cnt(n->1) \geq k) { // "n->val>= k" for lower_bound(k)
      auto pa = split(n->1, k);
     n->1 = pa.second;
      n->pull();
      return {pa.first, n};
      auto pa = split(n->r, k - cnt(n->1)); // and just "k"
      n->r = pa.first;
      n->pull();
      return {n, pa.second};
  } // 2753f4a9
  Node* merge(Node* 1, Node* r) {
    if (!1) return r;
    if (!r) return 1;
    if (1->v > r->v) {
     1->r = merge(1->r, r);
     1->pull();
      return 1;
    } else {
      r->1 = merge(1, r->1);
      r->pull();
      return r:
 } // 5e0342f4
 Node *ins(Node *t, Node *n, int pos) {
    auto pa = split(t, pos);
    return merge (merge (pa.first, n), pa.second);
  } // 9a5c6249
  // Example application: move the range (l, r) to index k
  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 \le 1) t = merge(ins(a, b, k), c);
    else t = merge(a, ins(c, b, k - r));
 } // 9c4818f6
};
FenwickTree.h
Description: Computes partial sums a[0] + a[1] + ... + a[pos - 1], and
```

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

```
Time: Both operations are \mathcal{O}(\log N).
```

int val, v, c = 1;

if (1) c += 1->c;

void pull() {

c = 1:

Node(int val): val(val), y(rand()) {}

e77418e0, 22 lines

```
template<class T> struct FT {
 vc < T > s;
 FT(int n): s(n) {}
 void update(int p, T v) {
    for (; p < sz(s); p |= p + 1) s[p] += v;
 T query(int p) {
   T res = 0;
    for (; p > 0; p \&= p - 1) res += s[p - 1];
    return res;
```

```
}
int lower_bound(T sum) {// min pos st sum of [0, pos] >= sum
// Returns n if no sum is >= sum, or -1 if empty sum is.
if (sum <= 0) return -1;
int p = 0;
for (int pw = 1 << 25; pw; pw >>= 1) {
   if (p + pw <= sz(s) && s[p + pw - 1] < sum)
        p += pw, sum -= s[p - 1];
}
return p;
} // 18c85520;
</pre>
```

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

"FenvickTree.h" f4066

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

RMQ.h

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

Usage: RMQ rmq(values); rmq.query(inclusive, exclusive); **Time:** $\mathcal{O}(|V|\log|V|+Q)$

```
template < class T > struct RMQ {
    vc<vc<T>> jmp;
    RMQ(vc<T> &a): jmp(1, a) {
        for (int pw = 1, k = 1; pw * 2 <= sz(a); pw *= 2, ++k) {
             jmp.emplace_back(sz(a) - pw * 2 + 1);
             for (int j = 0; j < sz(jmp[k]); ++j) {
                 jmp[k][j] = min(jmp[k - 1][j], jmp[k - 1][j + pw]);
             }
        }
    }
    T query(int a, int b) {
        assert(a < b); // or return inf if a == b
        int dep = 31 - _builtin_clz(b - a);
    return min(jmp[dep][a], jmp[dep][b - (1 << dep)]);
    }
};</pre>
```

MoQueries.h

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

```
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
vc<int> mo(vc<array<int, 2>> 0) {
 int L = 0, R = 0, blk = 350; // \sim N/sqrt(Q)
 vc < int > s(sz(Q)), res = s;
#define K(x) make_pair(x[0] / blk, x[1] ^ -(x[0] / blk & 1))
  iota(all(s), 0);
  sort(all(s), [\&](int s, int t) \{ return K(O[s]) < K(O[t]); \});
  for (int qi: s) {
    arrav < int, 2 > g = 0[gi];
    while (L > g.first) add(--L, 0);
    while (R < q.second) add(R++, 1);</pre>
    while (L < q.first) del(L++, 0);
    while (R > g.second) del(--R, 1);
    res[gi] = calc();
 return res;
} //844328d6
vc<int> moTree(vc<arrav<int, 2>> 0.vc<vc<int>> &ed.int root=0){
 int N = sz(ed), pos[2] = {}, blk = 350; // \sim N/sqrt(Q)
  vc < int > 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) make_pair(I[x[0]]/blk, I[x[1]] ^ -(I[x[0]]/blk&1))
  iota(all(s), 0);
 sort(all(s), [\&](int s, int t) { return K(Q[s]) < K(Q[t]); });
 for (int qi : s) for(int end = 0; end < 2; ++end) {</pre>
   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;
} //01876993
```

Numerical (4)

4.1 Polynomials and recurrences

Polynomial.h

c9b7b07a, 17 lines

```
struct Poly {
  vector<double> a;
  double operator()(double x) const {
   double val = 0;
   for (int i = sz(a); i--;) (val *= x) += a[i];
   return val;
```

```
void diff() {
    rep(i, 1, sz(a)) a[i-1] = i*a[i];
    a.pop_back();
  void divroot(double x0) {
    double b = a.back(), c; a.back() = 0;
    for(int i=sz(a)-1; i--;) c = a[i], a[i] = a[i+1]*x0+b, b=c;
    a.pop_back();
};
PolyRoots.h
Description: Finds the real roots to a polynomial.
Usage: polyRoots(\{\{2, -3, 1\}\}, -1e9, 1e9\}) // solve x^2-3x+2=0
Time: \mathcal{O}\left(n^2\log(1/\epsilon)\right)
"Polynomial.h"
                                                     088a8d2d, 23 lines
vc<double> polyRoots(Poly p, double xmin, double xmax) {
 if (sz(p.a) == 2) { return {-p.a[0]/p.a[1]}; }
 vc<double> ret;
 Poly der = p;
  der.diff();
  auto dr = polyRoots(der, xmin, xmax);
  dr.emplace_back(xmin-1);
 dr.emplace_back(xmax+1);
  sort(all(dr));
  for (int i = 0; i < sz(dr) - 1; ++i) {
    double l = dr[i], h = dr[i+1];
    bool sign = p(1) > 0;
    if (sign ^{(p(h) > 0)}) {
      for (int it = 0; it < 60; ++it) { // while (h - l > 1e-8)
        double m = (1 + h) / 2, f = p(m);
        if ((f \le 0) ^ sign) 1 = m;
        else h = m;
      ret.emplace_back((1 + h) / 2);
  return ret;
```

PolyInterpolate.h

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

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

BerlekampMassev.h

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

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

```
Time: \mathcal{O}(N^2)
"../number-theory/ModPow.h"
                                                     9ce628c7, 22 lines
vc<ll> berlekampMassev(vc<ll> s) {
 int n = sz(s), L = 0, m = 0;
 vc<11> C(n), B(n), T;
 C[0] = B[0] = 1;
 11 b = 1;
  for (int i = 0; ++m, i < n; ++i) {</pre>
   ll d = s[i] % mod;
   for (int j = 1; j <= L; ++j)</pre>
     d = (d + C[j] * s[i - j]) % mod;
    if (!d) continue;
    T = C; 11 coef = d * modpow(b, mod-2) % mod;
   for (int j = m; j < n; ++j)
     C[j] = (C[j] - coef * B[j - m]) % mod;
   if (2 * L > i) continue;
   L = i + 1 - L; B = T; b = d; m = 0;
 C.resize(L + 1); C.erase(C.begin());
  for (11 &x : C) x = (mod - x) % mod;
 return C:
```

LinearRecurrence.h

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

```
using Poly = vc<11>;
ll linearRec(Poly S, Poly tr, ll k) {
 int n = sz(tr);
  auto combine = [&] (Poly a, Poly b) {
   Poly res(n \star 2 + 1);
   for (int i = 0; i <= n; ++i)</pre>
     for (int j = 0; j <= n; ++j)
       res[i + j] = (res[i + j] + a[i] * b[j]) % mod;
    for (int i = 2 * n; i > n; --i) for (int j = 0; j < n; ++j)
     res[i - 1 - j] = (res[i - 1 - j] + res[i] * tr[j]) % mod;
   res.resize(n + 1);
   return res;
  Poly pol(n + 1), e(pol);
 pol[0] = e[1] = 1;
  for (++k; k; k /= 2) {
   if (k % 2) pol = combine(pol, e);
   e = combine(e, e);
  11 \text{ res} = 0;
  for (int i = 0; i < n; ++i)
   res = (res + pol[i + 1] * S[i]) % mod;
 return res;
```

4.2 Optimization

GoldenSectionSearch.h

Description: Finds the argument minimizing the function f in the interval [a,b] assuming f is unimodal on the interval, i.e. has only one local minimum. The maximum error in the result is eps. Works equally well for maximization with a small change in the code. See TernarySearch.h in the Various chapter for a discrete version.

```
Usage: double func(double x) { return 4+x+.3*x*x; } double xmin = gss(-1000,1000,func);

Time: \mathcal{O}(\log((b-a)/\epsilon)) 31d45b51, 14 lines

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

Integrate.h

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

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

IntegrateAdaptive.h

```
Description: Fast integration using an adaptive Simpson's rule.

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

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

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

Simplex.h

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

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

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

```
using vvd = vc<vd>;
const T eps = 1e-8, inf = 1/.0;
#define MP make pair
#define ltj(X) if(s == -1 || MP(X[j], N[j]) < MP(X[s], N[s])) s=j
struct LPSolver {
 int m, n;
 vi N, B;
 vvd D:
 LPSolver(const vvd& A, const vd& b, const vd& c) :
    m(sz(b)), n(sz(c)), N(n+1), B(m), D(m+2), vd(n+2)) {
      for (int i = 0; i < m; ++i)
        for (int j = 0; j < n; ++j) D[i][j] = A[i][j];</pre>
      for (int i = 0; i < m; ++i)
       B[i] = n+i, D[i][n] = -1, D[i][n+1] = b[i];
      for (int j = 0; j < n; ++j)
       N[j] = j, D[m][j] = -c[j];
      N[n] = -1, D[m+1][n] = 1;
  void pivot(int r, int s) {
   T \star a = D[r].data(), inv = 1 / a[s];
    for (int i = 0; i < m + 2; ++i)
      if (i != r && abs(D[i][s]) > eps) {
       T *b = D[i].data(), inv2 = b[s] * inv;
        for (int j = 0; j < n + 2; ++j) b[j] -= a[j] * inv2;</pre>
        b[s] = a[s] * inv2;
    for (int j = 0; j < n + 2; ++j)
     if (j != s) D[r][j] *= inv;
    for (int i = 0; i < m + 2; ++i)
     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;
      for (int j = 0; j \le n; ++j)
       if (N[j] != -phase) ltj(D[x]);
      if (D[x][s] >= -eps) return true;
      int r = -1;
      for (int i = 0; i < m; ++i) {</pre>
       if (D[i][s] <= eps) continue;</pre>
        if (r == -1 || MP(D[i][n+1] / D[i][s], B[i])
                     < MP(D[r][n+1] / D[r][s], B[r])) r = i;
      if (r == -1) return false;
      pivot(r, s);
 T solve(vd &x) {
    int r = 0;
    for (int i = 1; i < m; ++i)
     if (D[i][n+1] < D[r][n+1]) r = i;
    if (D[r][n+1] < -eps) {
      pivot(r, n);
      if (!simplex(2) || D[m+1][n+1] < -eps) return -inf;</pre>
      for (int i = 0; i < m; ++i) if (B[i] == -1) {
        int s = 0:
        for (int j = 1; j <= n; ++j) ltj(D[i]);</pre>
        pivot(i, s);
```

