

# Indian Institute of Technology KGP

# Aloo Paratha

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### Contest (1)

template.cpp #include <bits/stdc++.h> // #pragma GCC optimize("O3, unroll-loops") using namespace std; #define fr(i, a, b) for(int i = a; i < (b); ++i) **#define** rev(i, a, b) **for** (ll i = a; i >= b; i--) #define push\_back pb #define all(x) begin(x), end(x) #define sz(x) (int)(x).size() using 11 = long long; #define int long long typedef pair<int, int> pii; typedef vector<int> vi; **const** 11 mod1 = 1e9+7, mod2 = 998244353; const 11 INF = 1e9, INF2 = 1e18; #define ff first #define ss second void solve(){} int32 t main() ios\_base::sync\_with\_stdio(0); cin.tie(0); cout.tie(NULL); #ifndef ONLINE JUDGE freopen("input.txt", "r", stdin); freopen("output.txt", "w", stdout); #endif int t=1; cin >> t; while(t--) solve(); return 0;

### Mathematics (2)

### 2.1 Equations

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

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

$$ax + by = e$$

$$cx + dy = f \Rightarrow x = \frac{ed - bf}{ad - bc}$$

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

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

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

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

### 2.2 Recurrences

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

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

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

### 2.3 Trigonometry

30 lines

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

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

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

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

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

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

### 2.4 Geometry

### 2.4.1 Triangles

Side lengths: a, b, c

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

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

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

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

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

Length of bisector (divides angles in two):

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

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

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

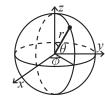
#### 2.4.2 Quadrilaterals

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

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

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

#### 2.4.3 Spherical coordinates



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

### 2.5 Derivatives/Integrals

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

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

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

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

Integration by parts:

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

#### template OrderStatisticTree

### 2.6 Sums

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

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

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

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

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

#### 2.7 Series

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

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

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

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

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

### 2.8 Probability theory

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

Expectation is linear:

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

For independent X and Y,

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

# 2.8.1 Discrete distributions Binomial distribution

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

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

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

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

#### First success distribution

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

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

#### Poisson distribution

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

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

# 2.8.2 Continuous distributions Uniform distribution

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

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

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

### Exponential distribution

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

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

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

#### Normal distribution

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

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

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

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

#### 2.9 Markov chains

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

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

### Data structures (3)

#### OrderStatisticTree.h

**Description:** A set (not multiset!) with support for finding the n'th element, and finding the index of an element. To get a map, change null-type. **Time:**  $\mathcal{O}(\log N)$ 

```
<ext/pb.ds/assoc.container.hpp>, <ext/pb.ds/tree.policy.hpp> 819d08, 15 lines
using namespace __gnu_pbds;
```

```
HashMap.h
```

```
Description: Hash map with mostly the same API as unordered_map, but
~3x faster. Uses 1.5x memory. Initial capacity must be a power of 2 (if
provided).
```

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

#### SegmentTree.h

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

Time:  $\mathcal{O}(\log N)$ 0f4bdb, 19 lines

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

#### LazySegmentTree.h

```
Description: LazySegmentTree.h
                                                     0666ae. 70 lines
class LazySegmentTree {
private:
    vector<int> t, lazy;
    int n;
    void build(vector<int>& a, int v, int tl, int tr) {
       if (t.1 == t.r) {
            t[v] = a[t1];
       } else {
            int tm = (tl + tr) / 2;
            build(a, v*2, tl, tm);
            build(a, v*2+1, tm+1, tr);
            t[v] = combine(t[v*2], t[v*2 + 1]);
    void push(int v) {
       t[v*2] += lazy[v];
        lazv[v*2] += lazv[v];
        t[v*2+1] += lazy[v];
        lazy[v*2+1] += lazy[v];
       lazy[v] = 0;
    void update(int v, int tl, int tr, int l, int r, int addend
        ) {
        if (1 > r)
            return;
        if (1 == t1 && tr == r) {
```

```
t[v] += addend;
            lazy[v] += addend;
        } else {
            push (v);
            int tm = (t1 + tr) / 2;
            update(v*2, t1, tm, 1, min(r, tm), addend);
            update(v*2+1, tm+1, tr, max(1, tm+1), r, addend);
            t[v] = combine(t[v*2], t[v*2+1]);
    int query(int v, int tl, int tr, int l, int r) {
        if (1 > r)
            return -INF;
        if (1 == t1 && tr == r)
            return t[v];
        push (v);
        int tm = (tl + tr) / 2;
        return combine (query (v*2, tl, tm, l, min(r, tm)),
                       query (v*2+1, tm+1, tr, max(1, tm+1), r))
    int combine(int a, int b) {
        return max(a, b); // Change this according to your
             requirement
public:
    LazySegmentTree(vector<int>& a) {
       n = a.size();
       t.assign(4*n, 0);
        lazy.assign(4*n, 0);
        build(a, 1, 0, n-1);
    void update(int 1, int r, int addend) {
        update(1, 0, n-1, 1, r, addend);
    int query(int 1, int r) {
        return query(1, 0, n-1, 1, r);
};
UnionFind.h
Description: UnionFind.h
                                                     3624b6, 17 lines
struct DSU
  vi par, size;
   DSU(int n) : par(n), size(n, 1) { iota(par.begin(), par.end
  int find(int x) {return x == par[x] ? x : par[x] = find(par[x
      ]);}
  void merge(int x, int y)
        int nx = find(x);
        int ny = find(y);
        if (nx!=ny)
            if(size[nx]<size[ny]) swap(nx,ny);</pre>
            par[ny] = nx;
            size[nx] += size[ny];
};
```

```
UnionFindRollback.h
```

