

**UDESC**

Universidade do Estado de Santa Catarina

# Não Treinamos o João

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## Contest (1)

```
.vimrc
7 lines

set nu ai si cindent et ts=4 sw=4 so=10 nosm undofile
```

```
inoremap {} {}<left><return><up><end><return>
" remap de chaves

au BufReadPost * if line("'\"") > 0 && line("'\"") <= line("$")
    | exe "normal! g'\"" | endif
" volta pro lugar onde estava quando saiu do arquivo
```

```
run.sh
2 lines

#!/bin/bash
g++ -std=c++20 -DBRUTE -O2 -Wall -Wextra -Wconversion -Wfatal-
errors -fsanitize=address,undefined $1 && ./a.out
```

```
hash.sh
3 lines

# 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
d41d8c, 49 lines

template <auto MOD, typename T = decltype(MOD)>
struct Mint {
    using U = long long;
    // se o modulo for long long, usar U = __int128
    using m = Mint<MOD, T>;
    T v;
    Mint(T val = 0) : v(val) {
        assert(sizeof(T) * 2 <= sizeof(U));
        if (v < -MOD || v >= 2 * MOD) v %= MOD;
        if (v < 0) v += MOD;
        if (v >= MOD) v -= MOD;
    }
    Mint(U val) : v(T(val % MOD)) {
        assert(sizeof(T) * 2 <= sizeof(U));
        if (v < 0) v += MOD;
```

```

    }
    bool operator==(m o) const { return v == o.v; }
    bool operator<(m o) const { return v < o.v; }
    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  $x = -b/2a$ .

$$ax + by = e \Rightarrow \begin{matrix} x = \frac{ed - bf}{ad - bc} \\ y = \frac{af - ec}{ad - bc} \end{matrix}$$

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 $\theta^\circ$

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

#### 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

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

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

$\vec{a}_\beta$  are the coordinates of vector  $\vec{a}$  in basis  $\beta$ .  
 $\vec{a}$  are the coordinates of vector  $\vec{a}$  in the canonical basis.  
 $\vec{b1}$  and  $\vec{b2}$  are the basis vectors for  $\beta$ .  
 $C$  is a matrix that changes from basis  $\beta$  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  $a_n = c_1 a_{n-1} + \dots + c_k a_{n-k}$ , and  $r_1, \dots, r_k$  are distinct roots of  $x^k - c_1 x^{k-1} - \dots - c_k$ , there are  $d_1, \dots, d_k$  s.t.

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

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

.vimrc run hash Mint troubleshoot

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

2.5 Geometry

2.5.1 Triangles

Side lengths:  $a, b, c$

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

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

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

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

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):

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

$$\text{Law of sines: } \frac{\sin \alpha}{a} = \frac{\sin \beta}{b} = \frac{\sin \gamma}{c} = \frac{1}{2R}$$

$$\text{Law of cosines: } a^2 = b^2 + c^2 - 2bc \cos \alpha$$

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

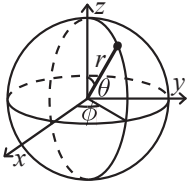
2.5.2 Quadrilaterals

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

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

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

2.5.3 Spherical coordinates



$$x = r \sin \theta \cos \phi \quad r = \sqrt{x^2 + y^2 + z^2}$$
$$y = r \sin \theta \sin \phi \quad \theta = \arccos(z / \sqrt{x^2 + y^2 + z^2})$$
$$z = r \cos \theta \quad \phi = \text{atan2}(y, x)$$

2.6 Derivatives/Integrals

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

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

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

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

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

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

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

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

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

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)  $\mu = \mathbb{E}(X) = \sum_x xp_X(x)$  and variance  $\sigma^2 = V(X) = \mathbb{E}(X^2) - (\mathbb{E}(X))^2 = \sum_x (x - \mathbb{E}(X))^2 p_X(x)$  where  $\sigma$  is the standard 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  $\text{Bin}(n, p)$ ,  $n = 1, 2, \dots$ ,  $0 \leq p \leq 1$ .

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

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

$\text{Bin}(n, p)$  is approximately  $\text{Po}(np)$  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  $\text{Fs}(p)$ ,  $0 \leq p \leq 1$ .

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

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

Poisson distribution

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

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

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

2.9.2 Continuous distributions

Uniform distribution

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

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

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

Exponential distribution

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

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

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

Normal distribution

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

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

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

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

2.10 Markov chains

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

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

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

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

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

**Description:** Set with operations to find the element by index - find\_by\_order(i) and to find the index of an element  
**Usage:** ordered\_set<int> seta;  
**Time:** All operations in  $\mathcal{O}(\log N)$ .  
<ext/pb\_ds/assoc.container.hpp>, <ext/pb\_ds/tree\_policy.hpp> d41d8c, 8 lines  
using namespace \_\_gnu\_pbds;

```
template <typename T>
using ordered_set =
    tree<T, null_type, less<T>, rb_tree_tag,
        tree_order_statistics_node_update>;
```

```
template <typename T, typename U>
using ordered_map = tree<T, U, less<T>, rb_tree_tag,
    tree_order_statistics_node_update>;
```

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  $\mathcal{O}(\log^2 N)$   
d41d8c, 151 lines

```
const ll INF = 1e18;
struct node {
    ll mi, smi, mx, smx, sum, lazy;
    int fmi, fmx;
    node() {
        mi = smi = INF;
        mx = smx = -INF;
        fmi = 0, fmx = 0, sum = 0, lazy = 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) {
        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);
    }
}
```

```

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;
    int n;
    void build(int _n) { // pra construir com tamanho, mas
        vazia
        n = _n;
        t.assign(n * 4, node());
    }
    void build(const vector<ll> &v) { // pra construir com
        vector
        n = (int)v.size();
        t.assign(n * 4, node());
        build(1, 0, n - 1, v);
    }
    void build(ll *bg, ll *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<ll> &a) {
        if (l == r) return t[p] = node(a[l]);
        int mid = (l + 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 - l + 1) * x;
        t[p].mi += x;
        t[p].mx += x;
        t[p].lazy += 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;
        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 (l == r) return;
        int mid = (l + 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 (l > R || r < L || x >= t[p].mx) return t[p];
    if (l >= L && r <= R && x > t[p].smx) {
        pushmin(p, x);
        return t[p];
    }
    pushdown(p, l, r);
    int mid = (l + 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 (l > R || r < L || x <= t[p].mi) return t[p];
    if (l >= L && r <= R && x < t[p].smi) {
        pushmax(p, x);
        return t[p];
    }
    pushdown(p, l, r);
    int mid = (l + 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 (l > R || r < L) return t[p];
    if (l >= L && r <= R) {
        pushsum(p, l, r, x);
        return t[p];
    }
    pushdown(p, l, r);
    int mid = (l + r) >> 1;
    return t[p] = updatesum(lc(p), l, mid, L, R, x) +
        updatesum(rc(p), mid + 1, r, L, R, x);
}
node query(int p, int l, int r, int L, int R) {
    if (l > R || r < L) return node();
    if (l >= L && r <= R) return t[p];
    pushdown(p, l, r);
    int mid = (l + r) >> 1;
    return query(lc(p), l, mid, L, R) + query(rc(p), mid +
        1, r, L, R);
}
ll query(int l, int r) { return query(1, 0, n - 1, l, r).
    sum; }
void updatemax(int l, int r, ll x) { updatemax(1, 0, n - 1,
    l, r, x); }
void updatemin(int l, int r, ll x) { updatemin(1, 0, n - 1,
    l, r, x); }
void updatesum(int l, int r, ll x) { updatesum(1, 0, n - 1,
    l, r, x); }
} seg;

```

## SegmentTree2D.cpp

**Description:** Segment Tree with point update and rectangular range query  
**Usage:** "seg" declared globally  
**Time:**  $\mathcal{O}(\log^2 N)$

d41d8c, 47 lines

```

struct SegTree2D {
    ll merge(ll a, ll b) { return a + b; }
    ll neutral = 0;

```

```

int n, m;
vector<vector<ll>> t;
void build(int _n, int _m) {
    n = _n, m = _m;
    t.assign(2 * n, vector<ll>(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]);
}
ll inner_query(int idx, int l, int r) {
    ll res = neutral;
    for (l += m, r += m + 1; l < r; l >= 1, r >= 1) {
        if (l & 1) res = merge(res, t[idx][l++]);
        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
ll query(int a, int b, int c, int d) {
    ll 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) {
        ll val = merge(t[i << 1][j + m], t[i << 1 | 1][j +
            m]);
        inner_update(i, j, val);
    }
}
} seg;

```

## ImplicitTreap.cpp

**Description:** Does everything

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

d41d8c, 124 lines

