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Lebenslangerschicksalsschatz

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Makefile hash hash-cpp

Contest (1)

Makefile

 $CXX = \alpha + +$ CXXFLAGS = -std=c++17 -O2 -Wall -Wextra -pedantic -Wshadow Wformat=2 -Wfloat-equal -Wconversion -Wlogical-op -Wshiftoverflow=2 -Wduplicated-cond -Wcast-qual -Wcast-align -Wno -unused-result -Wno-sign-conversion DEBUGFLAGS = -D_GLIBCXX_DEBUG -D_GLIBCXX_DEBUG_PEDANTIC -DLOCAL

-fsanitize=address -fsanitize=undefined -fno-sanitizerecover=all -fstack-protector -D FORTIFY SOURCE=2

DEBUG = falseifeq (\$(DEBUG),true) CXXFLAGS += \$ (DEBUGFLAGS)

hash.sh

Hashes a file, ignoring all whitespace and comments. Use for # verifying that code was correctly typed. cpp -dD -P -fpreprocessed | tr -d '[:space:]' | md5sum |cut -c-6

hash-cpp.sh

Hashes a file, ignoring all whitespace, comments and defines # verifying that code was correctly typed.

First do: chmod +x ./hash-cpp.sh

./hash-cpp.sh *.cpp start end

sed -n \$2','\$3' p' \$1 | sed '/^#w/d' | cpp -dD -P fpreprocessed | tr -d '[:space:]' | md5sum |cut -c-6

Mathematics (2)

2.1 Equations

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

The extremum is given by x = -b/2a.

$$ax + by = e$$

$$cx + dy = f$$

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

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

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

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

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

Trigonometry

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

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

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

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

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

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

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

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

2.3 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):

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

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

Pick's: A polygon on an integer grid strictly containing i lattice points and having b lattice points on the boundary has area $i + \frac{b}{2} - 1$. (Nothing similar in higher dimensions)

2.3.2 Quadrilaterals

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

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

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

2.3.3 Spherical coordinates



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

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

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

2.4 Derivatives/Integrals

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

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

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

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

Integration by parts:

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

Green's theorem:

Let C be a positive, smooth, simple curve. D is a region bounded by C.

$$\oint_C (Pdx + Qdy) = \int \int_D \left(\frac{\partial Q}{\partial x} - \frac{\partial P}{\partial y}\right)$$

To calculate area, $\frac{\partial Q}{\partial x} - \frac{\partial P}{\partial y} = 1$, usually, picking $Q = \frac{1}{2}x$ and $P = -\frac{1}{2}y$ suffice.

Then we have

$$\frac{1}{2} \oint_C x dy - \frac{1}{2} \oint_C y dx$$

Line integral:

C given by $x = x(t), y = y(t), t \in [a, b]$, then

$$\oint_C f(x,y)ds = \int_a^b f(x(t),y(t))ds$$

where,
$$ds = \sqrt{(\frac{dx}{dt})^2 + (\frac{dy}{dt})^2} dt$$
 or $\sqrt{(1 + (\frac{dy}{dx})^2} dx$

2.4.1 XOR sum

$$\bigoplus_{x=0}^{n-1} x = \{0, n-1, 1, n\}[n \operatorname{mod} 4]$$

$$\bigoplus_{x=l}^{r-1} x = \bigoplus_{a=0}^{r-1} a \oplus \bigoplus_{b=0}^{l-1} b$$

2.5 Series

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

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

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

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

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

2.6 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,

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

2.6.1 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 $Po(\lambda)$, $\lambda = t\kappa$.

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

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

2.6.2 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 $\text{Exp}(\lambda)$, $\lambda > 0$.

$$f(x) = \begin{cases} \lambda e^{-\lambda x} & x \ge 0\\ 0 & x < 0 \end{cases}$$

$$\mu = \frac{1}{\lambda}, \, \sigma^2 = \frac{1}{\lambda^2}$$

Normal distribution

Most real random values with mean 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

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

2.7 Markov chains

A Markov chain is a discrete random process with the property that the next state depends only on the current state. Let X_1, X_2, \ldots be a sequence of random variables generated by the Markov process. Then there is a transition matrix $\mathbf{P} = (p_{ij})$, with $p_{ij} = \Pr(X_n = i | X_{n-1} = j)$, 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 absorbing chain if

- 1. there is at least one absorbing state and
- **2**. it is possible to go from any state to at least one absorbing state in a finite number of steps.

A Markov chain is an A-chain if the states can be partitioned into two sets \mathbf{A} and \mathbf{G} , such that all states in \mathbf{A} are absorbing $(p_{ii}=1)$, and all states in \mathbf{G} leads to an absorbing state in \mathbf{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)

order-statistic-tree.h

Description: A set (not multiset!) with support for finding the n'th element, and finding the index of an element.

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

```
cbits/extc++.h> acfa21, 19 lines
template <typename K, typename V, typename Comp = std::less<K>>
using ordered_map = __gnu_pbds::tree
K, V, Comp,
__gnu_pbds::rb_tree_tag,
__gnu_pbds::tree_order_statistics_node_update
>;
template <typename K, typename Comp = std::less<K>>
using ordered_set = ordered_map<K, __gnu_pbds::null_type, Comp
>;
void example() {
    ordered_set<int> t, t2; t.insert(8);
```

```
auto it = t.insert(10).first;
assert(it == t.lower_bound(9));
assert(t.order_of_key(10) == 1); // num strictly smaller
assert(t.order_of_key(11) == 2);
assert(*t.find_by_order(0) == 8);
t.join(t2); // assuming T < T2 or T > T2, merge t2 into t
```

dsu.h

Description: Disjoint-set data structure. **Time:** $\mathcal{O}(\alpha(N))$

Time: $O(\alpha(N))$ 7d5db8, 14 lines struct UF { vector<int> e; UF (int n) : e(n, -1) {} bool same_set (int a, int b) { return find(a) == find(b); } int size(int x) { return -e[find(x)]; } int find(int x) { return e[x] < 0 ? x : e[x] = find(e[x]); } bool unite(int a, int b) { a = find(a), b = find(b); if (a == b) return 0; if (e[a] > e[b]) swap(a, b); e[a] += e[b]; e[b] = a; return 1; }

dsu-rollback.h

};

Description: Disjoint-set data structure with undo.

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

7ddf1d, 21 lines

da8783, 41 lines

```
struct RollbackUF {
  vector<int> e; vector<pair<int,int>> st;
  RollbackUF(int n) : e(n, -1) {}
  int size(int x) { return -e[find(x)]; }
  int find(int x) { return e[x] < 0 ? x : find(e[x]); }
  int time() { return st.size(); }
  void rollback(int t) {
   for (int i = time(); i --> t;)
     e[st[i].first] = st[i].second;
   st.resize(t);
  bool unite(int a, int b) {
   a = find(a), b = find(b);
   if (a == b) return false;
   if (e[a] > e[b]) swap(a, b);
   st.push back({a, e[a]});
   st.push_back({b, e[b]});
   e[a] += e[b]; e[b] = a;
    return true:
};
```

monotonic-queue.h

Description: Structure that supports all operations of a queue and get the minimum/maximum active value in the queue. Useful for sliding window 1D and 2D. For 2D problems, you will need to pre-compute another matrix, by making a row-wise traversal, and calculating the min/max value beginning in each cell. Then you just make a column-wise traverse as they were each an independent array.

Time: $\mathcal{O}(1)$

template<typename T> struct monotonic_queue {
 vector<T> as, aas;
 vector<T> bs, bbs;
 void reserve(int N) {
 as.reserve(N); aas.reserve(N);
 bs.reserve(N); bbs.reserve(N);

```
void reduce() {
   while (!bs.empty()) {
     as.push_back(bs.back());
     aas.push_back(aas.empty() ? bs.back() : (bs.back() * aas.
          back()));
     bs.pop_back(); bbs.pop_back();
 T get() {
   if (as.empty()) reduce();
    return (bbs.empty() ? aas.back() : (aas.back() * bbs.back()
 bool empty() const { return (as.empty() && bs.empty()); }
 int size() const { return int(as.size()) + int(bs.size()); }
    if (as.empty()) reduce();
    return as.back();
 void push(const T& val) {
   bs.push_back(val);
   bbs.push back(bbs.empty() ? val : (bbs.back() * val));
 void pop() {
   if (as.empty()) reduce();
   as.pop_back();
   aas.pop_back();
};
struct affine_t {
 int64 t b, c;
 affine_t operator*(affine_t rhs) {
    return { (rhs.b * b) % M, (rhs.b * c + rhs.c) % M};
};
```

point-context.h

Description: Examples of Segment Tree

3f5b19, 36 lines

```
struct seg node { // bbfc07
 int val:
 int mi, ma;
 seg_node() : mi(INT_MAX), ma(INT_MIN), val(0) {}
 seg_node(int x) : mi(x), ma(x), val(x) {}
 void merge(const seg_node& 1, const seg_node& r) {
   val = 1.val + r.val;
   mi = min(l.mi, r.mi);
   ma = max(1.ma, r.ma);
 void update(int x) {
   mi = ma = val = x;
 bool acc min(int& acc, int x) const {
   if (x >= mi) return true;
   if (acc > mi) acc = mi;
   return false;
 bool acc_max(int& acc, int x) const {
   if (x <= ma) return true;
   if (acc < ma) acc = ma;
   return false:
};
// 1 + min of (a, N) \le x
auto find_min_right = [&](seqtree<seq_node>& sq, int a, int x)
    -> int {
```

rec-lazy-segtree.h

Description: Segment Tree with Lazy update (half-open interval). **Time:** $\mathcal{O}(\lg(N) * Q)$

```
6f5fdf, 71 lines
template<class T> struct segtree_range {
 int N:
 vector<T> ts;
  segtree_range() {}
  explicit segtree_range(int N_{-}) : N(1 << _{-}1g(2*N_{-}1)) {
    ts.resize(2*N); build();
 template<class O> explicit segtree range(const vector<O>& A) {
    const int N_ = int(A.size());
    N = (1 << __lq(2*N_-1));
    ts.resize(2*N);
    for (int i = 0; i < N_{i} ++i) at (i) = T(A[i]);
    build();
 T& at(int a) { return ts[a + N]; }
  void build() { for (int a = N; --a; ) merge(a); }
  inline void push(int a) { ts[a].push(ts[2*a], ts[2*a+1]); }
  inline void merge(int a) { ts[a].merge(ts[2*a], ts[2*a+1]); }
  template < class Op, class E, class F, class... Args>
  auto query (int v, int 1, int r, int a, int b, Op op, E e, F f
       , Args&&... args) {
    if (1 >= b || r <= a) return e();
    if (1 >= a && r <= b) {
      return (ts[v].*f) (args...);
    int m = (1 + r)/2;
    push(v);
    return op (query (2*v, 1, m, a, b, op, args...), query (2*v+1, m, a, b, op, args...)
          m, r, a, b, op, args...));
 template < class Op, class E, class F, class... Args>
  auto query(int a, int b, Op op, E e, F f, Args&&... args) {
    return query(1, 0, N, a, b, op, e, f, args...);
 T query(int v, int 1, int r, int a, int b) {
   if (1 \ge b \mid | r \le a) return T();
    if (1 >= a && r <= b) return ts[v];</pre>
    int m = (1 + r)/2;
    push(v); T t;
    t.merge(query(2*v, 1, m, a, b), query(2*v+1, m, r, a, b));
    return t;
 T query(int a, int b) { return query(1, 0, N, a, b); }
 template < class F, class... Args > void update (int v, int 1,
       int r, int a, int b, F f, Args&&... args) {
    if (1 >= b || r <= a) return;
    if (1 >= a && r <= b) {
      (ts[v].*f) (args...);
      return;
    int m = (1 + r)/2;
    push (v);
    update(2*v, 1, m, a, b, f, args...);
    update(2*v+1, m, r, a, b, f, args...);
```

```
merge(v);
  template<class F, class... Args>
  void update(int a, int b, F f, Args&&... args) {
   update(1, 0, N, a, b, f, args...);
  template < class F, class... Args > int find first (int v, int 1,
       int r, int a, int b, F f, Args&&... args) {
    if (1 \ge b \mid | r \le a \mid | !(ts[v].*f)(args...)) return -1;
    if (1 + 1 == r) return 1;
    int m = (1 + r)/2;
   push(v);
    int cur = find_first(2*v, 1, m, a, b, f, args...);
    if (cur == -1) cur = find first(2*v+1, m, r, a, b, f, args
    return cur;
  template<class F, class... Args>
  int find_first(int a, int b, F f, Args&&... args) {
    return find_first(1, 0, N, a, b, f, args...);
};
```

lazy-context.h

Description: Examples of Segment Tree with Lazy update 44304b, 146 lines

```
struct seq_node {
  int sz, lz; int64 t sum;
  seg node() : sz(1), sum(0), lz(-1) {}
  seg node(int64 t val) : sz(1), sum(val), lz(-1) {}
  void push(const seg node& 1, const seg node& r) {
   if (1z == 2) {
     l.flip(lz);
     r.flip(lz);
    } else if (lz != -1) {
     l.assign(lz);
     r.assign(lz);
   1z = -1;
  void merge(const seq_node& 1, const seq_node& r) {
   sz = 1.sz + r.sz;
   sum = 1.sum + r.sum;
  void assign(int val) { // range \ a[i] <- val
   sum = sz * val;
   1z = val:
  void flip(int val) { // range \ a[i] <- !a[i]
   sum = sz - sum;
   if (1z == -1) 1z = 2;
   else if (1z == 0) 1z = 1;
   else if (1z == 1) 1z = 0;
   else lz = -1:
  int64_t get_sum() const { return sum; } // sum a[l, r]
template<typename T = int64 t> struct seg node {
 T val, lz_add, lz_set;
 int sz;
 bool to_set;
  seq_node(T n = 0) : val(n), lz_add(0), lz_set(0), sz(1),
      to set(0) {}
  void push(seq_node& 1, seq_node& r) {
   if (to_set) {
     l.assign(lz set);
     r.assign(lz_set);
     1z set = 0;
```

```
to set = false;
   if (lz add != 0) {
     1.add(lz add);
     r.add(lz_add);
     1z add = 0;
 void merge(const seg_node& 1, const seg_node& r) {
   sz = 1.sz + r.sz;
   val = 1.val + r.val;
 void add(T v) { // update \ range \ a[i] \leftarrow a[i] + v
   val += v * sz;
   lz_add += v;
 void assign(T v) { //update\ range\ a[i] <-v
   val = v * sz;
   1z add = 0:
   lz set = v;
   to_set = true;
 T get sum() const { return val; } // sum a/l, r)
// update range a[i] \leftarrow a[i] + b * (i - s) + c
// assuming b and c are non zero, be careful
// get sum a[l, r]
template<typename T = int64_t> struct seq_node {
T sum, lzB, lzC;
 int sz, idx;
 seq_node(int id = 0, T v = 0, int s = 0, T b = 0, T c = 0):
   sum(v), lzB(b), lzC(c - s * b), idx(id), sz(1) {}
 void push(seg_node& 1, seg_node& r) {
   l.add(lzB, lzC);
   r.add(lzB, lzC);
   lzB = lzC = 0;
 void merge(const seg node& 1, const seg node& r) {
   idx = min(l.idx, r.idx);
   sz = 1.sz + r.sz;
   sum = 1.sum + r.sum;
 T sum idx(T n) const { return n * (n + 1) / 2; }
 void add(T b, T c) {
   sum += b * (sum_idx(idx + sz) - sum_idx(idx)) + sz * c;
   1zB += b:
   1zC += c;
 T get_sum() const { return sum; }
// update range a[i] \leftarrow b * a[i] + c
// get sum a[l. r]
struct seg node {
 int sz: i64 sum, lzB, lzC;
 seq_node() : sz(1), sum(0), lzB(1), lzC(0) {}
 seg node(i64 v) : sz(1), sum(v), lzB(1), lzC(0) {}
 void push(seq_node& 1, seq_node& r) {
   1.add(lzB, lzC);
   r.add(1zB, 1zC):
   1zB = 1, 1zC = 0;
 void merge(const seg_node& 1, const seg_node& r) {
   sz = 1.sz + r.sz;
   sum = 1.sum + r.sum;
 void add(i64 b, i64 c) {
   sum = (b * sum + c * sz);
```

```
lzC = (lzC * b + c);
 i64 get_sum() const { return sum; }
// update range a[i] \leftarrow min(a[i], b);
// update range a[i] \leftarrow max(a[i], b):
// get val a[i]
struct seg node {
 int mn, mx;
 int 1z0, 1z1;
  seq_node() : mn(INT_MAX), mx(INT_MIN), lz0(INT_MAX), lz1(
       INT MIN) {}
  void push(seg_node& 1, seg_node& r) {
   1.minimize(1z0);
    1.maximize(lz1);
    r.minimize(lz0);
    r.maximize(lz1);
   lz0 = INT_MAX;
   lz1 = INT_MIN;
 void merge(const seg node& 1, const seg node& r) {
    mn = min(1.mn, r.mn);
    mx = max(1.mx, r.mx);
 void minimize(int val) {
    mn = lz0 = min(lz0, val);
    mx = 1z1 = min(1z0, 1z1);
 void maximize(int val) {
    mx = lz1 = max(lz1, val);
    mn = 1z0 = max(1z0, 1z1);
 pair<int, int> get() const { return {mx, mn}; }
segtree-2d.h
Description: 2D Segment Tree.
Time: \mathcal{O}(N \log^2 N) of memory, \mathcal{O}(\log^2 N) per query
"sparse_seg_tree.h"
                                                       09098e, 25 lines
template<class T> struct Node {
 node_t<T> seq; Node* c[2];
 Node() { c[0] = c[1] = nullptr; }
  void upd(int x, int y, T v, int L = 0, int R = SZ-1) { // add
    if (L == x \&\& R == x) \{ seg.upd(y,v); return; \}
    int M = (L+R) >> 1;
    if (x <= M) {
      if (!c[0]) c[0] = new Node();
      c[0] -> upd(x, y, v, L, M);
      if (!c[1]) c[1] = new Node();
      c[1] -> upd(x, y, v, M+1, R);
    seg.upd(y,v); // only for addition
            // seq.upd(y,c[0]?&c[0]->seq:nullptr,c[1]?&c[1]->
                 seq:nullptr);
 T query(int x1, int x2, int y1, int y2, int L = 0, int R = SZ
       -1) { // query sum of rectangle
    if (x1 <= L && R <= x2) return seg.query(y1,y2);
    if (x2 < L || R < x1) return 0;
    int M = (L+R) >> 1; T res = 0;
    if (c[0]) res += c[0]->query(x1, x2, y1, y2, L, M);
    if (c[1]) res += c[1]->query(x1, x2, y1, y2, M+1, R);
    return res;
```

lzB = (lzB * b);

rmq.h

Description: Range Minimum/Maximum Queries on an array. Returns min(V[a], V[a + 1], ... V[b]) in constant time. Returns a pair that holds the answer, first element is the value and the second is the index.

Usage: rmq_t<pair<int, int>> rmq(values); // values is a vector of pairs {val(i), index(i)} rmq.query(inclusive, exclusive); rmg_t<pair<int, int>, greater<pair<int, int>>> rmg(values) //max query Time: $\mathcal{O}(|V|\log|V|+Q)$

template<typename T, typename Cmp=less<T>> struct rmq_t : private Cmp { int N = 0; vector<vector<T>> table; const T& min(const T& a, const T& b) const { return Cmp:: operator()(a, b) ? a : b; } rmq_t() {} rmg t(const vector<T>& values) : N(int(values.size())), table (lq(N) + 1) { table[0] = values; for (int a = 1; a < int(table.size()); ++a) {</pre>

table[a][b] = min(table[a-1][b], table[a-1][b + (1 << (

table[a].resize(N - $(1 \ll a) + 1)$;

for (int b = 0; b + (1 << a) <= N; ++b)

fenwick-tree.h

};

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

return min(table[lg][a], table[lg][b - (1 << lg)]);</pre>

a-1))1):

