

Universidade do Estado de Santa Catarina

Não Treinamos o João

Enzo Rodrigues, Eric Grochowicz, João Oliveira

1	Contest	1						
2	Mathematics	1						
3	Data structures	3						
4	Numerical	5						
5	Number theory	8						
6	Combinatorial	9						
7	Graph	10						
8	Geometry	14						
9	Strings	17						
10) Various	18						
Contest (1)								
.vi	imre							
set nu ai si cindent et ts=4 sw=4 so=10 nosm undofile								
<pre>inoremap {} {}<left><return><up><end><return> " remap de chaves</return></end></up></return></left></pre>								
<pre>au BufReadPost * if line("'\"") > 0 && line("'\"") <= line("\$")</pre>								
ru	n.sh	nes						
#!/bin/bash g++ -std=c++20 -DBRUTE -O2 -Wall -Wextra -Wconversion -Wfatal- errors -fsanitize=address,undefined \$1 && ./a.out								
ha	sh.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								
Mint.cpp								
	mplate <auto mod,="" t="decltype(MOD)" typename=""></auto>							
st:	<pre>ruct Mint { using U = long long;</pre>							
// se o modulo for long long, usar U =int128 using m = Mint <mod, t="">;</mod,>								
	T v;							
	<pre>Mint(T val = 0) : v(val) { assert(sizeof(T) * 2 <= sizeof(U));</pre>							
	if $(v < -MOD \mid \mid v >= 2 * MOD) v %= MOD;$							
	if $(v < 0) v += MOD;$ if $(v >= MOD) v -= MOD;$							

Mint(U val) : v(T(val % MOD)) {

if (v < 0) v += MOD;

assert(sizeof(T) * 2 <= sizeof(U));</pre>

```
bool operator==(m o) const { return v == o.v; }
    bool operator<(m o) const { return v < o.v; }</pre>
   bool operator!=(m o) const { return v != o.v; }
    m pwr(m b, U e) {
       m res = 1;
        while (e > 0) {
           if (e & 1) res *= b;
           b *= b, e /= 2;
        return res;
    m &operator+=(m o) {
       v -= MOD - o.v;
        if (v < 0) v += MOD;
       return *this;
    m &operator-=(m o) {
       v -= o.v;
        if (v < 0) v += MOD;
        return *this;
    m &operator*=(m o) {
       v = (T) ((U) v * o.v % MOD);
        return *this:
    m &operator/=(m o) { return *this *= o.pwr(o, MOD - 2); }
    m &operator^=(U e) { return *this = pwr(*this, e); }
    friend m operator-(m a, m b) { return a -= b; }
    friend m operator+(m a, m b) { return a += b;
    friend m operator* (m a, m b) { return a *= b;
    friend m operator/(m a, m b) { return a /= b; }
    friend m operator^(m a, U e) { return a.pwr(a, e); }
};
troubleshoot.txt
                                                          52 lines
Pre-submit:
Write a few simple test cases if sample is not enough.
Are time limits close? If so, generate max cases.
Is the memory usage fine?
Could anything overflow?
Make sure to submit the right file.
Wrong answer:
Print your solution! Print debug output, as well.
Are you clearing all data structures between test cases?
Can your algorithm handle the whole range of input?
Read the full problem statement again.
Do you handle all corner cases correctly?
Have you understood the problem correctly?
Any uninitialized variables?
Any overflows?
Confusing N and M, i and j, etc.?
Are you sure your algorithm works?
What special cases have you not thought of?
Are you sure the STL functions you use work as you think?
Add some assertions, maybe resubmit.
Create some testcases to run your algorithm on.
Go through the algorithm for a simple case.
Go through this list again.
Explain your algorithm to a teammate.
Ask the teammate to look at your code.
Go for a small walk, e.g. to the toilet.
Is your output format correct? (including whitespace)
Rewrite your solution from the start or let a teammate do it.
```

Runtime error:

Have you tested all corner cases locally?

```
Any uninitialized variables?

Are you reading or writing outside the range of any vector?

Any assertions that might fail?

Any possible division by 0? (mod 0 for example)

Any possible infinite recursion?

Invalidated pointers or iterators?

Are you using too much memory?
```

Debug with resubmits (e.g. remapped signals, see Various).

Time limit exceeded:

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

Memory limit exceeded:

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

Mathematics (2)

2.1 Equations

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

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

$$ax + by = e$$

$$cx + dy = f$$

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

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

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.

2.2 Linear Operators

2.2.1 Rotate counter-clockwise by θ°

$$\begin{bmatrix} \cos\theta & -\sin\theta \\ \sin\theta & \cos\theta \end{bmatrix}$$

2.2.2 Reflect about the line y = mx

$$\frac{1}{m^2+1} \begin{bmatrix} 1-m^2 & 2m \\ 2m & m^2-1 \end{bmatrix}$$

2.2.3 Inverse of a 2x2 matrix A

$$\begin{bmatrix} a & b \\ c & d \end{bmatrix}^{-1} = \frac{1}{\det(A)} \begin{bmatrix} d & -b \\ -c & a \end{bmatrix}$$

.vimrc run hash Mint troubleshoot

2.2.4 Horizontal shear by K

$$\begin{bmatrix} 1 & K \\ 0 & 1 \end{bmatrix}$$

2.2.5 Vertical shear by K

$$\begin{bmatrix} 1 & 0 \\ K & 1 \end{bmatrix}$$

2.2.6 Change of basis

 \vec{a}_{β} are the coordinates of vector \vec{a} in basis β . \vec{a} are the coordinates of vector \vec{a} in the canonical basis. $\vec{b1}$ and $\vec{b2}$ are the basis vectors for β .

C is a matrix that changes from basis β to the canonical basis.

$$C\vec{a}_{\beta} = \vec{a}$$

$$C^{-1}\vec{a} = \vec{a}_{\beta}$$

$$C = \begin{bmatrix} b1_x & b2_x \\ b1_y & b2_y \end{bmatrix}$$

2.2.7 Properties of matrix operations

$$(AB)^{-1} = A^{-1}B^{-1}$$

$$(AB)^T = B^T A^T$$

$$(A^{-1})^T = (A^T)^{-1}$$

$$(A+B)^T = A^T + B^T$$

$$det(A) = det(A^T)$$

$$\det(AB) = \det(A)\det(B)$$

Let A be an NxN matrix:

$$det(kA) = K^N det(A)$$

2.3 Recurrences

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

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

Non-distinct roots r become polynomial factors, e.g. $a_n = (d_1n + d_2)r^n$.

2.4 Trigonometry

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

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

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

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

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

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

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

2.5 Geometry

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

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

Length of bisector (divides angles in two):

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

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

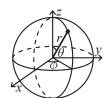
2.5.2 Quadrilaterals

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

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

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

2.5.3 Spherical coordinates



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

2.6 Derivatives/Integrals

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

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

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

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

Integration by parts:

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

2.7 Sums

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

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

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

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

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

2.8 Series

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

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

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

Não Treinamos o João - UDESC

$$\sin x = x - \frac{x^3}{3!} + \frac{x^5}{5!} - \frac{x^7}{7!} + \dots, (-\infty < x < \infty)$$
$$\cos x = 1 - \frac{x^2}{2!} + \frac{x^4}{4!} - \frac{x^6}{6!} + \dots, (-\infty < x < \infty)$$

2.9 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,

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

2.9.1 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 which yields success with probability p is Fs(p), $0 \le p \le 1$.

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

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

Poisson distribution

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

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

2.9.2 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

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

2.10 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 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)

OrderedSet.cpp

Usage: ordered_set<int> seta;

Time: All operations in $\mathcal{O}(\log N)$.

SegmentTreeBeats.cpp

Description: Updates of max, min and sum in range with queries of sum in range

Usage: "seg" is already declared globally, just call build Time: If sum update is not used: $\mathcal{O}(\log N)$, otherwise $\mathcal{O}\left(\log^2 N\right)$.

```
const 11 INF = 1e18;
struct node {
    11 mi, smi, mx, smx, sum, lazy;
    int fmi, fmx;
    node() {
        mi = smi = INF;
        mx = smx = -INF;
        fmi = 0, fmx = 0, sum = 0, lazv = 0;
    node(ll val) {
        mi = mx = sum = val;
        smi = INF, smx = -INF;
        fmx = fmi = 1;
        lazy = 0;
};
node operator+(node a, node b) {
    node ret;
    ret.sum = a.sum + b.sum;
    if (a.mi == b.mi) {
        ret.mi = a.mi;
        ret.fmi = a.fmi + b.fmi;
        ret.smi = min(a.smi, b.smi);
    } else if (a.mi < b.mi) {</pre>
        ret.mi = a.mi;
        ret.fmi = a.fmi;
        ret.smi = min(a.smi, b.mi);
    } else {
        ret.mi = b.mi;
        ret.fmi = b.fmi;
        ret.smi = min(b.smi, a.mi);
```