```

mt19937 rng((uint32_t)chrono::steady_clock::now().
    time_since_epoch().count());
namespace imp_treap {
    using T = ll; // mudar pra int se nao precisar pra melhorar
        a performance
    T merge(T a, T b) { return a + b; }
    T neutral = 0;
    struct node_info {
        node_info *l, *r;
        int y, size;
        T val, acc, add;
        bool rev;
        node_info() { }
    }

```

```

        node_info(T_val)
            : l(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->l, t->r);
        if (t->l) t->l->rev ^= 1;
        if (t->r) t->r->rev ^= 1;
    }
    t->acc += t->add * size(t);
    // t->acc += t->add se for RMQ
    t->val += t->add;
    if (t->l) t->l->add += t->add;
    if (t->r) t->r->add += t->add;
    t->add = 0;
}
inline void pull(node t) {
    if (t) {
        push(t->l), push(t->r);
        t->size = size(t->l) + size(t->r) + 1;
        t->acc = merge(t->val, merge(acc(t->l), acc(t->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->l, L, R->l);
        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->l);
        if (pos <= imp_key) {
            split(t->l, pos, L, t->l, 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->l);
        if (pos == imp_key) {
            node me = t;
            merge(t, t->l, t->r);
            delete me;
            return true;
        }
        bool ok;
        if (pos < imp_key) ok = remove(t->l, pos, add);
        else ok = remove(t->r, pos, imp_key + 1);
        pull(t);
        return ok;
    }
    inline T query(int l, int r) {
        if (l > r) return neutral;
        node L1, L2, R1, R2;
        split(root, r + 1, L1, R1);
        split(L1, l, L2, R2);
        T ans = acc(R2);
        merge(L1, L2, R2);
        merge(root, L1, R1);
        return ans;
    }
    inline void update_sum(int l, int r, T val) {
        if (l > r) return;
        node L1, L2, R1, R2;
        split(root, r + 1, L1, R1);
        split(L1, l, L2, R2);
        assert(R2);
        R2->add += val;
        merge(L1, L2, R2);
        merge(root, L1, R1);
    }
    inline void reverse(int l, int r) {
        if (l > r) return;
        node L1, L2, R1, R2;
        split(root, r + 1, L1, R1);
        split(L1, l, 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); }
}

```

### LichaoLazy.cpp

**Description:** Sendo  $N = MA - MI$ : insert((a, b)) minimiza tudo com  $ax + b$   $O(\log N)$  insert((a, b), l, r) minimiza com  $ax + b$  no range  $[l, r]$   $O(\log^2 N)$  shift((a, b)) soma  $ax + b$  em tudo  $O(1)$  shift((a, b), l, r) soma  $ax + b$  no range  $[l, r]$   $O(\log^2 N)$  query(x) retorna o valor da posicao x  $O(\log N)$  No inicio eh tudo LINF, se inserir (0, 0) fica tudo 0  
**Time:**  $O(n \log N)$  de memoria,  $O(n)$  de memoria se nao usar as operacoes de range

```

d41d8c.77 lines
template<int MI = int(-1e9), int MA = int(1e9)> struct lichao {
    struct line {
        ll a, b;
        ll la, lb; // lazy
        array<int, 2> ch;
        line(ll a_ = 0, ll b_ = LINF) :
            a(a_), b(b_), la(0), lb(0), ch({-1, -1}) {}
        ll operator ()(ll x) { return a*x + b; }
    };
    vector<line> ln;

```

```

int ch(int p, int d) {
    if (ln[p].ch[d] == -1) {
        ln[p].ch[d] = ln.size();
        ln.emplace_back();
    }
    return ln[p].ch[d];
}
lichao() { ln.emplace_back(); }

void prop(int p, int l, int r) {
    if (ln[p].la == 0 and ln[p].lb == 0) return;
    ln[p].a += ln[p].la, ln[p].b += ln[p].lb;
    if (l != r) {
        int pl = ch(p, 0), pr = ch(p, 1);
        ln[pl].la += ln[p].la, ln[pl].lb += ln[p].lb;
        ln[pr].la += ln[p].la, ln[pr].lb += ln[p].lb;
    }
    ln[p].la = ln[p].lb = 0;
}

ll query(int x, int p=0, int l=MI, int r=MA) {
    prop(p, l, r);
    ll ret = ln[p](x);
    if (ln[p].ch[0] == -1 and ln[p].ch[1] == -1) return ret;
    int m = l + (r-l)/2;
    if (x <= m) return min(ret, query(x, ch(p, 0), l, m));
    return min(ret, query(x, ch(p, 1), m+1, r));
}

void push(line s, int p, int l, int r) {
    prop(p, l, r);
    int m = l + (r-l)/2;
    bool L = s(l) < ln[p](l);
    bool M = s(m) < ln[p](m);
    bool R = s(r) < ln[p](r);
    if (M) swap(ln[p].a, s.a), swap(ln[p].b, s.b);
    if (s.b == LINF) return;
    if (L != M) push(s, ch(p, 0), l, m);
    else if (R != M) push(s, ch(p, 1), m+1, r);
}

void insert(line s, int a=MI, int b=MA, int p=0, int l=MI,
    int r=MA) {
    prop(p, l, r);
    if (a <= l and r <= b) return push(s, p, l, r);
    if (b < l or r < a) return;
    int m = l + (r-l)/2;
    insert(s, a, b, ch(p, 0), l, m);
    insert(s, a, b, ch(p, 1), m+1, r);
}

void shift(line s, int a=MI, int b=MA, int p=0, int l=MI, int
    r=MA) {
    prop(p, l, r);
    int m = l + (r-l)/2;
    if (a <= l and r <= b) {
        ln[p].la += s.a, ln[p].lb += s.b;
        return;
    }
    if (b < l or r < a) return;
    if (ln[p].b != LINF) {
        push(ln[p], ch(p, 0), l, m);
        push(ln[p], ch(p, 1), m+1, r);
        ln[p].a = 0, ln[p].b = LINF;
    }
    shift(s, a, b, ch(p, 0), l, m);
    shift(s, a, b, ch(p, 1), m+1, r);
}
};

```

# Numerical (4)

## 4.1 Factorization

PollardRhoMillerRabin.cpp

Description: Pollard Rho and Miller Rabin

```
d41d8c, 87 lines

namespace MillerRabin {
    inline ll mul_mod(ll a, ll b, ll m) { return (ll)((__int128
        )a * b % m); }
    inline ll power(ll b, ll e, ll m) {
        ll r = 1;
        b = b % m;
        while (e > 0) {
            if (e & 1) r = mul_mod(r, b, m);
            b = mul_mod(b, b, m), e >>= 1;
        }
        return r;
    }
    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 <=
    2^64
    int primes[] = {2, 325, 9375, 28178, 450775, 9780504,
        1795265022};

    // funciona para n <= 3*10^24 com os primos ate 41, mas tem
    que cuidar com overflow
    // int primes[] = {2, 3, 5, 7, 11, 13, 17, 19, 23, 29, 31,
        37, 41};

    bool prime(ll n) {
        if (n <= 2 || (n % 2 == 0)) return n == 2;
        ll 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 ll P = 1e6 + 1;
    ll seq[P];
    inline ll add_mod(ll x, ll y, ll m) { return (x += y) < m ?
        x : x - m; }
    inline ll mul_mod(ll a, ll b, ll m) { return (ll)((__int128
        )a * b % m); }

    ll rho(ll n) {
        if (n % 2 == 0) return 2;
        if (n % 3 == 0) return 3;
        ll x0 = rng() % n, c = rng() % n;
        while (1) {
            ll x = x0++, y = x, u = 1, v, t = 0;
            ll *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;
            }
        }
    }
}
```

```
v = u;
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(ll)> dfs = [&](ll x) {
        if (x == 1) return;
        if (x < Sieve::P) {
            auto fs = Sieve::factorize(x);
            f.insert(f.end(), fs.begin(), fs.end());
        } else if (MillerRabin::prime(x)) {
            f.push_back(x);
        } else {
            ll 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];
        pk *= k;
    }
    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;
}
```

EvaluateInterpolation.cpp

Description: Dado 'n' pontos (i, y[i]), i ∈ [0, n), avalia o polinomio de grau n-1 que passa por esses pontos em 'x' Tudo modular, precisa do mint

Time: O(N)