```
}
bool ok = simplex(1); x = vd(n);
for (int i = 0; i < m; ++i)
    if (B[i] < n) x[B[i]] = D[i][n+1];
return ok ? D[m][n+1] : inf;
}
};</pre>
```

4.3 Matrices

Determinant.h

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

IntDeterminant.h

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

```
Time: \mathcal{O}\left(N^3\right)
                                                        88ad2892, 18 lines
const 11 md = 12345;
auto det = [md] (vc<vc<ll>> &a) { //2612f2d6
  int n = sz(a); ll ans = 1;
  for (int i = 0; i < n; ++i) {</pre>
    for (int j = i + 1; j < n; ++j) {
      while (a[j][i] != 0) { // gcd step}
        11 t = a[i][i] / a[j][i];
        if (t) for (int k = i; k < n; ++k)</pre>
           a[i][k] = (a[i][k] - a[j][k] * t) % md;
        swap(a[i], a[j]);
         ans \star = -1;
    ans = ans * a[i][i] % md;
    if (!ans) return 0;
  return (ans + md) % md;
};
```

SolveLinear.h

Description: Solves A*x=b. If there are multiple solutions, an arbitrary one is returned. Returns rank, or -1 if no solutions. Data in A and b is lost. **Time:** $\mathcal{O}\left(n^2m\right)$

```
using vd = vc<double>;
constexpr double eps = 1e-12;

auto solveLinear = [](vc<vd> &A, vd &b, vd &x) { //790aae09
  int n = sz(A), m = sz(x), rank = 0, br, bc;
  if (n) assert(sz(A[0]) == m);
  vc<int> col(m); iota(all(col), 0);
  for (int i = 0; i < n; ++i) {
      double v, bv = 0;
  }
}</pre>
```

```
for (int r = i; r < n; ++r)</pre>
    for (int c = i; c < m; ++c)</pre>
      if ((v = fabs(A[r][c])) > bv) br = r, bc = c, bv = v;
  if (bv <= eps) {
    for (int j = i; j < n; ++j)
      if (fabs(b[j]) > eps) return -1;
    break:
  swap(A[i], A[br]);
  swap(b[i], b[br]);
  swap(col[i], col[bc]);
  for (int j = 0; j < n; ++j) swap(A[j][i], A[j][bc]);</pre>
  bv = 1/A[i][i];
  for (int j = i + 1; j < n; ++j) {
    double fac = A[j][i] * bv;
    b[j] -= fac * b[i];
    for (int k = i + 1; k < m; ++k) A[j][k] -= fac * A[i][k];
  ++rank;
x.assign(m, 0);
for (int i = rank; i--;) {
 b[i] /= A[i][i];
  x[col[i]] = b[i];
  for (int j = 0; j < i; ++j) b[j] -= A[j][i] * b[i];</pre>
return rank; // (multiple solutions if rank < m)
```

SolveLinear2.h

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

SolveLinearBinarv.h

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

```
using bs = bitset<1000>;
auto solveLinear = [](vc<bs> &A, vc<int> &b, bs &x, int m) {
 int n = sz(A), rank = 0, br;
 assert(m \le sz(x));
  vc<int> col(m); iota(all(col), 0);
  for (int i = 0; i < n; ++i) {</pre>
    for (br=i; br<n; ++br) if (A[br].any()) break;</pre>
    if (br == n) {
      for (int j = i; j < n; ++j) if (b[j]) return -1;
      break:
    int bc = int(A[br]._Find_next(i - 1));
    swap(A[i], A[br]);
    swap(b[i], b[br]);
    swap(col[i], col[bc]);
    for (int j = 0; j < n; ++j) if (A[j][i] != A[j][bc]) {</pre>
     A[j].flip(i); A[j].flip(bc);
```

```
for (int j = i + 1; j < n; ++j) 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;
    for (int j = 0; j < i; ++j) b[j] ^= A[j][i];
}
return rank; // (multiple solutions if rank < m)
}; //f6c54262</pre>
```

MatrixInverse.h

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

```
Time: \mathcal{O}\left(n^3\right)
                                                     453d5d7a, 35 lines
int matInv(vc<vc<double>>& A) {
 int n = sz(A); vi col(n);
 vc<vc<double>> tmp(n, vc<double>(n));
 rep(i, 0, n) tmp[i][i] = 1, col[i] = i;
  rep(i,0,n) {
    int r = i, c = i;
    rep(j,i,n) rep(k,i,n)
      if (fabs(A[j][k]) > fabs(A[r][c]))
        r = j, c = k;
    if (fabs(A[r][c]) < 1e-12) return i;</pre>
    A[i].swap(A[r]); tmp[i].swap(tmp[r]);
    rep(j,0,n)
      swap(A[j][i], A[j][c]), swap(tmp[j][i], tmp[j][c]);
    swap(col[i], col[c]);
    double v = A[i][i];
    rep(j,i+1,n) {
      double f = A[j][i] / v;
      A[j][i] = 0;
      rep(k, i+1, n) A[j][k] -= f*A[i][k];
      rep(k,0,n) tmp[j][k] -= f*tmp[i][k];
    rep(j, i+1, n) A[i][j] /= v;
    rep(j,0,n) tmp[i][j] /= v;
    A[i][i] = 1;
  for (int i = n-1; i > 0; --i) rep(j,0,i) {
    double v = A[j][i];
    rep(k,0,n) tmp[j][k] -= v*tmp[i][k];
  rep(i,0,n) rep(j,0,n) A[col[i]][col[j]] = tmp[i][j];
  return n;
```

Tridiagonal.h

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

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

This is useful for solving problems on the type

```
a_i = b_i a_{i-1} + c_i a_{i+1} + d_i, \ 1 \le i \le n,
```

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

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

Fails if the solution is not unique.

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

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

```
7a2f067a, 26 lines
using T = double;
vc<T> tridiagonal(vc<T> diag, const vc<T>& super,
    const vc<T>& sub, vc<T> b) {
  int n = sz(b); vi tr(n);
  rep(i, 0, n-1) {
    if (abs(diag[i]) < 1e-9 * abs(super[i])) { // diag[i] == 0
     b[i+1] = b[i] * diag[i+1] / super[i];
     if (i+2 < n) b[i+2] -= b[i] * sub[i+1] / super[i];</pre>
     diag[i+1] = sub[i]; tr[++i] = 1;
     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^{16}); higher for random inputs). Otherwise, use NTT/FFTMod. **Time:** $\mathcal{O}(N \log N)$ with $N = |A| + |B| \ (\sim 1s \text{ for } N = 2^{22})$

```
using C = complex<double>;
using vd = vc<double>;
auto fft = [&] (vc<C> &a) {
  int n = sz(a), L = 31 - __builtin_clz(n);
  static vc<C> R(2, 1), rt(2, 1); // R can be long double
  for (static int k = 2; k < n; k *= 2) {
   R.resize(n), rt.resize(n);
    auto x = polar(1., acos(-1) / k);
   for (int i = k; i < k * 2; ++i)</pre>
      rt[i] = R[i] = i & 1 ? R[i / 2] * x : R[i / 2];
  vc<int> rev(n);
  for (int i = 0; i < n; ++i)
   rev[i] = (rev[i / 2] | (i & 1) << L) / 2;
  for (int i = 0; i < n; ++i)</pre>
   if (i < rev[i]) swap(a[i], a[rev[i]]);</pre>
  for (int k = 1; k < n; k *= 2)
    for (int i = 0; i < n; i += k * 2) {
```

```
auto it1 = &a[i], it2 = it1 + k;
      for (int j = 0; j < k; ++j, ++it1, ++it2) {</pre>
        auto x = (double *) &rt[j + k], y = (double *) it2;
       C z(x[0]*y[0] - x[1]*y[1], x[0]*y[1] + x[1] * y[0]);
        *it2 = *it1 - z;
        *it1 += z;
}; //bfe3257c
auto conv = [&](vd &a,vd &b) {
 if (a.empty() || b.empty()) return vd();
 vd res(sz(a) + sz(b) - 1);
 int L = 32 - __builtin_clz(sz(res) - 1), n = 1 << L;</pre>
 vc<C> in(n), out(n);
 copy(all(a), begin(in));
 for (int i = 0; i < sz(b); ++i) in[i].imag(b[i]);</pre>
 fft(in);
 for (C &x: in) x \star = x;
 for (int i = 0; i < n; ++i)</pre>
   out[i] = in[-i & (n - 1)] - conj(in[i]);
 for (int i = 0; i < sz(res); ++i)
   res[i] = imag(out[i]) / (n * 4);
 return res;
}; //13bd14b6
```

FastFourierTransformMod.h

Description: Higher precision FFT, can be used for convolutions modulo arbitrary integers as long as $N \log_2 N \cdot \text{mod} < 8.6 \cdot 10^{14}$ (in practice 10^{16} or higher). Inputs must be in [0, mod).

Time: $\mathcal{O}(N \log N)$, where N = |A| + |B| (twice as slow as NTT or FFT) "FastFourierTransform.h"

```
auto convMod = [&](vc<int> &a, vc<int> &b, int M) {
 if (a.empty() || b.empty()) return vc<11>();
 vc<11> res(sz(a) + sz(b) - 1);
 int B=32-__builtin_clz(sz(res)), n=1<<B, cut = int(sqrt(M));</pre>
 vc<C> L(n), R(n), outs(n), outl(n);
 for (int i = 0; i < sz(a); ++i) L[i] = C(a[i]/cut, a[i]%cut);</pre>
 for (int i = 0; i < sz(b); ++i) R[i] = C(b[i]/cut, b[i]%cut);</pre>
 fft(L), fft(R);
 for (int i = 0; i < n; ++i) {</pre>
    int j = -i \& (n - 1);
    outl[j] = (L[i] + conj(L[j])) * R[i] / (n * 2.);
    outs[j] = (L[i] - conj(L[j])) * R[i] / (n * 2.) / 1i;
 fft (outl), fft (outs);
 for (int i = 0; i < sz(res); ++i) {
   11 \text{ av} = 11(\text{real}(\text{outl}[i]) + .5), \text{cv} = 11(\text{imag}(\text{outs}[i]) + .5),
       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 $g = \sum_{x} a[x]g^{xk}$ $\operatorname{root}^{(mod-1)/N}$. N must be a power of 2. Useful for convolution modulo specific nice primes of the form $2^a b + 1$, where the convolution result has size at most 2^a . For arbitrary modulo, see FFTMod. conv(a, b) = c, where $c[x] = \sum a[i]b[x-i]$. For manual convolution: NTT the inputs, multiply pointwise, divide by n, reverse(start+1, end), NTT back. Inputs must be in [0, mod).

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

```
"../number-theory/ModPow.h"
                                                      0bbe7f0e, 39 lines
constexpr 11 md = 998244353, root = 62;
// For p < 2^30 there is also e.g. 5 << 25, 7 << 26, 479 << 21
// and 483 \ll 21 (same root). The last two are > 10^9.
auto ntt = [&](vc<Mod> &a) {
```