Description: UnionFindRollback.h

6c5dd9, 34 lines

```
class DSU {
 private:
  vector<int> p, sz;
  // stores previous unites
  vector<pair<int &, int>> history;
  public:
  DSU(int n) : p(n), sz(n, 1) { iota(p.beqin(), p.end(), 0); }
  int get(int x) { return x == p[x] ? x : get(p[x]); }
  void unite(int a, int b) {
    a = get(a);
    b = qet(b);
    if (a == b) { return; }
    if (sz[a] < sz[b]) { swap(a, b); }</pre>
    // save this unite operation
    history.push_back({sz[a], sz[a]});
    history.push_back({p[b], p[b]});
    p[b] = a;
    sz[a] += sz[b];
  int snapshot() { return history.size(); }
  void rollback(int until) {
    while (snapshot() > until) {
      history.back().first = history.back().second;
      history.pop_back();
};
SubMatrix.h
```

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

```
Usage: SubMatrix<int> m(matrix);
m.sum(0, 0, 2, 2); // top left 4 elements
Time: \mathcal{O}(N^2+Q)
```

c59ada, 13 lines

```
template < class T >
struct SubMatrix {
 vector<vector<T>> p;
 SubMatrix(vector<vector<T>>& v) {
    int R = sz(v), C = sz(v[0]);
    p.assign(R+1, vector<T>(C+1));
    rep(r, 0, R) rep(c, 0, C)
      p[r+1][c+1] = v[r][c] + p[r][c+1] + p[r+1][c] - p[r][c];
 T sum(int u, int 1, int d, int r) {
    return p[d][r] - p[d][l] - p[u][r] + p[u][l];
};
```

#### Matrix.h

Description: Matrix.h

5742e0, 32 lines

```
template<class T> struct Matrix {
 typedef Matrix M;
 vector<vector<T>> d;
   Matrix(int n) {
        d.resize(n, vector < T > (n, 0));
 M operator*(const M& m) const {
   M a(m.d.size());
```

```
int N = m.d.size();
    rep(i,0,N) rep(j,0,N)
      rep(k, 0, N) \{a.d[i][j] += (d[i][k]*m.d[k][j]) mod1; a.d[i][
           j]%=mod1;}
    return a;
  vector<T> operator*(const vector<T>& vec) const {
        int N = this->d.size();
    vector<T> ret(N);
    rep(i, 0, N) rep(j, 0, N) {ret[i] += (d[i][j] * vec[j]) %mod1;}
        ret[i]%=mod1;}
    return ret;
  M operator^(ll p) const {
    assert (p >= 0);
   M a(this->d.size()), b(*this);
        int N = this->d.size();
    rep(i, 0, N) \ a.d[i][i] = 1;
    while (p) {
     if (p&1) a = a*b;
     b = b*b;
     p >>= 1;
    return a;
};
```

#### LineContainer.h

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

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

#### Treap.h

with log-time splits/joins, and is easy to augment with additional data.

```
Description: A short self-balancing tree. It acts as a sequential container
Time: \mathcal{O}(\log N)
                                                                    9556fc, 55 lines
```

```
struct Node {
 Node *1 = 0, *r = 0;
  int val, y, c = 1;
 Node(int val) : val(val), y(rand()) {}
 void recalc();
int cnt(Node* n) { return n ? n->c : 0; }
void Node::recalc() { c = cnt(1) + cnt(r) + 1; }
template<class F> void each(Node* n, F f) {
 if (n) { each(n->1, f); f(n->val); each(n->r, f); }
pair<Node*, Node*> split(Node* n, int k) {
 if (!n) return {};
 if (cnt(n->1) \Rightarrow k) { // "n->val>= k" for lower_bound(k)
    auto pa = split(n->1, k);
   n->1 = pa.second;
   n->recalc();
   return {pa.first, n};
    auto pa = split(n->r, k - cnt(n->1) - 1); // and just "k"
    n->r = pa.first;
   n->recalc();
   return {n, pa.second};
Node* merge(Node* 1, Node* r) {
 if (!1) return r;
  if (!r) return 1;
 if (1->y > r->y) {
   1->r = merge(1->r, r);
    1->recalc();
    return 1;
  } else {
    r->1 = merge(1, r->1);
    r->recalc();
    return r;
Node* ins(Node* t, Node* n, int pos) {
 auto pa = split(t, pos);
 return merge(merge(pa.first, n), pa.second);
// Example application: move the range (l, r) to index k
void move(Node*& t, int 1, int r, int k) {
 Node *a, *b, *c;
 tie(a,b) = split(t, 1); tie(b,c) = split(b, r - 1);
 if (k \le 1) t = merge(ins(a, b, k), c);
 else t = merge(a, ins(c, b, k - r));
```