```
d41d8c, 20 lines

mint evaluate_interpolation(int x, vector<mint> y) {
    int n = y.size();

    vector<mint> sulf(n+1, 1), fat(n, 1), ifat(n);
    for (int i = n-1; i >= 0; i--) sulf[i] = sulf[i+1] * (x - i);
    for (int i = 1; i < n; i++) fat[i] = fat[i-1] * i;
    ifat[n-1] = 1/fat[n-1];
    for (int i = n-2; i >= 0; i--) ifat[i] = ifat[i+1] * (i + 1);
}
```

```
mint pref = 1, ans = 0;
for (int i = 0; i < n; pref *= (x - i++)) {
    mint num = pref * sulf[i+1];

    mint den = ifat[i] * ifat[n-1 - i];
    if ((n-1 - i)%2) den *= -1;

    ans += y[i] * num * den;
}
return ans;
}
```

## 4.3 MIT Polynomials

GoldenSectionSearch.h

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

Usage: double func(double x) { return 4+x+.3\*x\*x; }

double xmin = gss(-1000,1000,func);

Time: O(log((b - a)/ε))

```
d41d8c, 14 lines

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

Polynomial.h

```
d41d8c, 17 lines

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

PolyRoots.h

Description: Finds the real roots to a polynomial.

Usage: poly\_roots({{-3,1}},-1e9,1e9) // solve x^2-3x+2 = 0

Time: O(n^2 log(1/ε))

```
"Polynomial.h"
d41d8c, 23 lines

vector<double> poly_roots(Poly p, double xmin, double xmax) {
    if (sz(p.a) == 2) { return {-p.a[0]/p.a[1]}; }
    vector<double> ret;
    Poly der = p;
```



```
der.diff();
auto dr = poly_roots(der, xmin, xmax);
dr.push_back(xmin-1);
dr.push_back(xmax+1);
sort(all(dr));
rep(i,0,sz(dr)-1) {
    double l = dr[i], h = dr[i+1];
    bool sign = p(l) > 0;
    if (sign ^ (p(h) > 0)) {
        rep(it,0,60) { // while (h - l > 1e-8)
            double m = (l + h) / 2, f = p(m);
            if ((f <= 0) ^ sign) l = m;
            else h = m;
        }
        ret.push_back((l + h) / 2);
    }
}
return ret;
}
```

PolyInterpolate.h

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

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

BerlekampMassey.h

**Description:** Recovers any  $n$ -order linear recurrence relation from the first  $2n$  terms of the recurrence. Useful for guessing linear recurrences after brute-forcing the first terms. Should work on any field, but numerical stability for floats is not guaranteed. Output will have size  $\leq n$ .  
**Usage:** BerlekampMassey({0, 1, 1, 3, 5, 11}) // {1, 2}  
"../number-theory/ModPow.h"

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

    ll b = 1;
    rep(i,0,n) { ++m;
        ll d = s[i] % mod;
        rep(j,1,L+1) d = (d + C[j] * s[i - j]) % mod;
        if (!d) continue;
        T = C; ll coef = d * modpow(b, mod-2) % mod;
        rep(j,m,n) C[j] = (C[j] - coef * B[j - m]) % mod;
        if (2 * L > i) continue;
        L = i + 1 - L; B = T; b = d; m = 0;
    }

    C.resize(L + 1); C.erase(C.begin());
    trav(x, C) x = (mod - x) % mod;
    return C;
}
```

LinearRecurrence.h

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

```
typedef vector<ll> Poly;
ll linearRec(Poly S, Poly tr, ll k) {
    int n = sz(S);

    auto combine = [&](Poly a, Poly b) {
        Poly res(n * 2 + 1);
        rep(i,0,n+1) rep(j,0,n+1)
            res[i + j] = (res[i + j] + a[i] * b[j]) % mod;
        for (int i = 2 * n; i > n; --i) rep(j,0,n)
            res[i - 1 - j] = (res[i - 1 - j] + res[i] * tr[j]) % mod;
        res.resize(n + 1);
        return res;
    };

    Poly pol(n + 1), e(pol);
    pol[0] = e[1] = 1;

    for (++k; k; k /= 2) {
        if (k % 2) pol = combine(pol, e);
        e = combine(e, e);
    }

    ll res = 0;
    rep(i,0,n) res = (res + pol[i + 1] * S[i]) % mod;
    return res;
}
```

Integrate.h

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

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

IntegrateAdaptive.h

**Description:** Fast integration using an adaptive Simpson's rule.  
**Usage:** double z, y; double h(double x) { return x\*x + y\*y + z\*z <= 1; } double g(double y) { ::y = y; return quad(h, -1, 1); } double f(double z) { ::z = z; return quad(g, -1, 1); } double sphereVol = quad(f, -1, 1), pi = sphereVol\*3/4;

```
typedef double d;
d simpson(d (*f)(d), d a, d b) {
    d c = (a+b) / 2;
    return (f(a) + 4*f(c) + f(b)) * (b-a) / 6;
}
d rec(d (*f)(d), d a, d b, d eps, d S) {
    d c = (a+b) / 2;
    d S1 = simpson(f, a, c);
    d S2 = simpson(f, c, b), T = S1 + S2;
    if (abs (T - S) <= 15*eps || b-a < 1e-10)
        return T + (T - S) / 15;
    return rec(f, a, c, eps/2, S1) + rec(f, c, b, eps/2, S2);
}
```

```
}
d quad(d (*f)(d), d a, d b, d eps = 1e-8) {
    return rec(f, a, b, eps, simpson(f, a, b));
}
```

Determinant.h

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

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

IntDeterminant.h

**Description:** Calculates determinant using modular arithmetics. Modulos can also be removed to get a pure-integer version.  
**Time:**  $\mathcal{O}(N^3)$

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

Simplex.h

**Description:** Solves a general linear maximization problem: maximize  $c^T x$  subject to  $Ax \leq b, x \geq 0$ . Returns -inf if there is no solution, inf if there are arbitrarily good solutions, or the maximum value of  $c^T x$  otherwise. The input vector is set to an optimal  $x$  (or in the unbounded case, an arbitrary solution fulfilling the constraints). Numerical stability is not guaranteed. For better performance, define variables such that  $x = 0$  is viable.  
**Usage:** vvd A = {{1,-1}, {-1,1}, {-1,-2}}; vd b = {1,1,-4}, c = {-1,-1}, x; T val = LPSolver(A, b, c).solve(x);  
**Time:**  $\mathcal{O}(NM * \#pivots)$ , where a pivot may be e.g. an edge relaxation.  $\mathcal{O}(2^n)$  in the general case.

```
typedef double T; // long double, Rational, double + mod<P>...
typedef vector<T> vd;
typedef vector<vd> vvd;

const T eps = 1e-8, inf = 1/.0;
#define MP make_pair
#define ltj(X) if(s == -1 || MP(X[j],N[j]) < MP(X[s],N[s])) s=j
```



```
struct LPSolver {
    int m, n;
    vi N, B;
    vvd D;

    LPSolver(const vvd& A, const vd& b, const vd& c) :
        m(sz(b)), n(sz(c)), N(n+1), B(m), D(m+2, vd(n+2)) {
        rep(i,0,m) rep(j,0,n) D[i][j] = A[i][j];
        rep(i,0,m) { B[i] = n+i; D[i][n] = -1; D[i][n+1] = b[i]; }
        rep(j,0,n) { N[j] = j; D[m][j] = -c[j]; }
        N[n] = -1; D[m+1][n] = 1;
    }

    void pivot(int r, int s) {
        T *a = D[r].data(), inv = 1 / a[s];
        rep(i,0,m+2) if (i != r && abs(D[i][s]) > eps) {
            T *b = D[i].data(), inv2 = b[s] * inv;
            rep(j,0,n+2) b[j] -= a[j] * inv2;
            b[s] = a[s] * inv2;
        }
        rep(j,0,n+2) if (j != s) D[r][j] *= inv;
        rep(i,0,m+2) if (i != r) D[i][s] *= -inv;
        D[r][s] = inv;
        swap(B[r], N[s]);
    }

    bool simplex(int phase) {
        int x = m + phase - 1;
        for (;;) {
            int s = -1;
            rep(j,0,n+1) if (N[j] != -phase) ltj(D[x]);
            if (D[x][s] >= -eps) return true;
            int r = -1;
            rep(i,0,m) {
                if (D[i][s] <= eps) continue;
                if (r == -1 || MP(D[i][n+1] / D[i][s], B[i])
                    < MP(D[r][n+1] / D[r][s], B[r])) r = i;
            }
            if (r == -1) return false;
            pivot(r, s);
        }
    }

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

math-simplex.cpp

**Description:** Simplex algorithm. WARNING- segfaults on empty (size 0) max cx st Ax<=b, x>=0 do 2 phases; 1st check feasibility; 2nd check bound- edness and ans

d41d8c, 40 lines