```
int n = sz(a), L = 31 - __builtin_clz(n);
  static vc<Mod> rt(2, 1);
  for (static int k = 2, s = 2; k < n; k *= 2, ++s) {
    rt.resize(n):
    array<Mod, 2> z{1, mdPow(root, md >> s)};
    for (int i = k; i < k * 2; ++i)
      rt[i] = rt[i / 2] * z[i & 1];
 vc<int> rev(n);
  for (int i = 0; i < n; ++i)
    rev[i] = (rev[i / 2] | (i & 1) << L) / 2;
  for (int i = 0; i < n; ++i)</pre>
   if (i < rev[i]) swap(a[i], a[rev[i]]);</pre>
  for (int k = 1; k < n; k *= 2)
    for (int i = 0; i < n; i += k * 2) {
      auto it1 = &a[i], it2 = it1 + k;
      for (int j = 0; j < k; ++j, ++it1, ++it2) {
       Mod z = rt[j + k] * *it2;
        *it2 = *it1 - z, *it1 += z;
}; //7c5a0cf5
auto conv = [&](vc<Mod> a, vc<Mod> b) -> vc<Mod> {
 if (a.empty() || b.empty()) return {};
 int s = sz(a) + sz(b) - 1,
     n = 1 \ll (32 - \underline{builtin_clz(s - 1)});
 Mod inv = md - (md - 1) / n;
  vc<Mod> L(a), R(b), out(n);
 L.resize(n), R.resize(n);
 ntt(L), ntt(R);
  for (int i = 0; i < n; ++i)</pre>
    out [-i \& (n - 1)] = L[i] * R[i] * inv;
  return {out.begin(), out.begin() + s};
1: //75d60f42
```

PolvInverse.h

Description: Compute the inversion of the Polynomial.

return {b.begin(), b.begin() + sz(a)};

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

```
"NumberTheoreticTransform.h"
                                                    ad42700a, 26 lines
auto invIter = [&](vc<Mod> &a, vc<Mod> &in, vc<Mod> &b) {
 int n = sz(in);
 vc<Mod> out(n);
  copy(a.begin(), a.begin() + min(sz(a), n), out.begin());
  auto conv = [&] {
    ntt(out);
    for (int i = 0; i < n; ++i) out[i] *= in[i];</pre>
    ntt(out), reverse(out.begin() + 1, out.end());
 conv(), fill(out.begin(), out.begin() + sz(b), 0), conv();
 b.resize(n);
  Mod inv = md - (md - 1) / n; inv *= inv;
  for (int i = n / 2; i < n; ++i)
    b[i] = out[i].x ? inv * (md - out[i].x) : 0;
}; //787b29bc
auto polyInv = [&] (vc<Mod> &a) -> vc<Mod> {
 if (a.empty()) return {};
  vc<Mod> b{mdPow(a[0], md - 2)};
 b.reserve(sz(a));
  while (sz(b) < sz(a)) {
    vc < Mod > in(sz(b) * 2);
    copy(all(b), in.begin()), ntt(in);
    invIter(a, in, b);
```

```
PolySart.h
```

Description: Compute the sqrt of a, where a is a polynomial and a[0] = 1(if not, ModSqrt is required).

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

a0bcdc24, 24 lines "PolyInverse.h"

```
auto polySqrt = [&](vc<Mod> &a) -> vc<Mod> {
 if (a.empty()) return {};
 vc<Mod> b{1}, ib{1};
 b.reserve(sz(a)), ib.reserve(sz(a));
  auto conv = [&](vc<Mod> &a, vc<Mod> &b) {
   for (int i = 0; i < sz(a); ++i) a[i] *= b[i];</pre>
   ntt(a), reverse(a.begin() + 1, a.end());
  while (sz(b) < sz(a)) {
   int h = sz(b), n = h * 2;
   vc<Mod> in(n), out(n);
   copy(all(ib), in.begin()), ntt(in);
   copy(all(b), out.begin());
   conv(out, out), fill(out.begin(), out.begin() + h, 0);
   Mod inv = md - (md - 1) / n;
   for (int i = h; i < min(n, sz(a)); ++i)
     out[i] = out[i] * inv - a[i];
    conv(out, in), b.resize(n), inv *= (md / 2);
   for (int i = h; i < n; ++i) b[i] = out[i] * inv;</pre>
   if (sz(b) < sz(a)) invIter(b, in, ib);</pre>
  return {b.begin(), b.begin() + sz(a)};
```

PolyExp.h

Description: Compute the Exp of a, where a is a polynomial and a[0] = 0. Time: $\mathcal{O}(n \log n)$

```
"PolyInverse.h", "../number-theory/ModInverse.h"
                                                    b11b7ae3, 36 lines
auto deri = [&](vc<Mod> a) {
  for (int i = 1; i < sz(a); ++i) a[i - 1] = a[i] * i;</pre>
  a.pop_back();
  return a;
auto polyExp = [&] (vc<Mod> &a) -> vc<Mod> {
  if (a.empty()) return {};
  vc<Mod> b{1}, ib{1};
  b.reserve(sz(a)), ib.reserve(sz(a));
  auto conv = [&] (vc<Mod> &a, vc<Mod> &b) {
    for (int i = 0; i < sz(a); ++i) a[i] *= b[i];</pre>
   ntt(a), reverse(a.begin() + 1, a.end());
  while (sz(b) < sz(a)) {
    int h = sz(b), n = h * 2;
    Mod inv = md - (md - 1) / n;
    vc<Mod> db(n), dib(n), A(deri(b)), B(n);
    copy(all(ib), dib.begin()), ntt(dib);
    copy(all(b), db.begin()), ntt(db);
    A.resize(n), conv(A, dib);
    for (int i = 0; i < n; ++i) B[i] = db[i] * dib[i];</pre>
    ntt(B), reverse(B.begin() + 1, B.end());
    fill(B.begin(), B.begin() + h, 0);
    vc<Mod> da(deri(vc<Mod>(a.begin(), a.begin() + h)));
    da.resize(n), ntt(da), conv(B, da);
    for (int i = min(n, sz(a)) - 1; i >= h; --i)
     A[i] = (A[i-1] - B[i-1] * inv) * iv[i] * inv - a[i];
    fill(A.begin(), A.begin() + h, 0), conv(A, db);
    b.resize(n):
    for (int i = h; i < n; ++i)
     b[i] = A[i].x ? inv * (md - A[i].x) : 0;
    if (sz(b) < sz(a)) invIter(b, dib, ib);</pre>
```

```
return {b.begin(), b.begin() + sz(a)};
```

PolvLn.h

Description: Compute the Ln of a, where a is a polynomial and a[0] = 1. Time: $\mathcal{O}(n \log n)$

```
"PolyInverse.h", "../number-theory/ModInverse.h"
auto deri = [&](vc<Mod> a) {
 for (int i = 1; i < sz(a); ++i) a[i - 1] = a[i] * i;
 a.pop back();
 return a;
auto inte = [&](vc<Mod> a) {
 for (int i = sz(a) - 1; i >= 1; --i)
   a[i] = a[i - 1] * iv[i];
 a[0] = 0;
 return a;
auto polyLn = [&] (vc<Mod> &a) -> vc<Mod> {
  if (a.empty()) return {};
  int n = 1 \ll (32 - \underline{\ }builtin_clz(2 * sz(a) - 2));
  Mod inv = md - (md - 1) / n;
  vc<Mod> b = polyInv(a), c = deri(a);
  b.resize(n), c.resize(n);
  ntt(b), ntt(c);
  for (int i = 0; i < n; ++i) b[i] = b[i] * c[i] * inv;
 ntt(b), reverse(b.begin() + 1, b.end());
 b = inte(b);
 return {b.begin(), b.begin() + sz(a)};
```

PolyPow.h

Description: Compute the $A^m \pmod{x^n}$, where the A is a polynomial and n is the size of A. Notice that you should change the constant term of the

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

```
"PolyLn.h", "PolyExp.h"
                                                     0528559c, 14 lines
auto polyPow = [&](vc<Mod> &a, ll k) {
 int n = sz(a), t = n;
 for (int i = 0; i < n; ++i) if (a[i].x) { t = i; break; }</pre>
 if (t * k \ge n) return vc < Mod > (n, 0);
 vc<Mod> b(a.begin() + t, a.end());
 b.resize(n - t * k);
 Mod a0 = b[0];
 b = polyLn(b);
 11 k1 = k % md;
 for (Mod &e: b) e *= k1;
 b = polyExp(b, mdPow(a0, k % (md - 1)));
 b.insert(b.begin(), t * k, 0);
 return b;
```

PolyDivision.h

Description: Polynomial floor division. No leading 0's.

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

"NumberTheoreticTransform.h"

```
auto polyDiv = [&](vc<Mod> a, vc<Mod> b)
               -> pair<vc<Mod>, vc<Mod>> {
 if (sz(a) < sz(b)) return {vc<Mod>(), a};
 int n = sz(a) - sz(b) + 1;
 vc<Mod> da(a.rbegin(), a.rend()), db(b.rbegin(), b.rend());
 da.resize(n), db.resize(n);
 da = conv(da, polyInv(db));
 da.resize(n), reverse(all(da));
 auto c = conv(da, b);
 a.resize(sz(b) - 1);
 for (int i = 0; i < sz(a); ++i) a[i] -= c[i];</pre>
```

```
return {da, a};
};
```

FastSubsetTransform.h

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

464cf360, 16 lines

```
void FST(vi& a, bool inv) {
 for (int n = sz(a), step = 1; step < n; step *= 2) {
   for (int i = 0; i < n; i += 2 * step) rep(j,i,i+step) {</pre>
     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

Modular Arithmetic.h

Description: Operators for modular arithmetic. You need to set mod to some number first and then you can use the structure.

```
constexpr 11 md = 998244353; // change to something else
struct Mod {
 11 x;
 Mod(ll x = 0): x(x) \{ \}
  Mod operator+(Mod b) {11 y=x+b.x;return y<md ? y : y - md; }</pre>
  Mod operator-(Mod b) { return x - b.x + (x < b.x ? md : 0); }
  Mod operator * (Mod b) { return x * b.x % md; }
 Mod operator / (Mod b) { return *this * inv(b); }
  void operator += (Mod b) { x += b.x; x < md ?: x -= md; }</pre>
  void operator *= (Mod b) { (x *= b.x) %= md; }
  void operator -= (Mod b) { x -= b.x; -x < 0 ?: x += md; }</pre>
  void operator /= (Mod b) { *this *= inv(b); }
  Mod inv(Mod a) {
   ll x, y, g = euclid(a.x, md, x, y);
   assert(q == 1);
    return (x + md) % md;
 Mod operator ^ (11 e) {
    if (!e) return 1;
    Mod r = *this ^ (e / 2); r = r * r;
    return e & 1 ? *this * r : r;
};
```