#### FenwickTree.h

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

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

```
struct FT {
  vector<ll> s;
 FT(int n) : s(n) {}
 void update(int pos, 11 dif) { // a[pos] \leftarrow= dif
    for (; pos < sz(s); pos |= pos + 1) s[pos] += dif;</pre>
 11 query (int pos) { // sum of values in [0, pos)
```

```
11 \text{ res} = 0;
    for (; pos > 0; pos &= pos - 1) res += s[pos-1];
    return res;
 int lower_bound(ll sum) \{// min \ pos \ st \ sum \ of \ [0, \ pos] >= sum
    // Returns n if no sum is \geq sum, or -1 if empty sum is.
    if (sum \leq 0) return -1;
    int pos = 0:
    for (int pw = 1 << 25; pw; pw >>= 1) {
      if (pos + pw <= sz(s) && s[pos + pw-1] < sum)
        pos += pw, sum -= s[pos-1];
    return pos;
};
```

#### FenwickTree2d.h

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

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

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

#### RMQ.h

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

```
rmg.guery(inclusive, exclusive);
Time: \mathcal{O}(|V|\log|V|+Q)
```

510c32, 16 lines

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

#### MoQueries.h

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

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

### Numerical (4)

### 4.1 Polynomials and recurrences

Polynomial.h

c9b7b0, 17 lines

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

```
void diff() {
    rep(i, 1, sz(a)) a[i-1] = i*a[i];
    a.pop_back();
  void divroot(double x0) {
    double b = a.back(), c; a.back() = 0;
    for(int i=sz(a)-1; i--;) c = a[i], a[i] = a[i+1]*x0+b, b=c;
    a.pop_back();
};
PolvRoots.h
Description: Finds the real roots to a polynomial.
Usage: polyRoots(\{\{2,-3,1\}\},-1e9,1e9) // solve x^2-3x+2=0
Time: \mathcal{O}\left(n^2\log(1/\epsilon)\right)
"Polynomial.h"
                                                        b00bfe, 23 lines
vector<double> polyRoots(Poly p, double xmin, double xmax) {
 if (sz(p.a) == 2) { return {-p.a[0]/p.a[1]}; }
 vector<double> ret;
 Polv der = p:
 der.diff();
  auto dr = polyRoots(der, xmin, xmax);
  dr.push back(xmin-1);
  dr.push back(xmax+1);
  sort (all (dr));
  rep(i, 0, sz(dr) - 1) {
    double l = dr[i], h = dr[i+1];
    bool sign = p(1) > 0;
    if (sign ^{(p(h) > 0)}) {
      rep(it,0,60) { // while (h - l > 1e-8)
        double m = (1 + h) / 2, f = p(m);
        if ((f <= 0) ^ sign) l = m;
        else h = m;
      ret.push_back((1 + h) / 2);
 return ret;
```

#### PolvInterpolate.h

**Description:** 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;
```

#### BerlekampMassev.h

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

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

```
Time: \mathcal{O}(N^2)
"../number-theory/ModPow.h"
```

```
96548b, 20 lines
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;
 11 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; 11 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());
  for (11& x : C) x = (mod - x) % mod;
 return C:
```

#### LinearRecurrence.h

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

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

```
typedef vector<11> Poly;
11 linearRec(Poly S, Poly tr, 11 k) {
 int n = sz(tr);
  auto combine = [&] (Poly a, Poly b) {
    Poly res(n \star 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);
 rep(i, 0, n) res = (res + pol[i + 1] * S[i]) % mod;
 return res;
```

### 4.2 Optimization

#### 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. 4756fc, 7 lines

```
template < class F >
double quad(double a, double b, F f, const int n = 1000) {
 double h = (b - a) / 2 / n, v = f(a) + f(b);
 rep(i,1,n*2)
```

```
v += f(a + i*h) * (i&1 ? 4 : 2);
return v * h / 3;
```

#### Simplex.h

T solve(vd &x) {

**Description:** Solves a general linear maximization problem: maximize  $c^T x$ subject to Ax < b, x > 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;</pre>
       if (r == -1 \mid | MP(D[i][n+1] / D[i][s], B[i])
                     < MP(D[r][n+1] / D[r][s], B[r])) r = i;
     if (r == -1) return false;
     pivot(r, s);
```

```
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;</pre>
  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;
```

#### Matrices

#### 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[i][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:  $\mathcal{O}(N^3)$ 

```
3313dc, 18 lines
const 11 mod = 12345;
ll det(vector<vector<ll>>& a) {
 int n = sz(a); ll ans = 1;
 rep(i,0,n) {
   rep(j,i+1,n) {
      while (a[j][i] != 0) { // gcd step}
       11 t = a[i][i] / a[j][i];
       if (t) rep(k,i,n)
         a[i][k] = (a[i][k] - a[j][k] * t) % mod;
        swap(a[i], a[j]);
        ans \star = -1;
   ans = ans * a[i][i] % mod;
   if (!ans) return 0;
 return (ans + mod) % mod;
```