```
vector<double> simplex(vector<vector<double>> A, vector<double>
    > b, vector<double> c) {
```

```
    int n = (int) A.size(), m = (int) A[0].size()+1, r = n, s = m
        -1;
    vector<vector<double>> D = vector<vector<double>> (n+2,
        vector<double>(m+1));
    vector<int> ix = vector<int> (n+m);
    for (int i=0; i<n+m; i++) ix[i] = i;
    for (int i=0; i<n; i++) {
        for (int j=0; j<m-1; j++) D[i][j]=-A[i][j];
        D[i][m-1] = 1;
        D[i][m] = b[i];
        if (D[r][m] > D[i][m]) r = i;
    }
    for (int j=0; j<m-1; j++) D[n][j]=c[j];
    D[n+1][m-1] = -1; int z = 0;
    for (double d;;) {
        if (r < n) {
            swap(ix[s], ix[r+m]);
            D[r][s] = 1.0/D[r][s];
            for (int j=0; j<m; j++) if (j!=s) D[r][j] *= -D[r][s];
            for(int i=0; i<=n+1; i++)if(i!=r) {
                for (int j=0; j<=m; j++) if(j!=s) D[i][j] += D[r][j] *
                    D[i][s];
                D[i][s] *= D[r][s];
            }
        }
        r = -1; s = -1;
        for (int j=0; j <m; j++) if (s<0 || ix[s]>ix[j]) {
            if (D[n+1][j]>eps || D[n+1][j]>-eps && D[n][j]>eps) s = j
                ;
        }
        if (s < 0) break;
        for (int i=0; i<n; i++) if(D[i][s]<-eps) {
            if (r < 0 || (d = D[r][m]/D[r][s]-D[i][m]/D[i][s]) < -eps
                || d < eps && ix[r+m] > ix[i+m]) r=i;
        }
        if (r < 0) return vector<double>(); // unbounded
    }
    if (D[n+1][m] < -eps) return vector<double>(); // infeasible
    vector<double> x(m-1);
    for (int i = m; i < n+m; i ++) if (ix[i] < m-1) x[ix[i]] = D[
        i-m][m];
    printf("%.21f\n", D[n][m]);
    return x; // ans: D[n][m]
}
```

SolveLinear.h

**Description:** Solves  $A * x = b$ . If there are multiple solutions, an arbitrary one is returned. Returns rank, or -1 if no solutions. Data in  $A$  and  $b$  is lost.

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

d41d8c, 38 lines

```
typedef vector<double> vd;
const double eps = 1e-12;

int solveLinear(vector<vd>& A, vd& b, vd& x) {
    int n = sz(A), m = sz(x), rank = 0, br, bc;
    if (n) assert(sz(A[0]) == m);
    vi col(m); iota(all(col), 0);

    rep(i,0,n) {
        double v, bv = 0;
        rep(r,i,n) rep(c,i,m)
            if ((v = fabs(A[r][c])) > bv)
                br = r, bc = c, bv = v;
        if (bv <= eps) {
            rep(j,i,n) if (fabs(b[j]) > eps) return -1;
            break;
        }
        swap(A[i], A[br]);
        swap(b[i], b[br]);
```

```
        swap(col[i], col[bc]);
        rep(j,0,n) swap(A[j][i], A[j][bc]);
        bv = 1/A[i][i];
        rep(j,i+1,n) {
            double fac = A[j][i] * bv;
            b[j] -= fac * b[i];
            rep(k,i+1,m) A[j][k] -= fac*A[i][k];
        }
        rank++;
    }

    x.assign(m, 0);
    for (int i = rank; i--;) {
        b[i] /= A[i][i];
        x[col[i]] = b[i];
        rep(j,0,i) b[j] -= A[j][i] * b[i];
    }
    return rank; // (multiple solutions if rank < m)
}
```

SolveLinear2.h

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

"SolveLinear.h" d41d8c, 7 lines

```
rep(j,0,n) if (j != i) // instead of rep(j,i+1,n)
// ... then at the end:
x.assign(m, undefined);
rep(i,0,rank) {
    rep(j,rank,m) if (fabs(A[i][j]) > eps) goto fail;
    x[col[i]] = b[i] / A[i][i];
fail;; }
```

SolveLinearBinary.h

**Description:** Solves  $Ax = b$  over  $\mathbb{F}_2$ . If there are multiple solutions, one is returned arbitrarily. Returns rank, or -1 if no solutions. Destroys  $A$  and  $b$ .

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

d41d8c, 34 lines

```
typedef bitset<1000> bs;

int solveLinear(vector<bs>& A, vi& b, bs& x, int m) {
    int n = sz(A), rank = 0, br;
    assert(m <= sz(x));
    vi col(m); iota(all(col), 0);
    rep(i,0,n) {
        for (br=i; br<n; ++br) if (A[br].any()) break;
        if (br == n) {
            rep(j,i,n) if(b[j]) return -1;
            break;
        }
        int bc = (int)A[br]._Find_next(i-1);
        swap(A[i], A[br]);
        swap(b[i], b[br]);
        swap(col[i], col[bc]);
        rep(j,0,n) if (A[j][i] != A[j][bc]) {
            A[j].flip(i); A[j].flip(bc);
        }
        rep(j,i+1,n) if (A[j][i]) {
            b[j] ^= b[i];
            A[j] ^= A[i];
        }
        rank++;
    }

    x = bs();
    for (int i = rank; i--;) {
        if (!b[i]) continue;
        x[col[i]] = 1;
        rep(j,0,i) b[j] ^= A[j][i];
    }
```

```
    return rank; // (multiple solutions if rank < m)
}
```

MatrixInverse.h

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

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

d41d8c, 35 lines

```
int matInv(vector<vector<double>>& A) {
    int n = sz(A); vi col(n);
    vector<vector<double>> tmp(n, vector<double>(n));
    rep(i,0,n) tmp[i][i] = 1, col[i] = i;

    rep(i,0,n) {
        int r = i, c = i;
        rep(j,i,n) rep(k,i,n)
            if (fabs(A[j][k]) > fabs(A[r][c]))
                r = j, c = k;
        if (fabs(A[r][c]) < 1e-12) return i;
        A[i].swap(A[r]); tmp[i].swap(tmp[r]);
        rep(j,0,n)
            swap(A[j][i], A[j][c]), swap(tmp[j][i], tmp[j][c]);
        swap(col[i], col[c]);
        double v = A[i][i];
        rep(j,i+1,n) {
            double f = A[j][i] / v;
            A[j][i] = 0;
            rep(k,i+1,n) A[j][k] -= f*A[i][k];
            rep(k,0,n) tmp[j][k] -= f*tmp[i][k];
        }
        rep(j,i+1,n) A[i][j] /= v;
        rep(j,0,n) tmp[i][j] /= v;
        A[i][i] = 1;
    }

    for (int i = n-1; i > 0; --i) rep(j,0,i) {
        double v = A[j][i];
        rep(k,0,n) tmp[j][k] -= v*tmp[i][k];
    }

    rep(i,0,n) rep(j,0,n) A[col[i]][col[j]] = tmp[i][j];
    return n;
}
```

Tridiagonal.h

**Description:**  $x = \text{tridiagonal}(d, p, q, b)$  solves the equation system

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

This is useful for solving problems on the type

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

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

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

Fails if the solution is not unique.

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

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

d41d8c, 26 lines

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

4.4 Optimization

4.5 Matrices

4.6 Convolutions

FFT.cpp

**Description:** Computa convolução (multiplicação) de polinômios em  $\mathcal{O}(N \log N)$ , sendo  $N$  a soma dos graus dos polinômios. Testado e sem erros de precisão com polinômios de grau até  $3 \cdot 10^5$  e constantes até  $10^6$ .

**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]);
    }

    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) {
            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;
    T g = 1;
    while ((g ^ (MOD / 2)) == 1) g += 1;
    if (inv) g = T(1) / g;
    for (int step = n / 2; step; step /= 2) {
        T w = g ^ (MOD / (n / step)), wn = 1;
        for (int i = 0; i < n / 2; i += step) {
            for (int j = 0; j < step; j++) {
                auto u = a[2 * i + j], v = wn * a[2 * i + j + step];
                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;
}