T query(int a, int b) const {

int $lq = __lq(b - a);$

```
Time: Both operations are \mathcal{O}(\log N).
                                                       f79f29, 28 lines
template<typename T> struct FT { // 8b7639
 vector<T> s;
  FT(int n) : s(n) {}
  FT(const vector<T>& A) : s(A) {
   const int N = int(s.size());
    for (int a = 0; a < N; ++a) {
     if ((a | (a + 1)) < N) s[a | (a + 1)] += s[a];
  void update(int pos, T dif) { // a[pos] += dif
    for (; pos < (int)s.size(); pos |= pos + 1) s[pos] += dif;
  T query(int pos) { // sum of values in [0, pos)
   T res = 0:
    for (; pos > 0; pos &= pos - 1) res += s[pos-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 pos = 0;
    for (int pw = 1 << 25; pw; pw >>= 1) {
     if (pos + pw <= (int)s.size() && s[pos + pw-1] < sum)
        pos += pw, sum -= s[pos-1];
    return pos;
};
```

fenwick-tree-2d.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: $\mathcal{O}(\log^2 N)$. (Use persistent segment trees for $\mathcal{O}(\log N)$.) "fenwick-tree.h"

```
aebbdc, 25 lines
template<typename T> struct FT2 {
 vector<vector<int>> ys; vector<FT<T>> ft;
 FT2(int limx) : ys(limx) {}
 void fakeUpdate(int x, int y) {
    for (; x < (int)ys.size(); x |= x + 1) ys[x].push_back(y);
 void init() {
   for(auto &v : ys){
     sort(v.begin(), v.end());
     v.resize(unique(v.begin(), v.end()) - v.begin());
     ft.emplace_back(v.size());
 int ind(int x, int y) {
   return (int) (lower_bound(ys[x].begin(), ys[x].end(), y) -
        vs[x].begin()); }
 void update(int x, int y, T dif) {
    for (; x < ys.size(); x |= x + 1)
      ft[x].update(ind(x, y), dif);
 T query(int x, int y) {
   T sum = 0;
    for (; x; x \&= x - 1) sum += ft[x-1].query(ind(x-1, y));
    return sum;
};
```

mo.h

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

```
5ef29d, 49 lines
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
vector<int> mo(vector<pair<int, int>> Q) { // d9247c
 int L = 0, R = 0, blk = 350; // \sim N/sqrt(Q)
 vector<int> s(int(Q.size())), res = s;
#define K(x) pair<int, int>(x.first/blk, x.second ^ -(x.first/
    blk & 1))
 iota(s.begin(), s.end(), 0);
 sort(s.begin(), s.end(), [&](int s, int t){ return K(Q[s]) < }
      K(Q[t]); });
 for (int qi : s) {
   auto q = Q[qi];
   while (L > q.first) add(--L, 0);
   while (R < q.second) add (R++, 1);
   while (L < q.first) del(L++, 0);
   while (R > q.second) del(--R, 1);
    res[gi] = calc();
 return res;
vector<int> moTree(vector<array<int, 2>> 0, vector<vector<int
    >>& ed, int root=0) { // b\bar{b}f891
 int N = int(ed.size()), pos[2] = {}, blk = 350; // \sim N/sqrt(Q)
 vector < int > s(int(Q.size())), res = s, I(N), L(N), R(N), in(N)
      ), par(N);
 add(0, 0), in[0] = 1;
```

```
auto dfs = [\&] (int x, int p, int dep, auto& f) -> void {
   par[x] = p;
   L[x] = N;
   if (dep) I[x] = N++;
   for (int y : ed[x]) if (y != p) f(y, x, !dep, f);
   if (!dep) I[x] = N++;
   R[x] = N;
 };
 dfs(root, -1, 0, dfs);
#define K(x) pii(I[x[0]] / blk, I[x[1]] ^ -(I[x[0]] / blk & 1))
 iota(s.begin(), s.end(), 0);
 sort(s.begin(), s.end(), [&](int s, int t){ return K(Q[s]) <</pre>
      K(Q[t]); });
 for (int qi : s) for (int end = 0; end < 2; ++end) {
   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;
```

line-container.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: $\mathcal{O}(\log N)$

```
8b2ace, 29 lines
struct Line {
 mutable lint k, m, p;
  bool operator<(const Line& o) const { return k < o.k; }</pre>
  bool operator<(lint x) const { return p < x; }</pre>
struct LineContainer : multiset<Line, less<>>> {
  // (for doubles, use inf = 1/.0, div(a,b) = a/b)
  static const lint inf = LLONG_MAX;
  lint div(lint a, lint b) { // floored division
    return a / b - ((a ^ b) < 0 && a % b); }
  bool isect(iterator x, iterator v) {
    if (y == end()) \{ x \rightarrow p = inf; return false; \}
    if (x->k == y->k) x->p = x->m > y->m ? inf : -inf;
    else x - > p = div(v - > m - x - > m, x - > k - v - > k);
    return x->p >= y->p;
  void add(lint k, lint m) {
    auto z = insert(\{k, m, 0\}), y = z++, x = y;
    while (isect(y, z)) z = erase(z);
    if (x != begin() \&\& isect(--x, y)) isect(x, y = erase(y));
    while ((y = x) != begin() \&\& (--x)->p >= y->p)
      isect(x, erase(y));
 lint query(lint x) {
    assert(!empty());
    auto 1 = *lower_bound(x);
    return 1.k * x + 1.m;
};
```

matrix.h

Description: Basic operations on square matrices. Usage: Matrix<int> A(N, vector<int>(N));

```
template <typename T> struct Matrix : vector<vector<T>>> {
 using vector<vector<T>>::vector;
 using vector<vector<T>>::size;
```

```
int h() const { return int(size()); }
  int w() const { return int((*this)[0].size()); }
  Matrix operator* (const Matrix& r) const {
    assert(w() == r.h());
   Matrix res(h(), vector<T>(r.w()));
   for (int i = 0; i < h(); ++i) {
     for (int j = 0; j < r.w(); ++j)
        for (int k = 0; k < w(); ++k) {
         res[i][j] += (*this)[i][k] * r[k][j];
    return res:
  friend vector<T> operator*(const Matrix<T>& A, const vector<T
    int N = int(A.size()), M = int(A[0].size());
   vector<T> y(N);
    for (int i = 0; i < N; ++i)
     for (int j = 0; j < M; ++j) y[i] += A[i][j] * b[j];
  Matrix& operator *= (const Matrix& r) { return *this = *this *
  Matrix pow(int n) const {
   assert(h() == w());
   Matrix x = *this, r(h(), vector<T>(w()));
    for (int i = 0; i < h(); ++i) r[i][i] = T(1);
   while (n) { if (n & 1) r *= x; x *= x; n >>= 1; }
};
```

range-color.h

Description: RangeColor structure, supports point queries and range updates, if C isn't int32_t change freq to map

```
Time: \mathcal{O}\left(\lg(L)*Q\right)
```

```
template<class T = int64_t, class C = int32_t> struct
    RangeColor{
  struct Node {
   T left, right; C color;
   bool operator < (const Node &n) const{ return right < n.
        right; }
  };
  C minInf;
  set < Node > st;
  vector<T> freq;
  RangeColor(T first, T last, C maxColor, C iniColor = C(0)):
      minInf(first - T(1)), freq(maxColor + 1) {
    freq[iniColor] = last - first + T(1);
    st.insert({first, last, iniColor});
  C query(T i) { //get\ color\ in\ position\ i}
   auto p = st.upper_bound({T(0), i - T(1), minInf});
    return p->color;
  void upd(T a, T b, C newColor) { //set newColor in [a, b]
   auto p = st.upper_bound({T(0), a - T(1), minInf});
   assert(p != st.end());
   T left = p->left, right = p->right;
   C old = p->color;
    freq[old] -= (right - left + T(1));
    p = st.erase(p);
    if (left < a) {</pre>
     freq[old] += (a - left);
     st.insert({left, a - T(1), old});
    if (b < right) {</pre>
```

```
freq[old] += (right - b);
    st.insert({b + T(1), right, old});
}
while ((p != st.end()) && (p->left <= b)){
    left = p->left, right = p->right;
    old = p->color;
    freq[old] -= (right - left + T(1));
    if (b < right){
        freq[old] += (right - b);
        st.erase(p);
        st.insert({b + T(1), right, old});
        break;
    } else p = st.erase(p);
}
freq[newColor] += (b - a + T(1));
st.insert({a, b, newColor});
}
T countColor(C x){ return freq[x]; }
};</pre>
```

implicit-treap.h

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

```
Time: \mathcal{O}(\log N) 2d0d97, 96 lines
```

```
mt19937 rng(chrono::steady_clock::now().time_since_epoch().
    count());
struct node {
 int v, p, sz;
 node *1, *r;
 bool rev;
 node(int k) : v(k), p(rng()), l(nullptr), rev(0), r(nullptr),
       sz(0) {}
int sz(node *t) {
 if (t == nullptr) return 0;
 return t->sz;
void push(node *t) {
 if (t == nullptr) return;
 if (t->rev) {
    swap(t->1, t->r);
   if (t->1 != nullptr) t->1->rev ^= t->rev;
   if (t->r != nullptr) t->r->rev ^= t->rev;
   t->rev = 0;
void updsz(node *t) {
 if (t == nullptr) return;
 push(t); push(t->1); push(t->r);
 t->sz = sz(t->1) + sz(t->r) + 1;
void split(node *t, node *&l, node *&r, int k) { //k on left
 push(t);
 if (t == nullptr) l = r = nullptr;
  else if (k \le sz(t->1)) {
    split (t->1, 1, t->1, k);
    r = t;
 else {
    split(t->r, t->r, r, k-1-sz(t->1));
   1 = t;
 updsz(t);
void merge(node *&t, node *1, node *r) {
 push(1); push(r);
 if (1 == nullptr) t = r;
 else if (r == nullptr) t = 1;
```

```
else if (1->p <= r->p) {
   merge(1->r, 1->r, r);
   t = 1;
 else (
   merge(r->1, 1, r->1);
   t = r;
 updsz(t);
void add(node *&t, node *c, int k) {
 push(t);
 if (t == nullptr) t = c;
 else if (c->p>=t->p) {
   split(t, c->1, c->r, k);
   t = c;
 else if (sz(t->1) >= k) add(t->1, c, k);
 else add(t->r, c, k-1-sz(t->1));
 updsz(t);
void del(node *&t, int k) {
 push(t);
 if (t == nullptr) return;
 if (sz(t->1) == k) merge(t, t->1, t->r);
 else if (sz(t->1) > k) del(t->1, k);
 else del(t->r, k);
 updsz(t);
void print(node *t) {
 if (r == nullptr) return;
 print(t->1);
 cout << t->v << ' ';
 print(t->r);
int main() {
 node *treap = nullptr;
 while(1) {
   int a:
   cin >> a;
   if (a == 1)
     int c, d;
     cin >> c >> d;
     node *r = new node(d);
     add(treap, r, c);
   } else if (a == 2) {
     int d;
     cin >> d;
     del(treap, d);
   print(treap);
```

Numerical (4)

polynomial.h

```
84593c, 17 lines
```

```
struct Poly {
  vector<double> a;
  double operator() (double x) const {
    double val = 0;
    for(int i = a.size(); i--;) (val *= x) += a[i];
    return val;
}
void diff() {
  for(int i = 1; i < a.size(); ++i) a[i-1] = i*a[i];</pre>
```

```
a.pop_back();
  void divroot(double x0) {
    double b = a.back(), c; a.back() = 0;
    for (int i = a.size()-1; i--; ) c = a[i], a[i] = a[i+1] *x0+b,
         b=c;
    a.pop_back();
};
```

poly-roots.h

Description: Finds the real roots to a polynomial.

Usage: poly_roots($\{\{2,-3,1\}\},-1e9,1e9$) // solve $x^2-3x+2=0$ Time: $\mathcal{O}\left(n^2\log(1/\epsilon)\right)$

"Polynomial.h" vector<double> poly_roots(Poly p, double xmin, double xmax) { if ((p.a).size() == 2) { return {-p.a[0]/p.a[1]}; } vector<double> ret; Poly der = p; der.diff(); auto dr = poly_roots(der, xmin, xmax); dr.push_back(xmin-1); dr.push_back(xmax+1); sort(dr.begin(), dr.end()); for(int i = 0; i < dr.size()-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.push_back((1 + h) / 2); return ret;

poly-interpolate.h

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

```
97a266, 13 lines
typedef vector<double> vd;
vd interpolate(vd x, vd y, int n) {
  vd 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;
```

lagrange.h

Description: Lagrange interpolation over a finite field and some combo stuff Time: $\mathcal{O}(N)$

```
"../number-theory/modular-arithmetic.h", "../number-theory/preparator.h" 84e6f3, 14 lines
template<typename T> struct interpolator_t {
 vector<T> pref, suff;
  interpolator_t(int N): pref(N), suff(N) {}
 T interpolate(const vector<T>& y, T x) {
   int N = int(y.size()); T res = 0;
```

```
pref[0] = suff[N - 1] = 1;
    for (int i = 0; i+1 < N; ++i) pref[i+1] = pref[i] * (x-i);
    for (int i = N-1; i > 0; --i) suff[i-1] = suff[i] * (x-i);
    for (int i = 0, sgn = (N \& 1 ? +1 : -1); i < N; ++i, <math>sgn *=
     res += y[i] * sgn * pref[i] * suff[i] * invFac[i] *
          invFac[N-1-i];
   return res;
};
```

berlekamp-massev.h

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

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

```
"ModularArithmetic.h"
                                                     4c4a48, 19 lines
template <typename num>
vector<num> BerlekampMassey(const vector<num>& s) {
 int n = int(s.size()), L = 0, m = 0;
 vector<num> C(n), B(n), T;
 C[0] = B[0] = 1;
 num b = 1;
 for(int i = 0; i < n; i++) { ++m;
   num d = s[i];
    for (int j = 1; j \le L; j++) d += C[j] * s[i - j];
   if (d == 0) continue;
   T = C; num coef = d / b;
    for (int j = m; j < n; j++) C[j] -= coef * B[j - m];
   if (2 * L > i) continue;
   L = i + 1 - L; B = T; b = d; m = 0;
 C.resize(L + 1); C.erase(C.begin());
 for (auto& x : C) x = -x;
 return C;
```

linear-recurrence.h

Description: Generates the k'th term of an n-order linear recurrence $S[i] = \sum_{j} S[i-j-1]tr[j]$, given $S[0 \dots n-1]$ and $tr[0 \dots n-1]$. Faster than matrix multiplication. Useful together with Berlekamp-Massey.

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

```
"ModularArithmetic.h"
                                                     Obaa7b, 22 lines
template <typename num>
num linearRec(const vector<num>& S, const vector<num>& tr, lint
     k) {
 int n = int(tr.size());
 assert(S.size() >= tr.size());
 auto combine = [&](vector<num> a, vector<num> b) {
   vector<num> res(n * 2 + 1);
   for (int i = 0; i \le n; i++) for (int j = 0; j \le n; j++)
        res[i + j] += a[i] \star b[j];
    for (int i = 2 * n; i > n; --i) for (int j = 0; j < n; j++)
     res[i - 1 - j] += res[i] * tr[j];
    res.resize(n + 1);
   return res;
 vector<num> 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);
 num res = 0;
```

```
for (int i = 0; i < n; i++) res += pol[i + 1] * S[i];
return res;
```

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. 7bb98e, 7 lines

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

integrate-adaptive.h

Description: Fast integration using an adaptive Simpson's rule. Usage: double sphereVolume = quad(-1, 1, [] (double x) { return quad(-1, 1, [&](double v)

```
return quad(-1, 1, [\&](double z) {
return x*x + y*y + z*z < 1; {);});});
                                                       92dd79, 15 lines
typedef double d;
```

```
#define S(a,b) (f(a) + 4*f((a+b) / 2) + f(b)) * (b-a) / 6
template <class F>
d rec(F& f, d a, d b, d eps, d S) {
  dc = (a + b) / 2;
  d S1 = S(a, c), S2 = S(c, b), T = S1 + S2;
  if (abs(T - S) \le 15 * eps | | b - a < 1e-10)
    return T + (T - S) / 15;
  return rec(f, a, c, eps / 2, S1) + rec(f, c, b, eps / 2, S2);
template<class F>
d \text{ quad}(d \text{ a, } d \text{ b, } F \text{ f, } d \text{ eps} = 1e-8)  {
  return rec(f, a, b, eps, S(a, b));
```

gaussian-elimination.h

Time: $\mathcal{O}(\min(N, M)NM)$

"../data-structures/matrix.h" b33fca, 75 lines

```
template<typename T> struct gaussian_elimination {
 int N, M;
 Matrix<T> A, E;
  vector<int> pivot;
 int rank, nullity, sgn;
  gaussian_elimination(const Matrix<T>& A_) : A(A_) {
    N = A.size(), M = A[0].size();
    E = Matrix<T>(N, vector<T>(N));
    for (int i = 0; i < N; ++i) E[i][i] = 1;
    rank = 0, nullity = M, sgn = 0;
    pivot.assign(M, -1);
    for (int col = 0, row = 0; col < M && row < N; ++col) {
      int sel = -1;
      for (int i = row; i < N; ++i) {</pre>
       if (A[i][col] != 0) {
          sel = i;
          break;
      if (sel == -1) continue;
     if (sel != row) {
        sgn += 1;
        swap(A[sel], A[row]);
        swap(E[sel], E[row]);
```

```
for (int i = 0; i < N; ++i) {
       if (i == row) continue;
       T c = A[i][col] / A[row][col];
       for (int j = col; j < M; ++j)
         A[i][j] = c * A[row][j];
       for (int j = 0; j < N; ++j)
         E[i][j] = c * E[row][j];
     pivot[col] = row++;
     ++rank, --nullity;
 pair<bool, vector<T>> solve(vector<T> b, bool reduced = false
     ) const {
   if (reduced == false) b = E * b;
   vector<T> x(M);
   for (int j = 0; j < M; ++j) {
     if (pivot[j] == -1) continue;
     x[j] = b[pivot[j]] / A[pivot[j]][j];
     b[pivot[j]] = 0;
    for (int i = 0; i < N; ++i)
     if (b[i] != 0) return {false, x};
   return {true, x};
  vector<vector<T>> kernel_basis() const {
   vector<vector<T>> basis;
   vector<T> e(M);
   for (int j = 0; j < M; ++j) {
     if (pivot[j] != -1) continue;
     auto y = solve(A * e, true).second;
     e[j] = 0, y[j] = -1;
     basis.push_back(y);
   return basis;
 Matrix<T> inverse() const {
   assert (N == M); assert (rank == N);
   Matrix<T> res(N, vector<T>(N));
   vector<T> e(N);
    for (int i = 0; i < N; ++i) {
     e[i] = 1;
     auto x = solve(e).second;
     for (int j = 0; j < N; ++j)
       res[j][i] = x[j];
     e[i] = 0;
   return res;
};
```

linear-solver-z2.h

Description: Solves Ax = b over \mathbb{F}_2 . If there are multiple solutions, one is returned arbitrarily. Returns true, or false if no solutions. Last column of a is b. c is the rank.

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

7a24e1, 26 lines

```
typedef bitset<2010> bs;
bool gauss(vector<bs> a, bs& ans, int n) {
  int m = int(a.size()), c = 0;
  bs pos; pos.set();
  for (int j = n-1, i; j >= 0; --j) {
    for (i = c; i < m; ++i)
      if (a[i][j]) break;
    if (i == m) continue;
    swap(a[c], a[i]);
    i = c++; pos[j] = 0;
    for (int k = 0; k < m; ++k)</pre>
```

```
if (a[k][j] && k != i)
    a[k] ^= a[i];
}
ans = pos;
for(int i = 0; i < m; ++i) {
    int ac = 0;
    for (int j = 0; j < n; ++j) {
        if (!a[i][j]) continue;
        if (!pos[j]) pos[j] = 1, ans[j] = ac^a[i][n];
        ac ^= ans[j];
}
if (ac != a[i][n]) return false;
}
return true;
}</pre>
```

simplex.h

Description: Solves a general linear maximization problem: maximize $c^T x$ subject to Ax < b, x > 0.

Time: $\mathcal{O}(N\overline{M}*\#pivots)$, where a pivot may be e.g. an edge relaxation. $\mathcal{O}(2^n)$ in the general case. WARNING- segfaults on empty (size 0) max cx st Ax <= b, x >= 0 do 2 phases; 1st check feasibility; 2nd check boundedness and ans

vector<double> simplex(vector<vector<double>> A, vector<double> b, vector<double> c) { int n = A.size(), m = A[0].size() + 1, r = n, s = m-1; vector<vector<double>> D = vector<vector<double>> (n+2, vector <double>(m+1));vector<int> ix = vector<int>(n + m); for (int i = 0; i < n + m; ++i) ix[i] = i; for (int i = 0; i < n; ++i) { for (int j = 0; j < m-1; ++j) D[i][j] = -A[i][j]; D[i][m - 1] = 1;D[i][m] = b[i];if (D[r][m] > D[i][m]) r = i;for (int j = 0; j < m-1; ++j) D[n][j] = c[j]; D[n + 1][m - 1] = -1; int z = 0;for (double d;;) { if (r < n) { swap(ix[s], ix[r + m]);D[r][s] = 1.0/D[r][s];for (int j = 0; $j \le m$; ++j) if (j != s) D[r][j] *= -D[r]][s]; for (int i = 0; $i \le n+1$; ++i) if (i != r) { for (int j = 0; $j \le m$; ++j) if (j != s) D[i][j] += D[r]][j] * D[i][s]; $D[i][s] \star= D[r][s];$ r = -1; s = -1; for (int j = 0; j < m; ++j) if (s < 0 || ix[s] > ix[j]) if (D[n+1][j] > eps || D[n+1][j] > -eps && D[n][j] > eps) s = i;if (s < 0) break; for (int i = 0; i < n; ++i) if (D[i][s] < -eps) { if $(r < 0 \mid | (d = D[r][m]/D[r][s]-D[i][m]/D[i][s]) < -eps$ | | d < eps && ix[r+m] > ix[i+m]) r = i;if (r < 0) return vector<double>(); // unbounded if (D[n+1][m] < -eps) return vector<double>(); // infeasible vector<double> x (m-1); for (int i = m; i < n+m; ++i) if (ix[i] < m-1) x[ix[i]] = D[i -m][m]; double result = D[n][m]; return x; // ans: D[n][m]

| tridiagonal

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

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

This is useful for solving problems on the type

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

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

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

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.

diag[1] == 0 is needed. **Time:** $\mathcal{O}(N)$ d0855f, 26 lines

```
typedef double T;
vector<T> tridiagonal(vector<T> diag, const vector<T> &super,
    const vector<T> &sub, vector<T> b) {
  int n = b.size(); vector<int> tr(n);
  for (int i = 0; i < n-1; ++i) {
    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;
```

4.1 Fourier transforms

fast-fourier-transform.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 $c[x] = \sum_i a[i]b[x-i]$. 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 $(\sum_i a_i^2 + \sum_i b_i^2) \log_2 N < 9 \cdot 10^{14}$ (in practice 10^{16} ; higher for random inputs). Otherwise, use NTT/FFTMod.