#### SolveLinear.h

**Description:** Solves A \* x = b. If there are multiple solutions, an arbitrary one is returned. Returns rank, or -1 if no solutions. Data in A and b is lost. Time:  $\mathcal{O}\left(n^2m\right)$ 44c9ab, 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"
                                                        08e495, 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:: }
```

#### MatrixInverse.h

**Description:** Invert matrix A. Returns rank; result is stored in A unless singular (rank < n). Can easily be extended to prime moduli; for prime powers, repeatedly set  $A^{-1} = A^{-1}(2I - AA^{-1}) \pmod{p^k}$  where  $A^{-1}$  starts as the inverse of A mod p, and k is doubled in each step. Time:  $\mathcal{O}\left(n^3\right)$ 

```
ebfff6, 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;</pre>
  A[i].swap(A[r]); tmp[i].swap(tmp[r]);
  rep(j,0,n)
   swap(A[j][i], A[j][c]), swap(tmp[j][i], tmp[j][c]);
  swap(col[i], col[c]);
  double v = A[i][i];
  rep(j,i+1,n) {
   double f = A[j][i] / v;
   A[j][i] = 0;
   rep(k,i+1,n) A[j][k] -= f*A[i][k];
   rep(k,0,n) tmp[j][k] \rightarrow 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] \rightarrow v*tmp[i][k];
rep(i,0,n) rep(j,0,n) A[col[i]][col[j]] = tmp[i][j];
return n;
```

#### 4.4 Fourier transforms

FastFourierTransform.h

**Description:** fft(a) computes  $\hat{f}(k) = \sum_x a[x] \exp(2\pi i \cdot kx/N)$  for all k. N must be a power of 2. Useful for convolution: conv(a, b) = c, where  $c[x] = \sum_x a[i]b[x-i]$ . For convolution of complex numbers or more than two vectors: FFT, multiply pointwise, divide by n, reverse(start+1, end), FFT back. Rounding is safe if  $(\sum_i a_i^2 + \sum_i b_i^2) \log_2 N < 9 \cdot 10^{14}$  (in practice  $10^{16}$ ; higher for random inputs). Otherwise, use NTT/FFTMod. **Time:**  $\mathcal{O}(N \log N)$  with N = |A| + |B| (~1s for  $N = 2^{22}$ )

```
typedef complex<double> C;
typedef vector<double> vd;
void fft(vector<C>& a) {
  int n = sz(a), L = 31 - __builtin_clz(n);
  static vector<complex<long double>> R(2, 1);
  static vector<C> rt(2, 1); // (^ 10% faster if double)
  for (static int k = 2; k < n; k \neq 2) {
   R.resize(n); rt.resize(n);
   auto x = polar(1.0L, acos(-1.0L) / k);
   rep(i,k,2*k) rt[i] = R[i] = i&1 ? R[i/2] * x : R[i/2];
  rep(i,0,n) \ rev[i] = (rev[i / 2] | (i \& 1) << L) / 2;
  rep(i,0,n) if (i < rev[i]) swap(a[i], a[rev[i]]);
  for (int k = 1; k < n; k *= 2)
   for (int i = 0; i < n; i += 2 * k) rep(j,0,k) {
     Cz = rt[j+k] * a[i+j+k]; // (25\% faster if hand-rolled)
     a[i + j + k] = a[i + j] - z;
     a[i + j] += z;
vd conv(const vd& a, const vd& b) {
 if (a.empty() || b.empty()) return {};
  vd res(sz(a) + sz(b) - 1);
  int L = 32 - \underline{\text{builtin\_clz}(\text{sz(res)})}, n = 1 << L;
  vector<C> in(n), out(n);
  copy(all(a), begin(in));
  rep(i,0,sz(b)) in[i].imag(b[i]);
  fft(in);
  for (C& x : in) x \star = x;
  rep(i,0,n) out[i] = in[-i & (n-1)] - conj(in[i]);
  rep(i, 0, sz(res)) res[i] = imag(out[i]) / (4 * n);
```

```
return res;
FastSubsetTransform.h
Description: Transform to a basis with fast convolutions of the form
c[z] \, = \, \sum_{z=x \oplus y} a[x] \cdot b[y], \text{ where } \oplus \text{ is one of AND, OR, XOR.} The size
of a must be a power of two.
Time: \mathcal{O}(N \log N)
                                                          464cf3, 16 lines
void FST(vi& a, bool inv) {
  for (int n = sz(a), step = 1; step < n; step *= 2) {
    for (int i = 0; i < n; i += 2 * step) rep(j,i,i+step) {</pre>
      int &u = a[j], &v = a[j + step]; tie(u, v) =
        inv ? pii(v - u, u) : pii(v, u + v); // AND
        inv ? pii(v, u - v) : pii(u + v, u); // OR
        pii(u + v, u - v);
 if (inv) for (int& x : a) x /= sz(a); // XOR only
vi conv(vi a, vi b) {
 FST(a, 0); FST(b, 0);
  rep(i,0,sz(a)) a[i] *= b[i];
 FST(a, 1); return a;
```

### Number theory (5)

#### 5.1 Modular arithmetic

\*this \*= other.inv(); }

Modular Arithmetic.h