SegmentTree2D ImplicitTreap

```
if (a.mx == b.mx) {
       ret.mx = a.mx;
       ret.fmx = a.fmx + b.fmx;
       ret.smx = max(a.smx, b.smx);
    } else if (a.mx > b.mx) {
       ret.mx = a.mx;
       ret.fmx = a.fmx;
       ret.smx = max(b.mx, a.smx);
   } else {
       ret.fmx = b.fmx;
       ret.mx = b.mx;
       ret.smx = max(a.mx, b.smx);
   return ret;
struct SegBeats {
   vector<node> t;
   void build(int _n) { // pra construir com tamanho, mas
        vazia
       n = _n;
       t.assign(n * 4, node());
   void build(const vector<11> &v) { // pra construir com
        vector
       n = (int)v.size();
       t.assign(n * 4, node());
       build(1, 0, n - 1, v);
   void build(11 *bg, 11 *en) { // pra construir com array de
        C
       build(vector<ll>(bg, en));
   inline int lc(int p) { return 2 * p; }
   inline int rc(int p) { return 2 * p + 1; }
   node build(int p, int l, int r, const vector<11> &a) {
       if (1 == r) return t[p] = node(a[1]);
       int mid = (1 + r) >> 1;
       return t[p] = build(lc(p), l, mid, a) + build(rc(p),
            mid + 1, r, a);
   void pushsum(int p, int l, int r, ll x) {
       t[p].sum += (r - 1 + 1) * x;
       t[p].mi += x;
       t[p].mx += x;
       t[p].lazv += x;
       if (t[p].smi != INF) t[p].smi += x;
       if (t[p].smx != -INF) t[p].smx += x;
   void pushmax(int p, ll x) {
       if (x <= t[p].mi) return;</pre>
       t[p].sum += t[p].fmi * (x - t[p].mi);
       if (t[p].mx == t[p].mi) t[p].mx = x;
       if (t[p].smx == t[p].mi) t[p].smx = x;
       t[p].mi = x;
   void pushmin(int p, ll x) {
       if (x >= t[p].mx) return;
       t[p].sum += t[p].fmx * (x - t[p].mx);
       if (t[p].mi == t[p].mx) t[p].mi = x;
       if (t[p].smi == t[p].mx) t[p].smi = x;
       t[p].mx = x;
   void pushdown(int p, int l, int r) {
       if (1 == r) return;
       int mid = (1 + r) >> 1;
       pushsum(lc(p), l, mid, t[p].lazy);
       pushsum(rc(p), mid + 1, r, t[p].lazy);
```

```
t[p].lazy = 0;
        pushmax(lc(p), t[p].mi);
        pushmax(rc(p), t[p].mi);
        pushmin(lc(p), t[p].mx);
        pushmin(rc(p), t[p].mx);
   node updatemin(int p, int l, int r, int L, int R, ll x) {
        if (1 > R \mid | r < L \mid | x >= t[p].mx) return t[p];
        if (1 >= L \&\& r <= R \&\& x > t[p].smx) {
            pushmin(p, x);
            return t[p];
        pushdown(p, l, r);
        int mid = (1 + r) >> 1;
        t[p] = updatemin(lc(p), l, mid, L, R, x) + updatemin(rc
             (p), mid + 1, r, L, R, x);
        return t[p];
   node updatemax(int p, int l, int r, int L, int R, ll x) {
        if (1 > R \mid | r < L \mid | x <= t[p].mi) return t[p];
        if (1 >= L \&\& r <= R \&\& x < t[p].smi) {
            pushmax(p, x);
            return t[p];
        pushdown(p, 1, r);
        int mid = (1 + r) >> 1;
        t[p] = updatemax(lc(p), l, mid, L, R, x) + updatemax(rc
             (p), mid + 1, r, L, R, x);
        return t[p];
    node updatesum(int p, int l, int r, int L, int R, ll x) {
        if (1 > R || r < L) return t[p];
        if (1 >= L && r <= R) {
            pushsum(p, 1, r, x);
            return t[p];
       pushdown(p, l, r);
        int mid = (1 + r) >> 1;
        return t[p] = updatesum(lc(p), 1, mid, L, R, x) +
                      updatesum(rc(p), mid + 1, r, L, R, x);
   node guery(int p, int l, int r, int L, int R) {
        if (1 > R || r < L) return node();
        if (1 >= L && r <= R) return t[p];
       pushdown(p, l, r);
        int mid = (1 + r) >> 1;
        return query(lc(p), 1, mid, L, R) + query(rc(p), mid +
             1, r, L, R);
   ll query (int 1, int r) { return query (1, 0, n - 1, 1, r).
    void updatemax(int 1, int r, 11 x) { updatemax(1, 0, n - 1,
          1, r, x); }
    void updatemin(int 1, int r, 11 x) { updatemin(1, 0, n - 1,
         1, r, x); }
   void updatesum(int 1, int r, 11 x) { updatesum(1, 0, n - 1,
          1, r, x); }
} seg;
SegmentTree2D.cpp
Description: Segment Tree with point update and rectangular range query
Usage: "seg" declared globally
Time: \mathcal{O}\left(\log^2 N\right)
                                                     d41d8c, 47 lines
struct SegTree2D {
   11 merge(ll a, ll b) { return a + b; }
```

11 neutral = 0;

```
int n, m;
    vector<vector<ll>> t;
    void build(int _n, int _m) {
        n = _n, m = _m;
        t.assign(2 * n, vector<11 > (2 * m, neutral));
        for (int i = 2 * n - 1; i >= n; i--)
            for (int j = m - 1; j > 0; j--)
                t[i][j] = merge(t[i][j << 1], t[i][j << 1 | 1])
        for (int i = n - 1; i > 0; i--)
            for (int j = 2 * m - 1; j > 0; j--)
                t[i][j] = merge(t[i << 1][j], t[i << 1 | 1][j])
    11 inner_query(int idx, int 1, int r) {
        11 res = neutral;
        for (1 += m, r += m + 1; 1 < r; 1 >>= 1, r >>= 1) {
            if (1 & 1) res = merge(res, t[idx][1++]);
            if (r \& 1) res = merge(res, t[idx][--r]);
        return res;
    // query do ponto (a, b) ate o ponto (c, d), retorna neutro
          se \ a > c \ ou \ b > d
    11 query(int a, int b, int c, int d) {
        11 res = neutral;
        for (a += n, c += n + 1; a < c; a >>= 1, c >>= 1) {
            if (a & 1) res = merge(res, inner_query(a++, b, d))
            if (c & 1) res = merge(res, inner_query(--c, b, d))
        return res;
    void inner_update(int idx, int i, ll x) {
        auto &c = t[idx];
        i += m;
        c[i] = x;
        for (i >>= 1; i > 0; i >>= 1) c[i] = merge(c[i << 1], c
             [i << 1 | 1]);
    void update(int i, int j, ll x) {
       i += n:
        inner update(i, j, x);
        for (i >>= 1; i > 0; i >>= 1) {
            11 \text{ val} = \text{merge}(t[i << 1][j + m], t[i << 1 | 1][j +
                m1):
            inner_update(i, j, val);
} seq;
```

Implicit Treap.cpp Description: Does everything Time: Expected $\mathcal{O}(\log N)$

d41d8c, 124 lines

5

```
node info(T val)
        : 1(0), r(0), y(rng()), size(0), val(_val), acc(0),
              add(0), rev(false) { }
};
using node = node_info *;
node root = 0;
inline int size(node t) { return t ? t->size : 0; }
inline T acc(node t) { return t ? t->acc : 0; }
inline bool rev(node t) { return t ? t->rev : false; }
inline void push (node t) {
   if (!t) return;
    if (rev(t)) {
        t->rev = false;
        swap(t->1, t->r);
        if (t->1) t->1->rev ^= 1;
        if (t->r) t->r->rev ^= 1;
    t->acc += t->add * size(t);
    // t \rightarrow acc += t \rightarrow add se for RMQ
    t->val += t->add;
    if (t->1) t->1->add += t->add;
    if (t->r) t->r->add += t->add;
    t->add = 0;
inline void pull(node t) {
    if (t) {
        push(t->1), push(t->r);
        t->size = size(t->1) + size(t->r) + 1;
        t\rightarrow acc = merge(t\rightarrow val, merge(acc(t\rightarrow l), acc(t\rightarrow r)))
void merge (node &t, node L, node R) {
    push(L), push(R);
    if (!L || !R) {
        t = L ? L : R;
    } else if (L->y > R->y) {
        merge(L->r, L->r, R);
        t = L;
    } else {
        merge(R->1, L, R->1);
        t = R;
    pull(t);
void split(node t, int pos, node &L, node &R, int add = 0)
    if (!t) {
        L = R = nullptr;
    } else {
        push(t);
        int imp_key = add + size(t->1);
        if (pos <= imp kev) {
            split(t->1, pos, L, t->1, add);
            R = t;
        } else {
            split(t->r, pos, t->r, R, imp_key + 1);
            L = t;
    pull(t);
inline void insert (node to, int pos) {
    node L, R;
    split(root, pos, L, R);
    merge(L, L, to);
    merge(root, L, R);
bool remove(node &t, int pos, int add = 0) {
```

```
if (!t) return false;
   push(t);
   int imp_key = add + size(t->1);
   if (pos == imp_key) {
        node me = t;
       merge(t, t->1, t->r);
       delete me;
       return true;
   bool ok:
   if (pos < imp_key) ok = remove(t->1, pos, add);
   else ok = remove(t->r, pos, imp_key + 1);
   pull(t);
   return ok;
inline T query(int 1, int r) {
   if (1 > r) return neutral;
   node L1, L2, R1, R2;
   split(root, r + 1, L1, R1);
   split(L1, 1, L2, R2);
   T ans = acc(R2);
   merge(L1, L2, R2);
   merge(root, L1, R1);
   return ans:
inline void update_sum(int 1, int r, T val) {
   if (1 > r) return;
   node L1, L2, R1, R2;
   split(root, r + 1, L1, R1);
   split(L1, 1, L2, R2);
   assert (R2);
   R2->add += val;
   merge(L1, L2, R2);
   merge(root, L1, R1);
inline void reverse(int 1, int r) {
   if (1 > r) return;
   node L1, L2, R1, R2;
   split(root, r + 1, L1, R1);
   split(L1, 1, L2, R2);
   R2 -> rev ^= 1;
   merge(L1, L2, R2);
   merge(root, L1, R1);
inline void insert(int pos, int val) { insert(new node_info
     (val), pos); }
inline bool remove(int pos) { return remove(root, pos); }
```

Numerical (4)

4.1 Factorization

PollardRhoMillerRabin.cpp

Description: Pollard Rho and Miller Rabin

```
inline bool composite(ll n, ll a, ll d, ll s) {
       ll x = power(a, d, n);
       if (x == 1 | | x == n - 1 | | a % n == 0) return false;
       for (int r = 1; r < s; r++) {
           x = mul\_mod(x, x, n);
           if (x == n - 1) return false;
       return true:
    // com esses "primos", o teste funciona garantido para n <=
    int primes[] = {2, 325, 9375, 28178, 450775, 9780504,
        1795265022};
   // funciona para n \le 3*10^24 com os primos ate 41, mas tem
         que cuidar com overflow
    37, 41};
   bool prime(ll n) {
       if (n <= 2 || (n % 2 == 0)) return n == 2;
       11 d = n - 1, r = 0;
       while (d \% 2 == 0) d /= 2, r++;
       for (int a : primes)
           if (composite(n, a, d, r)) return false;
       return true;
namespace PollardRho {
   mt19937 rng((uint32_t)chrono::steady_clock::now().
        time_since_epoch().count());
   const 11 P = 1e6 + 1;
   ll seg[P];
   inline 11 add_mod(11 x, 11 y, 11 m) { return (x += y) < m ?
         x : x - m; }
    inline 11 mul_mod(11 a, 11 b, 11 m) { return (11)((__int128
        )a * b % m); }
   ll rho(ll n) {
       if (n % 2 == 0) return 2;
       if (n % 3 == 0) return 3;
       11 \times 0 = rng() % n, c = rng() % n;
       while (1) {
           11 x = x0++, y = x, u = 1, v, t = 0;
           11 *px = seq, *py = seq;
           while (1) {
                *py++ = y = add_mod(mul_mod(y, y, n), c, n);
                *py++ = y = add mod(mul mod(y, y, n), c, n);
               if ((x = *px++) == y) break;
               u = mul_mod(u, abs(y - x), n);
               if (!u) return gcd(v, n);
               if (++t == 32) {
                   t = 0;
                   if ((u = gcd(u, n)) > 1 \&\& u < n) return u;
           if (t \&\& (u = gcd(u, n)) > 1 \&\& u < n) return u;
   }
vector<ll> factorize(ll x) {
   vector<ll> f;
   if (x == 1) return f;
    function < void(11) > dfs = [&](11 x) {
       if (x == 1) return;
       if (x < Sieve::P) {
```

TaylorShift FFT NTT NTTBigModulo

```
auto fs = Sieve::factorize(x);
        f.insert(f.end(), fs.begin(), fs.end());
   } else if (MillerRabin::prime(x)) {
        f.push_back(x);
   } else {
       11 d = PollardRho::rho(x);
       dfs(d);
       dfs(x / d);
};
dfs(x);
sort(f.begin(), f.end());
return f:
```