ModInverse.h

0c82d61c, 13 lines

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

```
Time: \mathcal{O}(n)
                                                          6cc69c62, 7 lines
constexpr 11 md = 1e9 + 7, LIM = 2e4 + 1; // change this
auto mdInv = [&] {
 vc<11> inv(LIM); inv[1] = 1;
  for (int i = 2; i < LIM; ++i)</pre>
```

```
CSU
    inv[i] = md - (md / i) * inv[md % i] % md;
  return inv:
ModPow.h
Time: \mathcal{O}(\log n)
                                                        b807a56c, 7 lines
"ModularArithmetic.h"
constexpr 11 md = 1000000007; // change this
auto mdPow = [&] (Mod b, ll e) {
  Mod res = 1:
  for (; e; b *= b, e /= 2)
   if (e & 1) res *= b;
  return res:
};
ModLog.h
Description: Returns the smallest x > 0 s.t. a^x = b \pmod{m}, or -1 if no
such x exists. modLog(a,1,m) can be used to calculate the order of a.
Time: \mathcal{O}\left(\sqrt{m}\right)
11 modLog(11 a, 11 b, 11 m) {
  11 n = (11)   sqrt(m) + 1, e = 1, f = 1, j = 1;
  unordered_map<11, 11> A;
  while (j \le n \&\& (e = f = e * a % m) != b % m)
   A[e * b % m] = j++;
  if (e == b % m) return j;
  if (__gcd(m, e) == __gcd(m, b))
    rep(i,2,n+2) if (A.count(e = e * f % m))
      return n * i - A[e];
  return -1;
ModSum.h
Description: Sums of mod'ed arithmetic progressions.
modsum(to, c, k, m) = \sum_{i=0}^{\text{to}-1} (ki+c)\%m. divsum is similar but for
floored division.
Time: \log(m), with a large constant.
                                                        5c5bc571, 16 lines
typedef unsigned long long ull;
ull sumsq(ull to) { return to / 2 * ((to-1) | 1); }
ull divsum(ull to, ull c, ull k, ull m) {
  ull res = k / m * sumsq(to) + c / m * to;
  k %= m; c %= m;
  if (!k) return res;
  ull to2 = (to * k + c) / m;
  return res + (to - 1) * to2 - divsum(to2, m-1 - c, m, k);
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, m);
ModMulLL.h
Description: Calculate a \cdot b \mod c (or a^b \mod c) for 0 \le a, b \le c \le 7.2 \cdot 10^{18}
Time: \mathcal{O}(1) for modmul, \mathcal{O}(\log b) for modpow
                                                         c36634fb, 6 lines
using ull = uint64 t;
constexpr 11 md = 1e9 + 7; // change this
auto mdMul = [&] (ull a, ull b) {
 ll ret = a * b - md * ull(1.L / md * a * b);
 return ret + md * (ret < 0) - md * (ret >= (11) md);
```

```
ModSart.h
Description: Tonelli-Shanks algorithm for modular square roots. Finds x
s.t. x^2 = a \pmod{p} (-x gives the other solution). p should be an odd prime
and 0 \le a < p.

Time: \mathcal{O}\left(\log^2 p\right) worst case, \mathcal{O}\left(\log p\right) for most p
                                                        3e061b4f, 18 lines
auto mdSgrt = [&](11 a) -> 11 {
 if (!a) return 0;
  if (mdPow(a, md / 2) != 1) return -1;
  if (md % 4 == 3) return mdPow(a, (md + 1) / 4);
  // a^{(n+3)/8} \text{ or } 2^{(n+3)/8} * 2^{(n-1)/4} \text{ works if } p \% 8 == 5
  11 s = md - 1, n = 2;
  int r = 0, m;
  while (s % 2 == 0) ++r, s /= 2;
  while (mdPow(n, md / 2) != md - 1) ++n;
  11 x = mdPow(a, (s + 1) / 2), b = mdPow(a, s), g = mdPow(n, s)
  for (; ; r = m) {
    11 t = b;
    for (m = 0; m < r && t != 1; ++m) (t *= t) %= md;
    if (!m) return x;
    11 \text{ qs} = \text{mdPow}(q, 111 << (r - m - 1));
    g = gs * gs % md, (x *= gs) %= md, (b *= g) %= md;
};
5.2 Primality
FastEratosthenes.h
Description: Prime sieve for generating all primes smaller than LIM.
Time: LIM=1e9 \approx 1.5s
                                                        22899edc, 24 lines
constexpr int LIM = 1e7;
bitset<LIM> is pr;
auto eratosthenes = [&] {
  const int S = round(sqrt(LIM)), R = LIM / 2;
  array<bool, S + 1> sieve{};
  vc<pair<int, int>> cp;
  for (int i = 3; i <= S; i += 2) if (!sieve[i]) {</pre>
    cp.emplace_back(i, i * i / 2);
    for (int j = i * i; j <= S; j += i * 2) sieve[j] = 1;</pre>
  vc < int > pr = {2};
  pr.reserve(int(LIM / log(LIM) * 1.1));
  for (int L = 1; L <= R; L += S) {
    array<bool, S> block{};
    for (auto &[p, idx]: cp)
      for (int i = idx; i < S + L; idx = (i += p))</pre>
        block[i - L] = 1;
    for (int i = 0; i < min(S, R - L); ++i) {
      if (!block[i]) pr.emplace_back((L + i) * 2 + 1);
  for (int i: pr) is_pr[i] = 1;
  return pr;
EulerSieve.h
Description: Calculate some arithmetic function in linear time.
Time: \mathcal{O}(n)
                                                        2d119e47, 27 lines
constexpr int LIM = 1e6 + 1; // change this
vc<int> pr, d(LIM), w(d), mu(d), phi(d);
bitset<LIM> n_pr;
auto eulerSieve = [&] {
  pr.reserve(int(LIM / log(LIM) * 1.1));
  phi[1] = d[1] = w[1] = mu[1] = n_pr[1] = 1;
```

for (int i = 2; i < LIM; ++i) {</pre>

pr.emplace_back(i);

if (!n_pr[i]) {

```
phi[i] = i - 1, mu[i] = -1, d[i] = 2, w[i] = i + 1;
    for (auto p: pr) {
     int t = p * i;
     if (t >= LIM) break;
     n_pr[t] = 1;
     if (i % p == 0) {
       phi[t] = phi[i] * p, mu[t] = 0;
       d[t] = d[i] * 2 - d[i / p];
       w[t] = w[i] + p * (w[i] - w[i / p]);
       phi[t] = phi[i] * (p - 1), mu[t] = -mu[i];
       d[t] = d[i] * 2;
       w[t] = w[i] * (p + 1);
};
```

MillerRabin.h

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

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

```
"ModMulLL.h"
                                                      60dcd132, 12 lines
bool isPrime(ull n) {
 if (n < 2 || n % 6 % 4 != 1) return (n | 1) == 3;</pre>
 ull A[] = \{2, 325, 9375, 28178, 450775, 9780504, 1795265022\},
      s = \underline{\quad} 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"
                                                    a33cf6ef, 18 lines
ull pollard(ull n) {
 auto f = [n](ull x) \{ return modmul(x, x, n) + 1; \};
 ull x = 0, y = 0, t = 30, prd = 2, i = 1, q;
  while (t++ % 40 || __gcd(prd, n) == 1) {
   if (x == y) x = ++i, y = f(x);
    if ((q = modmul(prd, max(x,y) - min(x,y), n))) prd = q;
   x = f(x), y = f(f(y));
 return __gcd(prd, n);
vector<ull> factor(ull n) {
 if (n == 1) return {};
 if (isPrime(n)) return {n};
 ull x = pollard(n);
  auto 1 = factor(x), r = factor(n / x);
 1.insert(1.end(), all(r));
 return 1:
```

5.3 Divisibility

Euclid.h

CRT ContinuedFractions FracBinarySearch IntPerm

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

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

CRT.h

Description: Chinese Remainder Theorem.

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

"euclid.h" 04d93a45, 7 lines 11 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 / q * m + a;return x < 0 ? x + m*n/q : 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 to

$$ax + by = d$$

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

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

5.4 Fractions

ContinuedFractions.h

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

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

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

FracBinarySearch.h

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

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

```
struct Frac { ll p, q; };
template < class F >
Frac fracBS(F f, ll N) {
 bool dir = 1, A = 1, B = 1;
 Frac lo{0, 1}, hi{1, 1}; // Set hi to 1/0 to search (0, N)
 if (f(lo)) return lo;
 assert(f(hi));
  while (A | | B)
    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 \star adv;
   dir = !dir:
   swap(lo, hi);
   A = B; B = !!adv;
 return dir ? hi : lo;
```

5.5 Pythagorean Triples

The Pythagorean triples are uniquely generated by

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

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

5.6 Primes

p = 962592769 is such that $2^{21} \mid p - 1$, which may be useful. For hashing use 970592641 (31-bit number), 31443539979727 (45-bit). 3006703054056749 (52-bit). There are 78498 primes less than 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 $\mathbb{Z}_{2^a}^{\times}$ is instead isomorphic to $\mathbb{Z}_2 \times \mathbb{Z}_{2^{a-2}}$.

5.7 Estimates

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

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

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:

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

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

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

Combinatorial (6)

6.1 Permutations

6.1.1 Factorial

n	1 2 3	4	5 6	7	8	3	9	10	
n!	1 2 6	24 1	20 72	0 504	0 403	$320 \ 362$	2880 36	528800	
n	11	12	13	1	4	15	16	17	
n!	4.0e7	′ 4.8e	8 6.2e	9 8.7	e10 1	.3e12 :	2.1e13	3.6e14	
n	20	25	30	40	50	100	150	171	
n!	2e18	2e25	3e32	8e47	3e64	9e157	6e262	>DBL_MA	ıΧ

IntPerm.h

Description: Permutation -> integer conversion. (Not order preserving.) Integer -> permutation can use a lookup table. Time: $\mathcal{O}(n)$ 044568e0, 6 lines

```
int permToInt(vi& v) {
  int use = 0, i = 0, r = 0;
  for(int x:v) r = r * ++i + \underline{\quad builtin_popcount(use & -(1<< x))},
    use |= 1 << x;
                                            // (note: minus, not \sim!)
  return r:
```

6.1.2 Cycles

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.

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

6.1.4 Burnside's lemma

Given a group G of symmetries and a set X, the number of elements of X up to symmetry equals

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

multinomial BellmanFord FloydWarshall

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

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

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

6.2 Partitions and subsets

6.2.1 Partition function

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

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

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

6.2.2 Lucas' Theorem

Let n, m be non-negative integers and p a prime. Write $n=n_kp^k+\ldots+n_1p+n_0$ and $m=m_kp^k+\ldots+m_1p+m_0$. Then $\binom{n}{m}\equiv\prod_{i=0}^k\binom{n_i}{m_i}\pmod{p}$.

6.2.3 Binomials

multinomial.h

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

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

6.3 General purpose numbers

6.3.1 Bernoulli numbers

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

Sums of powers:

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

Euler-Maclaurin formula for infinite sums:

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

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

6.3.2 Stirling numbers of the first kind

Number of permutations on n items with k cycles.

$$c(n,k) = c(n-1,k-1) + (n-1)c(n-1,k), \ c(0,0) = 1$$

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

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

6.3.3 Eulerian numbers

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

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

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

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

6.3.4 Stirling numbers of the second kind

Partitions of n distinct elements into exactly k groups.

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

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

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

6.3.5 Bell numbers

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

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

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

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

- sub-diagonal monotone paths in an $n \times n$ grid.
- strings with n pairs of parenthesis, correctly nested.
- binary trees with with n+1 leaves (0 or 2 children).

- ordered trees with n+1 vertices.
- ways a convex polygon with n + 2 sides can be cut into triangles by connecting vertices with straight lines.
- \bullet permutations of [n] with no 3-term increasing subseq.