```
Description: Modular Arithmetic.h
```

9f3a25, 43 lines

```
template <const int64 t MOD = mod1>
struct modint {
    int64 t value:
   modint() = default;
   modint(int64_t value_) : value(value_%MOD) {}
   modint<MOD> operator + (modint<MOD> other) const { int64_t
        c = this->value + other.value; return modint<MOD>(c >=
         MOD ? c - MOD : c); }
   modint<MOD> operator - (modint<MOD> other) const { int64_t
        c = this->value - other.value; return modint<MOD>(c <</pre>
           0 ? c + MOD : c); }
    modint<MOD> operator * (modint<MOD> other) const { int64_t
        c = (int64 t)this->value * other.value % MOD; return
        modint<MOD>(c < 0 ? c + MOD : c); }
    modint<MOD> & operator += (modint<MOD> other) { this->value
         += other.value; if (this->value >= MOD) this->value
         -= MOD; return *this; }
    modint<MOD> & operator -= (modint<MOD> other) { this->value
         -= other.value; if (this->value < 0) this->value
         += MOD; return *this; }
    modint<MOD> & operator *= (modint<MOD> other) { this->value
         = (int64_t)this->value * other.value % MOD; if (this
         ->value < 0) this->value += MOD; return *this; }
   modint<MOD> operator - () const { return modint<MOD>(this->
        value ? MOD - this->value : 0); }
   modint < MOD > pow(uint 64_t k) const { modint < MOD > x = *this,
        y = 1; for (; k; k >>= 1) { if (k & 1) y *= x; x *= x;
         } return y; }
   modint<MOD> inv() const { return pow(MOD - 2); } // MOD
         must be a prime*
    modint<MOD> operator / (modint<MOD> other) const { return
         *this * other.inv(); }
    modint<MOD> operator /= (modint<MOD> other)
                                                      { return
```

```
bool operator == (modint<MOD> other) const { return value
         == other.value; }
    bool operator != (modint<MOD> other) const { return value
         != other.value; }
    bool operator < (modint<MOD> other) const { return value <</pre>
         other.value; }
    bool operator > (modint<MOD> other) const { return value >
        other.value; }
    friend modint<MOD> operator * (int64_t value, modint<MOD> n
        ) { return modint<MOD>(value % MOD) * n; }
    friend istream & operator >> (istream & in, modint<MOD> &n)
          { return in >> n.value; }
    friend ostream & operator << (ostream & out, modint<MOD> n)
          { return out << n.value; }
using mint = modint<>;
template < const int64_t mod = mod1>
struct combi{
  int n; vector<mint> facts, finvs, invs;
  combi(int _n): n(_n), facts(_n), finvs(_n), invs(_n) {
    facts[0] = finvs[0] = 1;
    invs[1] = 1;
    for (int i = 2; i < n; i++) invs[i] = invs[mod % i] * (-</pre>
        mod / i);
    for(int i = 1; i < n; i++) {</pre>
      facts[i] = facts[i - 1] * i;
      finvs[i] = finvs[i - 1] * invs[i];
  inline mint fact(int n) { return facts[n]; }
  inline mint finv(int n) { return finvs[n]; }
  inline mint inv(int n) { return invs[n]; }
  inline mint ncr(int n, int k) { return n < k or k < 0 ? 0:
       facts[n] * finvs[k] * finvs[n-k]; }
```

#### ModInverse.h

**Description:** Pre-computation of modular inverses. Assumes LIM  $\leq$  mod and that mod is a prime.

```
const 11 mod = 1000000007, LIM = 200000;
l1* inv = new l1[LIM] - 1; inv[1] = 1;
rep(i,2,LIM) inv[i] = mod - (mod / i) * inv[mod % i] % mod;
```

#### ModPow.h

b83e45, 8 lines

```
const 11 mod = 1000000007; // faster if const

11 modpow(11 b, 11 e) {
    11 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.  $\operatorname{modLog}(a,1,m)$  can be used to calculate the order of a.

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

```
11 modLog(l1 a, l1 b, l1 m) {
    l1 n = (l1) sqrt(m) + 1, e = 1, f = 1, j = 1;
    unordered_map<l1, l1> 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))</pre>
```

```
return n * i - A[e];
return -1:
```

#### ModSum.h

Description: Sums of mod'ed arithmetic progressions.

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

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

5c5bc5, 16 lines

```
typedef unsigned long long ull;
ull sumsq(ull to) { return to / 2 * ((to-1) | 1); }
ull divsum(ull to, ull c, ull k, ull m) {
  ull res = k / m * sumsq(to) + c / m * to;
  k %= m; c %= m;
  if (!k) return res;
  ull to2 = (to * k + c) / m;
  return res + (to - 1) * to2 - divsum(to2, m-1 - c, m, k);
11 modsum(ull to, 11 c, 11 k, 11 m) {
  c = ((c % m) + m) % m;
  k = ((k % m) + m) % m;
  return to * c + k * sumsq(to) - m * divsum(to, c, k, m);
```

#### ModMulLL.h

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

```
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 \star (ret < 0) - M \star (ret >= (11)M);
ull modpow(ull b, ull e, ull mod) {
 ull ans = 1;
  for (; e; b = modmul(b, b, mod), e /= 2)
   if (e & 1) ans = modmul(ans, b, mod);
  return ans;
```