4.2 Polynomials and recurrences

TaylorShift.cpp

d41d8c, 18 lines

```
template <auto MOD, typename T = Mint<MOD>>
vector<T> shift(vector<T> a, int k) {
   int n = (int)a.size();
   vector<T> fat(n, 1), ifat(n), shifting(n);
   for (int i = 1; i < n; i++) fat[i] = fat[i - 1] * i;
   ifat[n - 1] = T(1) / fat[n - 1];
    for (int i = n - 1; i > 0; i--) ifat[i - 1] = ifat[i] * i;
   for (int i = 0; i < n; i++) a[i] *= fat[i];
   T pk = 1;
   for (int i = 0; i < n; i++) {
       shifting[n - i - 1] = pk * ifat[i];
   auto ans = multiply<MOD>(a, shifting);
   ans.erase(ans.begin(), ans.begin() + n - 1);
   for (int i = 0; i < n; i++) ans[i] *= ifat[i];
   return ans;
```

Optimization

Matrices

Convolutions

FFT.cpp

Description: Computa convolução (multiplicação) de polinômios em $O(N \log N)$, sendo N a soma dos graus dos polinômios. Testado e sem erros de precisão com polinômios de grau até 3.10⁵ e constantes até 10⁶. Time: $\mathcal{O}(N \log N)$.

d41d8c, 67 lines

```
struct base {
    double a. b:
   base (double _a = 0, double _b = 0) : a(_a), b(_b) { }
    const base operator+(const base &c) const { return base(a +
          c.a, b + c.b); }
    const base operator-(const base &c) const { return base(a -
         c.a, b - c.b); }
    const base operator*(const base &c) const {
        return base(a * c.a - b * c.b, a * c.b + b * c.a);
using poly = vector<base>;
const double PI = acos(-1);
void fft(poly &a, bool inv = 0) {
    int n = (int)a.size();
    for (int i = 0; i < n; i++) {
       int bit = n >> 1, j = 0, k = i;
        while (bit > 0) {
```

```
if (k & 1) j += bit;
           k >>= 1, bit >>= 1;
       if (i < j) swap(a[i], a[j]);</pre>
   double angle = 2 * PI / n * (inv ? -1 : 1);
   poly wn(n / 2);
    for (int i = 0; i < n / 2; i++) wn[i] = {cos(angle * i),}
        sin(angle * i) };
    for (int len = 2; len <= n; len <<= 1) {
        int aux = len / 2:
       int step = n / len;
        for (int i = 0; i < n; i += len) {</pre>
            for (int j = 0; j < aux; j++) {
                base v = a[i + j + aux] * wn[step * j];
               a[i + j + aux] = a[i + j] - v;
               a[i + j] = a[i + j] + v;
    for (int i = 0; inv && i < n; i++) a[i].a /= n, a[i].b /= n
vector<ll> multiply(vector<ll> &ta, vector<ll> &tb) {
    int n = (int)ta.size(), m = (int)tb.size();
    int t = n + m - 1, sz = 1;
   while (sz < t) sz <<= 1;
   poly a(sz), b(sz), c(sz);
    for (int i = 0; i < sz; i++) {
       a[i] = i < n ? base((double)ta[i]) : base(0);
       b[i] = i < m ? base((double)tb[i]) : base(0);
    fft(a, 0), fft(b, 0);
    for (int i = 0; i < sz; i++) c[i] = a[i] * b[i];
    fft(c, 1);
   vector<ll> res(sz);
    for (int i = 0; i < sz; i++) res[i] = ll(round(c[i].a));
    while ((int)res.size() > t && res.back() == 0) res.pop back
    return res;
```

NTT.cpp

d41d8c, 37 lines

```
template <auto MOD, typename T = Mint<MOD>>
void ntt(vector<T> &a, bool inv = 0) {
   int n = (int)a.size();
   auto b = a;
   Tq = 1;
   while ((q ^ (MOD / 2)) == 1) q += 1;
   if (inv) g = T(1) / g;
   for (int step = n / 2; step; step /= 2) {
       T w = q ^ (MOD / (n / step)), wn = 1;
       for (int i = 0; i < n / 2; i += step) {
           for (int j = 0; j < step; <math>j++) {
               auto u = a[2 * i + j], v = wn * a[2 * i + j +
               b[i + j] = u + v;
               b[i + n / 2 + j] = u - v;
```

```
wn = wn * w:
        swap(a, b);
    if (inv) {
        T invn = T(1) / n;
        for (int i = 0; i < n; i++) a[i] *= invn;
template <auto MOD, typename T = Mint<MOD>>
vector<T> multiply(vector<T> a, vector<T> b) {
    int n = (int)a.size(), m = (int)b.size();
    int t = n + m - 1, sz = 1;
    while (sz < t) sz <<= 1;
    a.resize(sz), b.resize(sz);
    ntt<MOD>(a, 0), ntt<MOD>(b, 0);
    for (int i = 0; i < sz; i++) a[i] *= b[i];
    ntt < MOD > (a, 1);
    while ((int)a.size() > t) a.pop_back();
    return a;
```

template <auto MOD, typename T = Mint<MOD>>

void ntt(vector<T> &a, bool inv = 0) {

int n = (int)a.size();

NTTBigModulo.cpp

d41d8c, 78 lines

```
auto b = a;
    Tq = 1;
    while ((q ^ (MOD / 2)) == 1) q += 1;
    if (inv) \alpha = T(1) / \alpha;
    for (int step = n / 2; step; step /= 2) {
        T w = q ^ (MOD / (n / step)), wn = 1;
        for (int i = 0; i < n / 2; i += step) {
            for (int j = 0; j < step; <math>j++) {
                auto u = a[2 * i + j], v = wn * a[2 * i + j +
                     stepl:
                b[i + j] = u + v;
                b[i + n / 2 + j] = u - v;
            wn = wn * w:
        swap(a, b);
    if (inv) {
        T invn = T(1) / n;
        for (int i = 0; i < n; i++) a[i] *= invn;
template <auto MOD, typename T = Mint<MOD>>
vector<T> multiply(vector<T> a, vector<T> b) {
    int n = (int)a.size(), m = (int)b.size();
    int t = n + m - 1, sz = 1;
    while (sz < t) sz <<= 1;
    a.resize(sz), b.resize(sz);
    ntt < MOD > (a, 0), ntt < MOD > (b, 0);
    for (int i = 0; i < sz; i++) a[i] *= b[i];
    ntt < MOD > (a, 1);
    while ((int)a.size() > t) a.pop_back();
    return a;
ll extended_gcd(ll a, ll b, ll &x, ll &y) {
    if (b == 0) {
        x = 1;
        y = 0;
```

```
return a:
    } else {
        11 g = extended_gcd(b, a % b, y, x);
        y = a / b * x;
        return q;
11 crt(array<int, 2> rem, array<int, 2> mod) {
    __int128 ans = rem[0], m = mod[0];
   11 g = extended_gcd(mod[1], (11)m, x, y);
    if ((ans - rem[1]) % g != 0) return -1;
    ans = ans + (_int128)1 * (rem[1] - ans) * (m / g) * y;
    m = (\underline{1}nt128) (mod[1] / g) * (m / g) * g;
    ans = (ans % m + m) % m;
    return (11) ans;
template <auto MOD1, auto MOD2, typename T = Mint<MOD1>,
     typename U = Mint<MOD2>>
vector<ll> big_multiply(vector<ll> ta, vector<ll> tb) {
    vector<T> al(ta.size()), bl(tb.size());
    vector<U> a2(ta.size()), b2(tb.size());
    for (int i = 0; i < (int)ta.size(); i++) a1[i] = ta[i];</pre>
    for (int i = 0; i < (int)tb.size(); i++) b1[i] = tb[i];</pre>
    for (int i = 0; i < (int)ta.size(); i++) a2[i] = ta[i];</pre>
    for (int i = 0; i < (int)tb.size(); i++) b2[i] = tb[i];</pre>
    auto c1 = multiply<MOD1>(a1, b1);
    vector<ll> res(c1.size());
    for (int i = 0; i < (int)res.size(); i++)</pre>
        res[i] = crt({c1[i].v, c2[i].v}, {MOD1, MOD2});
    return res;
const int MOD1 = 1004535809;
const int MOD2 = 1092616193;
```

Convolutions.cpp

```
d41d8c, 165 lines
vector<mint> and_convolution(vector<mint> A, vector<mint> B) {
    int n = (int)max(A.size(), B.size());
   int N = 0:
   while ((1 << N) < n) N++;
   A.resize(1 << N);
   B.resize(1 << N);
   vector<mint> C(1 << N);</pre>
    for (int j = 0; j < N; j++) {
        for (int i = (1 << N) - 1; i >= 0; i--) {
           if (~i >> j & 1) {
               A[i] += A[i | (1 << j)];
                B[i] += B[i | (1 << j)];
    for (int i = 0; i < 1 << N; i++) C[i] = A[i] * B[i];
    for (int j = 0; j < N; j++) {
        for (int i = 0; i < 1 << N; i++)
           if (~i >> j & 1) C[i] -= C[i | (1 << j)];
    return C;
vector<mint> gcd_convolution(vector<mint> A, vector<mint> B) {
    int N = (int)max(A.size(), B.size());
   A.resize(N + 1);
   B.resize(N + 1);
   vector<mint> C(N + 1);
    for (int i = 1; i \le N; i++) {
```

```
mint a = 0;
       mint b = 0:
        for (int j = i; j \le N; j += i) {
           a += A[j];
           b += B[j];
       C[i] = a * b;
    for (int i = N; i >= 1; i--)
        for (int j = 2 * i; j \le N; j += i) C[i] -= C[j];
    return C;
vector<mint> lcm_convolution(vector<mint> A, vector<mint> B) {
    int N = (int)max(A.size(), B.size());
   A.resize(N + 1);
   B.resize(N + 1);
    vector<mint> C(N + 1), a(N + 1), b(N + 1);
    for (int i = 1; i <= N; i++) {
        for (int j = i; j \le N; j += i) {
           a[j] += A[i];
           b[j] += B[i];
       C[i] = a[i] * b[i];
    for (int i = 1; i <= N; i++)
        for (int j = 2 * i; j \le N; j += i) C[j] -= C[i];
vector<mint> or convolution(vector<mint> A, vector<mint> B) {
    int n = (int)max(A.size(), B.size());
    int N = 0;
    while ((1 << N) < n) N++;
   A.resize(1 \ll N);
   B.resize(1 << N);
   vector<mint> C(1 << N);</pre>
    for (int j = 0; j < N; j++) {
        for (int i = 0; i < 1 << N; i++) {
            if (i >> j & 1) {
               A[i] += A[i ^ (1 << j)];
                B[i] += B[i ^ (1 << j)];
    for (int i = 0; i < 1 << N; i++) C[i] = A[i] * B[i];
    for (int j = N - 1; j >= 0; j--) {
        for (int i = (1 << N) - 1; i >= 0; i--)
           if (i >> j & 1) C[i] -= C[i ^ (1 << j)];
    return C;
vector<mint> subset convolution(vector<mint> A, vector<mint> B)
    int n = int(max(A.size(), B.size()));
    int N = 0:
    while ((1 << N) < n) N++;
   A.resize(1 << N), B.resize(1 << N);
    vector a(1 << N, vector < mint > (N + 1)), b(1 << N, vector <
        mint > (N + 1));
    for (int i = 0; i < 1 << N; i++) {
       int popcnt = __builtin_popcount(i);
       a[i][popcnt] = A[i];
       b[i][popcnt] = B[i];
    for (int j = 0; j < N; j++) {
        for (int i = 0; i < 1 << N; i++) {
           if (~i >> j & 1) continue;
```

```
for (int popcnt = 0; popcnt <= N; popcnt++) {</pre>
                a[i][popcnt] += a[i ^ (1 << j)][popcnt];
                b[i][popcnt] += b[i ^ (1 << j)][popcnt];
    vector c(1 << N, vector<mint>(N + 1));
    for (int i = 0; i < 1 << N; i++) {
        for (int j = 0; j \le N; j++)
            for (int k = 0; k + j \le N; k++) c[i][j + k] += a[i]
                 ][j] * b[i][k];
    for (int j = N - 1; j >= 0; j--) {
        for (int i = (1 << N) - 1; i >= 0; i--) {
            if (~i >> j & 1) continue;
            for (int popent = 0; popent <= N; popent++)</pre>
                c[i][popent] = c[i ^ (1 << j)][popent];
    vector<mint> ans(1 << N);</pre>
    for (int i = 0; i < 1 << N; i++) {
        int popcnt = __builtin_popcount(i);
        ans[i] = c[i][popcnt];
    return ans;
vector<mint> xor convolution(vector<mint> A, vector<mint> B) {
    int n = int(A.size());
    for (int rep = 0; rep < 2; rep++) {</pre>
        for (int len = n >> 1; len; len >>= 1) {
            for (int i = 0; i < n; i += len << 1) {
                for (int j = 0; j < len; j++) {
                    int id = i + j;
                    mint x = A[id];
                    mint y = A[id + len];
                    A[id] = x + y;
                    A[id + len] = x - y;
        swap(A, B);
    vector<mint> ans(n);
    for (int i = 0; i < n; i++) ans[i] = A[i] * B[i];
    for (int len = 1; len < n; len <<= 1) {
        for (int i = 0; i < n; i += len << 1) {
            for (int j = 0; j < len; j++) {
                int id = i + j;
                mint x = ans[id];
                mint y = ans[id + len];
                ans[id] = x + y;
                ans[id + len] = x - y;
    return ans:
vector<mint> xor multiply(vector<mint> A, vector<mint> B) {
    int N = 1:
    int n = int(max(A.size(), B.size()));
    while (N < n) N <<= 1;
    A resize(N):
    B.resize(N);
    auto ans = xor_convolution(A, B);
    for (int i = 0; i < N; i++) ans[i] /= N;
    return ans;
```