Graph (7)

7.1 Fundamentals

BellmanFord.h

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

```
7a4810de, 27 lines
constexpr ll INF = numeric_limits<ll>::max();
 int a, b, w, s() { return a < b ? a : -a; }</pre>
 Ed(int a, int b, int w): a(a), b(b), w(w) {}
struct Node { ll dist = INF; int prev = -1; };
auto bellmanFord = [](vc<Node> &nodes, vc<Ed> &eds, int s) {
 nodes[s].dist = 0;
  sort(all(eds), [](Ed a, Ed b) { return a.s() < b.s(); });</pre>
  int lim = sz(nodes) / 2 + 2; // /3+100 with shuffled vertices
  for (int i = 0; i < lim; ++i) for (Ed ed: eds) {
    Node cur = nodes[ed.a], &dest = nodes[ed.b];
    if (abs(cur.dist) == INF) continue;
    11 d = cur.dist + ed.w;
    if (d < dest.dist) {</pre>
      dest.prev = ed.a;
      dest.dist = (i < lim - 1 ? d : -INF);
  for (int i = 0; i < lim; ++i) for (Ed e : eds) {</pre>
    if (nodes[e.a].dist == -INF)
      nodes[e.b].dist = -INF;
}; // 96199983
```

FlovdWarshall.h

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

constexpr ll INF = numeric limits<ll>::max();

Time: $O(N^3)$ 379909b1, 14 lines

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

18d53978, 62 lines

```
struct PushRelabel {
  struct Edge {
   int v, rev;
   11 f, c;
  vc<vc<Edge>> g;
  vc<int> h, gap;
  vc<ll> ec;
  vc<vc<int>> hv;
  vc<Edge*> cur;
  PushRelabel(int n):
   q(n), ec(n), h(n), cur(n), hv(n * 2), gap(n * 2) {}
  void addEdge(int x, int y, ll c, ll rc = 0) {
   if (x == y) return;
   Edge a\{y, sz(g[y]), 0, c\}, b\{x, sz(g[x]), 0, rc\};
   q[x].emplace_back(a);
   g[y].emplace_back(b);
  void addFlow(Edge &e, ll f) {
   Edge &re = g[e.v][e.rev];
   if (!ec[e.v] && f) hv[h[e.v]].emplace_back(e.v);
   e.f += f, e.c -= f, ec[e.v] += f;
   re.f -= f, re.c += f, ec[re.v] -= f;
  11 maxFlow(int S, int T) {
   int n = sz(q);
   h[S] = n, ec[T] = 1, gap[0] = n - 1;
   for (int i = 0; i < n; ++i) cur[i] = g[i].data();</pre>
   for (auto &e: g[S]) addFlow(e, e.c);
    if (hv[0].empty()) return 0;
    for (int h now = 0; h now >= 0; ) {
     int u = hv[h_now].back();
     hv[h_now].pop_back();
     while (ec[u] > 0) {
       if (cur[u] == g[u].data() + sz(g[u])) {
          h[u] = numeric_limits<int>::max();
          for (auto &e : g[u]) {
            if (e.c && h[u] > h[e.v] + 1) {
             h[u] = h[e.v] + 1;
              cur[u] = &e;
          ++gap[h[u]];
          if(!--gap[h_now] && h_now < n) {</pre>
            for (int i = 0; i < n; ++i) {</pre>
              if(h_now < h[i] && h[i] < n) {</pre>
                --gap[h[i]];
                h[i] = n + 1;
          h now = h[u];
        } else if (cur[u] -> c && h[u] == h[cur[u] -> v] + 1) {
          addFlow(*cur[u], min(ec[u], cur[u]->c));
        } else ++cur[u];
     while (h_now >= 0 \&\& hv[h_now].empty()) --h_now;
    return -ec[S];
 bool leftOfMinCut (int a) { return h[a] >= sz(q); }
```

MinCostMaxFlow.h

 $\bf Description:$ Min-cost max-flow. Note that negative cost cycles are not supported.

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

c7cdbcb7, 91 lines

```
#include <bits/extc++.h>
constexpr 11 INF = numeric_limits<11>::max() / 2;
struct MCMF {
 struct Edge {
   int v. back:
   11 f, c;
 };
 int s, t;
 vc<vc<Edge>> q;
 vc<ll> dis, p;
 vc<int> vis;
 MCMF(int n): q(n), dis(n), p(n), vis(n) {}
 void add_edge(int x, int y, ll f, ll c) {
   Edge a\{y, sz(g[y]), f, c\}, b\{x, sz(g[x]), 0, -c\};
   g[x].emplace_back(a);
   g[y].emplace_back(b);
 bool setpi() {
   queue<int> 0;
   fill(all(dis), INF);
   dis[t] = 0;
   O.emplace(t);
    while (!Q.empty()) {
     int u = 0.front();
     () goq. ();
     vis[u] = 0;
     for (auto e: g[u]) {
       auto re = g[e.v][e.back];
       if (re.f && dis[e.v] > dis[u] + re.c) {
         dis[e.v] = dis[u] + re.c;
         if (!vis[e.v]) {
           vis[e.v] = 1;
           Q.emplace(e.v);
   return dis[s] != INF;
 bool path() {
   fill(all(dis), INF);
    __qnu_pbds::priority_queue<pair<11, int>> Q;
   vc<decltype(Q)::point_iterator> it(sz(p));
   dis[t] = 0;
   it[t] = Q.push({-dis[t], t});
    while (!Q.empty()) {
     int u = Q.top().second;
     Q.pop();
      for (auto e: g[u]) {
       auto re = g[e.v][e.back];
       if (re.f && dis[e.v] > dis[u] + re.c) {
         dis[e.v] = dis[u] + re.c;
         if (it[e.v] == Q.end())
           it[e.v] = Q.push({-dis[e.v], e.v});
          else Q.modify(it[e.v], {-dis[e.v], e.v});
    return dis[s] != INF;
```

```
ll dfs(int u, ll flow) {
   if (u == t || !flow) return flow;
   vis[u] = 1;
   11 w = flow;
    for (auto &e: g[u])
     if (e.f && !vis[e.v] && !e.c) {
       ll tmp = dfs(e.v, min(e.f, w));
       w -= tmp, e.f -= tmp, g[e.v][e.back].f += tmp;
       if (!w) return flow;
   return flow - w;
 pair<11, 11> mincost_flow(int S, int T) {
   s = S, t = T;
   if (!setpi()) return {};
   11 flow = 0, cost = 0, delta = 0, tmp;
      for (int u = 0; u < sz(g); ++u)
       for (auto &e: g[u]) e.c += dis[e.v] - dis[u];
      delta += dis[s];
     do {
       fill(all(vis), 0);
       tmp = dfs(s, INF);
       flow += tmp, cost += tmp * delta;
      }while (tmp);
    }while (path());
    return {flow, cost};
};
```

MinCut.h

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

GlobalMinCut.h

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

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

8b0e19d6, 21 lines

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

Gomory Hu.l

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

0418b3cd, 13 lines

76b5c93f, 24 lines

```
typedef array<11, 3> Edge;
vector<Edge> gomoryHu(int N, vector<Edge> ed) {
 vector<Edge> tree;
 vi par(N);
  rep(i,1,N) {
   PushRelabel D(N); // Dinic also works
   for (Edge t : ed) D.addEdge(t[0], t[1], t[2], t[2]);
   tree.push_back({i, par[i], D.calc(i, par[i])});
   rep(j,i+1,N)
     if (par[j] == par[i] && D.leftOfMinCut(j)) par[j] = i;
 return tree:
```

7.3 Matching

HopcroftKarp.h

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

```
Usage: vi btoa(m, -1); hopcroftKarp(g, btoa);
Time: \mathcal{O}\left(\sqrt{V}E\right)
```

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

MinimumVertexCover.h

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

```
"HopcroftKarp.h"
                                                  da419632, 20 lines
vi cover(vector<vi>& g, int n, int m) {
 vi match(m, -1);
 int res = dfsMatching(g, match);
 vector<bool> lfound(n, true), seen(m);
 for (int it : match) if (it != -1) lfound[it] = false;
 vi q, cover;
 rep(i,0,n) if (lfound[i]) q.push_back(i);
 while (!q.empty()) {
   int i = q.back(); q.pop_back();
   lfound[i] = 1;
   for (int e : g[i]) if (!seen[e] && match[e] != -1) {
     seen[e] = true;
     g.push_back(match[e]);
 rep(i,0,n) if (!lfound[i]) cover.push_back(i);
 rep(i,0,m) if (seen[i]) cover.push back(n+i);
 assert (sz(cover) == res);
 return cover;
```

Weighted Matching.h

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

```
Time: \mathcal{O}(N^2M)
                                                                              1e0fe93e, 31 lines
```

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

GeneralMatching.h

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

```
"../numerical/MatrixInverse-mod.h"
                                                    cb191272, 40 lines
vector<pii> generalMatching(int N, vector<pii>& ed) {
 vector<vector<ll>> mat(N, vector<ll>(N)), A;
 for (pii pa : ed) {
    int a = pa.first, b = pa.second, r = rand() % mod;
    mat[a][b] = r, mat[b][a] = (mod - r) % mod;
  int r = matInv(A = mat), M = 2*N - r, fi, fj;
  assert (r % 2 == 0);
  if (M != N) do {
    mat.resize(M, vector<ll>(M));
    rep(i,0,N) {
     mat[i].resize(M);
      rep(j,N,M) {
        int r = rand() % mod;
        mat[i][j] = r, mat[j][i] = (mod - r) % mod;
 } while (matInv(A = mat) != M);
  vi has (M, 1); vector<pii> ret;
  rep(it, 0, M/2) {
    rep(i,0,M) if (has[i])
      rep(j,i+1,M) if (A[i][j] && mat[i][j]) {
        fi = i; fj = j; goto done;
    } assert(0); done:
    if (fj < N) ret.emplace_back(fi, fj);</pre>
    has[fi] = has[fj] = 0;
    rep(sw, 0, 2) {
      11 a = modpow(A[fi][fj], mod-2);
      rep(i,0,M) if (has[i] && A[i][fj]) {
        ll b = A[i][fj] * a % mod;
        rep(j, 0, M) A[i][j] = (A[i][j] - A[fi][j] * b) % mod;
      swap(fi,fj);
 return ret;
```

7.4 DFS algorithms

SCC.h

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

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

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

Usage: TwoSat ts(number of boolean variables);

```
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;
  rep(i,0,n) if (comp[i] < 0) dfs(i, g, f);
}</pre>
```

BiconnectedComponents.h

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

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

2965e537, 33 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, e, y, top = me;
  for (auto pa : ed[at]) if (pa.second != par) {
    tie(y, e) = pa;
    if (num[y]) {
     top = min(top, num[y]);
     if (num[y] < me)
       st.push back(e);
    } else {
     int si = sz(st);
     int up = dfs(y, e, f);
      top = min(top, up);
     if (up == me) {
       st.push_back(e);
        f(vi(st.begin() + si, st.end()));
       st.resize(si);
     else if (up < me) st.push_back(e);</pre>
      else { /* e is a bridge */ }
  return top;
template<class F>
void bicomps(F f) {
  num.assign(sz(ed), 0);
  rep(i,0,sz(ed)) if (!num[i]) dfs(i, -1, f);
```

2SAT.h

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

```
ts.either(0, ~3); // Var 0 is true or var 3 is false
ts.setValue(2); // Var 2 is true
ts.atMostOne(\{0, \sim 1, 2\}); // <= 1 of vars 0, \sim 1 and 2 are true
ts.solve(); // Returns true iff it is solvable
ts.values[0..N-1] holds the assigned values to the vars
Time: \mathcal{O}(N+E), where N is the number of boolean variables, and E is the
number of clauses.
                                                    5f970691, 56 lines
struct TwoSat {
 int N;
 vector<vi> gr;
 vi values; // 0 = false, 1 = true
  TwoSat(int n = 0) : N(n), gr(2*n) {}
  int addVar() { // (optional)
    gr.emplace_back();
    gr.emplace back();
    return N++;
  void either(int f, int j) {
    f = max(2*f, -1-2*f);
    j = \max(2*j, -1-2*j);
    gr[f].push back(j^1);
    gr[j].push_back(f^1);
  void setValue(int x) { either(x, x); }
  void atMostOne(const vi& li) { // (optional)
    if (sz(li) <= 1) return;</pre>
    int cur = ~li[0];
    rep(i,2,sz(li)) {
     int next = addVar();
     either(cur, ~li[i]);
     either(cur, next);
     either(~li[i], next);
     cur = ~next;
    either(cur, ~li[1]);
 vi val, comp, z; int time = 0;
 int dfs(int i) {
    int low = val[i] = ++time, x; z.push_back(i);
    for(int e : gr[i]) if (!comp[e])
     low = min(low, val[e] ?: dfs(e));
    if (low == val[i]) do {
     x = z.back(); z.pop_back();
     comp[x] = low;
     if (values[x>>1] == -1)
       values[x>>1] = x&1;
    } while (x != i);
    return val[i] = low;
 bool solve() {
   values.assign(N, -1);
   val.assign(2*N, 0); comp = val;
    rep(i,0,2*N) if (!comp[i]) dfs(i);
    rep(i,0,N) if (comp[2*i] == comp[2*i+1]) return 0;
    return 1;
```

 ${\bf Euler Walk.h}$

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

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

7.5 Coloring

EdgeColoring.h

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

Time: $\mathcal{O}(NM)$ e210e25f, 31 lines

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

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.

16

9775a02f, 21 lines

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