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

```
inline int nxt_pow2(int s) { return 1 << (s > 1 ? 32 -
    __builtin_clz(s-1) : 0); }
template <typename dbl> struct cplx {
    dbl x, y;
    cplx(dbl x_ = 0, dbl y_ = 0) : x(x_), y(y_) { }
    friend cplx operator+(cplx a, cplx b) { return cplx(a.x + b.x , a.y + b.y); }
    friend cplx operator-(cplx a, cplx b) { return cplx(a.x - b.x , a.y - b.y); }
```

fast-subset-transform sum-of-powers

```
friend cplx operator*(cplx a, cplx b) { return cplx(a.x * b.x
       -a.y * b.y, a.x * b.y + a.y * b.x); }
  friend cplx conj(cplx a) { return cplx(a.x, -a.y); }
  friend cplx inv(cplx a) { dbl n = (a.x*a.x+a.y*a.y); return
      cplx(a.x/n,-a.y/n); }
template <typename T>
struct root_of_unity {};
template <typename dbl> struct root_of_unity<cplx<dbl>>> {
  static cplx<dbl> f(int k) {
   static const dbl PI = acos(-1);
   dbl a = 2*PI/k;
   return cplx<dbl>(cos(a), sin(a));
};
using M0 = modnum<998244353U>; // g = 3
using M1 = modnum<897581057U>; // q = 3
using M2 = modnum<880803841U>; // g = 26
using M3 = modnum<985661441U>; // q = 3
using M4 = modnum<943718401U>; // g = 7
using M5 = modnum<935329793U>; // g = 3
using M6 = modnum<918552577U>; // q = 5
constexpr unsigned primitive_root(unsigned M) {
 if (M == 880803841U) return 26U;
 else if (M == 943718401U) return 7U;
 else if (M == 918552577U) return 5U;
 else return 3U;
template<unsigned MOD> struct root_of_unity<modnum<MOD>> {
 static constexpr modnum<MOD> g0 = primitive root(MOD);
  static modnum<MOD> f(int K) {
   assert ((MOD-1)\%K == 0);
   return g0.pow((MOD-1)/K);
};
template<typename T> struct FFT {
 vector<T> rt; vector<int> rev;
 FFT(): rt(2, T(1)) {}
  void init(int N) {
   N = nxt pow2(N);
   if (N > int(rt.size())) {
     rev.resize(N); rt.reserve(N);
     for (int a = 0; a < N; ++a) {
       rev[a] = (rev[a/2] | ((a&1)*N)) >> 1;
      for (int k = int(rt.size()); k < N; k *= 2) {
       rt.resize(2*k);
       T z = root_of_unity < T > :: f(2*k);
       for (int a = k/2; a < k; ++a) {
         rt[2*a] = rt[a];
         rt[2*a+1] = rt[a] * z;
  void fft (vector<T>& xs, bool inverse) const {
   int N = int(xs.size());
   int s = __builtin_ctz(int(rev.size())/N);
   if (inverse) reverse(xs.begin() + 1, xs.end());
    for (int a = 0; a < N; ++a) {
     if (a < (rev[a] >> s))
       swap(xs[a], xs[rev[a] >> s]);
    for (int k = 1; k < N; k *= 2) {
     for (int a = 0; a < N; a += 2 * k) {
```

```
int u = a, v = u + k;
        for (int b = 0; b < k; ++b, ++u, ++v) {
          T z = rt[b + k] * xs[v];
          xs[v] = xs[u] - z;
          xs[u] = xs[u] + z;
    if (inverse) {
     for (int a = 0; a < N; ++a)
        xs[a] = xs[a] * inv(T(N));
 vector<T> convolve(vector<T> as, vector<T> bs) {
    int N = int(as.size()), M = int(bs.size());
    int K = N + M - 1, S = nxt_pow2(K); init(S);
    if (min(N, M) \le 64) {
     vector<T> res(K);
      for (int u = 0; u < N; ++u)
       for (int v = 0; v < M; ++v)
          res[u + v] = res[u + v] + as[u] * bs[v];
      return res;
    } else {
      as.resize(S), bs.resize(S);
      fft(as, false); fft(bs, false);
      for (int i = 0; i < S; ++i) as[i] = as[i] * bs[i];
      fft(as, true); as.resize(K);
      return as;
};
FFT<M0> FFT0; FFT<M1> FFT1;
FFT<M2> FFT2; FFT<M3> FFT3;
FFT<M4> FFT4; FFT<M5> FFT5; FFT<M6> FFT6;
// M0 M1 = 896005221510021121 (> 4.48 * 10^{17}, > 2^{58})
// M0 M1 M2 = 789204840662082423367925761 (> 7.892 * 10^26, > 2
// M0 M3 M4 M5 M6 =
     797766583174034668024539679147517452591562753 (> 7.977 *
//T = \{unsigned, unsigned long, modnum \in M \}
template<class T, unsigned M0, unsigned M1, unsigned M2>
T garner (modnum<M0> a0, modnum<M1> a1, modnum<M2> a2) {
 static const modnum<M1> INV M0 M1 = modnum<M1> (M0).inv();
  static const modnum<M2> INV_M0M1_M2 = (modnum < M2 > (M0) * M1).
  const modnum<M1> b1 = INV M0 M1 * (a1 - a0.x);
  const modnum<M2> b2 = INV_M0M1_M2 * (a2 - (modnum<M2> (b1.x) *
       M0 + a0.x));
  return (T(b2.x) * M1 + b1.x) * M0 + a0.x;
template < class T, unsigned M0, unsigned M1, unsigned M2,
    unsigned M3, unsigned M4>
T garner (modnum<M0> a0, modnum<M1> a1, modnum<M2> a2, modnum<M3
    > a3, modnum<M4> a4) {
  static const modnum<M1> INV M0 M1 = modnum<M1>(M0).inv();
  static const modnum<M2> INV M0M1 M2 = (modnum < M2 > (M0) * M1).
      inv():
  static const modnum<M3> INV_M0M1M2_M3 = (modnum<M3>(M0) * M1
      * M2).inv():
  static const modnum<M4> INV M0M1M2M3 M4 = (modnum<M4>(M0) *
      M1 * M2 * M3).inv();
  const modnum<M1> b1 = INV_M0_M1 \star (a1 - a0.x);
  const modnum<M2> b2 = INV_M0M1_M2 * (a2 - (modnum<M2> (b1.x) *
        M0 + a0.x));
```

```
const modnum<M3> b3 = INV M0M1M2 M3 * (a3 - ((modnum<M3>(b2.x
     ) * M1 + b1.x) * M0 + a0.x);
 const modnum<M4> b4 = INV M0M1M2M3 M4 * (a4 - (((modnum<M4>(
      b3.x) * M2 + b2.x) * M1 + b1.x) * M0 + a0.x));
 return (((T(b4.x) * M3 + b3.x) * M2 + b2.x) * M1 + b1.x) * M0
       + a0.x:
// results must be in [-448002610255888384, 448002611254132736]
vector<long long> convolve(const vector<long long>& as, const
    vector<long long>& bs) {
 static constexpr unsigned M0 = M0::M, M1 = M1::M;
 static const modnum<M1> INV_M0_M1 = modnum<M1>(M0).inv();
 if (as.empty() || bs.empty()) return {};
 const int len_as = int(as.size()), len_bs = int(bs.size());
 vector<modnum<M0>> as0(len_as), bs0(len_bs);
 for (int i = 0; i < len_as; ++i) as0[i] = as[i];</pre>
 for (int i = 0; i < len_bs; ++i) bs0[i] = bs[i];
 const vector<modnum<M0>> cs0 = FFT0.convolve(as0, bs0);
 vector<modnum<M1>> as1(len_as), bs1(len_bs);
 for (int i = 0; i < len_as; ++i) as1[i] = as[i];
 for (int i = 0; i < len_bs; ++i) bs1[i] = bs[i];</pre>
 const vector<modnum<M1>> cs1 = FFT1.convolve(as1, bs1);
 vector<long long> cs(len_as + len_bs - 1);
 for (int i = 0; i < len_as + len_bs - 1; ++i) {
    const modnum<M1> d1 = INV_M0_M1 * (cs1[i] - cs0[i].x);
   cs[i] = (d1.x > M1 - d1.x)
     ? (-1ULL - (static cast<unsigned long long>(M1 - 1U - d1.
          x) * M0 + (M0 - 1U - cs0[i].x))
      : (static_cast<unsigned long long>(d1.x) * M0 + cs0[i].x)
 return cs;
```

fast-subset-transform.h

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

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

5b9574, 16 lines

```
void FST(vector<int> &a, bool inv) {
  for (int n = a.size(), step = 1; step < n; step *= 2) {
    for (int i = 0; i < n; i += 2 * step) for(int j = i; j < i +
        step; ++j) {
    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(auto &x : a) x /= a.size(); // XOR only
}
vector<int> conv(vector<int> a, vector<int> b) {
  FST(a, 0); FST(b, 0);
  for(int i = 0; i < a.size(); ++i) a[i] *= b[i];
  FST(a, 1); return a;</pre>
```

sum-of-powers.h

Description: Computes monomials and sum of powers product certain polynomials. Check "General purpose numbers" section for more info. (Mono-

```
mials) pw(x) = x^d for a fixed d. \sum_{x=0}^{\infty} r^x f(x). (degree of f \le d). \sum_{x=0}^{N-1} r^x f(x).
```

(degree of $f \leq d$).

".../number-theory/modular-arithmetic.h", "/lagrange.h" 0b1060, 45 lines

vector<num> get monomials(int N, long long d) {

```
vector<int> pfac(N);
  for (int i = 2; i < N; ++i) pfac[i] = i;
  for (int p = 2; p < N; ++p) if (pfac[p] == p) {
   for (int m = 2*p; m < N; m += p) if (pfac[m] > p) pfac[m] =
  vector<num> pw(N);
  for (int i = 0; i < N; ++i) {
   if (i \le 1 \mid | pfac[i] == i) pw[i] = num(i).pow(d);
   else pw[i] = (pw[pfac[i]] * pw[i / pfac[i]]);
  return pw;
num sum_of_power_limit(num r, int d, const vector<num>& fs) {
 interpolator_t<num> M(d + 2);
  vector < num > qs(d + 1); qs[0] = 1;
  for (int x = 1; x \le d; ++x) qs[x] = qs[x - 1] * r;
  num ans = 0, cur_sum = 0;
  for (int x = 0; x \le d; ++x) {
   cur\_sum += qs[x] * fs[x];
   ans += cur_sum * invFac[d - x] * invFac[x + 1] * (((d - x)
        & 1) ? -1 : +1) * qs[d - x];
  // ans is equivalent to invFac(d + 1) * dp(d+1), where
  // for all x in [0, d], dp(x + 1) := E(d, d-x) + dp(x) * r,
      dp(0) = 0.
  // with E being the eulerian number. Works in O(d^2).
  ans *= (1 - r).pow(-(d + 1)) * fac[d + 1];
  return ans;
num sum_of_power(num r, int d, const vector<num>& fs, long long
  if (r == 0) return (0 < N) ? fs[0] : 0;
  interpolator_t<num> M(d + 10);
  vector < num > gs(d + 2); gs[0] = 0; num rr = 1;
  for (int x = 0; x \le d; ++x) {
   qs[x + 1] = qs[x] + rr * fs[x];
   rr *= r;
  if (r == 1) return M.interpolate(gs, N);
  const num c = sum_of_power_limit(r, d, fs);
  const num r inv = r.inv(); num rr inv = 1;
  for (int x = 0; x \le d + 1; ++x) {
   qs[x] = rr inv * (qs[x] - c);
   rr_inv *= r_inv;
  return c + r.pow(N) * M.interpolate(gs, N);
```

4.1.1 Duality

max $c^T x$ sjt to $Ax \leq b$. Dual problem is min $b^T x$ sjt to $A^T x \geq c$. By strong duality, min max value coincides.

4.1.2 Generating functions

A list of generating functions for useful sequences:

| $(1,1,1,1,1,1,\ldots)$ | $\frac{1}{1-z}$ |
|--|-------------------------|
| $(1,-1,1,-1,1,-1,\ldots)$ | $\frac{1}{1+z}$ |
| $(1,0,1,0,1,0,\ldots)$ | $\frac{1}{1-z^2}$ |
| $(1,0,\ldots,0,1,0,1,0,\ldots,0,1,0,\ldots)$ | $\frac{1}{1-z^2}$ |
| $(1, 2, 3, 4, 5, 6, \ldots)$ | $\frac{1}{(1-z)^2}$ |
| $\left(1, \binom{m+1}{m}, \binom{m+2}{m}, \binom{m+3}{m}, \ldots\right)$ | $\frac{1}{(1-z)^{m+1}}$ |
| $(1,c,\binom{c+1}{2},\binom{c+2}{3},\ldots)$ | $\frac{1}{(1-z)^c}$ |
| $(1,c,c^2,c^3,\ldots)$ | $\frac{1}{1-cz}$ |
| $(0,1,\frac{1}{2},\frac{1}{3},\frac{1}{4},\ldots)$ | $\ln \frac{1}{1-z}$ |

A neat manipulation trick is:

$$\frac{1}{1-z}G(z) = \sum_{n} \sum_{k \le n} g_k z^n$$

Number theory (5)

5.1 Modular arithmetic

modular-arithmetic.h

"mod-inv.h"

Description: Operators for modular arithmetic.

```
template<unsigned M > struct modnum {
 static constexpr unsigned M = M;
 using ll = int64_t; using ull = uint64_t; unsigned x;
 modnum& norm(unsigned a) { x = a < M ? a : a - M; return *</pre>
 constexpr modnum(11 a = 0U) : x(unsigned((a %= 11(M)) < 0 ? a)
       + 11(M) : a)) {}
 explicit operator int() const { return x; }
 modnum& operator+=(const modnum& a) { return norm(x + a.x); }
 modnum& operator-=(const modnum& a) { return norm(x - a.x + M
      ); }
 modnum\& operator*=(const modnum\& a) { x = unsigned(ull(x) * a) }
      .x % M); return *this; }
 modnum& operator/=(const modnum& a) { return (*this *= a.inv
 modnum operator+(const modnum& a) const { return (modnum(*
      this) += a); }
 modnum operator-(const modnum& a) const { return (modnum(*
      this) -= a); }
 modnum operator* (const modnum& a) const { return (modnum(*
      this) *= a); }
 modnum operator/(const modnum& a) const { return (modnum(*
      this) /= a); }
 template<typename T> friend modnum operator+(T a, const
      modnum& b) { return (modnum(a) += b); }
 template<typename T> friend modnum operator-(T a, const
      modnum& b) { return (modnum(a) -= b); }
 template<typename T> friend modnum operator*(T a, const
      modnum& b) { return (modnum(a) *= b); }
 template<typename T> friend modnum operator/(T a, const
      modnum& b) { return (modnum(a) /= b); }
 modnum operator+() const { return *this; }
 modnum operator-() const { return modnum() - *this; }
 modnum pow(ll e) const {
   if (e < 0) return inv().pow(-e);
   modnum b = x, xe = 1U;
   for (; e; e >>= 1) { if (e & 1) xe *= b; b *= b; }
   return xe;
 modnum inv() const { return minv(x, M); }
```

```
friend modnum inv(const modnum& a) { return a.inv(); }
explicit operator bool() const { return x; }
friend bool operator==(const modnum& a, const modnum& b) {
    return a.x == b.x; }
friend bool operator!=(const modnum& a, const modnum& b) {
    return a.x != b.x; }
friend ostream &operator<<(ostream& os, const modnum& a) {
    return os << a.x; }
friend istream &operator>>(istream& in, modnum& n) { ll v_;
    in >> v_; n = modnum(v_); return in; }
;
```

mod-inv.h

Description: Find x such that $ax \equiv 1 \pmod{m}$. The inverse only exist if a and m are coprimes.

```
int minv(int a, int m) {
   a %= m; assert(a);
   return a == 1 ? 1 : int(m - int64_t(minv(m, a)) * m / a);
}
```

mod-sum.h

62c1a3, 34 lines

Description: Sums of mod'ed arithmetic progressions.

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

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

decfb8, 17 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) {
    ull to2 = (to * k + c) / m;
    res += to * to2;
    res -= divsum(to2, m-1 - c, m, k) + to2;
  }
  return res;
}
lint modsum(ull to, lint c, lint k, lint m) {
  c = ((c % m) + m) % m;
  k = ((k % m) + m) % m;
  return to * c + k * sumsq(to) - m * divsum(to, c, k, m);
}
```

mod-mul h

Description: Calculate $a \cdot b \mod c$ (or $a^b \mod c$) for $0 \le a, b \le c \le 7.2 \cdot 10^{18}$. **Time:** $\mathcal{O}(1)$ for modmul, $\mathcal{O}(\log b)$ for modpow

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

mod-sart.h

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

```
int jacobi(int64_t a, int64_t m) { // Jacobi symbol (a/m)
int s = 1;
if (a < 0) a = a % m + m;</pre>
```

```
for (; m > 1; ) {
   a %= m;
   if (a == 0) return 0;
   const int r = __builtin_ctzll(a);
   if ((r \& 1) \&\& ((m + 2) \& 4)) s = -s;
   a >>= r;
   if (a & m & 2) s = -s;
   swap(a, m);
 return s;
vector<int64_t> mod_sqrt(int64_t a, int64_t p) {
 if (p == 2) return {a & 1};
 const int j = jacobi(a, p);
 if (j == 0) return {0};
 if (j == -1) return {};
 int64_t b, d;
  while (true) {
   b = xrand() % p;
   d = (b * b - a) % p;
   if (d < 0) d += p;
   if (jacobi(d, p) == -1) break;
 int64_t f0 = b, f1 = 1, g0 = 1, g1 = 0, tmp;
  for (int64_t e = (p + 1) >> 1; e; e >>= 1) {
   if (e & 1) {
     tmp = (q0 * f0 + d * ((q1 * f1) % p)) % p;
     q1 = (q0 * f1 + q1 * f0) % p;
     q0 = tmp;
   tmp = (f0 * f0 + d * ((f1 * f1) % p)) % p;
   f1 = (2 * f0 * f1) % p;
   f0 = tmp;
  return (q0 {q0, p - q0} : vector<
      int64_t>{p - g0, g0};
```

mod-range.h

Description: min $x \ge 0$ s.t. $l \le ((ax) \mod m) \le r$, m > 0, $a \ge 0$, $a \ge 0$, a

```
template<typename T> T mod_range(T m, T a, T 1, T r) {
    l = max(l, T(0));
    r = min(r, m - 1);
    if (l > r) return -1;
    a %= m;
    if (a == 0) return (l > 0) ? -1 : 0;
    const T k = (l + a - 1) / a;
    if (a * k <= r) return k;
    const T y = mod_range(a, m, a * k - r, a * k - l);
    return (y == -1) ? -1 : ((m * y + r) / a);
}</pre>
```

5.2 Primality

sieve.h

Description: Prime sieve for generating all primes up to a certain limit. pfac[i] is the lowest prime factor of i. Also useful if you need to compute any multiplicative function.

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

a76cb9, 24 lines

```
vector<int> run_sieve(int N) {
  vector<int> pfac(N + 1);
  vector<int> primes; primes.reserve(N+1);
  vector<int> mu(N + 1, -1); mu[1] = 1;
  vector<int> phi(N + 1); phi[1] = 1;
  for (int i = 2; i <= N; ++i) {
    if (!pfac[i]) {
      pfac[i] = i; primes.push_back(i);
      phi[i] = i - 1;</pre>
```

```
for (int p : primes) {
   if (p > N/i) break;
   pfac[p * i] = p;
   mu[p * i] *= mu[i];
   phi[p * i] = phi[i] * phi[p];
   if (i % p == 0) {
      mu[p * i] = 0;
      phi[p * i] = phi[i] * p;
      break;
   }
}
return primes;
```

miller-rabin.h

Description: Deterministic Miller-Rabin primality test. Guaranteed to work for numbers up to 2^{64} ; for larger numbers, extend A randomly.

pollard-rho.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.

```
"mod-mul.h", "extended-euclid.h", "miller-rabin.h"
                                                      6bf31f, 18 lines
ull pollard(ull n) {
 auto f = [n] (ull x, ull k) { return modmul(x, x, n) + k; };
 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, i);
   if ((q = modmul(prd, max(x,y) - min(x,y), n))) prd = q;
   x = f(x, i), y = f(f(y, i), i);
 return gcd(prd, n);
vector<ull> factor(ull n) {
 if (n == 1) return {};
 if (isPrime(n)) return {n};
 ull x = pollard(n);
 auto l = factor(x), r = factor(n / x);
 l.insert(l.end(), r.begin(), r.end());
 return 1;
```

5.3 Divisibility

extended-euclid.h

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

```
template<typename T>
T egcd(T a, T b, T &x, T &y) {
  if (!a) { x = 0, y = 1; return b; }
```

```
T g = egcd(b % a, a, y, x);
x -= y * (b/a); return g;
}
```

division-lemma.h

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

Description: This lemma let us exploit the fact tha the sequence (harmonic on integer division) has at most $2\sqrt{N}$ distinct elements, so we can iterate through every possible value of $\left\lfloor \frac{N}{i} \right\rfloor$, using the fact that the greatest integer j satisfying $\left\lfloor \frac{N}{i} \right\rfloor = \left\lfloor \frac{N}{j} \right\rfloor$ is $\left\lfloor \frac{N}{N} \right\rfloor$. This one computes the $\sum_{i=1}^{N} \left\lfloor \frac{N}{i} \right\rfloor i$.

```
int res = 0;
for (int a = 1, b; a <= N; a = b + 1) { // floor
   b = N / (N / a);
   // quotient (N/a) and there are (b - a + 1) elements
   int 1 = b - a + 1, r = a + b; // l * r / 2 = sum(i, j)
   if (1 & 1) r /= 2;
   else 1 /= 2;
   res += 1 * r * (N / a);
}
// [1, N), need to deal with case where a = N separately
for (int a = 1, b; a < N; a = b + 1) { // ceil
   const int k = (N - 1) / a + 1; // quotient k
   b = (N - 1) / (k - 1);
   int cnt = b - a + 1; // occur cnt times on interval [a, b]
}
```

divisors.h

Description: Generate all factors of n given it's prime factorization.

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

phi-function.h

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

 $\sum_{d|n} \phi(d) = n, \sum_{1 \le k \le n, \gcd(k,n) = 1} k = n\phi(n)/2, n >$ **Euler's thm**: $a, n \text{ coprime} \Rightarrow a^{\phi(n)} \equiv 1 \pmod{n}$.

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

da7671, 7 lines

```
const int n = int(1e5)*5;
vector<int> phi(n);
void calculatePhi() {
  for(int i = 0; i < n; ++i) phi[i] = i&1 ? i : i/2;
  for(int i = 3; i < n; i += 2) if (phi[i] == i)
      for(int j = i; j < n; j += i) phi[j] -= phi[j]/i;
}</pre>
```

discrete-log.h

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

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

6c6eb0, 18 lines "extended-euclid.h" template<typename T> T modLog(T a, T b, T m) { T k = 1, it = 0, g;while ((q = gcd(a, m)) != 1) { if (b == k) return it; if (b % g) return -1; b /= g; m /= g; ++it; k = k * a / g % m;T n = sqrtl(m) + 1, f = 1, j = 1;unordered_map<T, T> A; while $(j \le n)$ { f = f * a % m;A[f * b % m] = j++;for(int i = 1; $i \le n$; ++i) if (A.count(k = k * f % m) return n * i - A[k] + it; return -1:

prime-counting.h

Description: Count the number of primes up to N. Also useful for sum of primes.

Time: $\mathcal{O}(N^{3/4}/\log N)$

05ef7c, 22 lines

```
struct primes_t {
  vector<11> dp, w;
  ll pi(ll N) {
   const int sqrtN = int(sqrt(N));
    for (11 a = 1, b; a \le N; a = b+1) {
     b = N / (N / a);
     w.push_back(N/a);
    auto get = [\&](ll x) {
     if (x \le sartN) return int (x-1);
     return int(w.size() - N/x);
    reverse(w.begin(), w.end()); dp.reserve(w.size());
    for (auto& x : w) dp.push back(x-1);
    for (11 i = 2; i*i <= N; ++i) {
     if (dp[i-1] == dp[i-2]) continue;
     for (int j = int(w.size())-1; w[j] >= i*i; --j)
       dp[j] = dp[qet(w[j]/i)] - dp[i-2];
    return dp.back();
};
```

5.4 Chinese remainder theorem

chinese-remainder.h

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

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

5.5 Fractions

fractions.h

Description: Template that helps deal with fractions.

df1f1d, 31 lines

```
template<typename num = long long>
struct frac {
 num n, d;
 frac() : n(0), d(1) { }
 frac(num _n, num _d = 1): n(_n), d(_d){
   num g = gcd(n, d); n \neq g, d \neq g;
   if (d < 0) n *= -1, d *= -1;
   assert (d != 0);
 friend bool operator<(const frac& 1, const frac& r) { return
      1.n * r.d < r.n * 1.d; }
 friend bool operator == (const frac& 1, const frac& r) { return
       1.n == r.n && 1.d == r.d; }
 friend bool operator!=(const frac& 1, const frac& r) { return
       !(1 == r);
 friend frac operator+(const frac& 1, const frac& r) {
   num q = \gcd(1.d, r.d);
    return frac( r.d / g * 1.n + 1.d / g * r.n, 1.d / g * r.d);
 friend frac operator-(const frac& 1, const frac& r) {
   num g = gcd(1.d, r.d);
    return frac( r.d / g * 1.n - 1.d / g * r.n, 1.d / g * r.d);
  friend frac operator*(const frac& 1, const frac& r) { return
      frac(l.n * r.n, l.d * r.d); }
  friend frac operator/(const frac& 1, const frac& r) { return
      1 * frac(r.d, r.n); }
  friend frac& operator+=(frac& 1, const frac& r) { return 1 =
  friend frac& operator==(frac& 1, const frac& r) { return l =
      1-r; }
 template < class T > friend frac& operator *= (frac& 1, const T& r
      ) { return 1 = 1*r; }
 template < class T > friend frac& operator /= (frac& 1, const T& r
      ) { return 1 = 1/r; }
  friend ostream& operator<<(ostream& strm, const frac& a) {</pre>
   strm << a.n << "/" << a.d;
    return strm;
};
```

continued-fractions.h

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

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

Time: $\mathcal{O}(\log N)$ 61608b, 21 lines

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

```
if (abs(y = 1/(y - (dbl)a)) > 3*N) {
    return {NP, NQ};
}
LP = P; P = NP;
LQ = Q; Q = NQ;
}
```

frac-binary-search.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 { lint p, q; };
template<class F>
Frac fracBS(F f, lint N) {
 bool dir = 1, A = 1, B = 1;
 Frac left{0, 1}, right{1, 1}; // Set right to 1/0 to search
       (0, N]
  assert(!f(left)); assert(f(right));
  while (A | | B) {
   lint adv = 0, step = 1; // move right if dir, else left
    for (int si = 0; step; (step *= 2) >>= si) {
      adv += step;
      Frac mid{left.p * adv + right.p, left.q * adv + right.q};
      if (abs(mid.p) > N || mid.q > N || dir == !f(mid)) {
        adv -= step; si = 2;
    right.p += left.p * adv;
    right.q += left.q * adv;
    dir = !dir;
    swap(left, right);
    A = B; B = !!adv;
  return dir ? right : left;
```

5.5.1 Bézout's identity

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

$$ax + by = d$$

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

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

5.5.2 Pythagorean Triples

The Pythagorean triples are uniquely generated by

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

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

5.6 Primes

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

dirichlet-convolution int-perm lucas rolling-binomial

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 $\mathbb{Z}_2 \times \mathbb{Z}_{2^{a-2}}$.