d41d8c, 20 lines

Number theory (5)

5.1 Modular arithmetic

Modular Arithmetic.h.

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

```
d41d8c, 18 lines
const 11 mod = 17; // change to something else
struct Mod {
 11 x;
  Mod(ll xx) : x(xx) \{ \}
  Mod operator+(Mod b) { return Mod((x + b.x) % mod); }
  Mod operator-(Mod b) { return Mod((x - b.x + mod) % mod); }
  Mod operator*(Mod b) { return Mod((x * b.x) % mod); }
  Mod operator/(Mod b) { return *this * invert(b); }
  Mod invert (Mod a) {
   11 x, y, q = euclid(a.x, mod, x, y);
    assert(g == 1); return Mod((x + mod) % mod);
  Mod operator^(ll e) {
   if (!e) return Mod(1);
   Mod r = *this ^ (e / 2); r = r * r;
   return e&1 ? *this * r : r;
};
```

ModInverse.h

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

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

ModPow.h

d41d8c, 8 lines

d41d8c, 11 lines

```
const 11 mod = 1000000007; // faster if const
11 modpow(ll b, ll e) {
 11 \text{ ans} = 1:
  for (; e; b = b * b % mod, e /= 2)
   if (e & 1) ans = ans * b % mod;
  return ans:
```

ModLog.h

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

```
11 modLog(ll a, ll b, ll m) {
 unordered_map<11, 11> A;
 while (j \le n \&\& (e = f = e * a % m) != b % m)
  A[e * b % m] = j++;
 if (e == b % m) return j;
 if (__gcd(m, e) == __gcd(m, b))
   rep(i,2,n+2) if (A.count(e = e * f % m))
    return n * i - A[e];
 return -1;
```

ModSum.h

Description: Sums of mod'ed arithmetic progressions. modsum(to, c, k, m) = $\sum_{i=0}^{\text{to}-1} (ki+c)\%m$. divsum is similar but for

floored division.

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

```
typedef unsigned long long ull;
ull sumsq(ull to) { return to / 2 * ((to-1) | 1); }
ull divsum(ull to, ull c, ull k, ull m) {
 ull res = k / m * sumsq(to) + c / m * to;
 k %= m; c %= m;
 if (!k) return res;
 ull to2 = (to * k + c) / m;
 return res + (to - 1) * to2 - divsum(to2, m-1 - c, m, k);
ll modsum(ull to, ll c, ll k, ll m) {
 c = ((c % m) + m) % m;
 k = ((k \% m) + m) \% m;
 return to * c + k * sumsq(to) - m * divsum(to, c, k, m);
```

ModMulLL.h

Description: Calculate $a \cdot b \mod c$ (or $a^b \mod c$) for $0 < a, b < c < 7.2 \cdot 10^{18}$. **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) {
 ll ret = a * b - M * ull(1.L / M * a * b);
 return ret + M * (ret < 0) - M * (ret >= (11)M);
ull modpow(ull b, ull e, ull mod) {
 ull ans = 1;
 for (; e; b = modmul(b, b, mod), e /= 2)
   if (e & 1) ans = modmul(ans, b, mod);
 return ans;
```

ModSart.h

"ModPow.h"

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

Time: $\mathcal{O}(\log^2 p)$ worst case, $\mathcal{O}(\log p)$ for most p

11 sqrt(ll a, ll p) { a % = p; if (a < 0) a += p;if (a == 0) return 0; assert (modpow(a, (p-1)/2, p) == 1); // else no solution if (p % 4 == 3) return modpow(a, (p+1)/4, p); $// a^{(n+3)/8} \text{ or } 2^{(n+3)/8} * 2^{(n-1)/4} \text{ works if } p \% 8 == 5$ 11 s = p - 1, n = 2;int r = 0, m; while (s % 2 == 0)++r, s /= 2; while (modpow(n, (p-1) / 2, p) != p-1) ++n;11 x = modpow(a, (s + 1) / 2, p);11 b = modpow(a, s, p), g = modpow(n, s, p); for (;; r = m) { 11 t = b; for (m = 0; m < r && t != 1; ++m) t = t * t % p;if (m == 0) return x; 11 gs = modpow(g, 1LL << (r - m - 1), p);g = gs * gs % p;x = x * qs % p;b = b * g % p;

5.2 Primality

FastEratosthenes.h

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

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

MillerRabin.h

return pr;

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

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

```
"ModMulLL.h"
                                                      d41d8c, 12 lines
bool isPrime(ull n) {
 if (n < 2 || n % 6 % 4 != 1) return (n | 1) == 3;
 ull A[] = \{2, 325, 9375, 28178, 450775, 9780504, 1795265022\},
      s = \underline{builtin\_ctzll(n-1)}, d = n >> s;
  for (ull a : A) { // ^ count trailing zeroes
    ull p = modpow(a%n, d, n), i = s;
    while (p != 1 && p != n - 1 && a % n && i--)
      p = modmul(p, p, n);
    if (p != n-1 && i != s) return 0;
 return 1;
```

d41d8c, 24 lines

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

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

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

Divisibility

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

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

CRT.h

Description: Chinese Remainder Theorem.

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

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

5.3.1 Bézout's identity

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

$$ax + by = d$$

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

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

phiFunction.h

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

Euler's thm: $a, n \text{ coprime } \Rightarrow a^{\phi(n)} \equiv 1 \pmod{n}$.

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

d41d8c, 8 lines

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

5.4 Fractions

ContinuedFractions.h

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

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

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

```
typedef double d; // for N \sim 1e7; long double for N \sim 1e9
pair<11, 11> approximate(d x, 11 N) {
 11 LP = 0, LQ = 1, P = 1, Q = 0, inf = LLONG_MAX; d y = x;
    ll lim = min(P ? (N-LP) / P : inf, O ? (N-LO) / O : inf),
       a = (11) floor(y), b = min(a, lim),
       NP = b*P + LP, NQ = b*Q + LQ;
      // If b > a/2, we have a semi-convergent that gives us a
      // better approximation; if b = a/2, we *may* have one.
      // Return {P, Q} here for a more canonical approximation.
      return (abs(x - (d)NP / (d)NQ) < abs(x - (d)P / (d)Q)) ?
        make pair (NP, NO) : make pair (P, O);
    if (abs(y = 1/(y - (d)a)) > 3*N) {
      return {NP, NQ};
    LP = P; P = NP;
    LQ = Q; Q = NQ;
```

FracBinarySearch.h

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

Usage: fracBS([](Frac f) { return f.p>=3*f.q; }, 10); // {1,3} Time: $\mathcal{O}(\log(N))$ d41d8c, 25 lines

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

5.5 Pythagorean Triples

The Pythagorean triples are uniquely generated by

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

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

5.6 Primes

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

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

5.7 Estimates

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

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

Mobius Function

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

Mobius Inversion:

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

Other useful formulas/forms:

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

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

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

Combinatorial (6)

int use = 0, i = 0, r = 0;

use |= 1 << x;

6.1 Permutations

6.1.1 Factorial

		_		•	8	-	10	
n!	1 2 6	24 1	20 720	5040	40320	362880	3628800	
n	11	12	13	14	15	16	17	
							3 3.6e14	
n	20	25	30	40	50 10	00 - 150	171	
n!	2e18	2e25	3e32	$8e47 \ 3$	e64 9e	157 6e26	$62 > DBL_M$	AX

IntPerm.h

return r;

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

Time: $\mathcal{O}(n)$ int permToInt(vi& v) {

for (int x:v) $r = r * ++i + \underline{\quad}$ builtin_popcount (use & -(1<<x)), // (note: minus, not \sim !)

d41d8c, 6 lines

multinomial Dinitz Não Treinamos o João - UDESC 10

6.1.2 Cycles

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

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

6.1.3 Derangements

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

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

6.1.4 Burnside's lemma

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

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

where X^g are the elements fixed by q (q.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

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

Partitions and subsets

6.2.1 Partition function

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

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

6.2.2 Lucas' Theorem

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

6.2.3 Binomials

multinomial.h

rep(i, 1, sz(v)) rep(j, 0, v[i])c = c * ++m / (j+1);return c;

General purpose numbers

6.3.1 Bernoulli numbers

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

Sums of powers:

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

Euler-Maclaurin formula for infinite sums:

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

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

6.3.2 Stirling numbers of the first kind

Number of permutations on n items with k cycles.

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

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

6.3.3 Eulerian numbers

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

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

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

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

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) = $1, 1, 2, 5, 15, 52, 203, 877, 4140, 21147, \dots$ For p prime,

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

6.3.6 Labeled unrooted trees