```

## NTTBigModulo.cpp

d41d8c, 78 lines

```

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

```

```

ll crt(array<int, 2> rem, array<int, 2> mod) {
    __int128 ans = rem[0], m = mod[0];
    ll x, y;
    ll g = extended_gcd(mod[1], (ll)m, x, y);
    if ((ans - rem[1]) % g != 0) return -1;
    ans = ans + (__int128)1 * (rem[1] - ans) * (m / g) * y;
    m = (__int128)(mod[1] / g) * (m / g) * g;
    ans = (ans % m + m) % m;
    return (ll)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> a1(ta.size()), b1(tb.size());
    vector<U> a2(ta.size()), b2(tb.size());
    for (int i = 0; i < (int)ta.size(); i++) a1[i] = ta[i];
    for (int i = 0; i < (int)tb.size(); i++) b1[i] = tb[i];
    for (int i = 0; i < (int)ta.size(); i++) a2[i] = ta[i];
    for (int i = 0; i < (int)tb.size(); i++) b2[i] = tb[i];
    auto c1 = multiply<MOD1>(a1, b1);
    vector<ll> res(c1.size());
    for (int i = 0; i < (int)res.size(); i++)
        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);
    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 <= N; i++) {
        mint a = 0;
        mint b = 0;
        for (int j = i; j <= 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 <= 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 <= 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 <= N; j += i) C[j] -= C[i];
    return C;
}

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 << N);
    B.resize(1 << N);
    vector<mint> C(1 << N);
    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++) {
                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 <= N; j++)
            for (int k = 0; k + j <= 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 popcnt = 0; popcnt <= N; popcnt++)
                c[i][popcnt] -= c[i ^ (1 << j)][popcnt];
        }
    }
    vector<mint> ans(1 << N);
    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++) {
        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;
}
```

# Number theory (5)

## 5.1 Modular arithmetic

ModularArithmetic.h

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

```
"euclid.h"
d41d8c, 18 lines

const ll mod = 17; // change to something else
struct Mod {
    ll 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) {
        ll x, y, g = 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.

```
d41d8c, 3 lines

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

ModPow.h

```
d41d8c, 8 lines

const ll mod = 1000000007; // faster if const
```

```
ll modpow(ll b, ll e) {
    ll ans = 1;
    for (; e; b = b * b % mod, e /= 2)
        if (e & 1) ans = ans * b % mod;
    return ans;
}
```

ModLog.h

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

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

```
d41d8c, 11 lines

ll modLog(ll a, ll b, ll m) {
    ll n = (ll) sqrt(m) + 1, e = 1, f = 1, j = 1;
    unordered_map<ll, ll> A;
    while (j <= n && (e = f = e * a % m) != b % m)
        A[e * b % m] = j++;
    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.

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

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

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

```
d41d8c, 11 lines

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 >= (ll)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;
}
```

ModSqrt.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  $p$

```
"ModPow.h"
d41d8c, 24 lines

ll 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 or 2^(n+3)/8 * 2^(n-1)/4 works if p % 8 == 5
    ll 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;
    ll x = modpow(a, (s + 1) / 2, p);
    ll b = modpow(a, s, p), g = modpow(n, s, p);
    for (; r = m) {
        ll t = b;
        for (m = 0; m < r && t != 1; ++m)
            t = t * t % p;
        if (m == 0) return x;
        ll gs = modpow(g, 1LL << (r - m - 1), p);
        g = gs * gs % p;
        x = x * gs % p;
        b = b * g % p;
    }
}
```

## 5.2 Primality

FastEratosthenes.h

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

**Time:** LIM=1e9 ≈ 1.5s

d41d8c, 20 lines

```
const int LIM = 1e6;
bitset<LIM> isPrime;
vi eratosthenes() {
    const int S = (int)round(sqrt(LIM)), R = LIM / 2;
    vi pr = {2}, sieve(S+1); pr.reserve((int)(LIM/log(LIM)*1.1));
    vector<pii> cp;
    for (int i = 3; i <= S; i += 2) if (!sieve[i]) {
        cp.push_back({i, i * i / 2});
        for (int j = i * i; j <= 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;
        rep(i,0,min(S, R - L))
            if (!block[i]) pr.push_back((L + i) * 2 + 1);
    }
    for (int i : pr) isPrime[i] = 1;
    return pr;
}
```

**MillerRabin.h**  
**Description:** Deterministic Miller-Rabin primality test. Guaranteed to work for numbers up to  $7 \cdot 10^{18}$ ; for larger numbers, use Python and extend A randomly.  
**Time:** 7 times the complexity of  $a^b \bmod c$ .

```
"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 = __builtin_ctzll(n-1), d = n >> s;
    for (ull a : A) { // ^ count trailing zeroes
        ull p = modpow(a%n, d, n), i = s;
        while (p != 1 && p != n - 1 && a % n && i--)
            p = modmul(p, p, n);
        if (p != n-1 && i != s) return 0;
    }
    return 1;
}
```

**Factor.h**  
**Description:** Pollard-rho randomized factorization algorithm. Returns prime factors of a number, in arbitrary order (e.g. 2299 -> {11, 19, 11}).  
**Time:**  $\mathcal{O}\left(n^{1/4}\right)$ , less for numbers with small factors.

```
"ModMulLL.h", "MillerRabin.h" d41d8c, 18 lines
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 l = factor(x), r = factor(n / x);
    l.insert(l.end(), all(r));
    return l;
}
```

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

```
d41d8c, 5 lines
ll euclid(ll a, ll b, ll &x, ll &y) {
    if (!b) return x = 1, y = 0, a;
    ll d = euclid(b, a % b, y, x);
    return y -= a/b * x, d;
}
```

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

```
"euclid.h" d41d8c, 7 lines
ll crt(ll a, ll m, ll b, ll n) {
    if (n > m) swap(a, b), swap(m, n);
    ll x, y, g = euclid(m, n, x, y);
    assert((a - b) % g == 0); // else no solution
    x = (b - a) % n * x % n / g * m + a;
    return x < 0 ? x + m*n/g : x;
}
```

5.3.1 Bézout’s identity

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

$$ax + by = d$$

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

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

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

```
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  $x \geq 0$ , finds the closest rational approximation  $p/q$  with  $p, q \leq N$ . It will obey  $|p/q - x| \leq 1/qN$ .  
For consecutive convergents,  $p_{k+1}q_k - q_{k+1}p_k = (-1)^k$ .  $(p_k/q_k)$  alternates between  $> x$  and  $< x$ . If  $x$  is rational,  $y$  eventually becomes  $\infty$ ; 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 ~ 1e7; long double for N ~ 1e9
pair<ll, ll> approximate(d x, ll N) {
    ll LP = 0, LQ = 1, P = 1, Q = 0, inf = LLONG_MAX; d y = x;
    for (;;) {
        ll lim = min(P ? (N-LP) / P : inf, Q ? (N-LQ) / Q : inf),
            a = (ll)floor(y), b = min(a, lim),
            NP = b*P + LP, NQ = b*Q + LQ;
        if (a > b) {
            // If b > a/2, we have a semi-convergent that gives us a
            // better approximation; if b = a/2, we *may* have one.
            // Return {P, Q} here for a more canonical approximation.
            return (abs(x - (d)NP / (d)NQ) < abs(x - (d)P / (d)Q)) ?
                make_pair(NP, NQ) : make_pair(P, Q);
        }
        if (abs(y = 1/(y - (d)a)) > 3*N) {
            return {NP, NQ};
        }
        LP = P; P = NP;
        LQ = Q; Q = NQ;
    }
}
```

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

```
d41d8c, 25 lines
struct Frac { ll p, q; };

template<class F>
Frac fracBS(F f, ll N) {
    bool dir = 1, A = 1, B = 1;
    Frac lo{0, 1}, hi{1, 1}; // Set hi to 1/0 to search (0, N]
    if (f(lo)) return lo;
    assert(f(hi));
    while (A || B) {
        ll 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

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

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  $n < 1e19$ .

5.8 Mobius Function

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

Mobius Inversion:

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

Other useful formulas/forms:

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

6.1 Permutations

6.1.1 Factorial

$n$	1	2	3	4	5	6	7	8	9	10
$n!$	1	2	6	24	120	720	5040	40320	362880	3628800
$n$	11	12	13	14	15	16	17			
$n!$	4.0e7	4.8e8	6.2e9	8.7e10	1.3e12	2.1e13	3.6e14			
$n$	20	25	30	40	50	100	150	171		
$n!$	2e18	2e25	3e32	8e47	3e64	9e157	6e262	>DBL_MAX		

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

```
int permToInt(vi& v) {
    int use = 0, i = 0, r = 0;
    for(int x:v) r = r * ++i + __builtin_popcount(use & ~(1<<x)),
        use |= 1 << x;
    return r;
}
```

IntPerm multinomial

6.1.2 Cycles

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

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

6.1.3 Derangements

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

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

6.1.4 Burnside’s lemma

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

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

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

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

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

6.2 Partitions and subsets

6.2.1 Partition function

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

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

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

$n$	0	1	2	3	4	5	6	7	8	9	20	50	100
$p(n)$	1	1	2	3	5	7	11	15	22	30	627	~2e5	~2e8

6.2.2 Lucas’ Theorem

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

6.2.3 Binomials

multinomial.h

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

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

6.3 General purpose numbers

6.3.1 Bernoulli numbers

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

Sums of powers:

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

Euler-Maclaurin formula for infinite sums:

$$\begin{aligned} \sum_{i=m}^\infty f(i) &= \int_m^\infty f(x) dx - \sum_{k=1}^\infty \frac{B_k}{k!} f^{(k-1)}(m) \\ &\approx \int_m^\infty f(x) dx + \frac{f(m)}{2} - \frac{f'(m)}{12} + \frac{f'''(m)}{720} + O(f^{(5)}(m)) \end{aligned}$$

6.3.2 Stirling numbers of the first kind

Number of permutations on  $n$  items with  $k$  cycles.