5.6.1 Prime counting function $(\pi(x))$

The prime counting function is asymptotic to $\frac{x}{\log x}$, by the prime number theorem.

| | x | 10 | 10^{2} | 10^{3} | 10^{4} | 10^{5} | 10^{6} | 10^{7} | 10^{8} |
|-------|-----|----|----------|----------|----------|----------|----------|----------|-----------|
| π | (x) | 4 | 25 | 168 | 1.229 | 9.592 | 78.498 | 664.579 | 5.761.455 |

5.6.2 Sum of primes

For any multiplicative f:

$$S(n,p) = S(n,p-1) - f(p) \cdot (S(n/p,p-1) - S(p-1,p-1))$$

5.6.3 Moebius Function

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

Moebius Inversion:

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

Other useful formulas/forms:

$$\sum_{d|n} \phi(d) = n$$

$$\sum_{\substack{i < n \\ \gcd(i,n)=1}} i = n \frac{\phi(n)}{2}$$

$$\sum_{a=1}^{n} \sum_{b=1}^{n} [\gcd(a,b) = 1] = \sum_{d=1}^{n} \mu(d) \left| \frac{n}{d} \right|^{2}$$

$$\sum_{a=1}^{n} \sum_{b=1}^{n} \gcd(a,b) = \sum_{d=1}^{n} d \sum_{d|x}^{n} \left| \frac{n}{x} \right|^{2} \mu(\frac{x}{d})$$

$$\sum_{a=1}^{n} \sum_{b=a}^{n} \gcd(a,b) = \sum_{d=1}^{n} \sum_{d|x}^{n} \phi(\frac{x}{d})d$$

$$\sum_{a=1}^{n} \sum_{b=1}^{n} \text{lcm}(a,b) = \sum_{d=1}^{n} \mu(d) d \sum_{d \mid x}^{n} x \left(\frac{x}{2} \right)^{1/2} + 1$$

$$\sum_{a=1}^{n} \sum_{b=a+1}^{n} lcm(a,b) = \sum_{d=1}^{n} \sum_{d|x}^{n} \phi(\frac{x}{d}) \frac{x^{2}}{2d}$$

$$\sum_{a \in S} \sum_{b \in S} \gcd(a, b) = \sum_{d=1}^{n} \left(\sum_{x|d} \frac{d}{x} \mu(x)\right) \left(\sum_{d|v} \operatorname{freq}[v]\right)^{2}$$

$$\sum_{a \in S} \sum_{b \in S} \operatorname{lcm}(a, b) = \sum_{d=1}^{n} \left(\sum_{x \mid d} \frac{x}{d} \mu(x) \right) \left(\sum_{v \in S, d \mid v} v \right)^{2}$$

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

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

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

5.6.4 Dirichlet Convolution

Given a function f(x), let

$$(f * g)(x) = \sum_{d \mid x} g(d)f(x/d)$$

If the partial sums $s_{f*g}(n)$, $s_g(n)$ can be computed in O(1) and $s_f(1...n^{2/3})$ can be computed in $O\left(n^{2/3}\right)$ then all $s_f\left(\frac{n}{d}\right)$ can as well. Use

$$s_{f*g}(n) = \sum_{d=1}^{n} g(d)s_f(n/d).$$

$$\implies s_f(n) = \frac{s_{f*g}(n) - \sum_{d=2}^n g(d)s_f(n/d)}{g(1)}$$

- 1. If $f(x) = \mu(x)$ then g(x) = 1, (f * g)(x) = (x == 1), and $s_f(n) = 1 \sum_{i=2}^n s_f(n/i)$
- 2. If $f(x) = \phi(x)$ then g(x) = 1, (f * g)(x) = x, and $s_f(n) = \frac{n(n+1)}{2} \sum_{i=2}^n s_f(n/i)$

dirichlet-convolution.h

Description: Dirichlet convolution. Change $f,\,gs$ and fgs accordingly. This example calculates $\phi(N)$.

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

bd3050, 2

template<typename T, typename V> struct dirichlet_convolution V N; $// \sim N^{2/3}$ T inv; vector<V> fs; // can be any multiplicative function vector<T> psum; unordered_map<V, T> mapa; V f(V x) { return fs[x]; } T gs(V x) { return x; } T fqs(V x) { return T(x) * (x + 1) / 2; } dirichlet_convolution(V _N, const vector<V>& F) : N(_N + 1), fs(F), psum(N + 1) { inv = qs(1);for (V a = 0; a + 1 < N; ++a) psum[a + 1] = f(a + 1) + psum[a];T query(V x) { if (x < N) return psum[x];</pre> if (mapa.find(x) != mapa.end()) return mapa[x]; T ans = fqs(x); for $(V a = 2, b; a \le x; a = b + 1)$ { b = x / (x / a);ans -= (qs(b) - qs(a - 1)) * query(x / a); return mapa[x] = (ans / inv); };

Combinatorial (6)

6.1 Permutations

6.1.1 Factorial

| n | | | | 5 6 | | | | 9 | 10 | |
|----|-----|------------|------|-------|--------|--------|--------|-----------|--------|-----|
| n! | 1 2 | 6 | 24 1 | 20 72 | 0 504 | 10 40 | 320 36 | $2880\ 3$ | 628800 | |
| n | | | | | | | | 16 | | |
| n! | 4.0 | e 7 | 4.8e | 8 6.2 | e9 8.7 | 'e10 î | 1.3e12 | 2.1e13 | 3.6e14 | |
| n | 20 | | 25 | 30 | 40 | 50 | 100 | 150 | 171 | |
| n! | 2e1 | 8 | 2e25 | 3e32 | 8e47 | 3e64 | 9e157 | 7 6e262 | DBL_M | IAX |

```
int-perm.h
```

Description: Permutation -> integer conversion. (Not order preserving.) **Time:** $\mathcal{O}(n)$

6.1.2 Binomials

- Sum of every element in the n-th row of pascal triangle is 2^n .
- The product of the elements in each row is $\frac{(n+1)^n}{n!}$
- $\bullet \sum_{k=0}^{n} \binom{n}{k}^2 = \binom{2n}{n}$
- In a row p where p is a prime number, all the terms in that row except the 1s are multiples of p
- To count odd terms in row n, convert n to binary. Let x be the number of 1s in the binary representation. Then the number of odd terms will be 2^x
- Every entry in row $2^n 1$ is odd

lucas.h

Description: Lucas' thm: Let n, m be non-negative integers and p a prime. Write $n = n_k p^k + ... + n_1 p + n_0$ and $m = m_k p^k + ... + m_1 p + m_0$. Then $\binom{n}{m} \equiv \prod_{i=0}^k \binom{n_i}{m_i} \pmod{p}$. fact and ifact must hold pre-computed factorials / inverse factorials, e.g. from ModInv.h.

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

rolling-binomial.h

Description: $\binom{n}{k}$ (mod m) in time proportional to the difference between (n, k) and the previous (n, k).

"../number-theory/preparator.h" d087bf, 14 lines

```
using i64 = int64_t;
const int mod = int(1e9) + 7;
struct Bin {
  int N = 0,    K = 0; i64 r = 1;
  void m(int a, int b) { r = r * a % mod * invs[b] % mod; }
  i64 choose(int n, int k) {
   if (k > n || k < 0) return 0;
   while(N < n) ++N, m(N, N - K);
   while(K < k) ++K, m(N - K + 1, K);
   while(K > k) m(K, N - K + 1, K);
   while(N > n) m(N - K, N), --N;
   return r;
```

};

multinomial partitions

multinomial.h

Description: Computes
$$\binom{k_1 + \dots + k_n}{k_1, k_2, \dots, k_n} = \frac{(\sum k_i)!}{k_1! k_2! \dots k_n!}$$
. 864cdb, 7 lines

$$\binom{n}{k} = \frac{n}{k} \binom{n-1}{k-1} \qquad \qquad \binom{n}{k} \binom{n-k}{v} = \binom{n}{v} \binom{n-v}{k}$$

$$\sum_{k=0}^{n} \binom{n}{k} k = n2^{n-1} \qquad \qquad \sum_{j=0}^{k} \binom{m}{j} \binom{n-m}{k-j} = \binom{n}{k}$$

$$\sum_{m=0}^{n} \binom{m}{j} \binom{n-m}{k-j} = \binom{n+1}{k+1} \qquad \qquad \sum_{j=0}^{n} \binom{m}{j} \binom{n-k}{v} = \binom{n}{k}$$

$$(x+y)^n = \sum_{k=0}^{n} \binom{n}{k} x^{n-k} y^k \qquad \qquad \sum_{m=k}^{n} \binom{m}{r} = \binom{n+1}{k+1}$$

$$\sum_{j=0}^{m} \binom{m}{j}^2 = \binom{2m}{m}$$

$$2 \sum_{i=L}^{R} \binom{n}{i} - \binom{n}{L} - \binom{n}{R} = \sum_{i=L+1}^{R} \binom{n+1}{i}$$

6.1.3 Involutions

An involution is a permutation with maximum cycle length 2, and it is its own inverse.

$$a(n) = a(n-1) + (n-1)a(n-2), a(0) = a(1) = 1.$$

1, 1, 2, 4, 10, 26, 76, 232, 764, 2620, 9496, 35696, 140152

6.1.4 Cycles

Let the number of n-permutations whose cycle lenghts all belong to the set S be denoted by $g_S(n)$

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

6.1.5 The twelvefold way (from Stanley)

How many functions $f \colon N \to X$ are there?

| N | X | Any f | Injective | Surjective |
|---------|---------|--|---------------------|--------------------|
| dist. | dist. | x^n | $\frac{x!}{(x-n)!}$ | $x!\binom{n}{x}$ |
| indist. | dist. | $\binom{x+n-1}{n}$ | $\binom{x}{n}$ | $\binom{n-1}{n-x}$ |
| dist. | indist. | ${n \brace 1} + \ldots + {n \brace x}$ | $[n \le x]$ | $\binom{n}{k}$ |
| indist. | indist. | $p_1(n) + \dots p_x(n)$ | $[n \leq x]$ | $p_x(n)$ |

Where $\binom{a}{b} = \frac{1}{b!}(a)_b$, $p_x(n)$ is the number of ways to partition the integer n using x summand and $\binom{n}{x}$ is the number of ways to partition a set of n elements into x subsets (aka Stirling number of the second kind).

6.1.6 Burnside

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

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

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

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

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

6.1.7 Derangements

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

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

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.

$$p(0) = 1, \ p(n) = \sum_{k \in \mathbb{Z} \setminus \{0\}} (-1)^{k+1} p(n - k(3k - 1)/2)$$
$$p(n) \sim 0.145/n \cdot \exp(2.56\sqrt{n})$$
$$\frac{n}{p(n)} \begin{vmatrix} 0.1 & 2.3 & 4.5 & 6.7 & 8.9 & 20.50 & 100\\ 1 & 1.2 & 3.5 & 7.11 & 15.22 & 30.627 & \sim 2e5 & \sim 2e8 \end{vmatrix}$$

partitions.h

378a72, 16 lines

6.3 General purpose numbers

6.3.1 Bernoulli numbers

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

Sums of powers:

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

Euler-Maclaurin formula for infinite sums:

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

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

6.3.2 Stirling numbers of the first kind

Number of permutations on n items with k cycles.

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

c(8,k) = 8,0,5040,13068,13132,6769,1960,322,28,1c(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,

$$\mathcal{B}_{n+1} = \sum_{k=0}^{n} \binom{n}{k} \mathcal{B}_k$$

Also possible to calculate using Stirling numbers of the second kind,

$$B_n = \sum_{k=0}^n S(n,k)$$

If p is 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 : $(n-2)!/((d_1-1)!\cdots(d_n-1)!)$ # forests with exactly k rooted trees:

$$\binom{n}{k}k \cdot n^{n-k-1}$$

6.3.7 Catalan numbers

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

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

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

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

6.3.8 Super Catalan numbers

The number of monotonic lattice paths of a $n \times n$ grid that do not touch the diagonal.

$$S(n) = \frac{3(2n-3)S(n-1) - (n-3)S(n-2)}{n}$$
$$S(1) = S(2) = 1$$

1, 1, 3, 11, 45, 197, 903, 4279, 20793, 103049, 518859

6.3.9 Motzkin numbers

Number of ways of drawing any number of nonintersecting chords among n points on a circle. Number of lattice paths from (0, 0) to (n, 0) never going below the x-axis, using only steps NE, E, SE.

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

$$M(0) = M(1) = 1$$

 $1,\ 1,\ 2,\ 4,\ 9,\ 21,\ 51,\ 127,\ 323,\ 835,\ 2188,\ 5798,\ 15511,\ 41835,\ 113634$

6.3.10 Narayana numbers

Number of lattice paths from (0,0) to (2n,0) never going below the x-axis, using only steps NE and SE, and with k peaks.

$$N(n,k) = \frac{1}{n} \binom{n}{k} \binom{n}{k-1}$$
$$N(n,1) = N(n,n) = 1$$
$$\sum_{k=1}^{n} N(n,k) = C_n$$

1, 1, 1, 1, 3, 1, 1, 6, 6, 1, 1, 10, 20, 10, 1, 1, 15, 50

6.3.11 Schroder numbers

Number of lattice paths from (0, 0) to (n, n) using only steps N,NE,E, never going above the diagonal. Number of lattice paths from (0, 0) to (2n, 0) using only steps NE, SE and double east EE, never going below the x-axis. Twice the Super Catalan number, except for the first term.

1, 2, 6, 22, 90, 394, 1806, 8558, 41586, 206098

6.3.12 Triangles

Given rods of length 1, ..., n,

$$T(n) = \frac{1}{24} \left\{ \begin{array}{ll} n(n-2)(2n-5) & n \text{ even} \\ (n-1)(n-3)(2n-1) & n \text{ odd} \end{array} \right\}$$

is the number of distinct triangles (positive are) that can be constructed, i.e., the # of 3-subsets of [n] s.t. $x \le y \le z$ and $z \ne x + y$.

6.4 Fibonacci

$$Fib(x+y) = Fib(x+1)Fib(y) + Fib(x)Fib(y-1)$$

$$Fib(n+1)Fib(n-1) - Fib(n)^2 = (-1)^n$$

$$Fib(2n-1) = Fib(n)^2 - Fib(n-1)^2$$

$$\sum_{i=0}^{n} Fib(i) = Fib(n+2) - 1$$

$$\sum_{i=0}^{n} Fib(i)^{2} = Fib(n)Fib(n+1)$$

$$\sum_{i=0}^{n} Fib(i)^{3} = \frac{Fib(n)Fib(n+1)^{2} - (-1)^{n}Fib(n-1) + 1}{2}$$

6.5 Linear Recurrences

$$\begin{split} F_i &= \sum_{j=1}^K C_j F_{i-j} + D \\ \begin{bmatrix} 0 & 1 & 0 & 0 & & 0 & 0 \\ 0 & 0 & 1 & 0 & & 0 & 0 \\ 0 & 0 & 0 & 1 & & 0 & 0 \\ 0 & 0 & 0 & 1 & & & 0 & 0 \\ \vdots & \vdots & \ddots & \ddots & \ddots & \ddots & \ddots \\ C_K & C_{K-1} & C_{K-2} & C_{K-3} & & & \ddots & C_1 & 1 \\ 0 & 0 & 0 & 0 & & & & 0 & 1 \end{bmatrix} \begin{bmatrix} F_0 \\ F_1 \\ F_2 \\ F_3 \\ \vdots \\ F_{K-1} \\ D \end{bmatrix} = \begin{bmatrix} F_1 \\ F_2 \\ F_3 \\ \vdots \\ F_M \end{bmatrix} \\ \vdots \\ F_M \end{bmatrix}$$

6.6 Game Theory

A game can be reduced to Nim if it is a finite impartial game. Nim and its variants include:

6.6.1 Nim

Let $X = \bigoplus_{i=1}^n x_i$, then $(x_i)_{i=1}^n$ is a winning position iff $X \neq 0$. Find a move by picking k such that $x_k > x_k \oplus X$.

6.6.2 Misère Nim

Regular Nim, except that the last player to move *loses*. Play regular Nim until there is only one pile of size larger than 1, reduce it to 0 or 1 such that there is an odd number of piles. The second player wins (a_1, \ldots, a_n) if 1) there is a pile $a_i > 1$ and $\bigoplus_{i=1}^n a_i = 0$ or 2) all $a_i \le 1$ and $\bigoplus_{i=1}^n a_i = 1$.

6.6.3 Staircase Nim

Stones are moved down a staircase and only removed from the last pile. $(x_i)_{i=1}^n$ is an L-position if $(x_{2i-1})_{i=1}^{n/2}$ is (i.e. only look at odd-numbered piles).

nim-product.cpp

Description: Product of nimbers is associative, commutative, and distributive over addition (xor). Forms finite field of size 2^{2^k} . Application: Given 1D coin turning games $G_1, G_2, G_1 \times G_2$ is the 2D coin turning game defined as follows. If turning coins at x_1, x_2, \ldots, x_m is legal in G_1 and y_1, y_2, \ldots, y_n is legal in G_2 , then turning coins at all positions (x_i, y_j) is legal assuming that the coin at (x_m, y_n) goes from heads to tails. Then the grundy function g(x, y) of $G_1 \times G_2$ is $g_1(x) \times g_2(y)$.

Time: 64² xors per multiplication, memorize to speed up. 38fb87, 28 lines

```
using ull = uint64_t;
ull nim prod[64][64];
ull nim_prod2(int i, int j) {
 if (nim_prod[i][j]) return nim_prod[i][j];
 if ((i & j) == 0) return nim_prod[i][j] = 1ull << (i|j);</pre>
  int a = (i\&j) \& -(i\&j);
 return nim_prod[i][j] = nim_prod2(i ^ a, j) ^ nim_prod2((i ^
      a) | (a-1), (j^a) | (i & (a-1));
void all_nim_prod() {
 for (int i = 0; i < 64; i++) {
    for (int j = 0; j < 64; j++) {
      if ((i & j) == 0) nim_prod[i][j] = lull << (i|j);</pre>
        int a = (i&j) & -(i&j);
        nim_prod[i][j] = nim_prod[i ^ a][j] ^ nim_prod[(i ^ a)
             | (a-1)][(j ^ a) | (i & (a-1))];
ull get_nim_prod(ull x, ull y) {
 ull res = 0;
 for (int i = 0; i < 64 && (x >> i); ++i)
    if ((x >> i) & 1)
      for (int j = 0; j < 64 && (y >> j); ++j)
        if ((y >> j) & 1) res ^= nim_prod2(i, j);
 return res;
```

Graph (7)

7.1 Fundamentals

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

using pii = pair<int,int>;
vector<int> eulerWalk(vector<vector<pii>>>& gr, int nedges, int
 src=0) {
 int n = gr.size();
 vector<int> 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 = int(gr[x].size

```
if (it == end) { ret.push_back(x); s.pop_back(); continue; }
  tie(y, e) = qr[x][it++];
 if (!eu[e]) {
   D[x] --, D[y] ++;
   eu[e] = 1; s.push_back(y);
for (auto &x : D) if (x < 0 \mid \mid int(ret.size()) != nedges+1)
    return {};
return {ret.rbegin(), ret.rend()};
```

7.2 Network flow

push-relabel.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. id can be used to restore each edge and its amount of flow used.

Time: $\mathcal{O}\left(V^2\sqrt{E}\right)$ Better for dense graphs - Slower than Dinic (in practice)

```
template<typename flow_t = int> struct PushRelabel {
  struct edge_t { int dest, back; flow_t f, c; };
  vector<vector<edge_t>> g;
  vector<flow_t> ec;
  vector<edge_t*> cur;
  vector<vector<int>> hs; vector<int> H;
  PushRelabel(int n) : g(n), ec(n), cur(n), hs(2*n), H(n) {}
  void addEdge(int s, int t, flow_t cap, flow_t rcap = 0) { //
       d58501
    if (s == t) return;
    g[s].push_back({t, (int)g[t].size(), 0, cap});
    g[t].push_back({s, (int)g[s].size()-1, 0, rcap});
  void addFlow(edge_t& e, flow_t f) { // 2f7969
    edge_t &back = g[e.dest][e.back];
    if (!ec[e.dest] && f) hs[H[e.dest]].push_back(e.dest);
    e.f += f; e.c -= f; ec[e.dest] += f;
   back.f -= f; back.c += f; ec[back.dest] -= f;
  flow_t maxflow(int s, int t) { // 21100c
    int v = int(g.size()); H[s] = v; ec[t] = 1;
    vector<int> co(2*v); co[0] = v-1;
    for(int i = 0; i < v; ++i) cur[i] = q[i].data();</pre>
    for(auto& e : g[s]) addFlow(e, e.c);
    for (int hi = 0;;) {
     while (hs[hi].empty()) if (!hi--) return -ec[s];
     int u = hs[hi].back(); hs[hi].pop_back();
     while (ec[u] > 0) // discharge u
       if (cur[u] == g[u].data() + g[u].size()) {
          H[u] = 1e9;
          for(auto &e : q[u]) if (e.c && H[u] > H[e.dest]+1)
            H[u] = H[e.dest]+1, cur[u] = &e;
          if (++co[H[u]], !--co[hi] && hi < v)</pre>
            for (int i = 0; i < v; ++i) if (hi < H[i] && H[i] <
               --co[H[i]], H[i] = v + 1;
          hi = H[u];
        } else if (\operatorname{cur}[u] -> \operatorname{c \&\& H}[u] == \operatorname{H}[\operatorname{cur}[u] -> \operatorname{dest}] + 1)
          addFlow(*cur[u], min(ec[u], cur[u]->c));
        else ++cur[u];
 bool leftOfMinCut(int a) { return H[a] >= int(g.size()); }
```

dinitz.h

Description: Flow algorithm with complexity $O(VE \log U)$ where U =max |cap|. $O(\min(E^{1/2}, V^{2/3})E)$ if U = 1; $O(\sqrt{V}E)$ for bipartite matching. To obtain each partition A and B of the cut look at lvl, for $v \subset A$, |lvl[v]| > 0, for $u \subset B$, |lvl[u]| = 0. 7c5dcd, 75 lines

```
template<typename T = int> struct Dinitz {
 struct edge_t { int to, rev; T c, f; };
 vector<vector<edge_t>> adj;
 vector<int> lvl, ptr, q;
 Dinitz(int n) : lvl(n), ptr(n), q(n), adj(n) {}
 inline void addEdge(int a, int b, T c, T rcap = 0) { // 694
    adj[a].push_back({b, (int)adj[b].size(), c, 0});
   adj[b].push_back({a, (int)adj[a].size() - 1, rcap, 0});
 T dfs(int v, int t, T f) { // 8ffe6b
   if (v == t \mid \mid !f) return f;
    for (int &i = ptr[v]; i < int(adj[v].size()); ++i) {</pre>
      edge_t &e = adj[v][i];
     if (lvl[e.to] == lvl[v] + 1)
       if (T p = dfs(e.to, t, min(f, e.c - e.f))) {
         e.f += p, adj[e.to][e.rev].f -= p;
          return p;
    return 0;
 T maxflow(int s, int t) { // db2141
    T flow = 0; q[0] = s;
    for (int L = 0; L < 31; ++L) do { // 'int L=30' maybe
        faster for random data
      lvl = ptr = vector<int>(q.size());
      int qi = 0, qe = lvl[s] = 1;
      while (qi < qe && !lvl[t]) {
       int v = q[qi++];
        for (edge_t &e : adj[v])
         if (!lvl[e.to] && (e.c - e.f) >> (30 - L))
            q[qe++] = e.to, lvl[e.to] = lvl[v] + 1;
      while (T p = dfs(s, t, numeric_limits<T>::max()/4)) flow
          += p;
    } while (lvl[t]);
    return flow;
 bool leftOfMinCut(int v) { return bool(lvl[v] != 0); }
 pair<T, vector<pair<int,int>>> minCut(int s, int t) { // 727
      b22
   T cost = maxflow(s,t);
    vector<pair<int,int>> cut;
    for (int i = 0; i < int(adj.size()); i++) for(edge_t &e :</pre>
     if (lvl[i] && !lvl[e.to]) cut.push_back({i, e.to});
    return {cost, cut};
};
struct flow_demand_t {
 int src, sink;
 vector<int> d;
 Dinitz<int> flower;
 flow_demand_t(int N) : src(N + 1), sink(N + 2), d(N + 3),
      flower(N + 3) \{ \}
 void add_edge(int a, int b, int demand, int cap) {
   d[a] -= demand;
   d[b] += demand;
    flower.addEdge(a, b, cap - demand);
 int get_flow() {
```

```
const int INF = std::numeric limits<int>::max():
     int x = 0, y = 0;
    flower.add_edge(N, N-1, INF);
    for (int i = 0; i <= N; ++i) {
     if (d[i] < 0) {
       flower.add_edge(i, sink, -d[i]);
       x += -d[i];
     if (d[i] > 0) {
        flower.add_edge(src, i, d[i]);
       y += d[i];
    bool has circulation = (flower.maxflow(src, sink) == x && x
         == y);
    if (!has_circulation) return -1;
    return flower.maxflow(N-1, N);
};
```

min-cost-max-flow.h

Description: Min-cost max-flow. Assumes there is no negative cycle.