```
# on n vertices: n^{n-2}
# on k existing trees of size n_i: n_1 n_2 \cdots n_k n^{k-2}
# with degrees d_i: (n-2)!/((d_1-1)!\cdots(d_n-1)!)
```

6.3.7 Catalan numbers

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

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

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

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

Graph (7)

7.1 Fundamentals

7.2Network flow

Dinitz.cpp

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

```
struct dinitz {
 const bool scaling = false; // com \ scaling \rightarrow O(nm \ log(U)),
                               // com constante alta
 int lim:
 struct edge {
   int to, cap, rev, flow;
    edge(int to_, int cap_, int rev_, bool res_)
      : to(to_), cap(cap_), rev(rev_), flow(0), res(res_) {}
 };
 vector<vector<edge>> g;
 vector<int> lev, beg;
 11 F;
 dinitz(int n) : g(n), F(0) {}
 void add(int a, int b, int c) {
    g[a].emplace_back(b, c, g[b].size(), false);
   g[b].emplace_back(a, 0, g[a].size()-1, true);
 bool bfs(int s, int t) {
    lev = vector < int > (g.size(), -1); lev[s] = 0;
   beg = vector<int>(g.size(), 0);
    queue<int> q; q.push(s);
    while (q.size()) {
      int u = q.front(); q.pop();
      for (auto& i : g[u]) {
        if (lev[i.to] != -1 or (i.flow == i.cap)) continue;
        if (scaling and i.cap - i.flow < lim) continue;
       lev[i.to] = lev[u] + 1;
```

MinCostMaxFlow Hungarian

```
q.push(i.to);
   return lev[t] != -1;
  int dfs(int v, int s, int f = INF) {
   if (!f or v == s) return f;
    for (int& i = beg[v]; i < g[v].size(); i++) {</pre>
     auto& e = g[v][i];
     if (lev[e.to] != lev[v] + 1) continue;
     int foi = dfs(e.to, s, min(f, e.cap - e.flow));
     if (!foi) continue;
     e.flow += foi, q[e.to][e.rev].flow -= foi;
     return foi;
    return 0;
  11 max_flow(int s, int t) {
    for (\lim = \text{scaling} ? (1 << 30) : 1; \lim; \lim /= 2)
     while (bfs(s, t)) while (int ff = dfs(s, t)) F += ff;
    return F;
// Recupera as arestas do corte s-t
vector<pair<int, int>> get_cut(dinitz& g, int s, int t) {
 g.max_flow(s, t);
 vector<pair<int, int>> cut;
  vector < int > vis(q.q.size(), 0), st = {s};
  vis[s] = 1;
  while (st.size()) {
   int u = st.back(); st.pop_back();
   for (auto e : g.g[u]) if (!vis[e.to] and e.flow < e.cap)</pre>
     vis[e.to] = 1, st.push_back(e.to);
  for (int i = 0; i < g.g.size(); i++) for (auto e : g.g[i])
    if (vis[i] and !vis[e.to] and !e.res) cut.emplace_back(i, e
  return cut:
```

MinCostMaxFlow.cpp

Description: $min_cost_flow(s,t,f)$ computa o par (fluxo, custo) com $max(fluxo) \le f$ que tenha $min(custo) \ min_cost_flow(s,t)$: Fluxo maximo de custo minimo de s pra t Se for um dag, da pra substituir o SPFA por uma DP pra naopagar O(nm) no comeco Se nao tiver aresta com custo negativo, nao precisa do SPFA

Time: $\mathcal{O}(nm + f * mlogn)$

```
d41d8c, 115 lines
template<typename T> struct mcmf {
  struct edge {
    int to, rev, flow, cap; // para, id da reversa, fluxo,
        capacidade
   bool res; // se eh reversa
   T cost; // custo da unidade de fluxo
    edge(): to(0), rev(0), flow(0), cap(0), cost(0), res(false)
    edge(int to_, int rev_, int flow_, int cap_, T cost_, bool
      : to(to_), rev(rev_), flow(flow_), cap(cap_), res(res_),
          cost(cost_) {}
  };
  vector<vector<edge>> g;
  vector<int> par_idx, par;
  T inf;
  vector<T> dist;
```

```
mcmf(int n) : g(n), par_idx(n), par(n), inf(numeric_limits<T</pre>
    >::max()/3) {}
void add(int u, int v, int w, T cost) { // de u pra v com cap
      w e custo cost
  edge a = edge(v, g[v].size(), 0, w, cost, false);
  edge b = edge(u, g[u].size(), 0, 0, -cost, true);
  g[u].push_back(a);
  g[v].push_back(b);
vector<T> spfa(int s) { // nao precisa se nao tiver custo
     negativo
  deque<int> q;
  vector<bool> is_inside(q.size(), 0);
  dist = vector<T>(g.size(), inf);
  dist[s] = 0;
  q.push_back(s);
  is_inside[s] = true;
  while (!q.emptv()) {
    int v = q.front();
    q.pop_front();
    is_inside[v] = false;
    for (int i = 0; i < q[v].size(); i++) {
      auto [to, rev, flow, cap, res, cost] = g[v][i];
      if (flow < cap and dist[v] + cost < dist[to]) {</pre>
        dist[to] = dist[v] + cost;
        if (is_inside[to]) continue;
        if (!q.empty() and dist[to] > dist[q.front()]) q.
             push back (to);
        else q.push_front(to);
        is_inside[to] = true;
   }
  return dist;
bool dijkstra(int s, int t, vector<T>& pot) {
  priority queue<pair<T, int>, vector<pair<T, int>>, greater
  dist = vector<T>(q.size(), inf);
  dist[s] = 0;
  q.emplace(0, s);
  while (q.size()) {
    auto [d, v] = q.top();
    q.pop();
    if (dist[v] < d) continue;
    for (int i = 0; i < q[v].size(); i++) {
      auto [to, rev, flow, cap, res, cost] = q[v][i];
      cost += pot[v] - pot[to];
      if (flow < cap and dist[v] + cost < dist[to]) {</pre>
        dist[to] = dist[v] + cost;
        g.emplace(dist[to], to);
        par_idx[to] = i, par[to] = v;
  return dist[t] < inf;</pre>
pair<int, T> min_cost_flow(int s, int t, int flow = INF) {
  vector<T> pot(q.size(), 0);
  pot = spfa(s); // mudar algoritmo de caminho minimo aqui
```

```
int f = 0;
    T ret = 0:
    while (f < flow and dijkstra(s, t, pot)) {</pre>
      for (int i = 0; i < g.size(); i++)</pre>
        if (dist[i] < inf) pot[i] += dist[i];</pre>
      int mn_flow = flow - f, u = t;
      while (u != s) {
        mn_flow = min(mn_flow,
          g[par[u]][par_idx[u]].cap - g[par[u]][par_idx[u]].
               flow);
        u = par[u];
      ret += pot[t] * mn_flow;
      u = t;
      while (u != s) {
        g[par[u]][par_idx[u]].flow += mn_flow;
        g[u][g[par[u]][par_idx[u]].rev].flow -= mn_flow;
        u = par[u];
      f += mn_flow;
    return make_pair(f, ret);
  // Opcional: retorna as arestas originais por onde passa flow
  vector<pair<int,int>> recover() {
    vector<pair<int,int>> used;
    for (int i = 0; i < g.size(); i++) for (edge e : g[i])
      if(e.flow == e.cap && !e.res) used.push_back({i, e.to});
    return used;
};
```

7.3 Matching

Hungarian.cpp

Description: Resolve o problema de assignment (matriz n por n). Colocar os valores da matriz em 'a' (pode < 0). assignment() retorna um par com o valor do assignment minimo, e a coluna escolhida por cada linha Time: $\mathcal{O}\left(n^3\right)$.

```
d41d8c, 42 lines
template<typename T> struct hungarian {
 int n;
 vector<vector<T>> a;
 vector<T> u, v;
 vector<int> p, way;
 hungarian(int n_{-}): n(n_{-}), u(n+1), v(n+1), p(n+1), way(n+1) {
   a = vector<vector<T>>(n, vector<T>(n));
   inf = numeric_limits<T>::max();
 pair<T, vector<int>> assignment() {
   for (int i = 1; i <= n; i++) {
     p[0] = i;
     int j0 = 0;
     vector<T> minv(n+1, inf);
     vector<int> used(n+1, 0);
       used[j0] = true;
       int i0 = p[j0], j1 = -1;
       T delta = inf;
        for (int j = 1; j \le n; j++) if (!used[j]) {
         T cur = a[i0-1][j-1] - u[i0] - v[j];
```

```
if (cur < minv[j]) minv[j] = cur, way[j] = j0;</pre>
          if (minv[j] < delta) delta = minv[j], j1 = j;</pre>
        for (int j = 0; j \le n; j++)
          if (used[j]) u[p[j]] += delta, v[j] -= delta;
          else minv[j] -= delta;
        j0 = j1;
      } while (p[j0] != 0);
      do {
        int j1 = way[j0];
        p[j0] = p[j1];
        j0 = j1;
     } while (j0);
    vector<int> ans(n);
    for (int j = 1; j \le n; j++) ans[p[j]-1] = j-1;
    return make_pair(-v[0], ans);
};
```

Blossom.cpp

```
Description: Maximum matching in general graph.
Time: \mathcal{O}(n^3). If bipartite, \mathcal{O}(nm) and doesn't need 'contract' _{\text{d41d8c, 77 lines}}^{\text{times}}
vector<int> q[MAX];
int match[MAX]; // match[i] = com quem i esta matchzado ou -1
int n, pai[MAX], base[MAX], vis[MAX];
queue<int> q;
void contract(int u, int v, bool first = 1) {
  static vector<bool> bloss;
  static int 1;
  if (first) {
   bloss = vector<bool>(n, 0);
    vector<bool> teve(n, 0);
    int k = u; l = v;
    while (1) {
     teve[k = base[k]] = 1;
     if (match[k] == -1) break;
     k = pai[match[k]];
    while (!teve[l = base[l]]) l = pai[match[l]];
  while (base[u] != 1) {
   bloss[base[u]] = bloss[base[match[u]]] = 1;
   pai[u] = v;
   v = match[u];
    u = pai[match[u]];
  if (!first) return;
  contract(v, u, 0);
  for (int i = 0; i < n; i++) if (bloss[base[i]]) {</pre>
    base[i] = 1;
    if (!vis[i]) q.push(i);
    vis[i] = 1;
int getpath(int s) {
  for (int i = 0; i < n; i++) base[i] = i, pai[i] = -1, vis[i]
  vis[s] = 1; q = queue < int > (); q.push(s);
  while (q.size()) {
    int u = q.front(); q.pop();
    for (int i : q[u]) {
      if (base[i] == base[u] or match[u] == i) continue;
      if (i == s or (match[i] != -1 and pai[match[i]] != -1))
        contract(u, i);
      else if (pai[i] == -1) {
```