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

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

6.3.3 Eulerian numbers

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

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

6.3.4 Stirling numbers of the second kind

Partitions of  $n$  distinct elements into exactly  $k$  groups.

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

6.3.5 Bell numbers

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

$$C_n = \frac{1}{n+1} \binom{2n}{n} = \binom{2n}{n} - \binom{2n}{n+1} = \frac{(2n)!}{(n+1)!n!}$$
$$C_0 = 1, C_{n+1} = \frac{2(2n+1)}{n+2} C_n, C_{n+1} = \sum C_i C_{n-i}$$
$$C_n = 1, 1, 2, 5, 14, 42, 132, 429, 1430, 4862, 16796, 58786, \dots$$

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

Graph (7)

7.1 Fundamentals

7.2 Network flow

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

```
struct dinitz {
    const bool scaling = false; // com scaling -> O(nm log(U)),
    int lim; // com constante alta
    struct edge {
        int to, cap, rev, flow;
        bool res;
        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;
    ll 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;
```

```
                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++) {
            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, g[e.to][e.rev].flow -= foi;
            return foi;
        }
        return 0;
    }
    ll max_flow(int s, int t) {
        for (lim = 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(g.g.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)
            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.to);
    return cut;
}
```

MinCostMaxFlow.cpp

**Description:**  $min\_cost\_flow(s, t, f)$  computa o par  $(fluxo, custo)$  com  $max(fluxo) \leq f$  que tenha  $min(custo) min\_cost\_flow(s, t)$ : Fluxo máximo de custo mínimo de  $s$  pra  $t$  Se for um dag, da pra substituir o SPFA por uma DP pra naopagar  $O(nm)$  no começo Se nao tiver aresta com custo negativo, nao precisa do SPFA  
**Time:**  $O(nm + f * m \log n)$

```
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 res_)
            : 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>::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(g.size(), 0);
    dist = vector<T>(g.size(), inf);

    dist[s] = 0;
    q.push_back(s);
    is_inside[s] = true;

    while (!q.empty()) {
        int v = q.front();
        q.pop_front();
        is_inside[v] = false;

        for (int i = 0; i < g[v].size(); i++) {
            auto [to, rev, flow, cap, res, cost] = g[v][i];
            if (flow < cap and dist[v] + cost < dist[to]) {
                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<>> q;
    dist = vector<T>(g.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 < g[v].size(); i++) {
            auto [to, rev, flow, cap, res, cost] = g[v][i];
            cost += pot[v] - pot[to];
            if (flow < cap and dist[v] + cost < dist[to]) {
                dist[to] = dist[v] + cost;
                q.emplace(dist[to], to);
                par_idx[to] = i, par[to] = v;
            }
        }
    }
    return dist[t] < inf;
}

pair<int, T> min_cost_flow(int s, int t, int flow = INF) {
    vector<T> pot(g.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)) {
    for (int i = 0; i < g.size(); i++)
        if (dist[i] < inf) pot[i] += dist[i];

    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;
}

// Opcional: retorna as arestas originais por onde passa flow
= cap
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;
}
};
```

## LowerBoundMaxFlow.cpp

**Description:** precisa do dinitz add(a, b, l, r) adiciona aresta de a pra b, onde precisa passar f de fluxo,  $l \leq f \leq r$ . add(a, b, c) adiciona aresta de a pra b com capacidade c.

**Time:** Mesma complexidade do dinitz

d41d8c, 36 lines

```
struct lb_max_flow : dinitz {
    vector<int> d;
    lb_max_flow(int n) : dinitz(n + 2), d(n, 0) {}
    void add(int a, int b, int l, int r) {
        d[a] -= l;
        d[b] += l;
        dinitz::add(a, b, r - l);
    }
    void add(int a, int b, int c) {
        dinitz::add(a, b, c);
    }
    bool has_circulation() {
        int n = d.size();

        ll cost = 0;
        for (int i = 0; i < n; i++) {
            if (d[i] > 0) {
                cost += d[i];
                dinitz::add(n, i, d[i]);
            } else if (d[i] < 0) {
                dinitz::add(i, n+1, -d[i]);
            }
        }
    }
};
```

```
return (dinitz::max_flow(n, n+1) == cost);
}
bool has_flow(int src, int snk) {
    dinitz::add(snk, src, INF);
    return has_circulation();
}
}
ll max_flow(int src, int snk) {
    if (!has_flow(src, snk)) return -1;
    dinitz::F = 0;
    return dinitz::max_flow(src, snk);
}
};
```

## 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:**  $\mathcal{O}(n^3)$ .

d41d8c, 42 lines

```
template<typename T> struct hungarian {
    int n;
    vector<vector<T>> a;
    vector<T> u, v;
    vector<int> p, way;
    T inf;

    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);
            do {
                used[j0] = true;
                int i0 = p[j0], j1 = -1;
                T delta = inf;
                for (int j = 1; j <= 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;
                    if (minv[j] < delta) delta = minv[j], j1 = j;
                }
                for (int j = 0; j <= 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 <= 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'.

d41d8c, 77 lines

```
vector<int> g[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 l;
    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] != l) {
        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]]) {
        base[i] = l;
        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] = 0;
    vis[s] = 1; q = queue<int>(); q.push(s);
    while (q.size()) {
        int u = q.front(); q.pop();
        for (int i : g[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 -> 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; lazy = 0;
        ssz = 1; vsz = 0;
        subsum = x; vsum = 0;
    }
};

struct LCT {
    vector<node> t;
    LCT() {}
    LCT(int n) : t(n + 1) {}

    // <independant 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 &l = t[x].c[0], &r = t[x].c[1];
        push(l); push(r);
        t[x].sum = t[l].sum + t[r].sum + t[x].val;
        t[x].sz = t[l].sz + t[r].sz + 1;
        t[x].ssz = t[l].ssz + t[r].ssz + t[x].vsz + 1;
        t[x].subsum = t[l].subsum + t[r].subsum + t[x].vsum + t[x].
            val;
    }
    void push(int x) {
        if (!x) return;
        int &l = t[x].c[0], &r = t[x].c[1];
        if (t[x].flip) {
            swap(l, r);
            if (l) t[l].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 (l) t[l].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);
    }
}

// </independant splay tree code>

// making it a root in the rep. tree
void make_root(int u) {
    access(u);
    int l = t[u].c[0];
    t[l].flip ^= 1;
    swap(t[l].p, t[l].pp);
    t[u].vsz += t[l].ssz;
    t[u].vsum += t[l].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;
        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 -> 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] -> 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);
    }
    cut(u);
    int ans = 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 {
    int T;
    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() > 0) {
        id[i] = comp[i][0];
    }
}
};
```

## 7.8 Math

### 7.8.1 Number of Spanning Trees

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

### 7.8.2 Erdős–Gallai theorem

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

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

## 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.)

d41d8c, 28 lines

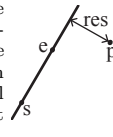
```
template <class T> int sgn(T x) { return (x > 0) - (x < 0); }
template<class T>
struct Point {
    typedef Point P;
    T x, y;
    explicit Point(T x=0, T y=0) : x(x), y(y) {}
    bool operator<(P p) const { return tie(x,y) < tie(p.x,p.y); }
    bool operator==(P p) const { return tie(x,y)==tie(p.x,p.y);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

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

#### lineDistance.h

**Description:**

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



"Point.h"

d41d8c, 4 lines

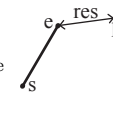
```
template<class P>
double lineDist(const P& a, const P& b, const P& p) {
    return (double)(b-a).cross(p-a)/(b-a).dist();
}
```

#### 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;



"Point.h"