#### ModSart.h

**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"
                                                        19a793, 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} \text{ or } 2^{(n+3)/8} * 2^{(n-1)/4} \text{ works if } p \% 8 == 5
 11 s = p - 1, n = 2;
  int r = 0, m;
  while (s % 2 == 0)
   ++r, s /= 2;
  while (modpow(n, (p-1) / 2, p) != p-1) ++n;
  11 x = modpow(a, (s + 1) / 2, p);
  11 b = modpow(a, s, p), g = modpow(n, s, p);
  for (;; r = m) {
   11 t = b;
    for (m = 0; m < r && t != 1; ++m)
     t = t * t % p;
    if (m == 0) return x;
    11 \text{ gs} = \text{modpow}(g, 1LL \ll (r - m - 1), p);
    q = qs * qs % p;
    x = x * gs % p;
```

```
b = b * q % p;
```

### 5.2 Primality

```
FastEratosthenes.h
Description: Prime sieve for generating all primes smaller than LIM.
Time: LIM=1e9 \approx 1.5s
int prime[1000000+1];
const int range = 1e6;
// Seive of Eratosthenes
void isPrime()
    for (int i = 2; i*i <= range; i++)</pre>
        if (prime[i] == 0)
             // its fine to start from i*i
             // 2*i,4*i,.. will be already marked by some
                  smaller prime numb
             for (int j = 111*i*i; j <= range; j+=i)</pre>
                 prime[j]=1;
    // if prime[i]==0 it means i is prime
// To find any n is prime or not
bool isPrime(ll n)
    if (n \le 1)
        return false:
    if (n \le 3)
        return true;
    if (n % 2 == 0 || n % 3 == 0)
        return false:
    for (11 i = 5; i * i <= n; i = i + 6)
        if (n \% i == 0 | | n \% (i + 2) == 0)
             return false:
    return true;
// code for finding number of divisors for all numbers from 2
     to N.
void divisors()
    // TC - O(NlogN)
    for (int i = 2; i < N; i++)
         for (int j = i; j < N; j += i)</pre>
             divis[j]++;
// Code for finding divisors of a number x
vi divisor(int x)
```

```
vi ans;
    int i = 1;
    while (i * i \le x)
        if (x % i == 0)
            ans.pb(i);
            if (x / i != i)
                ans.pb(x / i);
        i++;
    return ans;
// Code for finding factors of number x;
vector<pi> factor(int x)
    vector<pi> ans;
    for (int i = 2; i*i<=x; i++)</pre>
        if (x\%i == 0)
            int cnt=0;
            while (x\%i==0)
                cnt++;
                x/=i;
            ans.pb({i,cnt});
    if (x>1) ans.pb (\{x,1\});
    return ans;
// Fast Factorisation
// Code for finding all numbers upto X - (10^6)
// Normal Approach TG-O(N*sqrt(N))
// Store Lowest prime for each number using SIEVE IDEA
// Recursively call N/p till it reaches 1
// How it will calculate in log(X) becox max prime factors of X
      are log2(X)
void all_prime_factors(int X)
    // creating sp array
    int sp[1000000+1]; // initially zeroed
    int prime[1000000+1]; // initially zeroed
    const int range = 1e6;
    // Seive of Eratosthenes
    for (int i = 2; i <= range; i++)</pre>
        if(prime[i]==0)
            sp[i] = i;
            // its fine to start from i*i
            // 2*i,4*i,.. will be already marked by some
                 smaller prime numb
            for (int j = i*i; j \le range; j+=i)
                if(prime[j]==0)
                    prime[j]=1;
                    sp[j] = i;
```

```
// if prime[i]==0 it means i is prime
vector<int> factors[range+1];
for (int i = 2; i < X+1; i++)</pre>
    int num = i:
    while (num!=1)
        factors[i].push_back(sp[num]);
        num/=sp[num];
// final ans is stored in factors array
```

#### MillerRabin.h

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

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

```
"ModMulLL.h"
                                                       60dcd1, 12 lines
bool isPrime(ull n) {
  if (n < 2 || n % 6 % 4 != 1) return (n | 1) == 3;</pre>
  ull A[] = \{2, 325, 9375, 28178, 450775, 9780504, 1795265022\},
      s = \underline{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"
                                                     d8d98d, 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 1 = factor(x), r = factor(n / x);
 l.insert(l.end(), all(r));
  return 1;
```

### Divisibility

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

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

```
return v -= a/b * x, d;
CRT.h
Description: Chinese Remainder Theorem.
crt(a, m, b, n) computes x such that x \equiv a \pmod{m}, x \equiv b \pmod{n}. If
|a| < m and |b| < n, x will obey 0 \le x < \operatorname{lcm}(m, n). Assumes mn < 2^{62}.
Time: \log(n)
11 crt(ll a, ll m, ll b, ll n) {
  if (n > m) swap(a, b), swap(m, n);
  ll x, y, g = euclid(m, n, x, y);
  assert((a - b) % g == 0); // else no solution
  x = (b - a) % n * x % n / q * m + a;
  return x < 0 ? x + m*n/g : x;
```

#### 5.3.1 Bézout's identity

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

$$ax + by = d$$

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

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

#### phiFunction.h

**Description:** Euler's  $\phi$  function is defined as  $\phi(n) := \#$  of positive integers  $\leq n$  that are coprime with n.  $\phi(1) = 1$ , p prime  $\Rightarrow \phi(p^k) = (p-1)p^{k-1}$ ,  $m, n \text{ coprime } \Rightarrow \phi(mn) = \phi(m)\phi(n).$  If  $n = p_1^{k_1} p_2^{k_2} \dots p_r^{k_r}$  then  $\phi(n) =$  $(p_1-1)p_1^{k_1-1}...(p_r-1)p_r^{k_r-1}.$   $\phi(n)=n\cdot\prod_{n\mid n}(1-1/p).$  $\sum_{d|n} \phi(d) = n, \sum_{1 \le k \le n, \gcd(k,n)=1} k = n\phi(n)/2, n > 1$ Euler's thm: a, n 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)</pre>
    for (int j = i; j < LIM; j += i) phi[j] -= phi[j] / i;</pre>
```