Time: $\mathcal{O}(F(V+E)logV)$, being F the amount of flow.

```
e7ff44, 73 lines
template<class flow_t, class cost_t> struct min_cost {
 static constexpr flow t FLOW EPS = 1e-10L;
 static constexpr flow_t FLOW_INF = std::numeric_limits<flow_t</pre>
 static constexpr cost_t COST_EPS = 1e-10L;
 static constexpr cost t COST INF = std::numeric limits<cost t
      >::max();
 int n, m;
 vector<int> ptr, nxt, zu;
 vector<flow t> capa;
 vector<cost t> cost:
 void add_edge(int u, int v, flow_t w, cost_t c) { // d482f5
   nxt.push_back(ptr[u]); zu.push_back(v); capa.push_back(w);
        cost.push_back( c); ptr[u] = m++;
   nxt.push_back(ptr[v]); zu.push_back(u); capa.push_back(0);
        cost.push_back(-c); ptr[v] = m++;
 vector<cost_t> pot, dist;
 vector<bool> vis:
 vector<int> pari;
 // cost \ slopes[j] \ per \ flow \ when \ flows[j] <= flow <= flows[j] +
 vector<flow_t> flows; vector<cost_t> slopes;
 // The distance to a vertex might not be determined if it
      is >= dist[t].
     You can pass t = -1 to find a shortest path to each
 void shortest(int s, int t) { // e9bb0d
   using Entry = pair<cost_t, int>;
   priority_queue<Entry, vector<Entry>, std::greater<Entry>>
   for (int u = 0; u < n; ++u) { dist[u] = COST_INF; vis[u] =
        false; }
   for (que.emplace(dist[s] = 0, s); !que.empty(); ) {
     const cost_t c = que.top().first;
     const int u = que.top().second;
     que.pop();
     if (vis[u]) continue;
     vis[u] = true;
     if (u == t) return;
     for (int i = ptr[u]; \sim i; i = nxt[i]) if (capa[i] >
          FLOW_EPS) {
       const int v = zu[i];
       const cost_t cc = c + cost[i] + pot[u] - pot[v];
```

```
if (dist[v] > cc) { que.emplace(dist[v] = cc, v); pari[
          v] = i; 
pair<flow_t, cost_t> run(int s, int t, flow_t limFlow =
    FLOW INF) {
 assert(0 <= limFlow);
 pot.assign(n, 0);
 while (true) {
   bool upd = false;
   for (int i = 0; i < m; ++i) if (capa[i] > FLOW_EPS) {
     const int u = zu[i ^1], v = zu[i];
     const cost_t cc = pot[u] + cost[i];
     if (pot[v] > cc + COST_EPS) { pot[v] = cc; upd = true;
   if (!upd) break;
 dist.resize(n); vis.resize(n); pari.resize(n);
  flows.clear(); flows.push_back(0);
 slopes.clear();
  flow t flow = 0; cost t cost = 0;
 while (flow < limFlow) {</pre>
   shortest(s, t);
   if (!vis[t]) break;
   for (int u = 0; u < n; ++u) pot[u] += min(dist[u], dist[t])
   flow_t f = limFlow - flow;
   for (int v = t; v != s; ) {
     const int i = pari[v]; if (f > capa[i]) { f = capa[i];
          v = zu[i ^ 1];
   for (int v = t; v != s; ) {
     const int i = pari[v]; capa[i] -= f; capa[i ^ 1] += f;
          v = zu[i ^1;
   flow += f:
   cost += f * (pot[t] - pot[s]);
   flows.push_back(flow); slopes.push_back(pot[t] - pot[s]);
 return {flow, cost};
```

Matching

hopcroft-karp.h

};

Description: Fast bipartite matching algorithm. Graph q 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: vector<int> btoa(m, -1); hopcroftKarp(q, btoa);

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

d9a55d, 35 lines

```
using vi = vector<int>;
bool dfs(int a, int L, const vector<vi> &q, vi &btoa, vi &A, vi
  if (A[a] != L) return 0;
  A[a] = -1;
  for (auto &b : g[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 hopcroftKarp(const vector<vi> &g, vi &btoa) {
  int res = 0;
```

```
vector<int> A(g.size()), B(int(btoa.size())), cur, next;
for (;;) {
  fill(A.begin(), A.end(), 0), fill(B.begin(), B.end(), 0);
  cur.clear();
  for (auto &a : btoa) if (a != -1) A[a] = -1;
  for (int a = 0; a < g.size(); ++a) if (A[a] == 0) cur.
       push back(a);
  for (int lay = 1;; ++lay) {
   bool islast = 0; next.clear();
    for(auto &a : cur) for(auto &b : g[a]) {
      if (btoa[b] == -1) B[b] = lay, islast = 1;
      else if (btoa[b] != a && !B[b])
        B[b] = lay, next.push_back(btoa[b]);
    if (islast) break;
    if (next.empty()) return res;
    for(auto &a : next) A[a] = lay;
    cur.swap(next);
  for(int a = 0; a < int(g.size()); ++a)</pre>
    res += dfs(a, 0, q, btoa, A, B);
```

bipartite-matching.h

Description: Fast Kuhn! Simple maximum cardinality bipartite matching algorithm. Better than hopcroftKarp in practice. Worst case is O(VE) on an hairy tree. Shuffling the edges and vertices ordering should break some worst-case inputs.

Time: $\Omega(V\bar{E})$

19101e, 42 lines

```
struct bipartite_matching {
 int N. M. T:
 vector<vector<int>> adi;
 vector<int> match, seen;
 bipartite_matching(int a, int b) : N(a), M(a + b), adj(M),
 match(M, -1), seen(M, -1), T(0) {}
 void add_edge(int a, int b) +
   assert(0 \le a \&\& a \le N \&\& b + N \le M \&\& N \le b + N);
   adj[a].push_back(b + N);
 void shuffle_edges() { // useful to break some hairy tests
   mt19937 rng(chrono::steady_clock::now().time_since_epoch().
        count());
    for (auto& cur : adj)
     shuffle(cur.begin(), cur.end(), rng);
 bool dfs(int cur) {
   if (seen[cur] == T) return false;
   seen[cur] = T;
   for (int nxt : adj[cur])
     if (match[nxt] == -1) {
       match[nxt] = cur, match[cur] = nxt;
       return true;
    for (int nxt : adj[cur])
     if (dfs(match[nxt])) {
       match[nxt] = cur, match[cur] = nxt;
       return true;
    return false;
 int solve() {
   int res = 0;
   while (true) {
     int cur = 0; ++T;
     for (int i = 0; i < N; ++i)
```

if (match[i] == -1) cur += dfs(i);

if (cur == 0) break;

```
else res += cur;
   return res;
};
```

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}\left(N^2M\right)
                                                      7a2392, 31 lines
pair<int, vector<int>> hungarian(const vector<vector<int>> &a)
  if (a.empty()) return {0, {}};
 int n = a.size() + 1, m = a[0].size() + 1;
 vector < int > u(n), v(m), p(m), ans(n-1);
  for (int i = 1; i < n; ++i) {
    p[0] = i;
    int j0 = 0; // add "dummy" worker 0
    vector<int> dist(m, INT_MAX), pre(m, -1);
    vector<bool> done(m + 1);
      done[j0] = true;
      int i0 = p[j0], j1, delta = INT_MAX;
      for(int j = 1; j < m; ++j) if (!done[j]) {
        auto cur = a[i0-1][j-1] - u[i0] - v[j];
        if (cur < dist[j]) dist[j] = cur, pre[j] = j0;</pre>
        if (dist[j] < delta) delta = dist[j], j1 = j;</pre>
      for (int j = 0; j < m; ++j) {
        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;
  for (int j = 1; j < m; ++j) if (p[j]) ans [p[j]-1] = j-1;
 return {-v[0], ans}; // min cost
```

general-matching.h

Description: Maximum Matching for general graphs (undirected and non bipartite) using Edmond's Blossom Algorithm.

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

0b82ee, 68 lines

```
struct blossom_t {
 int t, n; // 1-based indexing!!
 vector<vector<int>> edges;
 vector<int> seen, parent, og, match, aux, Q;
 blossom_t(int _n) : n(_n), edges(n+1), seen(n+1),
 parent (n+1), og (n+1), match (n+1), aux (n+10), t(0) {}
 void addEdge(int u, int v) {
   edges[u].push_back(v);
   edges[v].push_back(u);
 void augment(int u, int v) {
   int pv = v, nv; // flip states of edges on u-v path
     pv = parent[v]; nv = match[pv];
     match[v] = pv; match[pv] = v;
     v = nv;
    } while(u != pv);
```

```
int lca(int v, int w) { // find LCA in O(dist)
   while (1) {
     if (v) {
       if (aux[v] == t) return v; aux[v] = t;
       v = og[parent[match[v]]];
     swap(v, w);
  void blossom(int v, int w, int a) {
    while (og[v] != a) {
     parent[v] = w; w = match[v]; // go other way around cycle
     if(seen[w] == 1) Q.push_back(w), seen[w] = 0;
                              // merge into supernode
     oq[v] = oq[w] = a;
     v = parent[w];
  bool bfs(int u) {
    for (int i = 1; i \le n; ++i) seen[i] = -1, oq[i] = i;
    Q = vector<int>(); Q.push_back(u); seen[u] = 0;
    for(int i = 0; i < 0.size(); ++i) {
     int v = Q[i];
     for(auto &x : edges[v]) {
       if (seen[x] == -1) {
          parent[x] = v; seen[x] = 1;
          if (!match[x]) return augment(u, x), true;
          Q.push_back(match[x]); seen[match[x]] = 0;
       } else if (seen[x] == 0 && oq[v] != oq[x]) {
          int a = lca(og[v], og[x]);
          blossom(x, v, a); blossom(v, x, a);
    return false;
  int solve() {
    int ans = 0; // find random matching (not necessary.
    vector < int > V(n-1); iota(V.beqin(), V.end(), 1); // constant
          improvement)
    shuffle(V.begin(), V.end(), mt19937(0x94949));
    for(auto &x : V) if(!match[x])
     for(auto &v : edges[x]) if (!match[v]) {
       match[x] = y, match[y] = x;
        ++ans; break;
    for (int i = 1; i \le n; ++i)
     if (!match[i] && bfs(i)) ++ans;
    return ans;
};
```

max-independent-set.h

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

7.4 DFS algorithms

centroid-decomposition.h

Description: Divide and Conquer on Trees.

dd21a1, 75 lines

```
template<typename T> struct centroid_t {
  int N;
  vector<vector<int>> adj;
  vector<vector<int>> dist; // dist to all ancestors
  vector<bool> blocked; // processed centroid
  vector<int> sz, depth, parent; // centroid parent
```

```
centroid_t(int _n) : N(_n), adj(_n), dist(32 - __builtin_clz(
     _n), vector<int>(_n)),
blocked(\underline{n}), sz(\underline{n}), depth(\underline{n}), parent(\underline{n}) {}
void add_edge(int a, int b) {
  adj[a].push_back(b);
  adj[b].push_back(a);
void dfs_sz(int cur, int prv) {
  sz[cur] = 1;
  for (int nxt : adj[cur]) {
    if (nxt == prv || blocked[nxt]) continue;
    dfs_sz(nxt, cur);
    sz[cur] += sz[nxt];
int find(int cur, int prv, int tsz) {
  for (int nxt : adj[cur])
    if (!blocked[nxt] && nxt != prv && 2*sz[nxt] > tsz)
      return find(nxt, cur, tsz);
  return cur:
void dfs_dist(int cur, int prv, int layer, int d) {
  dist[layer][cur] = d;
  for (int nxt : adj[cur]) {
    if (blocked[nxt] || nxt == prv) continue;
    dfs_dist(nxt, cur, layer, d + 1);
void get_path(int cur, int prv, int d, vector<int>& cur_path)
  cur_path.push_back(d);
  for (int nxt : adj[cur]) {
    if (nxt == prv || blocked[nxt]) continue;
    get_path(nxt, cur, d + 1, cur_path);
// solve for each subtree (cnt := \# of paths of length K
// that goes through vertex cur)
T solve subtree(int cur, int prv, int K) {
  vector<T> dp(sz[prv] + 1); dp[0] = 1;
  T cnt = 0;
  for (int nxt : adj[cur]) {
    if (blocked[nxt]) continue;
    vector<int> path;
    get_path(nxt, cur, 1, path);
    for (int d : path) {
      if (d > K || K - d > sz[prv]) continue;
      cnt += dp[K - d];
    for (int d : path) dp[d] += 1;
  return cnt;
T decompose(int cur, int K, int layer = 0, int prv_root = -1)
  dfs_sz(cur, -1);
  int root = find(cur, cur, sz[cur]);
  blocked[root] = true;
  depth[root] = layer;
  parent[root] = prv_root;
  dfs_dist(root, root, layer, 0);
  T res = solve_subtree(root, cur, K);
  for (int nxt : adj[root]) {
    if (blocked[nxt]) continue;
    res += decompose(nxt, K, layer + 1, root);
```

```
return res;
};
```

tarjan.h

Description: Finds all strongly connected components in a directed graph. **Usage:** sc.t s(g); $s.solve([\&](const vector < int > \& cc) {...}); visits all components in reverse topological order.$

```
Time: \mathcal{O}\left(E+V\right)
                                                      48ded7, 34 lines
struct scc_t {
    int n, t, scc_num;
    vector<vector<int>> adj;
    vector<int> low, id, stk, in_stk, cc_id;
    scc_t(const vector<vector<int>>& g) : n(int(g.size())), t
         (0), scc_num(0),
    adj(g), low(n,-1), id(n,-1), in_stk(n, false), cc_id(n) {}
    template < class F > void dfs (int cur, F f) {
        id[cur] = low[cur] = t++;
        stk.push_back(cur);
        in_stk[cur] = true;
        for (int nxt : adj[cur]) {
            if (id[nxt] == -1) {
                dfs(nxt, f);
                low[cur] = min(low[cur], low[nxt]);
            } else if (in stk[nxt])
                low[cur] = min(low[cur], id[nxt]);
        if (low[cur] == id[cur]) {
            vector<int> cc; cc.reserve(stk.size());
            while (true) {
                int v = stk.back(); stk.pop_back();
                in_stk[v] = false;
                cc.push back(v);
                cc_id[v] = scc_num;
                if (v == cur) break;
            } f(cc); scc_num ++;
    template < class F > void solve (F f) {
        stk.reserve(n);
        for (int r = 0; r < n; ++r)
            if (id[r] == -1) dfs(r, f);
};
```

bcc.h

Description: Finds all biconnected components in an undirected graph. In a biconnected component there are at least two distinct paths between any two nodes or the component is a bridge. Note that a node can be in several components. *blockcut* constructs the block cut tree of given graph. The first nodes represents the blocks, the others represents the articulation points.

```
nodes represents the blocks, the others represents the articulation points. Usage: int e.id = 0; vector<pair<int, int>> g(N); for (auto [a,b] : edges) { g[a].emplace_back(b, e.id); g[b].emplace_back(a, e.id++); } bcc_t b(g); b.solve([&] (const vector<int>& edges_id) {...}); Time: \mathcal{O}(E+V) 11efd8, 63 lines
```

```
struct bcc_t{
    using pii = pair<int, int>; // v, e_id
    int n, t;
    vector<vector<pii>> adj;
    vector<int> low, id, stk, is_art;
    bcc_t(const vector<vector<pii>> &g) : n(int(g.size())), t
        (0),
        adj(g), low(n,-1), id(n,-1), is_art(n) {}
    template<class F> void dfs(int cur, int e_par, F f){
        id[cur] = low[cur] = t++;
```

```
stk.push_back(e_par);
    int c = 0:
    for (auto [nxt, e_id] : adj[cur]) {
        if (id[nxt] == -1) {
            dfs(nxt, e_id, f);
            low[cur] = min(low[cur], low[nxt]);
            if (low[nxt] >= id[cur]) {
                is_art[cur] = true;
                auto top = find(stk.rbegin(), stk.rend(),
                vector<int> cc(stk.rbegin(), next(top));
                stk.resize(stk.size() - cc.size());
        else if (e_id != e_par) {
            low[cur] = min(low[cur], id[nxt]);
            if (id[nxt] < id[cur]) stk.push_back(e_id);</pre>
    if (e_par == -1) is_art[cur] = (c > 1) ? true : false;
template < class F > void solve (F f) {
    stk.reserve(n);
    for (int r = 0; r < n; ++r)
        if (id[r] == -1) dfs(r, -1, f);
auto blockcut(const vector<pii> &edges) {
   vector<vector<int>> cc;
   vector<int> cc_id(n);
    solve([&](const vector<int> &c) {
        set<int> vc;
        for(int e : c){
            auto [a, b] = edges[e];
            cc_id[a] = cc_id[b] = int(cc.size());
            vc.insert(a);
            vc.insert(b);
        cc.emplace_back(vc.begin(), vc.end());
    for (int a = 0; a < n; a++) if (is art[a]) {
        cc_id[a] = int(cc.size());
        cc.push back({a});
    int bcc_num = int(cc.size());
    vector<vector<int>> tree(bcc num);
    for(int c = 0; c < bcc_num && 1 < int(cc[c].size()); ++</pre>
        for(int a : cc[c]) if(is_art[a]) {
            tree[c].push_back(cc_id[a]);
            tree[cc_id[a]].push_back(c);
    return make_tuple(cc_id, cc, tree);
```

2sat.h

};

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

Usage: TwoSat ts(number of boolean variables); ts.either(0, \sim 3); // Var 0 is true or var 3 is false ts.set_value(2); // Var 2 is true ts.at_most_one($\{0, \sim 1, 2\}$); // <= 1 of vars 0, ~ 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.

```
"tarjan.h"
                                                      10ff6d, 42 lines
struct TwoSat {
    int N:
    vector<vector<int>> gr;
    vector<int> values; // 0 = false, 1 = true
    TwoSat(int n = 0) : N(n), gr(2*n) {}
    int add_var() { // (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 implies(int f, int j) { either(~f, j); }
    void set_value(int x) { either(x, x); }
    void at most one (const vector<int>& li) { // (optional)
        if (int(li.size()) <= 1) return;</pre>
        int cur = \simli[0];
        for (int i = 2; i < int(li.size()); ++i) {</pre>
            int next = add_var();
            either(cur, ~li[i]);
            either(cur, next);
            either(~li[i], next);
            cur = ~next;
        either(cur, ~li[1]);
    bool solve() {
        scc t s(gr);
        s.solve([](const vector<int> &v) { return; } );
        values.assign(N, -1);
        for (int i = 0; i < N; ++i) if (s.cc_id[2*i] == s.cc_id
             [2*i+1]) return 0;
        for (int i = 0; i < N; ++i) {
            if (s.cc id[2*i] < s.cc id[2*i+1]) values[i] =
            else values[i] = true;
        return 1;
```

7.5 Heuristics

maximal-cliques.h

Description: Runs a callback for all maximal cliques in a graph (given as a symmetric bitset matrix; self-edges not allowed). Possible optimization: on the top-most recursion level, ignore 'cands', and go through nodes in order of increasing degree, where degrees go down as nodes are removed.

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

```
57e107, 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 g = (P | X)._Find_first();
 auto cands = P & ~eds[q];
 for(int i = 0; i < eds.size(); ++i) if (cands[i]) {</pre>
   R[i] = 1;
   cliques(eds, f, P & eds[i], X & eds[i], R);
   R[i] = P[i] = 0; X[i] = 1;
```

chromatic-number.h

Description: Compute the chromatic number of a graph. Minimum number of colors needed to paint the graph in a way s.t. if two vertices share an edge, they must have distinct colors.