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;
}
```

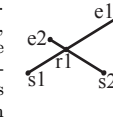
#### SegmentIntersection.h

**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<ll> and the intersection point does not have integer coordinates. Products of three coordinates are used in intermediate steps so watch out for overflow if using int or long long.

**Usage:** vector<P> inter = segInter(s1,e1,s2,e2);

if (sz(inter)==1)  
cout << "segments intersect at " << inter[0] << endl;



"Point.h", "OnSegment.h"

d41d8c, 13 lines

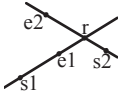
```
template<class P> vector<P> segInter(P a, P b, P c, P d) {
    auto oa = c.cross(d, a), ob = c.cross(d, b),
         oc = a.cross(b, c), od = a.cross(b, d);
    // Checks if intersection is single non-endpoint point.
    if (sgn(oa) * sgn(ob) < 0 && sgn(oc) * sgn(od) < 0)
        return {(a * ob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

```
    if (onSegment(c, d, b)) s.insert(b);
    if (onSegment(a, b, c)) s.insert(c);
    if (onSegment(a, b, d)) s.insert(d);
    return {all(s)};
}
```

lineIntersection.h

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

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



sideOf.h

**Description:** Returns where *p* is as seen from *s* towards *e*. 1/0/-1 ⇔ 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"
template<class P>
int sideOf(P s, P e, P p) { return sgn(s.cross(e, p)); }
```

```
template<class P>
int sideOf(const P& s, const P& e, const P& p, double eps) {
    auto a = (e-s).cross(p-s);
    double l = (e-s).dist()*eps;
    return (a > l) - (a < -l);
}
```

OnSegment.h

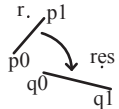
**Description:** Returns true iff p lies on the line segment from s to e. Use (segDist(s,e,p)<=epsilon) instead when using Point<double>.

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

linearTransformation.h

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

```
"Point.h"
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 || (y == 0 && x < 0);
    }
    Angle t90() const { return {-y, x, t + (half() && x >= 0)}; }
    Angle t180() const { return {-x, -y, t + half()}; }
    Angle t360() const { return {x, y, t + 1}; }
};
bool operator<(Angle a, Angle b) {
    // add a.dist2() and b.dist2() to also compare distances
    return make_tuple(a.t, a.half(), a.y * (1l)b.x) <
        make_tuple(b.t, b.half(), a.x * (1l)b.y);
}
```

```
// Given two points, this calculates the smallest angle between
// them, i.e., the angle that covers the defined line segment.
pair<Angle, Angle> segmentAngles(Angle a, Angle b) {
    if (b < a) swap(a, b);
    return (b < a.t180() ?
        make_pair(a, b) : make_pair(b, a.t360()));
}
Angle operator+(Angle a, Angle b) { // point a + vector b
    Angle r(a.x + b.x, a.y + b.y, a.t);
    if (a.t180() < r) r.t--;
    return r.t180() < a ? r.t360() : r;
}
Angle angleDiff(Angle a, Angle b) { // angle b - angle a
    int tu = b.t - a.t; a.t = b.t;
    return {a.x*b.x + a.y*b.y, a.x*b.y - a.y*b.x, tu - (b < a)};
}
```

8.2 Circles

CircleIntersection.h

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

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

CircleTangents.h

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

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

CirclePolygonIntersection.h

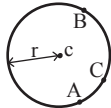
**Description:** Returns the area of the intersection of a circle with a ccw polygon.  
**Time:** O(n)

```
"../content/geometry/Point.h"
typedef Point<double> P;
#define arg(p, q) atan2(p.cross(q), p.dot(q))
double circlePoly(P c, double r, vector<P> ps) {
    auto tri = [&](P p, P q) {
        auto r2 = r * r / 2;
        P d = q - p;
        auto a = d.dot(p)/d.dist2(), b = (p.dist2()-r*r)/d.dist2();
        auto det = a * a - b;
        if (det <= 0) return arg(p, q) * r2;
        auto s = max(0., -a-sqrt(det)), t = min(1., -a+sqrt(det));
        if (t < 0 || 1 <= s) return arg(p, q) * r2;
        P u = p + d * s, v = p + d * t;
        return arg(p,u) * r2 + u.cross(v)/2 + arg(v,q) * r2;
    };
    auto sum = 0.0;
    rep(i,0,sz(ps))
        sum += tri(ps[i] - c, ps[(i + 1) % sz(ps)] - c);
    return sum;
}
```

circumcircle.h

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

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



MinimumEnclosingCircle.h

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



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

```
"circumcircle.h"
d41d8c, 17 lines

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

### 8.3 Polygons

**InsidePolygon.h**  
**Description:** Returns true if p lies within the polygon. If strict is true, it returns false for points on the boundary. The algorithm uses products in intermediate steps so watch out for overflow.  
**Usage:** vector<P> v = {P{4,4}, P{1,2}, P{2,1}};  
bool in = inPolygon(v, P{3, 3}, false);  
**Time:**  $\mathcal{O}(n)$

```
"Point.h", "OnSegment.h", "SegmentDistance.h"
d41d8c, 11 lines

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

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

```
"Point.h"
d41d8c, 6 lines

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

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

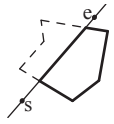
```
"Point.h"
d41d8c, 9 lines

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

**PolygonCut.h**  
**Description:** Returns a vector with the vertices of a polygon with everything to the left of the line going from s to e cut away.  
**Usage:** vector<P> p = ...;  
p = polygonCut(p, P(0,0), P(1,0));

```
"Point.h", "lineIntersection.h"
d41d8c, 13 lines


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



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

```
"Point.h"
d41d8c, 13 lines

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



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

**PointInsideHull.h**  
**Description:** Determine whether a point t lies inside a convex hull (CCW order, with no collinear points). Returns true if point lies within the hull. If strict is true, points on the boundary aren't included.  
**Time:**  $\mathcal{O}(\log N)$

```
"Point.h", "sideOf.h", "OnSegment.h"
d41d8c, 14 lines

typedef Point<ll> P;
```

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

**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) sgn(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 ls = cmp(lo + 1, lo), ms = cmp(m + 1, m);
        (ls < ms || (ls == ms && ls == cmp(lo, m)) ? hi : lo) = m;
    }
    return lo;
}

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

### kdTree.h

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

"Point.h"	d41d8c, 63 lines
<pre>typedef long long T; typedef Point&lt;T&gt; P; const T INF = numeric_limits&lt;T&gt;::max();</pre>	

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

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

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

    Node(vector<P>&& vp) : pt(vp[0]) {
        for (P p : vp) {
            x0 = min(x0, p.x); x1 = max(x1, p.x);
            y0 = min(y0, p.y); y1 = max(y1, p.y);
        }
        if (vp.size() > 1) {
            // split on x if width >= height (not ideal...)
            sort(all(vp), x1 - x0 >= y1 - y0 ? on_x : on_y);
            // divide by taking half the array for each child (not
            // best performance with many duplicates in the middle)
            int half = sz(vp)/2;
            first = new Node({vp.begin(), vp.begin() + half});
            second = new Node({vp.begin() + half, vp.end()});
        }
    }
};
```

```
struct KDTree {
    Node* root;
    KDTree(const vector<P>& vp) : root(new Node({all(vp)})) {}

    pair<T, P> search(Node *node, const P& p) {
        if (!node->first) {
            // uncomment if we should not find the point itself:
```

```
        // if (p == node->pt) return {INF, P()};
        return make_pair((p - node->pt).dist2(), node->pt);
    }

    Node *f = node->first, *s = node->second;
    T bfirst = f->distance(p), bsec = s->distance(p);
    if (bfirst > bsec) swap(bsec, bfirst), swap(f, s);

    // search closest side first, other side if needed
    auto best = search(f, p);
    if (bsec < best.first)
        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);
}
};
```