#### 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:  $\mathcal{O}(\log N)$ dd6c5e, 21 lines

```
typedef double d; // for N \sim 1e7; long double for N \sim 1e9
pair<11, 11> approximate(d x, 11 N) {
 11 LP = 0, LQ = 1, P = 1, Q = 0, inf = LLONG_MAX; d y = x;
    ll lim = min(P ? (N-LP) / P : inf, Q ? (N-LQ) / Q : inf),
       a = (11) 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)NO) < abs(x - (d)P / (d)O)) ?
   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

04d93a, 7 lines

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

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

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

### 5.5 Pythagorean Triples

The Pythagorean triples are uniquely generated by

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

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

#### 5.6 Primes

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

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

### 5.7 Estimates

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

#### IntPerm multinomial

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

#### Mobius Function

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

Mobius Inversion:

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

Other useful formulas/forms:

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

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

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

### Combinatorial (6)

#### 6.1 Permutations

#### 6.1.1 Factorial

n	1 2 3	4	5 6	7	8	9	10	
n!	1 2 6	24 1	20 720	5040	40320	362880	3628800	
n	11	12	13	14	1.5	16	17	
n!	4.0e7	4.8e	8 6.2e	9 8.7e	10 1.3e	12 2.1e1	3 3.6e14	
n	20	25	30	40	50 10	00 150	0 171	
n!	2e18	2e25	3e32	$8e47 \ 3$	e64 9e	157  6e26	$62 > DBL_M$	ΑX

#### IntPerm.h

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

044568, 6 lines int permToInt(vi& v) { **int** use = 0, i = 0, r = 0;

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

### 6.1.2 Cycles

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

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

### 6.1.3 Derangements

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

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

6.1.4 Burnside's lemma

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

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

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

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

#### Partitions and subsets 6.2

#### **6.2.1** Partition function

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

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

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

#### 6.2.2 Lucas' Theorem

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

#### 6.2.3 Binomials

multinomial.h

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

11 c = 1, m = v.emptv() ? 1 : v[0];rep(i, 1, sz(v)) rep(j, 0, v[i]) c = c \* ++m / (j+1);

### General purpose numbers

#### 6.3.1 Bernoulli numbers

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

Sums of powers:

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

Euler-Maclaurin formula for infinite sums:

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

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

10

6.3.2 Stirling numbers of the first kind

Number of permutations on n items with k cycles.

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

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

### $6(3,3) = 0.0, 1.3, 11, 50, 274, 1764, 13068, 109584, \dots$

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

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

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

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

#### 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

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

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

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

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

## Graph (7)

#### 7.1 Fundamentals

#### BellmanFord.h

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

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

#### FlovdWarshall.h

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

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

```
531245, 12 lines
const 11 inf = 1LL << 62;</pre>
void floydWarshall(vector<vector<11>>& m) {
  int n = sz(m);
  rep(i, 0, n) m[i][i] = min(m[i][i], OLL);
  rep(k, 0, n) rep(i, 0, n) rep(j, 0, n)
```

```
if (m[i][k] != inf && m[k][j] != inf) {
    auto newDist = max(m[i][k] + m[k][j], -inf);
    m[i][j] = min(m[i][j], newDist);
rep(k, 0, n) if (m[k][k] < 0) rep(i, 0, n) rep(j, 0, n)
  if (m[i][k] != inf && m[k][j] != inf) m[i][j] = -inf;
```

#### TopoSort.h

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

```
Time: \mathcal{O}(|V| + |E|)
                                                        d678d8, 8 lines
vi topoSort(const vector<vi>& qr) {
 vi indeg(sz(gr)), q;
 for (auto& li : gr) for (int x : li) indeg[x]++;
 rep(i, 0, sz(gr)) if (indeg[i] == 0) q.push_back(i);
 rep(j,0,sz(q)) for (int x : gr[q[j]])
    if (--indeq[x] == 0) q.push_back(x);
 return q;
```

#### 7.2 Network flow

#### PushRelabel.h

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

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