```
pai[i] = u;
       if (match[i] == -1) return i;
       i = match[i];
       vis[i] = 1; q.push(i);
 return -1;
int blossom() {
 int ans = 0;
 memset(match, -1, sizeof(match));
 for (int i = 0; i < n; i++) if (match[i] == -1)
   for (int j : g[i]) if (match[j] == -1) {
     match[i] = j;
     match[j] = i;
     ans++;
     break:
 for (int i = 0; i < n; i++) if (match[i] == -1) {
   int j = getpath(i);
   if (j == -1) continue;
   ans++;
   while (j != -1) {
     int p = pai[j], pp = match[p];
     match[p] = j;
     match[j] = p;
     j = pp;
 return ans;
```

7.4 DFS algorithms

7.5 Coloring

7.6 Heuristics

7.7 Trees

```
LinkCutTree.cpp
```

```
d41d8c, 208 lines
const int N = 1e5 + 9;
struct node {
 int p = 0, c[2] = \{0, 0\}, pp = 0;
 bool flip = 0;
 int sz = 0, ssz = 0, vsz = 0; // sz \Rightarrow aux tree size, ssz =
       subtree\ size\ in\ rep\ tree\ ,\ vsz=virtual\ tree\ size
 long long val = 0, sum = 0, lazy = 0, subsum = 0, vsum = 0;
 node() {}
 node(int x) {
   val = x; sum = x;
   sz = 1; lazv = 0;
   ssz = 1; vsz = 0;
   subsum = x; vsum = 0;
struct LCT {
 vector<node> t;
 LCT() {}
 LCT(int n) : t(n + 1) {}
 // <independent splay tree code>
 int dir(int x, int y) { return t[x].c[1] == y; }
 void set(int x, int d, int y) {
   if (x) t[x].c[d] = y, pull(x);
   if (y) t[y].p = x;
```

```
void pull(int x) {
 if (!x) return;
  int &1 = t[x].c[0], &r = t[x].c[1];
  push(1); push(r);
  t[x].sum = t[1].sum + t[r].sum + t[x].val;
  t[x].sz = t[1].sz + t[r].sz + 1;
  t[x].ssz = t[1].ssz + t[r].ssz + t[x].vsz + 1;
  t[x].subsum = t[1].subsum + t[r].subsum + t[x].vsum + t[x].
void push(int x) {
 if (!x) return;
  int &1 = t[x].c[0], &r = t[x].c[1];
  if (t[x].flip) {
   swap(1, r);
   if (1) t[1].flip ^= 1;
   if (r) t[r].flip ^= 1;
    t[x].flip = 0;
  if (t[x].lazy) {
   t[x].val += t[x].lazy;
    t[x].sum += t[x].lazy * t[x].sz;
    t[x].subsum += t[x].lazy * t[x].ssz;
    t[x].vsum += t[x].lazy * t[x].vsz;
    if (1) t[1].lazy += t[x].lazy;
    if (r) t[r].lazy += t[x].lazy;
    t[x].lazy = 0;
void rotate(int x, int d) {
 int y = t[x].p, z = t[y].p, w = t[x].c[d];
  swap(t[x].pp, t[y].pp);
  set(y, !d, w);
  set(x, d, y);
  set(z, dir(z, y), x);
void splay(int x) {
  for (push(x); t[x].p;) {
    int y = t[x].p, z = t[y].p;
    push(z); push(y); push(x);
    int dx = dir(y, x), dy = dir(z, y);
    if (!z) rotate(x, !dx);
    else if (dx == dy) rotate(y, !dx), rotate(x, !dx);
    else rotate(x, dy), rotate(x, dx);
// </independent splay tree code>
// making it a root in the rep. tree
void make_root(int u) {
  access(u);
  int 1 = t[u].c[0];
  t[1].flip ^= 1;
  swap(t[1].p, t[1].pp);
  t[u].vsz += t[l].ssz;
  t[u].vsum += t[1].subsum;
  set(u, 0, 0);
// make the path from root to u a preferred path
// returns last path-parent of a node as it moves up the tree
int access(int _u) {
 int last = _u;
  for (int v = 0, u = _u; u; u = t[v = u].pp) {
    splay(u); splay(v);
    t[u].vsz -= t[v].ssz;
    t[u].vsum -= t[v].subsum;
    int r = t[u].c[1];
    t[u].vsz += t[r].ssz;
```

12

cut(u);

int ans = component_size(u);

```
t[u].vsum += t[r].subsum;
   t[v].pp = 0;
   swap(t[r].p, t[r].pp);
   set(u, 1, v);
   last = u;
 splay(_u);
 return last;
void link(int u, int v) { // u \rightarrow v
  // assert (!connected(u, v));
 make root(v);
 access(u); splay(u);
 t[v].pp = u;
 t[u].vsz += t[v].ssz;
 t[u].vsum += t[v].subsum;
void cut(int u) { // cut par[u] \Rightarrow u, u is non root vertex
 access(u);
 assert(t[u].c[0] != 0);
 t[t[u].c[0]].p = 0;
 t[u].c[0] = 0;
 pull(u);
// parent of u in the rep. tree
int get_parent(int u) {
 access(u); splay(u); push(u);
 u = t[u].c[0]; push(u);
 while (t[u].c[1]) {
   u = t[u].c[1]; push(u);
 splay(u);
 return u;
// root of the rep. tree containing this node
int find_root(int u) {
 access(u); splay(u); push(u);
 while (t[u].c[0]) {
   u = t[u].c[0]; push(u);
 splay(u);
 return u;
bool connected(int u, int v) {
  return find_root(u) == find_root(v);
// depth in the rep. tree
int depth(int u) {
 access(u); splay(u);
 return t[u].sz;
int lca(int u, int v) {
  // assert(connected(u, v));
 if (u == v) return u;
 if (depth(u) > depth(v)) swap(u, v);
 access(v);
 return access(u);
int is_root(int u) {
 return get_parent(u) == 0;
int component_size(int u) {
 return t[find root(u)].ssz;
int subtree_size(int u) {
 int p = get_parent(u);
 if (p == 0) {
   return component_size(u);
```

```
link(p, u);
   return ans;
 long long component_sum(int u) {
   return t[find_root(u)].subsum;
 long long subtree_sum(int u) {
   int p = get_parent(u);
   if (p == 0) {
     return component_sum(u);
    cut(u);
    long long ans = component_sum(u);
   link(p, u);
    return ans;
  // sum of the subtree of u when root is specified
 long long subtree_query(int u, int root) {
    int cur = find_root(u);
   make_root(root);
   long long ans = subtree sum(u);
   make_root(cur);
   return ans;
 // path sum
 long long query(int u, int v) {
   int cur = find_root(u);
   make_root(u); access(v);
   long long ans = t[v].sum;
   make_root(cur);
    return ans;
 void upd(int u, int x) {
   access(u); splay(u);
    t[u].val += x;
 // add x to the nodes on the path from u to v
 void upd(int u, int v, int x) {
   int cur = find_root(u);
   make root(u); access(v);
   t[v].lazy += x;
   make root(cur);
}t[2];
Block-Cut-Tree.cpp
Description: Block cut tree
Time: \mathcal{O}(n+m)
                                                     d41d8c, 65 lines
struct Bct {
   vector<int> tin, low, stk, art, id, splits;
   vector<vector<int>> adj, g, comp, up;
   int n, sz, m;
   void build(int _n, int _m) {
       n = _n, m = _m;
       adj.resize(n);
   void add_edge(int u, int v) {
       adj[u].emplace_back(v);
       adj[v].emplace_back(u);
   void dfs(int u, int p) {
       low[u] = tin[u] = ++T;
       stk.emplace_back(u);
        for (auto v : adj[u])
            if (tin[v] == -1) {
```

```
dfs(v, u);
                low[u] = min(low[u], low[v]);
                if (low[v] >= tin[u]) {
                    int x;
                    sz++;
                    do {
                        assert(stk.size());
                        x = stk.back();
                        stk.pop_back();
                        comp[x].emplace_back(sz);
                    } while (x != v);
                    comp[u].emplace_back(sz);
            } else if (v != p) {
                low[u] = min(low[u], tin[v]);
    inline bool is_articulation_point(int u) { return art[id[u
    inline int number_of_splits(int u) { return splits[id[u]];
    void work() {
       T = sz = 0;
        stk.clear();
        tin.resize(n, -1);
        comp.resize(n);
       low.resize(n);
        for (int i = 0; i < n; i++)
            if (tin[i] == -1) dfs(i, 0);
        art.resize(sz + n + 1);
        splits.resize(n + sz + 1, 1);
        id.resize(n);
        g.resize(sz + n + 1);
        for (int i = 0; i < n; i++) {
            if ((int)comp[i].size() > 1) {
                id[i] = ++sz;
                art[id[i]] = 1;
                splits[id[i]] = (int)comp[i].size();
                for (auto u : comp[i]) {
                    g[id[i]].emplace_back(u);
                    g[u].emplace_back(id[i]);
            } else if (comp[i].size()) {
                id[i] = comp[i][0];
       }
};
```

13

7.8 Math

7.8.1 Number of Spanning Trees

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

7.8.2 Erdős–Gallai theorem

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

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

Geometry (8)

8.1 Geometric primitives

Point.h

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

```
template \langle class T \rangle int sgn(T x) \{ return (x > 0) - (x < 0); \}
template<class T>
struct Point {
 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); }</pre>
  bool operator==(P p) const { return tie(x,y)==tie(p.x,p.y); }
  P operator+(P p) const { return P(x+p.x, y+p.y); }
  P operator-(P p) const { return P(x-p.x, y-p.y); }
  P operator*(T d) const { return P(x*d, y*d); }
  P operator/(T d) const { return P(x/d, y/d); }
  T dot(P p) const { return x*p.x + y*p.y; }
  T cross(P p) const { return x*p.y - y*p.x; }
  T cross(P a, P b) const { return (a-*this).cross(b-*this); }
  T dist2() const { return x*x + v*v; }
  double dist() const { return sqrt((double)dist2()); }
  // angle to x-axis in interval [-pi, pi]
  double angle() const { return atan2(y, x); }
  P unit() const { return *this/dist(); } // makes dist()=1
  P perp() const { return P(-y, x); } // rotates +90 degrees
  P normal() const { return perp().unit(); }
  // returns point rotated 'a' radians ccw around the origin
  P rotate(double a) const {
   return P(x*cos(a)-y*sin(a),x*sin(a)+y*cos(a)); }
  friend ostream& operator<<(ostream& os, P p) {</pre>
   return os << "(" << p.x << "," << p.v << ")"; }
```

lineDistance.h

Description:

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



```
d41d8c, 4 lines
template<class P>
double lineDist(const P& a, const P& b, const P& p) {
  return (double) (b-a).cross(p-a)/(b-a).dist();
```

SegmentDistance.h Description:

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



```
Usage: Point < double > a, b(2,2), p(1,1);
bool onSegment = segDist(a,b,p) < 1e-10;
                                                       d41d8c, 6 lines
typedef Point < double > P;
double segDist(P& s, P& e, P& p) {
 if (s==e) return (p-s).dist();
 auto d = (e-s) . dist2(), t = min(d, max(.0, (p-s) . dot(e-s)));
 return ((p-s)*d-(e-s)*t).dist()/d;
```

SegmentIntersection.h

Description:

If a unique intersection point between the line segments going from s1 to e1 and from s2 to e2 exists then it is returned. If no intersection point exists an empty vector is returned. If infinitely many exist a vector with 2 elements is returned, containing the endpoints of the common line segment. The wrong position will be returned if P is Point<|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 = seqInter(s1,e1,s2,e2);
if (sz(inter) == 1)
```

```
cout << "segments intersect at " << inter[0] << endl;</pre>
"Point.h", "OnSegment.h"
                                                      d41d8c, 13 lines
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 {all(s)};
```

lineIntersection.h

Description:

If a unique intersection point of the lines going through s1,e1 and s2,e2 exists {1, point} is returned. If no intersection point exists $\{0, (0,0)\}$ is returned and if infinitely many exists $\{-1, e^2\}$ (0,0)} is returned. The wrong position will be returned if P is Point<ll> and the intersection point does not have integer coordinates. Products of three coordinates are used in inter- 1 mediate steps so watch out for overflow if using int or ll.

```
Usage: auto res = lineInter(s1,e1,s2,e2);
```

```
if (res.first == 1)
cout << "intersection point at " << res.second << endl;</pre>
"Point.h"
                                                         d41d8c, 8 lines
template<class P>
```

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

sideOf.h

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

```
Usage: bool left = sideOf(p1,p2,q)==1;
"Point.h"
                                                      d41d8c, 9 lines
template<class P>
int sideOf(P s, P e, P p) { return sqn(s.cross(e, p)); }
template<class P>
int sideOf(const P& s, const P& e, const P& p, double eps) {
 auto a = (e-s).cross(p-s);
 double l = (e-s).dist()*eps;
 return (a > 1) - (a < -1);
```

OnSegment.h

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

```
d41d8c, 3 lines
template < class P > bool on Segment (P s, P e, P p) {
 return p.cross(s, e) == 0 \&\& (s - p).dot(e - p) <= 0;
```

linearTransformation.h Description:

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



```
d41d8c, 6 lines
typedef Point < double > P:
P linearTransformation(const P& p0, const P& p1,
    const P& q0, const P& q1, const P& r) {
  P dp = p1-p0, dq = q1-q0, num(dp.cross(dq), dp.dot(dq));
  return q0 + P((r-p0).cross(num), (r-p0).dot(num))/dp.dist2();
```

Angle.h

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

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

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

8.2 Circles

CircleIntersection.h

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

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.

CirclePolygonIntersection.h

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

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

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

circumcircle.h

Description:

"Point.h"

The circumcirle of a triangle is the circle intersecting all three vertices. ccRadius returns the radius of the circle going through points A, B and C and ccCenter returns the center of the same circle.



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

MinimumEnclosingCircle.h

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

return A + (b*c.dist2()-c*b.dist2()).perp()/b.cross(c)/2;

8.3 Polygons

InsidePolygon.h

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

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

PolygonArea.

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

PolygonCenter.h

Description: Returns the center of mass for a polygon.

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

PolygonCut.h Description:

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



d41d8c, 13 lines

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

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

ConvexHull.h

return res;

Description:

"Point.h"

Returns a vector of the points of the convex hull in counterclockwise order. Points on the edge of the hull between two other points are not considered part of the hull. $\mathbf{Time}: \mathcal{O}(n\log n)$



d41d8c, 13 lines

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

HullDiameter.h

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

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

```
"Point.h"
                                                      d41d8c, 12 lines
typedef Point<11> P;
array<P, 2> hullDiameter(vector<P> S) {
 int n = sz(S), j = n < 2 ? 0 : 1;
  pair<11, array<P, 2>> res({0, {S[0], S[0]}});
  rep(i,0,j)
    for (;; j = (j + 1) % n) {
      res = \max(\text{res}, \{(S[i] - S[j]).dist2(), \{S[i], S[j]\}\});
      if ((S[(j+1) % n] - S[j]).cross(S[i+1] - S[i]) >= 0)
  return res.second;
```

PointInsideHull.h

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

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

```
d41d8c, 14 lines
"Point.h", "sideOf.h", "OnSegment.h"
typedef Point<11> P;
bool inHull(const vector<P>& 1, P p, bool strict = true) {
  int a = 1, b = sz(1) - 1, r = !strict;
  if (sz(1) < 3) return r && onSegment(1[0], 1.back(), p);</pre>
  if (sideOf(1[0], 1[a], 1[b]) > 0) swap(a, b);
  if (sideOf(1[0], 1[a], p) >= r || sideOf(1[0], 1[b], p) <= -r)
    return false:
  while (abs(a - b) > 1) {
    int c = (a + b) / 2;
    (sideOf(1[0], 1[c], p) > 0 ? b : a) = c;
  return sqn(l[a].cross(l[b], p)) < r;</pre>
```

LineHullIntersection.h

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

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

```
int endB = extrVertex(poly, (b - a).perp());
if (cmpL(endA) < 0 \mid | cmpL(endB) > 0)
  return {-1, -1};
array<int, 2> res;
rep(i, 0, 2) {
  int lo = endB, hi = endA, n = sz(poly);
  while ((lo + 1) % n != hi) {
    int m = ((lo + hi + (lo < hi ? 0 : n)) / 2) % n;
    (cmpL(m) == cmpL(endB) ? lo : hi) = m;
  res[i] = (lo + !cmpL(hi)) % n;
  swap (endA, endB);
if (res[0] == res[1]) return {res[0], -1};
if (!cmpL(res[0]) && !cmpL(res[1]))
  switch ((res[0] - res[1] + sz(poly) + 1) % sz(poly)) {
    case 0: return {res[0], res[0]};
    case 2: return {res[1], res[1]};
return res;
```

8.4 Misc. Point Set Problems

ClosestPair.h

Description: Finds the closest pair of points.

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

```
"Point.h"
                                                      d41d8c, 17 lines
typedef Point<ll> P;
pair<P, P> closest (vector<P> v) {
 assert(sz(v) > 1);
  set<P> S;
  sort(all(v), [](P a, P b) { return a.y < b.y; });</pre>
  pair<11, pair<P, P>> ret{LLONG_MAX, {P(), P()}};
  int j = 0;
  for (P p : v) {
   P d{1 + (ll)sgrt(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"
                                                      d41d8c, 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>
struct Node {
 P pt; // if this is a leaf, the single point in it
 T x0 = INF, x1 = -INF, y0 = INF, y1 = -INF; // bounds
 Node *first = 0, *second = 0;
 T distance (const P& p) { // min squared distance to a point
   T x = (p.x < x0 ? x0 : p.x > x1 ? x1 : p.x);
    T y = (p.y < y0 ? y0 : p.y > y1 ? y1 : p.y);
    return (P(x,y) - p).dist2();
  Node (vector < P > && vp) : pt (vp[0]) {
    for (P p : vp) {
```

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

FastDelaunav.h

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

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

```
"Point.h"
typedef Point<11> 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 {
  Q rot, o; P p = arb; bool mark;
  P& F() { return r()->p; }
  Q& r() { return rot->rot; }
  Q prev() { return rot->o->rot; }
  Q next() { return r()->prev(); }
bool circ(P p, P a, P b, P c) { // is p in the circumcircle?
  111 p2 = p.dist2(), A = a.dist2()-p2,
      B = b.dist2()-p2, C = c.dist2()-p2;
  return p.cross(a,b) *C + p.cross(b,c) *A + p.cross(c,a) *B > 0;
```

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

8.5 3D

PolyhedronVolume.h

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

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

Point3D.h

Description: Class to handle points in 3D space. T can be e.g. double or long long. $\frac{\text{d41d8c, 32 lines}}{\text{d42d8c, 32 lines}}$

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

3dHull.h

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

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

```
vector<F> hull3d(const vector<P3>& A) {
  assert(sz(A) >= 4);
  \label{eq:vector} \text{vector} < \text{PR} >> \ \texttt{E} \left( \text{sz} \left( \texttt{A} \right) \text{, vector} < \text{PR} > \left( \text{sz} \left( \texttt{A} \right) \text{, } \left\{ -1 \text{, } -1 \right\} \right) \right) \text{;}
#define E(x,y) E[f.x][f.y]
  vector<F> FS;
  auto mf = [\&] (int i, int j, int k, int l) {
    P3 q = (A[j] - A[i]).cross((A[k] - A[i]));
    if (q.dot(A[1]) > q.dot(A[i]))
       q = q * -1;
    F f{q, i, j, k};
    E(a,b).ins(k); E(a,c).ins(j); E(b,c).ins(i);
    FS.push_back(f);
  rep(i,0,4) rep(j,i+1,4) rep(k,j+1,4)
    mf(i, j, k, 6 - i - j - k);
  rep(i,4,sz(A)) {
     rep(j, 0, sz(FS)) {
      F f = FS[j];
       if(f.q.dot(A[i]) > f.q.dot(A[f.a])) {
         E(a,b).rem(f.c);
         E(a,c).rem(f.b);
         E(b,c).rem(f.a);
         swap(FS[j--], FS.back());
         FS.pop_back();
    int nw = sz(FS);
     rep(j,0,nw) {
       F f = FS[i];
\#define C(a, b, c) if (E(a,b).cnt() != 2) mf(f.a, f.b, i, f.c);
       C(a, b, c); C(a, c, b); C(b, c, a);
  for (F& it : FS) if ((A[it.b] - A[it.a]).cross(
    A[it.c] - A[it.a]).dot(it.q) <= 0) swap(it.c, it.b);
  return FS:
```

sphericalDistance.h

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

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

Strings (9)

AhoCorasick.cpp

Description: query retorna o somatorio do numero de matches de todas as stringuinhas na stringona

d41d8c, 40 lines

```
Time: insert in \mathcal{O}(|s| \log SIGMA).
```

```
namespace aho {
  map<char, int> to[MAX];
```

SuffixArray Manacher Lyndon IntervalContainer

```
int link[MAX], idx, term[MAX], exit[MAX], sobe[MAX];
  void insert(string& s) {
   int at = 0;
    for (char c : s) {
     auto it = to[at].find(c);
     if (it == to[at].end()) at = to[at][c] = ++idx;
     else at = it->second;
   term[at]++, sobe[at]++;
#warning nao esquece de chamar build() depois de inserir
 void build() {
   queue<int> q;
   q.push(0);
   link[0] = exit[0] = -1;
   while (q.size()) {
     int i = q.front(); q.pop();
     for (auto [c, j] : to[i]) {
       int 1 = link[i];
       while (l != -1 \text{ and } !to[l].count(c)) l = link[l];
       link[j] = 1 == -1 ? 0 : to[1][c];
       exit[j] = term[link[j]] ? link[j] : exit[link[j]];
       if (exit[j]+1) sobe[j] += sobe[exit[j]];
        q.push(j);
  int query(string& s) {
   int at = 0, ans = 0;
    for (char c : s) {
     while (at != -1 and !to[at].count(c)) at = link[at];
     at = at == -1 ? 0 : to[at][c];
     ans += sobe[at];
   return ans;
```