```
Time: \mathcal{O}\left(N2^N\right)
```

```
template<class T> int min_colors(int N, const T& gr) {
 vector<int> adj(N);
 for (int a = 0; a < N; ++a) {
    for (int b = a + 1; b < N; ++b) {
      if (!gr[a][b]) continue;
      adj[a] = (1 << b);
      adj[b] = (1 << a);
 static vector\langle unsigned \rangle dp(1 << N), buf(1 << N), w(1 << N);
 for (int mask = 0; mask < (1 << N); ++mask) {
   bool ok = true:
    for (int i = 0; i < N; ++i) if (mask & 1 << i) {
      if (adj[i] & mask) ok = false;
    if (ok) dp[mask]++;
   buf[mask] = 1;
    w[mask] = \underline{\quad} builtin_popcount(mask) % 2 == N % 2 ? 1 : -1;
 for (int i = 0; i < N; ++i) {
    for (int mask = 0; mask < (1 << N); ++mask) if (!(mask & 1
      dp[mask^{(1 << i)}] += dp[mask];
 for (int colors = 1; colors <= N; ++colors) {
   unsigned S = 0:
    for (int mask = 0; mask < (1 << N); ++mask)
      S += (buf[mask] *= dp[mask]) * w[mask];
   if (S) return colors:
 assert (false);
```

cycle-counting.cpp

Description: Counts 3 and 4 cycles

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

```
int count_cycles(const vector<vector<int>>& adj, const vector<</pre>
    int>& deg) +
 const int N = int(adj.size());
 vector<int> idx(N), loc(N);
 iota(idx.begin(), idx.end(), 0);
  sort(idx.begin(), idx.end(), [&](const int& a, const int& b)
       { return deg[a] < deg[b]; });
  for (int i = 0; i < N; ++i) loc[idx[i]] = i;
  vector<vector<int>> gr(N);
  for (int a = 0; a < N; ++a) {
    for (int b : adj[a]) {
      if (loc[a] < loc[b]) gr[a].push_back(b);</pre>
  int cycle3 = 0;
    vector<bool> seen(N, false);
    for (int a = 0; a < N; ++a) {
      for (int b : gr[a]) seen[b] = true;
      for (int b : gr[a]) {
        for (int c : gr[b]) {
          if (seen[c]) {
```

edge-coloring lca compress-tree heavylight

edge-coloring.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.)

vector<int> misra_gries(int N, vector<pair<int, int>> eds) {

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

const int M = int(eds.size());

for (int y : {fan[0], u, end})

for (int i = 0; i < M; ++i)

return ret;

```
vector < int > cc(N + 1), ret(M), fan(N), free(N), loc;
for (auto e : eds) ++cc[e.first], ++cc[e.second];
int u, v, ncols = *max element(cc.begin(), cc.end()) + 1;
vector<vector<int>> adj(N, vi(ncols, -1));
for (auto 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& z = free[y] = 0; adj[y][z] != -1; z++);

for (tie(u, v) = eds[i]; adj[u][ret[i]] != v;) ++ret[i];

7.6 Trees

lca.h

Description: Data structure for computing lowest common ancestors and build Euler Tour in a tree. Edges should be an adjacency list of the tree, either directed or undirected.

```
Time: \mathcal{O}\left(N\log N + Q + Q\log\right)
                                                       279920, 19 lines
struct lca t
 int T = 0;
  vector<int> time, path, walk;
  rmq_t<int> rmq;
  lca t(vector<vector<int>> &edges) : time(int(edges.size())),
  rmq((dfs(edges, 0, -1), walk)) {}
  void dfs(vector<vector<int>> &edges, int v, int p) {
    time[v] = T++;
    for(int u : edges[v]) if (u != p) {
      path.push_back(v), walk.push_back(time[v]);
      dfs(edges, u, v);
 int lca(int a, int b) {
    if (a == b) return a;
    tie(a, b) = minmax(time[a], time[b]);
    return path[rmq.query(a, b)];
```

compress-tree.h

};

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

```
Time: \mathcal{O}(|S| \log |S|)
"LCA.h"
                                                      ae0a91, 20 lines
vector<pair<int,int>> compressTree(lca_t &lca, const vector<int
    >& subset) {
 static vector<int> rev; rev.resize(lca.time.size());
 vector<int> li = subset, &T = lca.time;
 auto cmp = [&](int a, int b) { return T[a] < T[b]; };</pre>
 sort(li.begin(), li.end(), cmp);
 int m = li.size()-1;
 for (int i = 0; i < m; ++i) {
   int a = li[i], b = li[i+1];
   li.push_back(lca.lca(a, b));
 sort(li.begin(), li.end(), cmp);
 li.erase(unique(li.begin(), li.end()), li.end());
 for (int i = 0; i < int(li.size()); ++i) rev[li[i]] = i;</pre>
 vector<pair<int, int>> ret = {{0, li[0]}};
 for (int i = 0; i < li.size()-1; ++i) {
   int a = li[i], b = li[i+1];
    ret.emplace_back(rev[lca.lca(a, b)], b);
 return ret;
```

heavylight.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 supports commutative segtree modifications/queries on paths, edges and subtrees. Takes as input the full adjacency list with pairs of (vertex, value). USE_EDGES being true means that values are stored in the edges and are initialized with the adjacency list, otherwise values are stored in the nodes and are initialized to the T defaults value.

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

```
"../data-structures/lazy-segtree.h" 053b35, 70 lines template<br/>bool use_edges> struct HLD_t { int N, T{};
```

```
vector<vector<int>> adj;
vector<int> sz, depth, chain, par, in, out, preorder;
HLD t() {}
HLD_t(const \ vector < vector < int >> & G, int r = 0) : N(int(G.size))
     ())), adj(G),
sz(N), depth(N), chain(N), par(N, -1), in(N), out(N),
    preorder(N) {
  dfs_sz(r); chain[r] = r; dfs_hld(r); }
void dfs_sz(int cur) {
  if (~par[cur]) {
    adj[cur].erase(find(adj[cur].begin(), adj[cur].end(), par
  sz[cur] = 1;
  for (auto& nxt : adj[cur]) {
    par[nxt] = cur; depth[nxt] = 1 + depth[cur];
    dfs_sz(nxt); sz[cur] += sz[nxt];
    if (sz[nxt] > sz[adj[cur][0]]) swap(nxt, adj[cur][0]);
void dfs_hld(int cur) {
  in[cur] = T++;
  preorder[in[cur]] = cur;
  for (auto& nxt : adj[cur]) {
    chain[nxt] = (nxt == adj[cur][0] ? chain[cur] : nxt);
    dfs_hld(nxt);
  out[cur] = T;
int lca(int a, int b) {
  while (chain[a] != chain[b]) {
    if (in[a] < in[b]) swap(a, b);</pre>
    a = par[chain[a]];
  return (in[a] < in[b] ? a : b);</pre>
bool is_ancestor(int a, int b) { return in[a] <= in[b] && in[</pre>
    b] < out[a]; }</pre>
int climb(int a, int k) {
  if (depth[a] < k) return -1;
  int d = depth[a] - k;
  while (depth[chain[a]] > d) a = par[chain[a]];
  return preorder[in[a] - depth[a] + d];
int kth_on_path(int a, int b, int K) {
  int m = lca(a, b);
  int x = depth[a] - depth[m], y = depth[b] - depth[m];
  if (K > x + y) return -1;
  return (x > K ? climb(a, K) : climb(b, x + y - K));
// bool is true if path should be reversed (only for
     noncommutative operations)
const vector<tuple<bool, int, int>>& get path(int a, int b)
  static vector<tuple<bool, int, int>> L, R;
  L.clear(); R.clear();
  while (chain[a] != chain[b]) {
    if (depth[chain[a]] > depth[chain[b]]) {
      L.push_back({true, in[chain[a]], in[a] + 1});
      a = par[chain[a]];
      R.push_back({false, in[chain[b]], in[b] + 1});
      b = par[chain[b]];
  if (depth[a] > depth[b]) {
    L.push_back({true, in[b] + use_edges, in[a] + 1});
    R.push_back({false, in[a] + use_edges, in[b] + 1});
```

};

return L;

L.insert(L.end(), R.rbegin(), R.rend());

tree-isomorphism manhattan-mst directed-mst

- Rebuild the tree after every \sqrt{N} queries.
- Consider vertices with $> \text{or} < \sqrt{N}$ degree separately.
- For subtree updates, note that there are $O(\sqrt{N})$ distinct sizes among child subtrees of any vertex.

Block Tree: Use a DFS to split edges into contiguous groups of size \sqrt{N} to $2\sqrt{N}$.

7.7 Other

manhattan-mst.h

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

Time: O(Nlog N)

```
<dsu.h>
                                                     de8170, 28 lines
typedef Point<int> P;
pair<vector<array<int, 3>>, int> manhattanMST(vector<P> ps) {
 vector<int> id(ps.size());
 iota(id.begin(), id.end(), 0);
 vector<array<int, 3>> edges;
 for (int k = 0; k < 4; ++k) {
   sort(id.begin(), id.end(), [&](int i, int j) {
     return (ps[i]-ps[j]).x < (ps[j]-ps[i]).y;});
    map<int, int> sweep;
    for(auto& i : id) {
      for (auto it = sweep.lower_bound(-ps[i].y);
        it != sweep.end(); sweep.erase(it++)) {
        int j = it->second;
       P d = ps[i] - ps[j];
        if (d.y > d.x) break;
        edges.push_back({d.y + d.x, i, j});
      sweep[-ps[i].y] = i;
    if (k \& 1) for (auto\& p : ps) p.x = -p.x;
    else for(auto& p : ps) swap(p.x, p.y);
 sort(edges.begin(), edges.end());
 UF uf(ps.size());
 int cost = 0:
  for (auto e: edges) if (uf.unite(e[1], e[2])) cost += e[0];
  return {edges, cost};
```

directed-mst.h

Description: Edmonds' algorithm for finding the weight of the 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/dsu-rollback.h"

struct Edge { int a, b; ll w; };

struct Node {
   Edge key;
   Node *l, *r;
   ll delta;
   void prop() {
      key.w += delta;
      if (l) l->delta += delta;
      if (r) r->delta += delta;
      delta = 0;
   }
   Edge top() { prop(); return key; }
};
Node *merge(Node *a, Node *b) {
```

```
tree-isomorphism.h Time: \mathcal{O}(N \log(N))
```

92e59f, 51 lines

```
map<vector<int>, int> delta;
struct tree_t {
 int n;
  pair<int, int> centroid;
  vector<vector<int>> edges;
  vector<int> sz;
  tree_t (vector<vector<int>>& graph) :
   edges(graph), sz(edges.size()) {}
  int dfs_sz(int v, int p) {
   sz[v] = 1;
   for (int u : edges[v]) {
     if (u == p) continue;
     sz[v] += dfs_sz(u, v);
    return sz[v];
  int dfs(int tsz, int v, int p) {
    for (int u : edges[v]) {
     if (u == p) continue;
     if (2*sz[u] <= tsz) continue;
     return dfs(tsz, u, v);
    return centroid.first = v;
  pair<int,int> find centroid(int v) {
    int tsz = dfs sz(v, -1);
    centroid.second = dfs(tsz, v, -1);
    for (int u : edges[centroid.first]) {
     if (2*sz[u] == tsz)
       centroid.second = u;
    return centroid;
```

7.6.1 Sqrt Decomposition

HLD generally suffice; If not, here are some common strategies int u : edges[v]) {
 if (u == p) continue;
 offset.push_back(hash_it(u, v));

```
if (!a || !b) return a ?: b;
 a->prop(), b->prop();
 if (a->key.w > b->key.w) swap(a, b);
 swap(a->1, (a->r = merge(b, a->r)));
 return a:
void pop(Node * a) { a->prop(); a = merge(a->1, a->r); }
pair<11, vector<int>> dmst(int n, int r, vector<Edge>& g) {
 RollbackUF uf(n);
 vector<Node*> heap(n);
 for (Edge e : g) heap[e.b] = merge(heap[e.b], new Node(e));
 11 \text{ res} = 0;
 vector<int> 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;
  for (int s = 0; s < n; ++s) {
   int u = s, qi = 0, w;
    while (seen[u] < 0) {
     if (!heap[u]) return {-1,{}};
     Edge e = heap[u]->top();
     heap[u]->delta -= e.w, pop(heap[u]);
     Q[qi] = e, path[qi++] = u, seen[u] = s;
      res += e.w, u = uf.find(e.a);
     if (seen[u] == s) {
       Node \star cyc = 0;
       int end = qi, time = uf.time();
       do cyc = merge(cyc, heap[w = path[--qi]]);
       while (uf.unite(u, w));
       u = uf.find(u), heap[u] = cyc, seen[u] = -1;
        cycs.push_front({u, time, {&Q[qi], &Q[end]}});
    for (int i = 0; i < qi; ++i) 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;
 for (int i = 0; i < n; ++i) par[i] = in[i].a;
```

return {res, par}; 7.8 Theorems

7.8.1 Euler's theorem

Let V, A and F be the number of vertices, edges and faces of connected planar graph, V - A + F = 2

7.8.2 Eulerian Cycles

The number of Eulerian cycles in a directed graph G is: $t_w(G) \prod_{v \in G} (\deg v - 1)!$, where $t_w(G)$ is the number of arborescences ("directed spanning" tree) rooted at w (Check Number of Spanning Trees)

7.8.3 Dilworth's theorem

For any partially ordered set, the sizes of the max antichain and of the min chain decomposition are equal. Equivalent to Konig's theorem on the bipartite graph (U,V,E) where U=V=S and (u,v) is an edge when u < v. Those vertices outside the min vertex cover in both U and V form a max antichain

7.8.4 König-Egervary theorem

For Bipartite Graphs, the number of edges in the maximum matching is greater than or equal the number of vertices in the minimum cover

Maximum Weight Closure

Given a vertex-weighted directed graph G. Turn the graph into a flow network, adding weight ∞ to each edge. Add vertices S, T. For each vertex v of weight w, add edge (S, v, w) if w > 0, or edge (v, T, -w) if w < 0. Sum of positive weights minus minimum S-T cut is the answer. Vertices reachable from S are in the closure. The maximum-weight closure is the same as the complement of the minimum-weight closure on the graph with edges reversed.

7.8.5 Maximum Weighted Independent Set in a Bipartite Graph

This is the same as the minimum weighted vertex cover. Solve this by constructing a flow network with edges (S, u, w(u)) for $u \in L$, (v, T, w(v)) for $v \in R$ and (u, v, ∞) for $(u, v) \in E$. The minimum S, T-cut is the answer. Vertices adjacent to a cut edge are in the vertex cover. Num

7.8.6 Number of Spanning Trees

Define Laplacian Matrix as L = D - A, D being a Diagonal Matrix with $D_{i,i} = deq(i)$ and A an Adjacency Matrix. Create an $N \times N$ Laplacian matrix mat, and for each edge $a \to b \in G$, do mat[a][b]--, mat[b][b]++ (and mat[b][a]--, mat [a] [a] ++ if G is undirected). Remove the ith row and column and take the determinant; this yields the number of directed spanning trees rooted at i (if G is undirected, remove any row/column).

7.8.7 Tutte Matrix

- A graph has a perfect matching iff the *Tutte* matrix has a non-zero determinant.
- The rank of the *Tutte* matrix is equal to twice the size of the maximum matching. The maximum cost matching can be found by polynomial interpolation.

7.8.8 Menger's theorem

- Vertices: A graph is k-connected iff all pairwise vertices are connected to at least k internally disjoint paths.
- Edges: A graph is called k-edge-connected if the removal of at least k edges of the graph keeps it connected. A graph is k-edge-connected iff for all pairwise vertices u and v, exist kpaths which link u to v without sharing an edge.

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 \text{class T} \rangle int \text{sgn}(\text{T}|\mathbf{x}) { return (\mathbf{x} > 0) - (\mathbf{x} < 0); }
template<class T>
struct Point {
 typedef Point P;
 T x, y;
 explicit Point (T x=0, T y=0) : x(x), y(y) {}
 bool operator (P p) const \{ return tie(x,y) < tie(p.x,p.y) \}
 bool operator==(P p) const { return tie(x,y)==tie(p.x,p.y); }
 P operator+(P p) const { return P(x+p.x, y+p.y); }
 P operator-(P p) const { return P(x-p.x, y-p.y); }
 P operator*(T d) const { return P(x*d, y*d); }
 P operator/(T d) const { return P(x/d, y/d); }
 T dot(P p) const { return x*p.x + y*p.y; }
 T cross(P 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)); }
using P = Point < double >;
```

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.

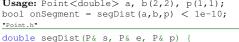


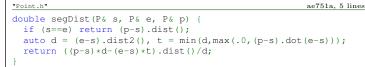
```
template<class P>
double lineDist(const P& a, const P& b, const P& p) {
 return (double) (b-a).cross(p-a)/(b-a).dist();
```

SegmentDistance.h

Description:

Returns the shortest distance between point p and the line segment from point s to e. **Usage:** Point < double > a, b(2,2), p(1,1);





SegmentIntersection.h

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;</pre>
"Point.h", "OnSegment.h"
template<class P> vector<P> segInter(P a, P b, P c, P d) {
 auto oa = c.cross(d, a), ob = c.cross(d, b),
     oc = a.cross(b, c), od = a.cross(b, d);
  // Checks if intersection is single non-endpoint point.
 if (sgn(oa) * sgn(ob) < 0 \&\& sgn(oc) * sgn(od) < 0)
    return { (a * ob - b * oa) / (ob - oa) };
 if (onSegment(c, d, a)) s.insert(a);
 if (onSegment(c, d, b)) s.insert(b);
 if (onSegment(a, b, c)) s.insert(c);
 if (onSegment(a, b, d)) s.insert(d);
 return {s.begin(), s.end()};
```

SegmentIntersectionQ.h

Description: Like segmentIntersection, but only returns true/false. Products of three coordinates are used in intermediate steps so watch out for overflow if using int or long long.

```
"Point.h"
                                                     1ff4ba, 16 lines
template<class P>
bool segmentIntersectionQ(P s1, P e1, P s2, P e2) {
 if (e1 == s1) {
    if (e2 == s2) return e1 == e2;
    swap(s1,s2); swap(e1,e2);
 P v1 = e1-s1, v2 = e2-s2, d = s2-s1;
  auto a = v1.cross(v2), a1 = d.cross(v1), a2 = d.cross(v2);
 if (a == 0) { // parallel
    auto b1 = s1.dot(v1), c1 = e1.dot(v1),
       b2 = s2.dot(v1), c2 = e2.dot(v1);
    return !a1 && max(b1,min(b2,c2)) <= min(c1,max(b2,c2));
 if (a < 0) { a = -a; a1 = -a1; a2 = -a2; }
  return (0 <= a1 && a1 <= a && 0 <= a2 && a2 <= a);
```

LineIntersection.h

Description:

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

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



```
if (res.first == 1)
cout << "intersection point at " << res.second << endl;</pre>
template<class P>
pair<int, P> lineInter(P s1, P e1, P s2, P e2) {
  auto d = (e1 - s1).cross(e2 - s2);
 if (d == 0) // if parallel
```

```
auto p = s2.cross(e1, e2), q = s2.cross(e2, s1);
return {1, (s1 * p + e1 * q) / d};
}
```

LineProjectionReflection.h

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

"Point.h"

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

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;
```

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 Description:

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



```
typedef Point<double> P;
P linearTransformation(const P& p0, const P& p1,
  const P& q0, const P& q1, const P& r) {
  P dp = p1-p0, dq = q1-q0, num(dp.cross(dq), dp.dot(dq));
  return q0 + P((r-p0).cross(num), (r-p0).dot(num))/dp.dist2();
}
```

Angle.h

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

```
Usage: vector<Angle> v = \{w[0], w[0].t360() ...\}; // sorted int j = 0; rep(i,0,n) \{ while (v[j] < v[i].t180()) ++j; \} // sweeps j such that (j-i) represents the number of positively // oriented triangles with vertices at 0 and i Of0602, 34 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) {
 // add a.dist2() and b.dist2() to also compare distances
  return make_tuple(a.t, a.half(), a.y * (11)b.x) <</pre>
    make_tuple(b.t, b.half(), a.x * (11)b.y);
// Given two points, this calculates the smallest angle between
// them, i.e., the angle that covers the defined line segment.
pair<Angle, Angle> segmentAngles(Angle a, Angle b) {
 if (b < a) swap(a, b);
 return (b < a.t180() ?
    make_pair(a, b) : make_pair(b, a.t360()));
Angle operator+(Angle a, Angle b) { // point a + vector b
 Angle r(a.x + b.x, a.y + b.y, a.t);
 if (a.t180() < r) r.t--;
 return r.t180() < a ? r.t360() : r;
Angle angleDiff(Angle a, Angle b) { // angle \ b - angle \ a}
 int tu = b.t - a.t; a.t = b.t;
 return \{a.x*b.x + a.y*b.y, a.x*b.y - a.y*b.x, tu - (b < a)\};
```

AngleCmp.h

Description: Useful utilities for dealing with angles of rays from origin. OK for integers, only uses cross product. Doesn't support (0,0).

```
template <class P>
bool sameDir(P s, P t) {
 return s.cross(t) == 0 \&\& s.dot(t) > 0;
// checks 180 \le s...t < 360?
template <class P>
bool isReflex(P s, P t) {
 auto c = s.cross(t);
 return c ? (c < 0) : (s.dot(t) < 0);
// operator < (s,t) for angles in [base,base+2pi)
template <class P>
bool angleCmp(P base, P s, P t) {
 int r = isReflex(base, s) - isReflex(base, t);
 return r ? (r < 0) : (0 < s.cross(t));
// is x in [s,t] taken ccw? 1/0/-1 for in/border/out
template <class P>
int angleBetween (P s, P t, P x) {
 if (sameDir(x, s) || sameDir(x, t)) return 0;
 return angleCmp(s, x, t) ? 1 : -1;
```

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" c64785, 10 lines
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.

```
rPoint.h" b0153d, 13 lines
template<class P>
vector<pair<P, P>> tangents(P c1, double r1, P c2, double r2) {
   P d = c2 - c1;
   double dr = r1 - r2, d2 = d.dist2(), h2 = d2 - dr * dr;
   if (d2 == 0 || h2 < 0) return {};
   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;
}
```

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.