```
b0d5b15b, 12 lines
```

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

MaximumClique.h

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

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

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

MaximumIndependentSet.h

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

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.

Usage: auto jmp = treeJump(par); lca(jmp, dep, a, b); **Time:** construction $\mathcal{O}(N \log N)$, queries $\mathcal{O}(\log N)$

9fb3804b, 22 lines

```
auto treeJump = [](vc<int> &a) {
  int d = 32 - __builtin_clz(sz(a));
  vc<vc<int>> jmp(d, a);
  for (int i = 1; i < d; ++i) for (int j = 0; j < sz(a); ++j)
    jmp[i][j] = jmp[i - 1][jmp[i - 1][j]];
  return jmp;
};
auto lca = [](vc<vc<int>> &tbl, vc<int> &dep, int a, int b) {
  if (dep[a] < dep[b]) swap(a, b);</pre>
  auto jmp = [&](int nod, int h) {
    for (int i = 0; i < sz(tbl); ++i) if (h & (1 << i))
      nod = tbl[i][nod];
    return nod;
  };
  a = jmp(a, dep[a] - dep[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.h

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

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

```
"../data-structures/RMQ.h"
struct LCA {
 int T = 0;
 vc<int> tim, pth, ret;
 RMO<int> rmg;
 LCA(vc<vc<int>> &g): tim(sz(g)), rmq((dfs(g, 0, -1), ret)) {}
 void dfs(vc<vc<int>> &q, int u, int p) {
   tim[u] = T++;
    for (int v: q[u]) if (v != p) {
     pth.emplace_back(u), ret.emplace_back(tim[u]);
     dfs(g, v, u);
 int lca(int a, int b) {
   if (a == b) return a;
   tie(a, b) = minmax(tim[a], tim[b]);
    return pth[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, origindex) representing a tree rooted at 0. The root points to itself.

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

"LCA.h"

```
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/SegmentTree.h"
```

```
4f3827a7, 46 lines
template<bool EDGE> struct HLD {
 int n, tim = 0;
  vc<vc<int>> q;
 vc<int> par, siz, dep, rt, pos;
  Sea tr:
  HLD(vc<vc<int>> g): n(sz(g)), g(g), par(n, -1), siz(n, 1),
                      dep(n), rt(n), pos(n), tr(n) {
                      dfsSiz(0); dfsHld(0); }
  void dfsSiz(int u) {
    if (~par[u]) g[u].erase(find(all(g[u]), par[u]));
    for (auto &v: g[u]) {
      par[v] = u, dep[v] = dep[u] + 1;
      dfsSiz(v);
      siz[u] += siz[v];
      if (siz[v] > siz[g[u][0]]) swap(v, g[u][0]);
  } //f3bc4162
 void dfsHld(int u) {
    pos[u] = tim++;
    for (auto v: q[u]) {
      rt[v] = (v == q[u][0] ? rt[u] : v);
      dfsHld(v);
  } //bac86e27
  template < class F > void process (int u, int v, F op) {
    for (; rt[u] != rt[v]; v = par[rt[v]]) {
      if (dep[rt[u]] > dep[rt[v]]) swap(u, v);
      op(pos[rt[v]], pos[v] + 1);
```

if (dep[u] > dep[v]) swap(u, v);

```
CSU
```

```
op(pos[u] + EDGE, pos[v] + 1);
  } //974632cb
  void modifyPath(int u, int v, int val) {
   process(u, v, [&](int 1, int r) { tr.modify(1, r, val); });
  int queryPath(int u, int v) {
   int res = 0;
   process(u, v, [&](int 1, int r) {
     (res += tr.query(1, r)) %= P;
   return res;
  int querySubtree(int u) { // modifySubtree is similar
    return tr.query(pos[u] + EDGE, pos[u] + siz[u]);
};
```

LinkCutTree.h

Description: link-cut Tree. Supports BST-like augmentations. (Can be used in place of HLD). Current implementation supports update value at a node, and query max on a path.

```
Time: All operations take amortized \mathcal{O}(\log N).
                                                      2534771f. 77 lines
struct Node {
  Node *p, *pp, *c[2];
 bool flip;
  int val, xval;
  void push() {
    if (flip) {
     if (c[0]) c[0]->flip ^= 1;
     if (c[1]) c[1]->flip ^= 1;
     swap(c[0], c[1]), flip = 0;
  void pull() {
   xval = val;
   if (c[0]) xval ^= c[0]->xval;
    if (c[1]) xval ^= c[1]->xval;
  bool dir() { return this == p->c[1]; }
  void rot() {
   bool t = dir();
   Node *y = p, *&z = c[!t];
   p = y->p;
   if (p) p->c[y->dir()] = this;
   if (z) z \rightarrow p = y;
   y - > c[t] = z;
   y->p = this;
    (z = y) \rightarrow pull();
  void g() { if (p) p->g(), pp = p->pp; push(); }
  void splay() {
    for (g(); p; rot()) if (p->p)
      (p->dir() == dir() ? p : this) ->rot();
    pull();
  Node *access() {
    for (Node *y = 0, *z = this; z; y = z, z = z->pp) {
     z->splav();
     if (z->c[1]) z->c[1]->pp = z, z->c[1]->p = 0;
     if (y) y->p=z;
     z -> c[1] = y;
     z->pull();
    splay();
    flip ^= 1;
    return this;
  } // 2060f82f
```

```
struct LinkCut {
 vc<Node> nodes;
 LinkCut(int n): nodes(n) {}
 bool cut(int u, int v) {
   Node *y = nodes[v].access();
   Node *x = nodes[u].access();
   if (x->c[0] != y || y->c[1]) return 0;
   x->c[0] = y->p = y->pp = 0;
   x \rightarrow pull():
   return 1;
 bool is_connected(int u, int v) {
   if (u == v) return 1;
   Node *x = nodes[u].access();
   Node *y = nodes[v].access();
   return x->p;
 bool link(int u, int v) {
   if (is_connected(u, v)) return 0;
   nodes[u].access()->pp = &nodes[v];
   return 1;
 } // 892ea56a
 void update(int u, int c) {
   nodes[u].access()->val = c;
 int query(int u, int v) {
   nodes[v].access();
   return nodes[u].access()->xval;
};
```

DirectedMST.h

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

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

```
"../data-structures/UnionFindRollback.h"
                                                     39e620f1, 60 lines
struct Edge { int a, b; ll w; };
struct Node {
 Edge key;
 Node *1, *r;
 11 delta;
  void prop()
   kev.w += delta:
   if (1) 1->delta += delta;
   if (r) r->delta += delta;
   delta = 0:
 Edge top() { prop(); return key; }
Node *merge(Node *a, Node *b) {
 if (!a || !b) return a ?: b;
 a->prop(), b->prop();
 if (a->key.w > b->key.w) swap(a, b);
  swap(a->1, (a->r = merge(b, a->r)));
  return a;
void pop(Node*& a) { a->prop(); a = merge(a->1, a->r); }
pair<ll, vi> dmst(int n, int r, vector<Edge>& g) {
 RollbackUF uf(n);
 vector<Node*> heap(n);
 for (Edge e : q) heap[e.b] = merge(heap[e.b], new Node{e});
 11 \text{ res} = 0;
 vi seen(n, -1), path(n), par(n);
 seen[r] = r;
 vector<Edge> Q(n), in(n, \{-1,-1\}), comp;
  deque<tuple<int, int, vector<Edge>>> cycs;
  rep(s,0,n) {
```

```
int u = s, qi = 0, w;
  while (seen[u] < 0) {</pre>
    if (!heap[u]) return {-1,{}};
    Edge e = heap[u] \rightarrow top();
    heap[u]->delta -= e.w, pop(heap[u]);
    Q[qi] = e, path[qi++] = u, seen[u] = s;
    res += e.w, u = uf.find(e.a);
    if (seen[u] == s) {
      Node * cvc = 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};
```

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 *i*th row and column and take the determinant; this yields the number of directed spanning trees rooted at i (if G is undirected, remove any row/column).

7.8.2 Erdős–Gallai theorem

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

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

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.)

```
template \langle class T \rangle int sgn(T x) \{ return (x > 0) - (x < 0); \}
template<class T>
struct Point {
  using P = Point;
  explicit Point (T x = 0, T y = 0): x(x), y(y) {}
 bool operator<(P p) const { return tie(x,y) < tie(p.x,p.y); }</pre>
 bool operator==(P p) const { return tie(x,y)==tie(p.x,p.y); }
 P operator + (P p) const { return P(x + p.x, y +p.y); }
```

```
P operator - (P p) const { return P(x - p.x, y -p.y); }
  P operator * (T d) const { return P(x * d, y *d); }
  P operator / (T d) const { return P(x / d, y /d); }
  T dot(P p) const { return x * p.x + y * p.y; }
 T cross(P p) const { return x * p.y - y * p.x; }
  T cross(P a, P b) const { return (a-*this).cross(b-*this); }
  T dist2() const { return x * x + y * y; }
  double dist() const { return sqrt((double)dist2()); }
  // angle to x-axis in interval [-pi, pi]
  double angle() const { return atan2(y, x); }
  P unit() const { return *this / dist(); } // makes dist()=1
  P perp() const { return P(-y, x); } // rotates +90 degrees
  P normal() const { return perp().unit(); }
  // returns point rotated 'a' radians ccw around the origin
  P rotate (double a) const {
   return P(x * cos(a) - y * sin(a), x * sin(a) + y * cos(a));
  friend ostream &operator<<(ostream &os, P p) {</pre>
    return os << "(" << p.x << "," << p.y << ")";
};
```

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" f6bf6b55, 4 lines
template<class P>
double lineDist(const P& a, const P& b, const P& p) {
 return (double) (b-a).cross(p-a)/(b-a).dist();

SegmentDistance.h

Description:

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

 $\begin{array}{ll} \textbf{Usage:} \; \texttt{Point} < \texttt{double} > \; \texttt{a, b(2,2), p(1,1);} \\ \texttt{bool onSegment = segDist(a,b,p)} \; < \; \texttt{1e-10;} \\ \end{array}$



nt.h" 5c88f46f, 6 lines

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

SegmentIntersection.h

Description:

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

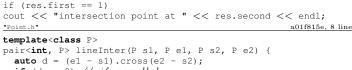


```
Usage: vector<P> inter = segInter(s1,e1,s2,e2);
if (sz(inter)==1)
cout << "segments intersect at " << inter[0] << endl;
"Point.h", "onSegment.h"
```

template<class P> vector<P> segInter(P a, P b, P c, P d) {

lineIntersection.h

Description:



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

sideOf.h

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

```
Usage: bool left = sideOf(p1,p2,q) ==1;
"Point.h"
```

OnSegment.h

Description: Returns true iff p lies on the line segment from s to e. Use $(segDist(s,e,p) \le psilon)$ instead when using Point double.

linearTransformation.h

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

```
"Point.h" 03a3061b, 6 lines
```

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

```
Usage: vector<Angle> v = \{w[0], w[0].t360() ...\}; // sorted int j = 0; rep(i,0,n) \{ while (v[j] < v[i].t180()) ++j; \} // sweeps j such that (j-i) represents the number of positively oriented triangles with vertices at 0 and i 0f0602c7, 35 lines
```

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

8.2 Circles

CircleIntersection.h

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

"Point.h"

84d6d345, 11 lines

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

CircleTangents.h

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

"Point.h" b0153d0e, 13 lines

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

CirclePolygonIntersection.h

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

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

```
"../../content/geometry/Point.h"
                                                     a1ee63d6, 19 lines
typedef Point < double > P;
#define arg(p, q) atan2(p.cross(q), p.dot(q))
double circlePoly(P c, double r, vector<P> ps) {
  auto tri = [&] (P p, P q) {
   auto r2 = r * r / 2;
   P d = q - p;
    auto a = d.dot(p)/d.dist2(), b = (p.dist2()-r*r)/d.dist2();
   auto det = a * a - b;
   if (det <= 0) return arg(p, q) * r2;</pre>
    auto s = max(0., -a-sqrt(det)), t = min(1., -a+sqrt(det));
   if (t < 0 || 1 <= s) return arg(p, g) * r2;</pre>
   P u = p + d * s, v = p + d * t;
   return arg(p,u) * r2 + u.cross(v)/2 + arg(v,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.