SuffixArray.cpp

Description: kasai recebe o suffix array e calcula lcp[i], o lcp entre s[sa[i], ..., n-1] e s[sa[i+1], ..., n-1].

```
Time: \mathcal{O}(N \log N). Kasai in \mathcal{O}(N).
                                                       d41d8c, 34 lines
vector<int> suffix_array(string s) {
 int n = s.size(), N = max(n, 260);
  vector<int> sa(n), ra(n);
  for (int i = 0; i < n; i++) sa[i] = i, ra[i] = s[i];
  for (int k = 0; k < n; k ? k *= 2 : k++) {
   vector<int> nsa(sa), nra(n), cnt(N);
    for (int i = 0; i < n; i++) nsa[i] = (nsa[i]-k+n)%n, cnt[ra[
        i]]++;
    for(int i = 1; i < N; i++) cnt[i] += cnt[i-1];</pre>
    for (int i = n-1; i+1; i--) sa[--cnt[ra[nsa[i]]]] = nsa[i];
    for(int i = 1, r = 0; i < n; i++) nra[sa[i]] = r += ra[sa[i]]
        ]] !=
     ra[sa[i-1]] or ra[(sa[i]+k)%n] != ra[(sa[i-1]+k)%n];
    ra = nra:
    if (ra[sa[n-1]] == n-1) break;
  return vector<int>(sa.begin()+1, sa.end());
vector<int> kasai(string s, vector<int> sa) {
 int n = s.size(), k = 0;
```

```
vector<int> ra(n), lcp(n);
for (int i = 0; i < n; i++) ra[sa[i]] = i;

for (int i = 0; i < n; i++, k -= !!k) {
   if (ra[i] == n-1) { k = 0; continue; }
   int j = sa[ra[i]+1];
   while (i+k < n and j+k < n and s[i+k] == s[j+k]) k++;
   lcp[ra[i]] = k;
}
return lcp;
}</pre>
```

Manacher.cpp

Description: manacher recebe um vetor de T e retorna o vetor com tamanho dos palindromos ret $[2^*i]$ = tamanho do maior palindromo centrado em i ret $[2^*i+1]$ = tamanho maior palindromo centrado em i e i+1 **Time:** all in $\mathcal{O}(N)$.

```
d41d8c, 43 lines
template<typename T> vector<int> manacher(const T& s) {
 int 1 = 0, r = -1, n = s.size();
 vector < int > d1(n), d2(n);
 for (int i = 0; i < n; i++) {
   int k = i > r ? 1 : min(d1[1+r-i], r-i);
   while (i+k < n \&\& i-k >= 0 \&\& s[i+k] == s[i-k]) k++;
   d1[i] = k--;
   if (i+k > r) l = i-k, r = i+k;
 1 = 0, r = -1;
 for (int i = 0; i < n; i++) {
    int k = i > r ? 0 : min(d2[1+r-i+1], r-i+1); k++;
    while (i+k \le n \&\& i-k >= 0 \&\& s[i+k-1] == s[i-k]) k++;
    d2[i] = --k;
   if (i+k-1 > r) 1 = i-k, r = i+k-1;
 vector<int> ret(2*n-1);
 for (int i = 0; i < n; i++) ret[2*i] = 2*d1[i]-1;
 for (int i = 0; i < n-1; i++) ret[2*i+1] = 2*d2[i+1];
 return ret;
// verifica se a string s[i..j] eh palindromo
template<typename T> struct palindrome {
 vector<int> man;
 palindrome(const T& s) : man(manacher(s)) {}
 bool query(int i, int j) {
   return man[i+j] >= j-i+1;
};
// tamanho do maior palindromo que termina em cada posicao
template<typename T> vector<int> pal end(const T& s) {
 vector<int> ret(s.size());
 palindrome<T> p(s);
 ret[0] = 1;
 for (int i = 1; i < s.size(); i++) {
   ret[i] = min(ret[i-1]+2, i+1);
   while (!p.query(i-ret[i]+1, i)) ret[i]--;
 return ret:
```

Lyndon.cpp

int n = s.size();

Description: Duval algorithm to find the Lyndon factorization of a string. Also contains a function to find the minimum cyclic shift of a string. **Time:** All $\mathcal{O}(n)$

```
vector<string> duval(string const &s) {
```

```
vector<string> factorization;
    while (i < n) {
        int j = i + 1, k = i;
        while (j < n \&\& s[k] \le s[j]) {
            if (s[k] < s[j]) k = i;
            else k++;
            j++;
        while (i \le k) {
            factorization.push_back(s.substr(i, j - k));
            i += j - k;
    return factorization;
string min_cyclic_shift(string s) {
    s += s;
    int n = s.size();
    int i = 0, ans = 0;
    while (i < n / 2) {
        ans = i;
        int j = i + 1, k = i;
        while (j < n \&\& s[k] <= s[j]) {
            if (s[k] < s[j]) k = i;
            else k++;
            j++;
        while (i \le k) i += j - k;
    return s.substr(ans, n / 2);
```

Various (10)

int i = 0;

10.1 Intervals

IntervalContainer.h

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

```
Time: \mathcal{O}(\log N)
                                                     d41d8c, 23 lines
set<pii>::iterator addInterval(set<pii>& is, int L, int R) {
 if (L == R) return is.end();
 auto it = is.lower_bound({L, R}), before = it;
  while (it != is.end() && it->first <= R) {
    R = max(R, it->second);
    before = it = is.erase(it);
 if (it != is.begin() && (--it)->second >= L) {
   L = min(L, it->first);
    R = max(R, it->second);
    is.erase(it);
 return is.insert(before, {L,R});
void removeInterval(set<pii>& is, int L, int R) {
 if (L == R) return;
  auto it = addInterval(is, L, R);
  auto r2 = it->second;
 if (it->first == L) is.erase(it);
  else (int&)it->second = L;
 if (R != r2) is.emplace (R, r2);
```

IntervalCover.h

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

Time: $\mathcal{O}(N \log N)$ d41d8c, 19 lines

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

ConstantIntervals.h

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

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

```
Time: \mathcal{O}\left(k\log\frac{n}{k}\right)
                                                                                 d41d8c, 19 lines
template<class F, class G, class T>
```

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

10.2 Misc. algorithms

TernarySearch.h

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

```
Usage: int ind = ternSearch(0, n-1, [&](int i){return a[i];});
Time: \mathcal{O}(\log(b-a))
                                                             d41d8c, 11 lines
```

```
template<class F>
int ternSearch(int a, int b, F f) {
  assert(a <= b);
 while (b - a >= 5) {
   int mid = (a + b) / 2;
```

```
if (f(mid) < f(mid+1)) a = mid; //(A)
  else b = mid+1;
rep(i,a+1,b+1) if (f(a) < f(i)) a = i; // (B)
```

LIS.h

Description: Compute indices for the longest increasing subsequence. Time: $\mathcal{O}(N \log N)$ d41d8c, 17 lines

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

FastKnapsack.h

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

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

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

10.3 Dynamic programming

KnuthDP.h

Description: When doing DP on intervals: $a[i][j] = \min_{i < k < j} (a[i][k] + a[i][k])$ 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 $f(b,c) \leq f(a,d)$ and $f(a,c) + f(b,d) \le f(a,d) + f(b,c)$ for all $a \le b \le c \le d$. Consider also: LineContainer (ch. Data structures), monotone queues, ternary search. Time: $\mathcal{O}\left(N^2\right)$

DivideAndConquerDP.h

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

```
d41d8c, 18 lines | FastMod.h
```

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

10.4 Debugging tricks

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

10.5 Optimization tricks

builtin ia32 ldmxcsr(40896); disables denormals (which make floats 20x slower near their minimum value).

10.5.1 Bit backs

d41d8c, 16 lines

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

10.5.2 Pragmas

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

Não Treinamos o João - UDESC

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

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

FastInput.h

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

Usage: ./a.out < input.txt

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

d41d8c, 17 lines

```
inline char gc() { // like getchar()
  static char buf[1 << 16];</pre>
  static size_t bc, be;
  if (bc >= be) {
   buf[0] = 0, bc = 0;
   be = fread(buf, 1, sizeof(buf), stdin);
  return buf[bc++]; // returns 0 on EOF
int readInt() {
  int a, c;
  while ((a = gc()) < 40);
  if (a == '-') return -readInt();
  while ((c = gc()) >= 48) a = a * 10 + c - 480;
  return a - 48;
```

BumpAllocator.h

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

d41d8c, 8 lines

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

FastInput BumpAllocator 20

Techniques (A)

techniques.txt

Combinatorics

159 lines

Recursion Divide and conquer Finding interesting points in N log N Algorithm analysis Master theorem Amortized time complexity Greedy algorithm Scheduling Max contiquous subvector sum Invariants Huffman encoding Graph theory Dynamic graphs (extra book-keeping) Breadth first search Depth first search * Normal trees / DFS trees Dijkstra's algorithm MST: Prim's algorithm Bellman-Ford Konig's theorem and vertex cover Min-cost max flow Lovasz toggle Matrix tree theorem Maximal matching, general graphs Hopcroft-Karp Hall's marriage theorem Graphical sequences Floyd-Warshall Euler cycles Flow networks * Augmenting paths * Edmonds-Karp Bipartite matching Min. path cover Topological sorting Strongly connected components Cut vertices, cut-edges and biconnected components Edge coloring * Trees Vertex coloring * Bipartite graphs (=> trees) * 3^n (special case of set cover) Diameter and centroid K'th shortest path Shortest cycle Dynamic programming Knapsack Coin change Longest common subsequence Longest increasing subsequence Number of paths in a dag Shortest path in a dag Dynprog over intervals Dynprog over subsets Dynprog over probabilities Dynprog over trees 3^n set cover Divide and conquer Knuth optimization Convex hull optimizations RMQ (sparse table a.k.a 2^k-jumps) Bitonic cycle Log partitioning (loop over most restricted)

Computation of binomial coefficients Pigeon-hole principle Inclusion/exclusion Catalan number Pick's theorem Number theory Integer parts Divisibility Euclidean algorithm Modular arithmetic * Modular multiplication * Modular inverses * Modular exponentiation by squaring Chinese remainder theorem Fermat's little theorem Euler's theorem Phi function Frobenius number Ouadratic reciprocity Pollard-Rho Miller-Rabin Hensel lifting Vieta root jumping Game theory Combinatorial games Game trees Mini-max Nim Games on graphs Games on graphs with loops Grundy numbers Bipartite games without repetition General games without repetition Alpha-beta pruning Probability theory Optimization Binary search Ternary search Unimodality and convex functions Binary search on derivative Numerical methods Numeric integration Newton's method Root-finding with binary/ternary search Golden section search Matrices Gaussian elimination Exponentiation by squaring Sorting Radix sort Geometry Coordinates and vectors * Cross product * Scalar product Convex hull Polygon cut Closest pair Coordinate-compression Ouadtrees KD-trees All segment-segment intersection Discretization (convert to events and sweep) Angle sweeping Line sweeping Discrete second derivatives Strings Longest common substring Palindrome subsequences

Knuth-Morris-Pratt Tries Rolling polynomial hashes Suffix array Suffix tree Aho-Corasick Manacher's algorithm Letter position lists Combinatorial search Meet in the middle Brute-force with pruning Best-first (A*) Bidirectional search Iterative deepening DFS / A* Data structures LCA (2^k-jumps in trees in general) Pull/push-technique on trees Heavy-light decomposition Centroid decomposition Lazy propagation Self-balancing trees Convex hull trick (wcipeg.com/wiki/Convex_hull_trick) Monotone queues / monotone stacks / sliding queues Sliding queue using 2 stacks Persistent segment tree

21