MinimumEnclosingCircle.h

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

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

CircleLine.h

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

```
"Point.h", "lineDistance.h", "LineProjectionReflection.h"
                                                         debf86, 8 lines
template<class P>
vector<P> circleLine(P c, double r, P a, P b) {
 double h2 = r*r - a.cross(b,c)*a.cross(b,c)/(b-a).dist2();
 if (h2 < 0) return {};
  P p = lineProj(a, b, c), h = (b-a).unit() * sqrt(h2);
 if (h2 == 0) return {p};
  return \{p - h, p + h\};
```

CircleCircleArea.h

Description: Calculates the area of the intersection of 2 circles 8bf2b6, 12 lines

```
template<class P>
double circleCircleArea(P c, double cr, P d, double dr) {
 if (cr < dr) swap(c, d), swap(cr, dr);
  auto A = [&] (double r, double h) {
   return r*r*acos(h/r)-h*sqrt(r*r-h*h);
  };
  auto 1 = (c - d) \cdot dist(), a = (1*1 + cr*cr - dr*dr)/(2*1);
 if (1 - cr - dr >= 0) return 0; // far away
 if (1 - cr + dr <= 0) return M PI*dr*dr;
 if (1 - cr >= 0) return A(cr, a) + A(dr, 1-a);
 else return A(cr, a) + M_PI*dr*dr - A(dr, a-1);
```

CirclePolygonIntersection.h

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

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

```
"Point.h"
                                                     cf9deb, 18 lines
#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;
    auto s = max(0., -a-sqrt(det)), t = min(1., -a+sqrt(det));
   if (t < 0 \mid | 1 \le s) return arg(p, q) * r2;
   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;
  for (int i = 0; i < ps.size(); ++i)
   sum += tri(ps[i] - c, ps[(i + 1) % ps.size()] - c);
  return sum;
```

8.3 Polygons

InsidePolygon.h

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

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

"Point.h", "OnSegment.h", "SegmentDistance.h" f9442d, 12 lines

```
template<class P>
bool inPolygon(vector<P> &p, P a, bool strict = true) {
 int cnt = 0, n = p.size();
 for (int i = 0; i < n; ++i) {
   P q = p[(i + 1) % n];
    if (onSegment(p[i], q, a)) return !strict; // change to
        // -1 if u need to detect points in the boundary
    //or: if (segDist(p[i], q, a) \le eps) return ! strict;
    cnt ^= ((a.y<p[i].y) - (a.y<q.y)) * a.cross(p[i], q) > 0;
 return cnt;
```

Polygon Area, h

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

```
"Point.h"
                                                     3794ee, 17 lines
template<class T>
T polygonArea(vector<Point<T>> &v) {
 T a = v.back().cross(v[0]);
 for(int i = 0; i < v.size()-1; ++i)
   a += v[i].cross(v[i+1]);
  return abs(a)/2.0;
Point<T> polygonCentroid(vector<Point<T>> &v) { // not tested
 Point<T> cent(0,0); T area = 0;
 for(int i = 0; i < v.size(); ++i) {</pre>
    int j = (i+1) % (v.size()); T a = cross(v[i], v[j]);
    cent += a * (v[i] + v[j]);
    area += a;
 return cent/area/(T)3;
```

PolygonCenter.h

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

```
"Point.h"
                                                       26a00f, 8 lines
P polygonCenter(const vector<P>& v) {
 P res(0, 0); double A = 0;
 for (int i = 0, j = v.size() - 1; i < v.size(); j = ++i) {
    res = res + (v[i] + v[j]) * v[j].cross(v[i]);
   A += v[j].cross(v[i]);
 return res / A / 3;
```

PolygonCut.h Description:

thing to the left of the line going from s to e cut away.

```
Returns a vector with the vertices of a polygon with every-
Usage: vector<P> p = ...;
p = polygonCut(p, P(0,0), P(1,0));
```

```
"Point.h", "lineIntersection.h"
                                                         7df36f, 11 lines
vector<P> polygonCut(const vector<P>& poly, P s, P e) {
 vector<P> res;
 for(int i = 0; i < poly.size(); ++i) {</pre>
   P cur = poly[i], prev = i ? poly[i-1] : poly.back();
    bool side = s.cross(e, cur) < 0;</pre>
    if (side != (s.cross(e, prev) < 0))</pre>
      res.push_back(lineInter(s, e, cur, prev).second);
    if (side) res.push_back(cur);
 return res:
```

ConvexHull.h

Description:

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



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

"Point.h"

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

HullDiameter.h

Description: Returns the two points with max distance on a convex hull (ccw, no duplicate/colinear points). 0e0c1f, 11 lines

```
array<P, 2> hullDiameter(vector<P> S) {
 int n = S.size(), j = n < 2 ? 0 : 1;
 pair<lint, array<P, 2>> res({0, {S[0], S[0]}});
 for (int i = 0; i < j; ++i)
   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 colinear points). Returns true if point lies within the hull. If strict is true, points on the boundary aren't included.

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

```
"Point.h", "sideOf.h", "OnSegment.h"
                                                        7b8514, 12 lines
bool inHull(const vector<P> &1, P p, bool strict = true) {
 int a = 1, b = 1.size() - 1, r = !strict;
 if (1.size() < 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 \mid \mid sideOf(1[0], 1[b], p) <= -r)
    return false;
 while (abs(a - b) > 1) {
    int c = (a + b) / 2;
    (sideOf(1[0], 1[c], p) > 0 ? b : a) = c;
 return sgn(l[a].cross(l[b], p)) < r;</pre>
```

PolvUnion.h

Description: Calculates the area of the union of n polygons (not necessarily convex). The points within each polygon must be given in CCW order. Guaranteed to be precise for integer coordinates up to 3e7. If epsilons are needed, add them in sideOf as well as the definition of sgn.

Time: $\mathcal{O}(N^2)$, where N is the total number of points

```
"Point.h", "sideOf.h"
double rat(P a, P b) { return sqn(b.x) ? a.x/b.x : a.y/b.y; }
double polyUnion(vector<vector<P>>& poly) {
 double ret = 0;
 for(int i = 0; i < poly.size(); ++i)
    for(int v = 0; v < poly[i].size(); ++v) {
```

LineHullIntersection HalfPlane ClosestPair KdTree

```
P A = poly[i][v], B = poly[i][(v + 1) % poly[i].size()];
   vector<pair<double, int>> segs = {{0, 0}, {1, 0}};
   for(int j = 0; j < poly.size(); ++j) if (i != j) {
     for(int u = 0; u < poly[j]; ++u) {
       P C = poly[j][u], D = poly[j][(u + 1) % poly[j].size
       int sc = sideOf(A, B, C), sd = sideOf(A, B, D);
       if (sc != sd) {
         double sa = C.cross(D, A), sb = C.cross(D, B);
         if (min(sc, sd) < 0)
           segs.emplace_back(sa / (sa - sb), sqn(sc - sd));
       } else if (!sc && !sd && j<i && sgn((B-A).dot(D-C))</pre>
         segs.emplace_back(rat(C - A, B - A), 1);
         segs.emplace_back(rat(D - A, B - A), -1);
     }
 sort(segs.begin(), segs.end());
 for(auto& s : segs) s.first = min(max(s.first, 0.0), 1.0);
 double sum = 0;
 int cnt = segs[0].second;
 for(int j = 1; j < segs.size(); ++j) {
   if (!cnt) sum += segs[j].first - segs[j - 1].first;
   cnt += segs[j].second;
 ret += A.cross(B) * sum;
return ret / 2;
```

LineHullIntersection.h

Description: Line-convex polygon intersection. The polygon must be ccw and have no colinear 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: $\mathcal{O}(N + Q \log n)$

```
"Point.h"
                                                      65ebb6, 39 lines
typedef array<P, 2> Line;
#define cmp(i,j) sgn(dir.perp().cross(poly[(i)%n]-poly[(j)%n]))
#define extr(i) cmp(i + 1, i) >= 0 && cmp(i, i - 1 + n) < 0
int extrVertex(vector<P>& poly, P dir) {
  int n = poly.size(), left = 0, right = n;
  if (extr(0)) return 0;
  while (left + 1 < right) {</pre>
   int m = (left + right) / 2;
    if (extr(m)) return m;
   int ls = cmp(left + 1, left), ms = cmp(m + 1, m);
    (ls < ms \mid \mid (ls == ms && ls == cmp(left, m)) ? right : left
        ) = m;
  return left;
#define cmpL(i) sgn(line[0].cross(poly[i], line[1]))
array<int, 2> lineHull(Line line, vector<P>& poly) {
  int endA = extrVertex(poly, (line[0] - line[1]).perp());
  int endB = extrVertex(poly, (line[1] - line[0]).perp());
  if (cmpL(endA) < 0 \mid \mid cmpL(endB) > 0)
   return {-1, -1};
  array<int, 2> res;
  for(int i = 0; i < 2; ++i) {
   int left = endB, right = endA, n = poly.size();
    while ((left + 1) % n != right) {
```

```
int m = ((left + right + (left < right ? 0 : n)) / 2) % n</pre>
    (cmpL(m) == cmpL(endB) ? left : right) = m;
  res[i] = (left + !cmpL(right)) % 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) % poly.size()) {
    case 0: return {res[0], res[0]};
    case 2: return {res[1], res[1]};
return res;
```

HalfPlane.h

Description: Halfplane intersection area

"Point.h", "lineIntersection.h"

```
#define eps 1e-8
typedef Point < double > P;
struct Line {
 P P1, P2;
 // Right hand side of the ray P1 -> P2
 explicit Line (P \ a = P(), P \ b = P()) : P1(a), P2(b) {};
 P intpo(Line y) {
   pair<int, P> r = lineInter(P1, P2, y.P1, y.P2);
   assert (r.first == 1);
   return r.second:
 P dir() { return P2 - P1; }
 bool contains (P x) {
   return (P2 - P1).cross(x - P1) < eps;
 bool out(P x) { return !contains(x); }
template<class T>
bool mycmp(Point<T> a, Point<T> b) {
 // return atan2(a.y, a.x) < atan2(b.y, b.x);
 if (a.x * b.x < 0) return a.x < 0;
 if (abs(a.x) < eps) {
   if (abs(b.x) < eps) return a.y > 0 && b.y < 0;
   if (b.x < 0) return a.y > 0;
   if (b.x > 0) return true;
 if (abs(b.x) < eps) {
   if (a.x < 0) return b.y < 0;
   if (a.x > 0) return false;
 return a.cross(b) > 0;
bool cmp(Line a, Line b) { return mycmp(a.dir(), b.dir()); }
double Intersection_Area(vector <Line> b) {
 sort(b.begin(), b.end(), cmp);
 int n = b.size();
 int q = 1, h = 0, i;
 vector<Line> c(b.size() + 10);
 for (i = 0; i < n; i++) {
    while (q < h \&\& b[i].out(c[h].intpo(c[h - 1]))) h--;
   while (q < h \&\& b[i].out(c[q].intpo(c[q + 1]))) q++;
   c[++h] = b[i];
   if (q < h \&\& abs(c[h].dir().cross(c[h-1].dir())) < eps) {
     if (c[h].dir().dot(c[h-1].dir()) > 0) {
```

if (b[i].out(c[h].P1)) c[h] = b[i];

```
}else {
      // The area is either 0 or infinite.
      // If you have a bounding box, then the area is
           definitely 0.
      return 0;
while (q < h-1 \&\& c[q].out(c[h].intpo(c[h-1]))) h--;
while (q < h-1 \&\& c[h].out(c[q].intpo(c[q + 1]))) q++;
// Intersection is empty. This is sometimes different from
     the case when
// the intersection area is 0.
if (h - q <= 1) return 0;
c[h + 1] = c[q];
vector<P> s;
for (i = q; i \le h; i++) s.push_back(c[i].intpo(c[i + 1]));
s.push_back(s[0]);
double ans = 0;
for (i = 0; i < (int) s.size()-1; i++) ans += s[i].cross(s[i])
    + 1]);
return ans / 2;
```

8.4 Misc. Point Set Problems

ClosestPair.h

c0a94b, 70 lines

Description: Finds the closest pair of points.

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

```
"Point.h"
                                                     32b14f, 16 lines
pair<P, P> closest (vector<P> v) {
  assert(v.size() > 1);
  sort(v.begin(), v.end(), [](Pa, Pb) { return a.y < b.y; });
  pair<int64_t, pair<P, P>> ret{LLONG_MAX, {P(), P()}};
  int j = 0;
  for(P &p : v) {
    P d{1 + (int64_t)sqrt(ret.first), 0};
    while (v[j].y \le p.y - d.x) S.erase(v[j++]);
    auto lo = S.lower_bound(p - d), hi = S.upper_bound(p + d);
    for (; lo != hi; ++lo)
      ret = min(ret, {(*lo - p).dist2(), {*lo, p}});
    S.insert(p);
  return ret.second;
```

KdTree.h

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

```
"Point.h"
                                                      915562, 63 lines
typedef long long T;
typedef Point<T> P;
const T INF = numeric_limits<T>::max();
bool on_x(const P& a, const P& b) { return a.x < b.x; }</pre>
bool on_y(const P& a, const P& b) { return a.y < b.y; }</pre>
  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]) {
```

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DelaunayTriangulation FastDelaunay RectangleUnionArea

```
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 the box is wider than high (not best
           heuristic...)
      sort(vp.begin(), vp.end(), 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 = vp.size()/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({vp.begin(), 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);
};
```

DelaunayTriangulation.h

Description: Computes the Delaunay triangulation of a set of points. Each circumcircle contains none of the input points. If any three points are colinear or any four are on the same circle, behavior is undefined.

```
Time: \mathcal{O}\left(n^2\right)
"Point.h", "3dHull.h" f6175a, 10 lines template < class P, class F> void delaunay (vector < P>& ps, F trifun) {
    if (ps.size() == 3) { int d = (ps[0].cross(ps[1], ps[2]) < 0) }
    ;
    trifun(0,1+d,2-d); }
    vector < P>> ps; ps. ps. emplace_back(p.x, p.y, p.dist2()); if (ps.size() > 3) for (auto &t: hull3d(p3)) if ((p3[t.b]-p3[t.a]).
        cross(p3[t.c]-p3[t.a]).dot(P3(0,0,1)) < 0)  
    trifun(t.a, t.c, t.b); }
```

FastDelaunav.h

Description: Fast Delaunay triangulation. There must be no duplicate points. If all points are on a line, no triangles will be returned. Should work for doubles as well, though there may be precision issues in 'circ'. Returns triangles in order $\{t[0][0], t[0][1], t[0][2], t[1][0], \ldots\}$, all counter-clockwise.

```
Time: \mathcal{O}(n \log n)
"Point.h"
                                                       a1f392, 89 lines
typedef Point<ll> P;
typedef struct Quad* Q;
typedef __int128_t lll; // (can be ll if coords are < 2e4)
P arb(LLONG_MAX, LLONG_MAX); // not equal to any other point
struct Quad {
 bool mark; Q o, rot; P p;
 P F() { return r()->p; }
 Q r() { return rot->rot; }
 Q prev() { return rot->o->rot; }
 Q next() { return rot->r()->o->rot; }
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 = \text{new Quad}\{0, 0, 0, \text{orig}\}, q1 = \text{new Quad}\{0, 0, 0, \text{arb}\},
    q2 = new Quad\{0, 0, 0, dest\}, q3 = new Quad\{0, 0, 0, arb\};
  q0 -> 0 = q0; q2 -> 0 = q2; // 0 - 0, 2 - 2
  q1->0 = q3; q3->0 = q1; // 1-3, 3-1
  q0 - rot = q1; q1 - rot = q2;
 q^2 - rot = q^3; q^3 - rot = q^0;
 return q0;
void splice(Q a, Q b) {
  swap(a->o->rot->o, b->o->rot->o); swap(a->o, b->o);
Q connect(Q a, Q b) {
 Q = makeEdge(a->F(), b->p);
  splice(q, a->next());
  splice(q->r(), b);
  return q;
pair<0,0> rec(const vector<P>& s) {
 if (sz(s) \le 3)  {
    0 = \text{makeEdge}(s[0], s[1]), b = \text{makeEdge}(s[1], s.back());
    if (sz(s) == 2) return { a, a->r() };
    splice(a->r(), b);
    auto side = s[0].cross(s[1], s[2]);
    Q c = side ? connect(b, a) : 0;
    return {side < 0 ? c->r() : a, side < 0 ? c : b->r() };
#define H(e) e->F(), e->p
#define valid(e) (e->F().cross(H(base)) > 0)
 O A, B, ra, rb;
 int half = (sz(s) + 1) / 2;
  tie(ra, A) = rec({s.begin(), s.begin() + half});
  tie(B, rb) = rec({s.begin() + half, s.end()});
  while ((B->p.cross(H(A)) < 0 \&& (A = A->next()))
    (A->p.cross(H(B)) > 0 && (B = B->r()->o)));
  O base = connect(B->r(), A);
  if (A->p == ra->p) ra = base->r();
  if (B->p == rb->p) rb = base;
#define DEL(e, init, dir) Q e = init->dir; if (valid(e)) \
 while (circ(e->dir->F(), H(base), e->F())) {
    Q t = e->dir; \
    splice(e, e->prev()); \
```

```
splice(e->r(), e->r()->prev()); \
   e = t; \
 for (;;) {
   DEL(LC, base->r(), o); DEL(RC, base, prev());
   if (!valid(LC) && !valid(RC)) break;
   if (!valid(LC) || (valid(RC) && circ(H(RC), H(LC))))
     base = connect(RC, base->r());
     base = connect(base->r(), LC->r());
 return { ra, rb };
vector<P> triangulate(vector<P> pts) {
 sort(pts.begin(), pts.end()); assert(unique(pts.begin(), pts
      .end()) == pts.end());
 if (pts.size() < 2) return {};</pre>
 Q e = rec(pts).first;
 vector<Q> q = {e};
 int qi = 0;
 while (e->o->F().cross(e->F(), e->p) < 0) e = e->o;
#define ADD { Q c = e; do { c->mark = 1; pts.push_back(c->p); \
 q.push_back(c->r()); c = c->next(); } while (c != e); }
 ADD; pts.clear();
 while (qi < sz(q)) if (!(e = q[qi++]) -> mark) ADD;
 return pts;
```

RectangleUnionArea.h

Description: Sweep line algorithm that calculates area of union of rectangles in the form $[x1,x2)\times[y1,y2)$

```
gles in the form [zt, z2) x [yt, y2)
Usage: vector<pair<int,int>, pair<int,int>> rectangles;
rectangles.push.back({{x1, x2}, {y1, y2}});
lint result = area(rectangles);

struct seg_node{
  int val, cnt, lz;
  seg_node(int n = INT_MAX, int c = 0): val(n), cnt(c), lz(0)
```

```
void push(seq_node& 1, seq_node& r) {
    if(lz){
      1.add(lz);
      r.add(lz);
      1z = 0;
  void merge(const seg_node& 1, const seg_node& r) {
    if(1.val < r.val) val = 1.val, cnt = 1.cnt;
    else if(1.val > r.val) val = r.val, cnt = r.cnt;
    else val = 1.val, cnt = 1.cnt + r.cnt;
  void add(int n){
    val += n;
    1z += n;
  int get_sum() { return (val ? 0 : cnt); }
// x1 y1 x2 y2
lint solve(const vector<array<int, 4>>&v){
  vector<int>vs;
  for(auto& [a, b, c, d] : v) {
    ys.push_back(b);
    ys.push_back(d);
  sort(ys.begin(), ys.end());
  ys.erase(unique(ys.begin(), ys.end()), ys.end());
  vector<array<int, 4>>e;
  for(auto [a, b, c, d] : v) {
```

```
b = int(lower_bound(ys.begin(), ys.end(), b) - ys.begin());
 d = int(lower_bound(ys.begin(), ys.end(), d) - ys.begin());
 e.push back({a, b, d, 1});
 e.push_back({c, b, d, -1});
sort(e.begin(), e.end());
int m = (int)ys.size();
segtree_range<seg_node>seg(m-1);
for(int i=0;i<m-1;i++) seg.at(i) = seg_node(0, ys[i+1] - ys[i
    ]);
seq.build();
int last = INT_MIN, total = ys[m-1] - ys[0];
lint ans = 0;
for(auto [x, y1, y2, c] : e) {
 ans += (lint) (total - seg.query(0, m-1).get_sum()) * (x -
     last);
 last = x;
 seq.update(y1, y2, &seq_node::add, c);
return ans;
```

$8.5 \quad 3D$

PolyhedronVolume.h

Description: Magic formula for the volume of a polyhedron. Faces should point outwards. 832599, 6 lines

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

Point3D.h

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

```
8058ae, 32 lines
template<class T> struct 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 {
   return tie(x, y, z) < tie(p.x, p.y, p.z); }
  bool operator == (R p) const {
   return tie(x, y, z) == tie(p.x, p.y, p.z); }
  P operator+(R p) const { return P(x+p.x, y+p.y, z+p.z); }
  P operator-(R p) const { return P(x-p.x, y-p.y, z-p.z); }
  P operator*(T d) const { return P(x*d, y*d, z*d); }
  P operator/(T d) const { return P(x/d, y/d, z/d); }
  T dot(R p) const { return x*p.x + y*p.y + z*p.z; }
  P cross(R p) const {
    return P(y*p.z - z*p.y, z*p.x - x*p.z, x*p.y - y*p.x);
  T dist2() const { return x*x + y*y + z*z; }
  double dist() const { return sqrt((double)dist2()); }
  //Azimuthal angle (longitude) to x-axis in interval [-pi, pi]
  double phi() const { return atan2(v, x); }
  //Zenith angle (latitude) to the z-axis in interval [0, pi]
  double theta() const { return atan2(sqrt(x*x+y*y),z); }
  P unit() const { return *this/(T)dist(); } //makes dist()=1
  //returns unit vector normal to *this and p
  P normal(P p) const { return cross(p).unit(); }
  //returns point rotated 'angle' radians ccw around axis
  P rotate(double angle, P axis) const {
   double s = sin(angle), c = cos(angle); P u = axis.unit();
    return u*dot(u)*(1-c) + (*this)*c - cross(u)*s;
```

};

3dHull.h

Description: Computes all faces of the 3-dimension hull of a point set. *No four points must be coplanar*, or else random results will be returned. All faces will point outwards.

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

```
"Point3D.h"
                                                      3ed613, 48 lines
typedef Point3D<double> P3;
struct PR {
 void ins(int x) { (a == -1 ? a : b) = x; }
 void rem(int x) { (a == x ? a : b) = -1; }
 int cnt() { return (a != -1) + (b != -1); }
 int a, b;
struct F { P3 q; int a, b, c; };
vector<F> hull3d(const vector<P3>& A) {
 assert(A.size() >= 4);
 vector<vector<PR>> E(A.size(), vector<PR>(A.size(), {-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);
  };
 for (int i=0; i<4; i++) for (int j=i+1; j<4; j++) for (k=j+1; k<4; k
   mf(i, j, k, 6 - i - j - k);
  for(int i=4; i<A.size();++i) {</pre>
    for(int j=0; j<FS.size(); ++j) {</pre>
     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 = FS.size();
    for(int j=0; j<nw; j++) {</pre>
     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 (auto &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. All angles measured in radians. The algorithm starts by converting the spherical coordinates to cartesian coordinates so if that is what you have you can use only the two last rows. dx*radius is then the difference between the two points in the x direction and d*radius is the total distance between the points.

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

$\underline{\text{Strings}}$ (9)

kmp.h

Description: failure[x] computes the length of the longest prefix of s that ends at x, other than s[0...x] itself (abacaba -> -1,0,0,1,0,1,2,3). Can be used to find all occurrences of a pattern in a text.

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

```
vector<int> prefix_function(const string& S) {
  vector<int> fail = {-1}; fail.reserve(S.size());
  for (int i = 0; i < int(S.size()); ++i) {
    int j = fail.back();
    while (j != -1 && S[i] != S[j]) j = fail[j];
    fail.push back(i+1);</pre>
```

duval.h

return fail:

Description: A string is called simple (or a Lyndon word), if it is strictly smaller than any of its own nontrivial suffixes.