"Point.h" 1caa3aea, 9 lines

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

MinimumEnclosingCircle.h

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

8.3 Polygons

InsidePolygon.h

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

```
Usage: vectorP> v = \{P\{4,4\}, P\{1,2\}, P\{2,1\}\};bool in = inPolygon(v, P\{3, 3\}, false);
Time: O(n)
```

"Point.h", "OnSegment.h", "SegmentDistance.h" 2bf504ba, 11 lines

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

PolygonArea.h

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

"Point.h"

f1230037, 6 lines

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

PolygonCenter.h

"Point.h"

Description: Returns the center of mass for a polygon. **Time:** $\mathcal{O}(n)$

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

PolygonCut.h

Description:

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

```
Usage: vector\langle P \rangle p = ...;

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



ConvexHull.h

Description:

Returns a vector of the points of the convex hull in counterclockwise order. Points on the edge of the hull between two other points are not considered part of the hull.



31095458, 13 lines

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

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

HullDiameter.h

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

"Point.h" c571b8ed, 12 lines

```
typedef Point<11> P;
array<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 convex hull (CCW order, with no collinear points). Returns true if point lies within the hull. If strict is true, points on the boundary aren't included.

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

9706dcc8, 9 lines

```
"Point.h", "sideOf.h", "OnSegment.h" 71446bda, 14 lines
```

typedef Point<11> P;

LineHullIntersection ClosestPair kdTree FastDelaunay

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

LineHullIntersection.h

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

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

```
"Point.h"
                                                    7cf45b1b, 39 lines
#define cmp(i,j) sqn(dir.perp().cross(poly[(i)%n]-poly[(j)%n]))
#define extr(i) cmp(i + 1, i) >= 0 && cmp(i, i - 1 + n) < 0
template <class P> int extrVertex(vector<P>& poly, P dir) {
 int n = sz(poly), lo = 0, hi = n;
  if (extr(0)) return 0;
  while (lo + 1 < hi) {
   int m = (lo + hi) / 2;
   if (extr(m)) return m;
   int 1s = cmp(1o + 1, 1o), ms = cmp(m + 1, m);
    (ls < ms \mid | (ls == ms \&\& ls == cmp(lo, m)) ? hi : lo) = m;
  return lo;
#define cmpL(i) sqn(a.cross(poly[i], b))
template <class P>
array<int, 2> lineHull(P a, P b, vector<P>& poly) {
 int endA = extrVertex(poly, (a - b).perp());
  int endB = extrVertex(poly, (b - a).perp());
  if (cmpL(endA) < 0 \mid | cmpL(endB) > 0)
   return {-1, -1};
  array<int, 2> res;
  rep(i, 0, 2) {
   int lo = endB, hi = endA, n = sz(poly);
    while ((lo + 1) % n != hi) {
     int m = ((lo + hi + (lo < hi ? 0 : n)) / 2) % n;</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;
```

8.4 Misc. Point Set Problems

ClosestPair.h

Description: Finds the closest pair of points.

```
Time: \mathcal{O}(n \log n)
"Point.h"
                                                      d58e8aa8, 17 lines
using P = Point<double>;
auto closest = [](vc<P> v) {
 assert (sz(v) > 1);
 set<P> S;
 sort(all(v), [](P a, P b) { return a.y < b.y; });</pre>
  pair<ll,pair<P, P>> ret{numeric_limits<ll>::max(),{P(),P()}};
  int j = 0;
  for (P p: v) {
    P d(1 + 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);
  return ret.second;
}; // 4e29a1c8
```

kdTree.h

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

bac5b040, 63 lines

```
typedef long long T;
```

```
typedef Point<T> P;
const T INF = numeric_limits<T>::max();
bool on x(const P& a, const P& b) { return a.x < b.x; }
bool on_y(const P& a, const P& b) { return a.y < b.y; }</pre>
struct Node {
 P pt; // if this is a leaf, the single point in it
 T x0 = INF, x1 = -INF, y0 = INF, y1 = -INF; // bounds
  Node *first = 0, *second = 0;
 T distance (const P& p) { // min squared distance to a point
    T x = (p.x < x0 ? x0 : p.x > x1 ? x1 : p.x);
    T y = (p.y < y0 ? y0 : p.y > y1 ? y1 : p.y);
    return (P(x,y) - p).dist2();
  Node (vector<P>&& vp) : pt(vp[0]) {
    for (P p : vp) {
      x0 = min(x0, p.x); x1 = max(x1, p.x);
      y0 = min(y0, p.y); y1 = max(y1, p.y);
    if (vp.size() > 1) {
      // split on x if width >= height (not ideal...)
      sort(all(vp), x1 - x0 >= y1 - y0 ? on_x : on_y);
      // divide by taking half the array for each child (not
      // best performance with many duplicates in the middle)
      int half = sz(vp)/2;
      first = new Node({vp.begin(), vp.begin() + half});
      second = new Node({vp.begin() + half, vp.end()});
};
struct KDTree {
  Node* root;
  KDTree(const vector<P>& vp) : root(new Node({all(vp)})) {}
  pair<T, P> search(Node *node, const P& p) {
    if (!node->first) {
      // uncomment if we should not find the point itself:
      // if (p = node \rightarrow pt) return \{INF, P()\};
      return make_pair((p - node->pt).dist2(), node->pt);
```

```
Node *f = node->first, *s = node->second;
    T bfirst = f->distance(p), bsec = s->distance(p);
    if (bfirst > bsec) swap(bsec, bfirst), swap(f, s);
    // search closest side first, other side if needed
    auto best = search(f, p);
    if (bsec < best.first)</pre>
      best = min(best, search(s, p));
    return best:
  // find nearest point to a point, and its squared distance
  // (requires an arbitrary operator< for Point)
 pair<T, P> nearest (const P& p) {
    return search(root, p);
};
```

FastDelaunav.h

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

```
Time: \mathcal{O}(n \log n)
"Point.h"
                                                      eefdf506, 88 lines
typedef Point<ll> P;
typedef struct Ouad* O;
typedef __int128_t 111; // (can be ll if coords are < 2e4)
P arb(LLONG_MAX, LLONG_MAX); // not equal to any other point
struct Ouad {
  O rot, o; P p = arb; bool mark;
  P& F() { return r()->p; }
  Q& r() { return rot->rot; }
  O 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 \rightarrow rot, r \rightarrow p = arb, r \rightarrow o = i & 1 ? <math>r : r \rightarrow r();
  r->p = orig; r->F() = dest;
  return r;
void splice(Q a, Q b) {
  swap(a->o->rot->o, b->o->rot->o); swap(a->o, b->o);
Q connect(Q a, Q b) {
  Q = makeEdge(a->F(), b->p);
  splice(q, a->next());
  splice(q->r(), b);
  return q;
pair<Q,Q> rec(const vector<P>& s) {
  if (sz(s) <= 3) {
    Q = makeEdge(s[0], s[1]), b = makeEdge(s[1], s.back());
    if (sz(s) == 2) return { a, a->r() };
    splice(a->r(), b);
    auto side = s[0].cross(s[1], s[2]);
```

0 c = side ? connect(b, a) : 0;

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

3D

PolyhedronVolume.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;
```

Description: Class to handle points in 3D space. T can be e.g. double or 8058aeda, 32 lines

```
template<class T> struct Point3D {
  typedef Point3D P;
  typedef const P& R;
  T x, y, z;
  explicit Point3D(T x=0, T y=0, T z=0) : x(x), y(y), z(z) {}
  bool operator<(R p) const {</pre>
```

```
return tie(x, y, z) < tie(p.x, p.y, p.z); }</pre>
 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: $\mathcal{O}\left(n^2\right)$

typedef Point3D<double> P3;

```
"Point3D.h"
                                                                          5b45fc3f, 49 lines
```

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

swap(FS[j--], FS.back());

FS.pop_back();

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

sphericalDistance.h

Description: Returns the shortest distance on the sphere with radius radius between the points with azimuthal angles (longitude) f1 (ϕ_1) and f2 (ϕ_2) from x axis and zenith angles (latitude) t1 (θ_1) and t2 (θ_2) from z axis (0 = north pole). All angles measured in radians. The algorithm starts by converting the spherical coordinates to cartesian coordinates so if that is what you have you can use only the two last rows. dx*radius is then the difference between the two points in the x direction and d*radius is the total distance between the points. 611f0797, 8 lines

```
double sphericalDistance(double f1, double t1,
    double f2, double t2, double radius) {
  double dx = \sin(t2) \cdot \cos(f2) - \sin(t1) \cdot \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: p[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: $\mathcal{O}(n)$

```
f84022ea, 17 lines
auto getBorder = [](string s) {
 vc<int> p(sz(s));
  for (int i = 1; i < sz(s); ++i) {
    int q = p[i - 1];
    while (g \&\& s[g] != s[i]) g = p[g - 1];
    p[i] = g + (s[g] == s[i]);
 return p;
}; // 43f874fa
auto KMP = [&getBorder](string s, string pat) {
 vc < int > p = getBorder(pat + ' \ 0' + s), ans;
 for (int i = sz(p) - sz(s); i < sz(p); ++i) {
   if (p[i] == sz(pat)) ans.emplace back(i - 2 * sz(pat));
 return ans;
}; // 8ffe484c
```

Zfunc.h

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

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

```
3ae5260c, 12 lines
vi Z(string S) {
 vi z(sz(S));
 int 1 = -1, r = -1;
 rep(i,1,sz(S)) {
```

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

Manacher.h

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

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

e7ad794a, 13 lines

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

MinRotation.h

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

Description: Finds the lexicographically smallest rotation of a string.

Usage: rotate(v.beqin(), v.beqin()+minRotation(v), v.end());

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

SuffixArrav.h

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

Time: $\mathcal{O}(n \log n)$ 195132ce, 24 lines