### FastDelaunay.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], ...}, all counter-clockwise.

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

"Point.h"	d41d8c, 88 lines
<pre>typedef Point&lt;ll&gt; P; typedef struct Quad* Q; typedef __int128_t lll; // (can be ll if coords are &lt; 2e4) P arb(LLONG_MAX, LLONG_MAX); // not equal to any other point  struct Quad {     Q rot, o; P p = arb; bool mark;     P&amp; F() { return r()-&gt;p; }     Q&amp; r() { return rot-&gt;rot; }     Q prev() { return rot-&gt;o-&gt;rot; }     Q next() { return r()-&gt;prev(); } } *H;  bool circ(P p, P a, P b, P c) { // is p in the circumcircle?     lll 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 &gt; 0; }  Q makeEdge(P orig, P dest) {     Q r = H ? H : new Quad{new Quad{new Quad{new Quad{0}}}};     H = r-&gt;o; r-&gt;r()-&gt;r() = r;     rep(i,0,4) r = r-&gt;rot, r-&gt;p = arb, r-&gt;o = i &amp; 1 ? r : r-&gt;r();     r-&gt;p = orig; r-&gt;F() = dest;     return r; }  void splice(Q a, Q b) {     swap(a-&gt;o-&gt;rot-&gt;o, b-&gt;o-&gt;rot-&gt;o); swap(a-&gt;o, b-&gt;o); }  Q connect(Q a, Q b) {     Q q = makeEdge(a-&gt;F(), b-&gt;p);     splice(q, a-&gt;next());     splice(q-&gt;r(), b);     return q; }  pair&lt;Q,Q&gt; rec(const vector&lt;P&gt;&amp; s) {     if (sz(s) &lt;= 3) {         Q a = makeEdge(s[0], s[1]), b = makeEdge(s[1], s.back());         if (sz(s) == 2) return { a, a-&gt;r() };</pre>	

```
        splice(a->r(), b);
        auto side = s[0].cross(s[1], s[2]);
        Q c = side ? connect(b, a) : 0;
        return {side < 0 ? c->r() : a, side < 0 ? c : b->r() };
    }

#define H(e) e->F(), e->p
#define valid(e) (e->F().cross(H(base)) > 0)
    Q A, B, ra, rb;
    int half = sz(s) / 2;
    tie(ra, A) = rec({all(s) - half});
    tie(B, rb) = rec({sz(s) - half + all(s)});
    while ((B->p.cross(H(A)) < 0 && (A = A->next())) ||
            (A->p.cross(H(B)) > 0 && (B = B->r()->o)));
    Q base = connect(B->r(), A);
    if (A->p == ra->p) ra = base->r();
    if (B->p == rb->p) rb = base;

#define DEL(e, init, dir) 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; \
    }
    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 };
}
```

## 8.5 3D

### PolyhedronVolume.h

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

d41d8c, 6 lines
<pre>template&lt;class V, class L&gt; double signedPolyVolume(const V&amp; p, const L&amp; trilst) {     double v = 0;     for (auto i : trilst) v += p[i.a].cross(p[i.b]).dot(p[i.c]);     return v / 6; }</pre>

### Point3D.h

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

d41d8c, 32 lines
<pre>template&lt;class T&gt; struct Point3D {     typedef Point3D P;     typedef const P&amp; R;</pre>

```
T x, y, z;
explicit Point3D(T x=0, T y=0, T z=0) : x(x), y(y), z(z) {}
bool operator<(R p) const {
    return tie(x, y, z) < tie(p.x, p.y, p.z); }
bool operator==(R p) const {
    return tie(x, y, z) == tie(p.x, p.y, p.z); }
P operator+(R p) const { return P(x+p.x, y+p.y, z+p.z); }
P operator-(R p) const { return P(x-p.x, y-p.y, z-p.z); }
P operator*(T d) const { return P(x*d, y*d, z*d); }
P operator/(T d) const { return P(x/d, y/d, z/d); }
T dot(R p) const { return x*p.x + y*p.y + z*p.z; }
P cross(R p) const {
    return P(y*p.z - z*p.y, z*p.x - x*p.z, x*p.y - y*p.x);
}
T dist2() const { return x*x + y*y + z*z; }
double dist() const { return sqrt((double)dist2()); }
//Azimuthal angle (longitude) to x-axis in interval [-pi, pi]
double phi() const { return atan2(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:**  $\mathcal{O}(n^2)$

```
"Point3D.h" d41d8c, 49 lines
```

```
typedef Point3D<double> P3;
```

```
struct PR {
    void ins(int x) { (a == -1 ? a : b) = x; }
    void rem(int x) { (a == x ? a : b) = -1; }
    int cnt() { return (a != -1) + (b != -1); }
    int a, b;
};
```

```
struct F { P3 q; int a, b, c; };
```

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

    rep(i,4,sz(A)) {
        rep(j,0,sz(FS)) {
            F f = FS[j];
            if(f.q.dot(A[i]) > f.q.dot(A[f.a])) {
                E(a,b).rem(f.c);
                E(a,c).rem(f.b);
```

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

sphericalDistance.h

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

```
d41d8c, 8 lines
```

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

Strings (9)

AhoCorasick.cpp

**Description:** query retorna o somatorio do numero de matches de todas as stringuinhas na stringona  
**Time:** insert in  $\mathcal{O}(|s| \log SIGMA)$ .

```
d41d8c, 40 lines
```

```
namespace aho {
    map<char, int> to[MAX];
    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 l = link[i];
                while (l != -1 and !to[l].count(c)) l = link[l];
                link[j] = l == -1 ? 0 : to[l][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) {
    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];
        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-1]] !=
            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;
}
```

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 l = 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[l+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;
}
l = 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 <= n && i-k >= 0 && s[i+k-1] == s[i-k]) k++;
    d2[i] = --k;
    if (i+k-1 > r) l = 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

**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  $\mathcal{O}(n)$

d41d8c, 35 lines

```
vector<string> duval(string const &s) {
    int n = s.size();
    int i = 0;
    vector<string> factorization;
    while (i < n) {
        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 <= 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 <= k) i += j - k;
return s.substr(ans, n / 2);
}
```

Various (10)

10.1 Intervals

IntervalContainer.h

**Description:** Add and remove intervals from a set of disjoint intervals. Will merge the added interval with any overlapping intervals in the set when adding. Intervals are [inclusive, exclusive).  
**Time:**  $\mathcal{O}(\log N)$

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]; });
    T cur = G.first;
    int at = 0;
    while (cur < G.second) { // (A)
        pair<T, int> mx = make_pair(cur, -1);
        while (at < sz(I) && I[S[at]].first <= cur) {
            mx = max(mx, make_pair(I[S[at]].second, S[at]));
            at++;
        }
        if (mx.second == -1) return {};
        cur = mx.first;
        R.push_back(mx.second);
    }
```

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){...});  
**Time:**  $\mathcal{O}(k \log \frac{n}{k})$

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) {
 g(i, to, p);
 i = to; p = q;
 } else {
 int mid = (from + to) >> 1;
 rec(from, mid, f, g, i, p, f(mid));
 rec(mid+1, to, f, g, i, p, q);
 }
}

template<class F, class G>
void constantIntervals(int from, int to, F f, G g) {
 if (to <= from) return;
 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) < ... < f(i) ≥ ... ≥ f(b). To reverse which of the sides allows non-strict inequalities, change the < marked with (A) to <=, and reverse the loop at (B). To minimize f, change it to >, also at (B).
Usage: int ind = ternSearch(0,n-1,[&](int i){return a[i];});
Time: O(log(b-a))

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)
 return a;
}

LIS.h
Description: Compute indices for the longest increasing subsequence.
Time: 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<p> res;
 rep(i,0,sz(S)) {
 // change 0 -> i for longest non-decreasing subsequence
 auto it = lower\_bound(all(res), p{S[i], 0});
 if (it == res.end()) res.emplace\_back(), it = res.end()-1;
 \*it = {S[i], i};
 prev[i] = it == res.begin() ? 0 : (it-1)->second;
 }

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

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

d41d8c, 16 lines

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

10.3 Dynamic programming

KnuthDP.h

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

d41d8c, 18 lines

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

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

10.4 Debugging tricks

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

10.5 Optimization tricks

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

10.5.1 Bit hacks

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

10.5.2 Pragmas

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

FastMod.h

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

d41d8c, 8 lines

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

FastInput.h

**Description:** Read an integer from stdin. Usage requires your program to pipe in input from file.  
**Usage:** `./a.out < input.txt`  
**Time:** About 5x as fast as `cin/scanf`.

d41d8c, 17 lines

```
inline char gc() { // like getchar()
    static char buf[1 << 16];
    static size_t bc, be;
    if (bc >= be) {
```

```
        buf[0] = 0, bc = 0;
        be = fread(buf, 1, sizeof(buf), stdin);
    }
    return buf[bc++]; // returns 0 on EOF
}

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

BumpAllocator.h

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

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*) {}
```

# Techniques (A)

techniques.txt	159 lines
Recursion	
Divide and conquer	
Finding interesting points in N log N	
Algorithm analysis	
Master theorem	
Amortized time complexity	
Greedy algorithm	
Scheduling	
Max contiguous 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	
2-SAT	
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)	
Combinatorics	

## techniques

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
Quadratic 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
Quadtrees
KD-trees
All segment-segment intersection
Sweeping
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