Time: $\mathcal{O}\left(N\right)$ d9b2cb, 27 lines

```
template <typename T>
pair<int, vector<string>> duval(int n, const T &s) {
 assert (n >= 1);
 //s += s //uncomment if you need to know the min cyclic
       string
 vector<string> factors; // strings here are simple and in non
      -inc order
 int i = 0, ans = 0;
 while (i < n) { // until n/2 to find min cyclic string
   ans = i;
   int j = i + 1, k = i;
   while (j < n + n \&\& !(s[j % n] < s[k % n])) {
     if (s[k % n] < s[j % n]) k = i;
     else k++;
     j++;
    while (i \le k) {
      factors.push_back(s.substr(i, j-k));
      i += j - k;
 return {ans, factors};
  // returns 0-indexed position of the least cyclic shift
 // min cyclic string will be s.substr(ans, n/2)
template <typename T>
pair<int, vector<string>> duval(const T &s) {
 return duval((int) s.size(), s);
```

z-algorithm.h

Description: z[x] computes the length of the longest common prefix of s[i:] and s, except z[0] = 0. (abacaba -> 0010301) **Time:** $\mathcal{O}(n)$

7c8c64, 16 lines

```
vector<int> Z(const string& S) {
 vector<int> z(S.size());
  int 1 = -1, r = -1;
  for(int i = 1; i < int(S.size()); ++i) {</pre>
   z[i] = i >= r ? 0 : min(r - i, z[i - 1]);
   while (i + z[i] < int(S.size()) \&\& S[i + z[i]] == S[z[i]])
   if (i + z[i] > r) l = i, r = i + z[i];
 return z;
vector<int> get_prefix(string a, string b) {
  string str = a + '0' + b;
 vector < int > k = z(str);
  return vector<int>(k.begin() + int(a.size())+1, k.end());
```

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}(N)
```

```
87e1f0, 13 lines
array<vector<int>, 2> manacher(const string &s) {
  int n = s.size();
  array<vector<int>, 2> p = {vector<int>(n+1), vector<int>(n)};
  for (int z = 0; z < 2; ++z) 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) 1 = L, r = R;
  return p;
```

min-rotation.h

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

```
2a08fd, 8 lines
int min_rotation(string s) {
  int a=0, N=s.size(); s += s;
  for (int b = 0; b < N; ++b) for (int i = 0; i < N; ++i) {
   if (a+i == b \mid \mid s[a+i] < s[b+i]) {b += max(0, i-1); break;}
   if (s[a+i] > s[b+i]) \{ a = b; break; \}
  return a;
```

aho-corasick.h

5159f5, 41 lines

```
const int sigma = 26;
array<int, sigma> init;
for (int i = 0; i < sigma; i++) init[i] = -1;
vector<array<int, sigma>> trie(1, init);
vector<int> out(1, -1), parent(n, -1), ids(n);
for (int i = 0; i < n; i++) {
  int cur = 0;
  for (char ch : s[i]) {
   int c = ch - 'a';
   if (trie[cur][c] == -1) {
     trie[cur][c] = (int)trie.size();
     trie.push_back(init);
     out.push_back(-1);
    cur = trie[cur][c];
```

```
if (out[cur] == -1) out[cur] = i;
 ids[i] = out[cur];
vector<int> bfs; bfs.reserve(trie.size());
vector<int> f(trie.size());
for (int c = 0; c < sigma; c++) {
 if (trie[0][c] == -1) trie[0][c] = 0;
 else bfs.push_back(trie[0][c]);
for (int z = 0; z < (int)bfs.size(); z++) {
 int cur = bfs[z];
 for (int c = 0; c < sigma; c++) {
   if (trie[cur][c] == -1) {
     trie[cur][c] = trie[f[cur]][c];
   else {
     int nxt = trie[cur][c];
     int fail = trie[f[cur]][c];
     if (out[nxt] == -1) out[nxt] = out[fail];
     else parent[out[nxt]] = out[fail];
     f[nxt] = fail;
     bfs.push back(nxt);
```

suffix-arrav.h

Description: Builds suffix array for a string, first element is the size of the string. The 1cp function calculates longest common prefixes for neighbouring strings in suffix array. The returned vector is of size n+1.

Time: $\mathcal{O}(N \log N)$ where N is the length of the string for creation of the SA. $\mathcal{O}(N)$ for longest common prefixes.

```
<../data-structures/rmq.h>
                                                     5393da, 47 lines
mt19937 rng(chrono::steady clock::now().time since epoch().
    count()):
struct suffix_array_t {
 int N, H; vector<int> sa, invsa, lcp;
  rmq_t<pair<int, int>> rmq;
 bool cmp(int a, int b) { return invsa[a + H] < invsa[b + H];</pre>
  void ternary_sort(int a, int b) {
   if (a == b) return;
    int md = sa[a+rng() % (b-a)], lo = a, hi = b;
    for (int i = a; i < b; ++i) if (cmp(sa[i], md)) swap(sa[i],
          sa[lo++]);
    for (int i = b-1; i \ge lo; --i) if (cmp(md, sa[i])) swap(sa
        [i], sa[--hi]);
    ternary_sort(a, lo);
    for (int i = lo; i < hi; ++i) invsa[sa[i]] = hi-1;</pre>
    if (hi-lo == 1) sa[lo] = -1;
    ternary_sort(hi, b);
  suffix_array_t() {}
  template<typename I>
  suffix_array_t(I begin, I end): N(int(end - begin)+1), sa(N)
    vector<int> v(begin, end); v.push_back(INT_MIN);
    invsa = v; iota(sa.begin(), sa.end(), 0);
   H = 0; ternary_sort(0, N);
    for (H = 1; H <= N; H *= 2)
     for (int j = 0, i = j; i != N; i = j)
        if (sa[i] < 0) {
          while (j < N \&\& sa[j] < 0) j += -sa[j];
          sa[i] = -(j - i);
        else { j = invsa[sa[i]] + 1; ternary_sort(i, j); }
    for (int i = 0; i < N; ++i) sa[invsa[i]] = i;
    lcp.resize(N-1); int K = 0;
```

```
for (int i = 0; i < N-1; ++i) {
      if (invsa[i] > 0) while (v[i + K] == v[sa[invsa[i] - 1] +
            K]) ++K;
      lcp[invsa[i]-1] = K; K = max(K - 1, 0);
    vector<pair<int, int>> lcp_index(N-1);
    for (int i = 0; i < N-1; ++i) lcp_index[i] = {lcp[i], 1 + i}
    rmq = rmq_t<pair<int, int>>(std::move(lcp_index));
 pair<int, int> rmq_query(int a, int b) const { return rmq.
      query(a, b); }
  pair<int, int> get_split(int a, int b) const { return rmq.
      query(a, b-1); }
 int get_lcp(int a, int b) const {
   if (a == b) return N - a;
    a = invsa[a], b = invsa[b];
   if (a > b) swap(a, b);
    return rmq_query(a + 1, b + 1).first;
};
```

suffix-tree.h

Description: Builds suffix-tree informations based by emulating it over the suffix-array and lcp, root of the tree represents the special character (size of string for suffix-array), can therefore be ignored when calculating stuff.

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

```
<../data-structures/rmq.h>, "suffix-array.h"
                                                     31aacf, 46 lines
struct suffix tree t {
 int N. V:
  vector<vector<int>> ch;
  vector<array<int, 2>> sa_range;
  vector<int> leaves, par, depth;
  vector<int> suff link;
  vector<bool> is_unique_link, has_unique_child;
  suffix_array_t us;
  suffix tree t() {}
  suffix_tree_t(string S) : N(int(S.size())), V(0), ch(2*N+1),
  sa\_range(2*N+1), leaves(N+1), par(2*N+1), depth(2*N+1),
  us(S.begin(), S.end()) { dfs(0, N+1, -1); }
  void dfs(int a, int b, int prv) {
    int cur = V++;
    par[cur] = prv;
    if (prv != -1) ch[prv].push_back(cur);
    sa_range[cur] = {a, b};
    if (b - a == 1) {
      leaves[us.sa[a]] = cur;
      depth[cur] = N - us.sa[a];
    } else {
      int d = us.get_split(a, b).first;
      depth[cur] = d;
      int mi = a;
      while (b - mi >= 2) {
        auto [nd, nmi] = us.get_split(mi, b);
        if (nd != d) break;
        dfs(mi, nmi, cur);
        mi = nmi;
      dfs(mi, b, cur);
 void build_links() {
    suff_link.resize(V, -1), is_unique_link.resize(V),
         has unique child.resize(V);
    for (int i = 0; i < N; ++i) {
      for (int cur = leaves[i], link = leaves[i+1];; cur = par[
        if (cur == 0 || suff_link[cur] != -1) break;
        suff link[cur] = link;
```

```
is_unique_link[cur] = (sa_range[cur][1] - sa_range[cur
            [0] == (sa_range[link][1] - sa_range[link][0]);
       if (is_unique_link[cur]) has_unique_child[link] = true;
       while (~link && depth[link] + 1 > depth[par[cur]]) link
             = par[link];
};
```

suffix-automaton.h

```
Description: Suffix automaton
                                                     c4406e, 38 lines
template<int offset = 'a'> struct array_state {
  array<int, 26> as;
  array_state() { fill(begin(as), end(as), ~0); }
  int& operator[](char c) { return as[c - offset]; }
  int count(char c) { return (~as[c - offset] ? 1 : 0); }
template<typename Char, typename state = map<Char, int>> struct
     suffix_automaton {
  struct node t {
   int len, link; int64_t cnt;
   state next;
  };
  int N, cur;
 vector<node_t> nodes;
  suffix_automaton() : N(1), cur(0), nodes{node_t{0, -1, 0}},
      {}}} {}
  node_t& operator[](int v) { return nodes[v]; };
  void append(Char c) {
   int v = cur; cur = N++;
   nodes.push\_back(node\_t\{nodes[v].len + 1, 0, 1, \{\}\});
    for (; ~v && !nodes[v].next.count(c); v = nodes[v].link) {
     nodes[v].next[c] = cur;
    if (~∨) {
     const int u = nodes[v].next[c];
     if (nodes[v].len + 1 == nodes[u].len) {
       nodes[cur].link = u;
      } else {
       const int clone = N++;
       nodes.push_back(nodes[u]);
       nodes[clone].len = nodes[v].len + 1;
       nodes[u].link = nodes[cur].link = clone;
        for (; ~v && nodes[v].next[c] == u; v = nodes[v].link)
          nodes[v].next[c] = clone;
```

9.1 Suffix Automaton

};

9.1.1 Number of different substrings

Is the number of paths in the automaton starting at the root.

$$d(v) = 1 + \sum_{v \to w} d(w)$$

9.1.2 Total length of different substrings

Is the sum of children answers and paths starting at each children.

$$ans(v) = \sum_{v \to w} d(w) + ans(w)$$

9.1.3 Lexicographically K-th substring

Is the K-th lexicographically path, so you can search using the number of paths from each state

9.1.4 Smallest cyclic shift

Construct for string S + S. Greedily search the minimal character.

9.1.5 Number of occurrences

For each state not created by cloning, initialize cnt(v) = 1. Then, just do a dfs to calculate cnt(v)

$$cnt(link(v)) + = cnt(v)$$

9.1.6 First occurrence position

When we create a new state cur do first(pos) = len(cur) - 1. When we clone q as clone do first(clone) = first(q). Answer is first(v) - size(P) + 1, where v is the state of string P

9.1.7 All occurrence positions

From first(v) do a dfs using suffix link, from link(u) go to u.

Various (10)

10.1 Intervals

interval-container.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)
                                                     f47dfb, 22 lines
set<pair<int,int>>::iterator addInterval(set<pair<int,int>> &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<pair<int,int>> &is, int L, int R) {
 if (L == R) return;
 auto it = addInterval(is, L, R);
 auto r2 = it->second;
 if (it->first == L) is.erase(it);
 else (int&)it->second = L;
 if (R != r2) is.emplace (R, r2);
```

interval-cover.h

Description: Compute indices of smallest set of intervals covering another interval. Intervals should be [inclusive, exclusive). To support [inclusive, inclusive, change (A) to add | | R.empty(). Returns empty set on failure (or if G is empty). Time: $\mathcal{O}(N \log N)$

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

constant-intervals.h

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

```
Usage: constantIntervals(0, sz(v), [&](int x){
return v[x]; {\( \) (int lo, int hi, T val) \( \) (... \( \));
Time: \mathcal{O}\left(k\log\frac{n}{k}\right)
                                                         753a4c, 19 lines
template<class F, class G, class T>
void rec(int from, int to, F& f, G& q, int& i, T& p, T q) {
 if (p == q) return;
 if (from == to) {
    g(i, to, p);
    i = to; p = q;
    int mid = (from + to) >> 1;
    rec(from, mid, f, q, i, p, f(mid));
    rec(mid+1, to, f, g, i, p, q);
template<class F, class G>
void constantIntervals(int from, int to, F f, G g) {
 if (to <= from) return;
 int i = from; auto p = f(i), q = f(to-1);
 rec(from, to-1, f, g, i, p, q);
 g(i, to, q);
```

10.2 Misc. algorithms

basis-manager.h

Description: A list of basis values sorted in decreasing order, where each value has a unique highest bit. d5bcd3, 41 lines

```
const int BITS = 60;
template<typename T> struct xor basis {
 int N = 0;
 array<T, BITS> basis;
 T min_value(T start) const {
    if (N == BITS) return 0;
```

```
for (int i = 0; i < N; ++i)
   start = min(start, start ^ basis[i]);
  return start:
T max_value(T start = 0) const {
 if (N == BITS) return ((T) 1 << BITS) - 1;
  for (int i = 0; i < N; ++i)
   start = max(start, start ^ basis[i]);
 return start;
bool add(T x) {
 x = \min value(x);
 if (x == 0) return false;
 basis[N++] = x;
  // Insertion sort.
  for (int k = N - 1; k > 0 && basis[k] > basis[k - 1]; k--)
   swap(basis[k], basis[k - 1]);
  return true;
void merge(const xor_basis<T>& other) {
  for (int i = 0; i < other.n && N < BITS; <math>i++)
   add(other.basis[i]);
void merge(const xor_basis<T>& a, const xor_basis<T>& b) {
 if (a.N > b.N) {
   *this = a;
   merge(b);
  } else {
    *this = b;
   merge(a);
```

ternary-search.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 nonstrict inequalities, change the < marked with (A) to <=, and reverse the loop at (B). To minimize f, change it to >, also at (B). If you are dealing with real numbers, you'll need to pick $m_1 = (2a+b)/3.0$ and $m_2 = (a+2b)/3.0$. Consider setting a constant number of iterations for the search, usually [200,300] iterations are sufficient for problems with error limit as 10^{-6} .

Usage: int ind = ternSearch(0, n-1, [&] (int i) {return a[i];});

```
Time: \mathcal{O}(\log(b-a)) 35ef73, 12 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; } for (int i = a+1; i <= b; ++i) if (f (a) < f (i)) a = i; // (B) return a; }
```

count-triangles.h

Description: Counts x, y >= 0 such that $Ax + By \le C$.

```
flovd-cvcle.h
```

Description: Detect loop in a list. Consider using mod template to avoid overflow. **Time:** $\mathcal{O}(n)$

template < class F >
pair < int, int > find (int x0, F f) {
 int t = f(x0), h = f(t), mu = 0, lam = 1;
 while (t != h) t = f(t), h = f(f(h));
 h = x0;
 while (t != h) t = f(t), h = f(h), ++mu;
 h = f(t);
 while (t != h) h = f(h), ++lam;
 return {mu, lam};

10.3 Dynamic programming

divide-and-conquer-dp.h

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

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

```
struct DP { // Modify at will:
 int lo(int ind) { return 0; }
 int hi(int ind) { return ind; }
 lint f(int ind, int k) { return dp[ind][k]; }
 void store(int ind, int k, lint v) { res[ind] = {k, v}; }
 void rec(int L, int R, int LO, int HI) {
   if (L >= R) return;
   int mid = (L + R) \gg 1;
   pair<lint, int> best(LLONG MAX, LO);
   for (int k = max(LO, lo(mid)); k \le min(HI, hi(mid)); ++k)
     best = min(best, make_pair(f(mid, k), k));
   store(mid, best.second, best.first);
   rec(L, mid, LO, best.second);
    rec(mid+1, R, best.second, HI);
 void solve(int L, int R) { rec(L, R, INT_MIN, INT_MAX); }
struct DP { // Modify at will:
 vector<int>a, freq;
 vector<lint>old, cur;
 lint cnt;
 int lcur, rcur;
 DP(const \ vector < int > \&_a, \ int \ n): \ a(\_a), \ freq(n), \ old(n+1,
      linf), cur(n+1, linf), cnt(0), lcur(0), rcur(0){}
 int lo(int ind) { return 0; }
 int hi(int ind) { return ind; }
 void add(int k, int c) { cnt += freq[a[k]]++; }
 void del(int k, int c) { cnt -= --freq[a[k]]; }
 lint C(int 1, int r) {
   while(lcur > 1) add(--lcur, 0);
   while(rcur < r) add(rcur++, 1);</pre>
   while(lcur < 1) del(lcur++, 0);</pre>
   while(rcur > r) del(--rcur, 1);
    return cnt;
 lint f(int ind, int k) { return old[k] + C(k, ind); }
 void store(int ind, int k, lint v) { cur[ind] = v; }
 void rec(int L, int R, int LO, int HI) {
   if (L >= R) return;
   int mid = (L + R) \gg 1;
   pair<lint, int> best(LLONG_MAX, LO);
```

for (int $k = max(LO, lo(mid)); k \le min(HI, hi(mid)); ++k)$

best = min(best, make_pair(f(mid, k), k));

```
store(mid, best.second, best.first);
rec(L, mid, LO, best.second);
rec(mid+1, R, best.second, HI);
};
```

knuth-dp.h

b456ab, 10 lines

2cef33, 50 lines

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:** $\mathcal{O}(N^2)$

digit-dp.h

Description: Compute how many # between 1 and N have K distinct digits in the base L without leading zeros;

```
Usage: auto hex_to_dec = [&] (char c) -> int { return ('A' <= c && c <= 'F' ? (10 + c - 'A') : (c - '0')); }; digit_dp<modnum<int(1e9) + 7>, hex_to_dec>(N, K); Time: \mathcal{O}(NK)
```

```
template<typename T, class F> T digit_dp(const string& S, int K
    , F& L) {
 const int base = 16;
 const int len = int(S.size());
 vector<bool> w(base);
 vector<vector<T>> dp(len + 1, vector<T>(base + 2));
 int cnt = 0:
 for (int d = 0; d < len; ++d) {
   // adding new digit to numbers with prefix < s
   for (int x = 0; x \le base; ++x) {
     dp[d + 1][x] += dp[d][x] * x;
     dp[d + 1][x + 1] += dp[d][x] * (base - x);
    // adding strings whith prefix only 0's and last digit != 0
   if (d) dp[d + 1][1] += (base - 1);
   // adding prefix equal to s and last digit < s, first digit
         cannot be 0
    for (int x = 0; x < L(S[d]); ++x)
     if (d == 0 \&\& x == 0) continue;
     if (w[x]) dp[d + 1][cnt] += 1;
     else dp[d + 1][cnt + 1] += 1;
    // marking if the last digit appears in the prefix of s
   if (w[L(S[d])] == false) {
     w[L(S[d])] = true;
     cnt++;
 // adding string k
 dp[len][cnt] += 1;
 return dp[len][K];
```

knapsack-bounded.h

Description: You are given n types of items, each items has a weight and a quantity. Is possible to fill a knapsack with capacity k using any subset of items?

```
Time: \mathcal{O}(Wn)
```

9bddad, 15 lines

```
vector<int> how_many(n+1), dp(k+1);
```

```
dp[0] = 1;
for (int i = 1; i <= n; ++i) cin >> how_many[i];
for (int i = 1; i <= n; ++i) {
  for (int j = k-items[i]; j >= 0; --j) {
    if (dp[j]) {
      int x = 1;
      while (x <= how_many[i] &&
        j + x*items[i] <= k && !dp[j + x*items[i]]) {
      dp[j + x*items[i]] = 1;
        ++x;
    }
    }
}</pre>
```

knapsack-bounded-costs.h

Description: You are given n types of items, you have e[i] items of i-th type, and each item of i-th type weight w[i] and cost c[i]. What is the minimal cost you can get by picking some items weighing at most W in total? **Time:** $\mathcal{O}(Wn)$

```
3ade3c, 28 lines
<MinQueue.h>
const int maxn = 1000;
const int maxm = 100000;
const int inf = 0x3f3f3f;
minQueue<int> g[maxm];
array<int, maxm> dp; // the minimum cost dp[i] I need to pay in
      order to fill the knapsack with total weight i
int w[maxn], e[maxn], c[maxn]; // weight, number, cost
int main() {
  int n, m;
  cin >> n >> m;
  for (int i = 1; i \le n; i++) cin >> w[i] >> c[i] >> e[i];
  for (int i = 1; i <= m; i++) dp[i] = inf;
  for (int i = 1; i <= n; i++) {
    for (int j = 0; j < w[i]; j++) q[j].clear();
    for (int j = 0; j <= m; j++) {
      minQueue<int> &mq = q[j % w[i]];
      if (mq.size() > e[i]) mq.pop();
      mq.add(c[i]);
      mq.push(dp[j]);
      dp[j] = mq.getMin();
  cout << "Minimum value i can pay putting a total weight " <<</pre>
      m \ll " is " \ll dp[m] \ll ' n';
  for (int i = 0; i <= m; i++) cout << dp[i] << " " << i << '\n
  cout << "\n";
```

two-max-equal-sum.h

Description: Two maximum equal sum disjoint subsets, s[i] = 0 if v[i] wasn't selected, s[i] = 1 if v[i] is in the first subset and s[i] = 2 if v[i] is in the second subset

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

```
rec[i][d] = newdp[d] == a ? 1 : newdp[d] == b ? 2 : 0;
}
for(int j = i-1, d = sum; j >= 0 ; j--)
  d += (s[j] = rec[j][d]) ? s[j] == 2 ? v[j] : -v[j] : 0;
return {dp[sum], s};
```

10.4 Optimization tricks

10.4.1 Bit hacks

- 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.

hashmap.h

Description: Faster/better hash maps, taken from CF

09a72f, 19 lines

```
#include < bits / extc++.h>
struct splitmix64_hash {
 static uint64 t splitmix64(uint64 t x) {
   x += 0x9e3779b97f4a7c15;
   x = (x^{(x)} >> 30) \times 0xbf58476d1ce4e5b9;
   x = (x^{(x)} > 27) \times 0x94d049bb133111eb;
    return x^(x >> 31);
 size_t operator()(uint64_t x) const {
   static const uint64 t FIXED RANDOM = std::chrono::
        steady_clock::now().time_since_epoch().count();
    return splitmix64(x + FIXED_RANDOM);
};
template <typename K, typename V, typename Hash =
    splitmix64 hash>
using hash_map = __gnu_pbds::gp_hash_table<K, V, Hash>;
template <typename K, typename Hash = splitmix64_hash>
using hash_set = hash_map<K, __gnu_pbds::null_type, Hash>;
```

10.5 Bit Twiddling Hack

```
hacks.h
                                                        829b7d, 22 lines
// Iterate on non-empty submasks of a bitmask.
for (int s = m; s > 0; s = (m & (s - 1)))
// Iterate on non-zero bits of a bitset B.
for (int j = B._Find_next(0); j < MAXV; j = B._Find_next(j))</pre>
ll next_perm(ll v) { // compute next perm i.e.
 ll t = v | (v-1); // 00111,01011,01101,10011 ...
  return (t + 1) \mid (((\sim t \& -\sim t) - 1) >> (\_builtin\_ctz(v) + 1))
template<typename F> // All subsets of size k of {0..N-1}
void iterate k subset(ll N, ll k, F f) {
  11 \text{ mask} = (111 << k) - 1;
  while (!(mask & 111<<N)) { f(mask);</pre>
    11 t = mask \mid (mask-1);
    mask = (t+1) \mid (((\sim t \& -\sim t) - 1) >> (\underline{builtin\_ctzll(mask)})
         +1));
template<typename F> // All subsets of set
void iterate_mask_subset(ll set, F f) { ll mask = set;
 do f(mask), mask = (mask-1) & set;
```

```
while (mask != set);
```

10.6 Random Numbers

random-numbers.h

Description: An example on the usage of generator and distribution. Use shuffle instead of random shuffle.

b28ebe. 9 lines