```
struct SA {
 vc<int> sa, ht, rk;
  SA(string s, int lim = 256) {
   int n = sz(s) + 1, k = 0, a, b, i, j, p;
   vc<int> x(all(s) + 1), y(n), ws(max(n, lim));
    rk = sa = ht = y, iota(all(sa), 0);
    for (j = 0, p = 0; p < n; j = max(1, j * 2), lim = p) {
     p = j, iota(all(v), n - j);
     for (i = 0; i < n; ++i) if (sa[i] >= j) y[p++] = sa[i]-j;
     fill(all(ws), 0);
     for (i = 0; i < n; ++i) ++ws[x[i]];
     for (i = 1; i < lim; ++i) ws[i] += ws[i - 1];</pre>
      for (i = n; i--; ) sa[--ws[x[y[i]]]] = y[i];
      swap(x, y), p = 1, x[sa[0]] = 0;
      for (i = 1; i < n; ++i) {
       a = sa[i - 1], b = sa[i];
       x[b] = (y[a] == y[b] && y[a+j] == y[b+j]) ? p - 1 : p++;
```

```
for (i = 1; i < n; ++i) rk[sa[i]] = i;
for (i = 0; i < n - 1; ht[rk[i++]] = k)
    for (k && --k, j = sa[rk[i] - 1]; s[i+k] == s[j+k]; ++k);
}
};</pre>
```

Hashing.h

Description: Self-explanatory methods for string hashing. 3f02d837, 44 lines

```
// Arithmetic mod 2^64-1. 2x slower than mod 2^64 and more
// code, but works on evil test data (e.g. Thue-Morse, where
// ABBA... and BAAB... of length 2^10 hash the same mod 2^64).
// "typedef ull H;" instead if you think test data is random,
// or work mod 10^9+7 if the Birthday paradox is not a problem.
struct H {
  typedef uint64_t ull;
 ull x; H(ull x=0) : x(x) {}
#define OP(O,A,B) H operator O(H o) { ull r = x; asm \
  (A "addq %%rdx, %0\n adcq $0,%0" : "+a"(r) : B); return r; }
  OP(+,,"d"(o.x)) OP(*,"mul %1\n", "r"(o.x) : "rdx")
 H operator-(H o) { return *this + ~o.x; }
 ull get() const { return x + !~x; }
 bool operator==(H o) const { return get() == o.get(); }
 bool operator<(H o) const { return get() < o.get(); }</pre>
static const H C = (11)1e11+3; // (order \sim 3e9; random \ also \ ok)
struct HashInterval {
  vector<H> ha, pw;
  HashInterval(string& str) : ha(sz(str)+1), pw(ha) {
    pw[0] = 1;
    rep(i, 0, sz(str))
     ha[i+1] = ha[i] * C + str[i],
      pw[i+1] = pw[i] * C;
 H hashInterval(int a, int b) { // hash [a, b)
    return ha[b] - ha[a] * pw[b - a];
vector<H> getHashes(string& str, int length) {
 if (sz(str) < length) return {};</pre>
 H h = 0, pw = 1;
 rep(i,0,length)
   h = h * C + str[i], pw = pw * C;
  vector<H> ret = {h};
  rep(i,length,sz(str)) {
    ret.push_back(h = h * C + str[i] - pw * str[i-length]);
 return ret;
H hashString(string& s){H h{}; for(char c:s) h=h*C+c;return h;}
```

AhoCorasick.h

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

Time: construction takes $\mathcal{O}(26N)$, where N= sum of length of patterns. find(x) is $\mathcal{O}(N)$, where N= length of x. findAll is $\mathcal{O}(NM)$. f35677c4, 66 lines

```
struct AhoCorasick {
  enum {alpha = 26, first = 'A'}; // change this!
```

```
struct Node {
  // (nmatches is optional)
  int back, next[alpha], start = -1, end = -1, nmatches = 0;
  Node(int v) { memset(next, v, sizeof(next)); }
vector<Node> N:
vi backp;
void insert(string& s, int j) {
  assert(!s.empty());
  int n = 0:
  for (char c : s) {
    int& m = N[n].next[c - first];
    if (m == -1) { n = m = sz(N); N.emplace_back(-1); }
    else n = m;
  if (N[n].end == -1) N[n].start = j;
  backp.push_back(N[n].end);
  N[n].end = j;
  N[n].nmatches++;
AhoCorasick(vector<string>& pat) : N(1, -1) {
  rep(i,0,sz(pat)) insert(pat[i], i);
  N[0].back = sz(N);
  N.emplace_back(0);
  queue<int> q;
  for (q.push(0); !q.empty(); q.pop()) {
    int n = q.front(), prev = N[n].back;
    rep(i,0,alpha) {
      int &ed = N[n].next[i], y = N[prev].next[i];
      if (ed == -1) ed = y;
      else {
       N[ed].back = v;
        (N[ed].end == -1 ? N[ed].end : backp[N[ed].start])
          = N[y].end;
        N[ed].nmatches += N[y].nmatches;
        q.push(ed);
vi find(string word) {
  int n = 0;
  vi res; // ll count = 0:
  for (char c : word) {
    n = N[n].next[c - first];
    res.push back(N[n].end);
    // count \neq 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: \mathcal{O}(\log N)
                                                    edce4766, 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;
```

IntervalCover.h

if (R != r2) is.emplace(R, r2);

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

Time: $\mathcal{O}(N \log N)$ 9e9d8de7, 19 lines

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

ConstantIntervals.h

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

```
 \begin{array}{ll} \textbf{Usage:} & \text{constantIntervals}(0, \text{sz(v), [\&](int x)}\{\text{return v[x];}\}, \\ \text{[\&](int lo, int hi, T val)}\{\ldots\}); \\ \textbf{Time:} & \mathcal{O}\left(k\log\frac{n}{k}\right) \\ & \text{753a4cd9, 19 lines} \end{array}
```

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

10.2 Misc. algorithms

TernarySearch.h

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

Usage: int ind = ternSearch(0,n-1,[&](int i){return a[i];}); Time: $\mathcal{O}(\log(b-a))$ 9155b4fb, 11 lines

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

LIS.h

Description: Compute indices for the longest increasing subsequence. **Time:** $\mathcal{O}\left(N\log N\right)$ 2932a052, 17 lines

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

10.3 Dynamic programming

KnuthDP.h

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

```
DivideAndConquerDP.h
```

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

```
d38d2b27, 18 lines
struct DP { // Modify at will:
 int lo(int ind) { return 0; }
 int hi(int ind) { return ind; }
 11 f(int ind, int k) { return dp[ind][k]; }
 void store(int ind, int k, ll v) { res[ind] = pii(k, v); }
 void rec(int L, int R, int LO, int HI) {
   if (L >= R) return;
   int mid = (L + R) \gg 1;
   pair<11, int> best(LLONG_MAX, LO);
    rep(k, max(LO,lo(mid)), min(HI,hi(mid)))
     best = min(best, make_pair(f(mid, k), k));
    store(mid, best.second, best.first);
   rec(L, mid, LO, best.second+1);
   rec(mid+1, R, best.second, HI);
 void solve(int L, int R) { rec(L, R, INT_MIN, INT_MAX); }
};
```

10.4 Debugging tricks

- signal (SIGSEGV, [] (int) { _Exit(0); }); converts segfaults into Wrong Answers. Similarly one can catch SIGABRT (assertion failures) and SIGFPE (zero divisions). _GLIBCXX_DEBUG failures generate SIGABRT (or SIGSEGV on gcc 5.4.0 apparently).
- feenableexcept (29); kills the program on NaNs (1), 0-divs (4), infinities (8) and denormals (16).

10.5 Optimization tricks

__builtin_ia32_ldmxcsr(40896); disables denormals (which make floats 20x slower near their minimum value).

10.5.1 Bit hacks

- x & -x is the least bit in x.
- for (int x = m; x;) { --x &= m; ... } loops over all subset masks of m (except m itself).
- c = x&-x, r = x+c; (((r^x) >> 2)/c) | r is the next number after x with the same number of bits set.
- rep(b,0,K) rep(i,0,(1 << K))
 if (i & 1 << b) D[i] += D[i^(1 << b)];
 computes all sums of subsets.</pre>

10.5.2 Pragmas

- #pragma GCC optimize ("Ofast") will make GCC auto-vectorize loops and optimizes floating points better.
- #pragma GCC target ("avx2") can double performance of vectorized code, but causes crashes on old machines.
- #pragma GCC optimize ("trapv") kills the program on integer overflows (but is really slow).

FastMod.h

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

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

FastInput.h

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.

7b3c7051, 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.

745db225, 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

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

"BumpAllocator.h"

2dd6c977, 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
```

Usage: vector<vector<int, small<int>>> ed(N); bb66d422, 14 lines char buf[450 << 20] alignas(16);

```
size_t buf_ind = sizeof buf;
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) {}
};
```

SIMD.h

Description: Cheat sheet of SSE/AVX intrinsics, for doing arithmetic on several numbers at once. Can provide a constant factor improvement of about 4, orthogonal to loop unrolling. Operations follow the pattern "_mm(256)?_name_(si(128|256)|epi(8|16|32|64)|pd|ps)". Not all are described here; grep for _mm_ in /usr/lib/gcc/*/4.9/include/ for more. If AVX is unsupported, try 128-bit operations, "emmintrin.h" and #define __SSE__ and __MMX__ before including it. For aligned memory use _mm_malloc(size, 32) or int buf[N] alignas(32), but prefer loadu/storeu.

```
#pragma GCC target ("avx2") // or sse4.1
#include "immintrin.h"
typedef __m256i mi;
#define L(x) _mm256_loadu_si256((mi*)&(x))
// High-level/specific methods:
// load(u)?\_si256, store(u)?\_si256, setzero\_si256, \_mm\_malloc
// blendv_(epi8|ps|pd) (z?y:x), movemask_epi8 (hibits of bytes)
// i32gather_epi32(addr, x, 4): map addr[] over 32-b parts of x
// sad_epu8: sum of absolute differences of u8, outputs 4xi64
// maddubs_epi16: dot product of unsigned i7's, outputs 16xi15
// madd_epi16: dot product of signed i16's, outputs 8xi32
// extractf128_si256(, i) (256->128), cvtsi128_si32 (128->lo32)
// permute2f128_si256(x,x,1) swaps 128-bit lanes
// shuffle_epi32(x, 3*64+2*16+1*4+0) == x for each lane
// shuffle_epi8(x, y) takes a vector instead of an imm
// Methods that work with most data types (append e.g. _epi32):
// set1. blend (i8?x:y), add, adds (sat.), mullo, sub, and/or,
// and not, abs, min, max, sign(1,x), cmp(gt|eq), unpack(lo|hi)
int sumi32(mi m) { union {int v[8]; mi m;} u; u.m = m;
 int ret = 0; rep(i,0,8) ret += u.v[i]; return ret; }
mi zero() { return _mm256_setzero_si256(); }
mi one() { return _mm256_set1_epi32(-1); }
bool all_zero(mi m) { return _mm256_testz_si256(m, m); }
bool all_one(mi m) { return _mm256_testc_si256(m, one()); }
11 example_filteredDotProduct(int n, short* a, short* b) {
```

```
int i = 0; 11 r = 0;
mi zero = _mm256_setzero_si256(), acc = zero;
while (i + 16 <= n) {
    mi va = L(a[i]), vb = L(b[i]); i += 16;
    va = _mm256_and_si256(_mm256_cmpgt_epi16(vb, va), va);
    mi vp = _mm256_madd_epi16(va, vb);
    acc = _mm256_add_epi64(_mm256_unpacklo_epi32(vp, zero),
        _mm256_add_epi64(acc, _mm256_unpacklo_epi32(vp, zero)));
}
union {11 v[4]; mi m;} u; u.m = acc; rep(i,0,4) r += u.v[i];
for (;i<n;++i) if (a[i] < b[i]) r += a[i]*b[i]; // <- equiv
return r;</pre>
```