```
0ae1d4, 48 lines
struct PushRelabel
 struct Edge {
   int dest, back;
   11 f, c;
 vector<vector<Edge>> q;
 vector<11> ec;
 vector<Edge*> cur;
 vector<vi> hs; vi H;
 PushRelabel(int n) : g(n), ec(n), cur(n), hs(2*n), H(n) {}
 void addEdge(int s, int t, ll cap, ll rcap=0) {
   if (s == t) return;
   g[s].push_back({t, sz(g[t]), 0, cap});
   g[t].push_back({s, sz(g[s])-1, 0, rcap});
 void addFlow(Edge& e, ll f) {
   Edge &back = g[e.dest][e.back];
   if (!ec[e.dest] && f) hs[H[e.dest]].push_back(e.dest);
   e.f += f; e.c -= f; ec[e.dest] += f;
   back.f -= f; back.c += f; ec[back.dest] -= f;
 11 calc(int s, int t) {
   int v = sz(q); H[s] = v; ec[t] = 1;
   vi co(2*v); co[0] = v-1;
   rep(i,0,v) cur[i] = g[i].data();
   for (Edge& e : q[s]) addFlow(e, e.c);
    for (int hi = 0;;) {
     while (hs[hi].empty()) if (!hi--) return -ec[s];
     int u = hs[hi].back(); hs[hi].pop_back();
     while (ec[u] > 0) // discharge u
       if (cur[u] == g[u].data() + sz(g[u])) {
         H[u] = 1e9;
         for (Edge& e : q[u]) if (e.c && H[u] > H[e.dest]+1)
           H[u] = H[e.dest]+1, cur[u] = &e;
```

```
if (++co[H[u]], !--co[hi] && hi < v)</pre>
          rep(i, 0, v) if (hi < H[i] && H[i] < v)
            --co[H[i]], H[i] = v + 1;
      } else if (cur[u]->c && H[u] == H[cur[u]->dest]+1)
        addFlow(*cur[u], min(ec[u], cur[u]->c));
      else ++cur[u];
bool leftOfMinCut(int a) { return H[a] >= sz(g); }
```

#### MinCostMaxFlow.h

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

**Time:**  $\mathcal{O}\left(FE\log(V)\right)$  where F is max flow.  $\mathcal{O}\left(VE\right)$  for setpi. <sub>58385b, 79 lines</sub>

```
#include <bits/extc++.h>
const 11 INF = numeric limits<11>::max() / 4;
struct MCMF {
 struct edge {
    int from, to, rev;
    11 cap, cost, flow;
  };
  vector<vector<edge>> ed;
  vi seen;
  vector<ll> dist, pi;
  vector<edge*> par;
  MCMF(int N) : N(N), ed(N), seen(N), dist(N), pi(N), par(N) {}
  void addEdge(int from, int to, ll cap, ll cost) {
    if (from == to) return;
    ed[from].push_back(edge{ from, to, sz(ed[to]), cap, cost, 0 });
    ed[to].push back(edge{ to,from,sz(ed[from])-1,0,-cost,0 });
  void path(int s) {
    fill(all(seen), 0);
    fill(all(dist), INF);
    dist[s] = 0; ll di;
    __gnu_pbds::priority_queue<pair<ll, int>> q;
    vector<decltype(q)::point_iterator> its(N);
    q.push({ 0, s });
    while (!q.empty()) {
      s = q.top().second; q.pop();
      seen[s] = 1; di = dist[s] + pi[s];
      for (edge& e : ed[s]) if (!seen[e.to]) {
       11 val = di - pi[e.to] + e.cost;
        if (e.cap - e.flow > 0 && val < dist[e.to]) {</pre>
          dist[e.to] = val;
          par[e.to] = &e;
          if (its[e.to] == q.end())
            its[e.to] = q.push({ -dist[e.to], e.to });
            q.modify(its[e.to], { -dist[e.to], e.to });
    rep(i, 0, N) pi[i] = min(pi[i] + dist[i], INF);
  pair<11, 11> maxflow(int s, int t) {
```

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

#### EdmondsKarp.h

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

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

# MinCut.h

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

#### GlobalMinCut.h

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

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

8b0e19, 21 lines

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

#### GomorvHu.h

**Description:** Given a list of edges representing an undirected flow graph, returns edges of the Gomory-Hu tree. The max flow between any pair of vertices is given by minimum edge weight along the Gomory-Hu tree path. **Time:**  $\mathcal{O}(V)$  Flow Computations

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

### 7.3 Matching

hopcroftKarp.h

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

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

f612e4, 42 line

```
bool dfs(int a, int L, vector<vi>& g, vi& btoa, vi& A, vi& B) {
   if (A[a] != L) return 0;
   A[a] = -1;
   for (int b : g[a]) if (B[b] == L + 1) {
```

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

## DFSMatching.h Description: DFSMatching.h

da1d4b, 70 lines

12

```
struct bipartite {
    int n, m;
    vector<vector<int>> q;
    vector<bool> paired;
    vector<int> match;
    bipartite(int n, int m): n(n), m(m), g(n), paired(n), match
         (m, -1) \{ \}
    void add(int a, int b) {
        g[a].push_back(b);
    vector<size_t> ptr;
    bool kuhn(int v) {
        for(size_t &i = ptr[v]; i < g[v].size(); i++) {</pre>
            int &u = match[q[v][i]];
            if(u == -1 \mid | (dist[u] == dist[v] + 1 && kuhn(u)))
                u = v;
                paired[v] = true;
                return true;
        return false;
```

```
vector<int> dist:
bool bfs() {
    dist.assign(n, n);
    int que[n];
    int st = 0, fi = 0;
    for (int v = 0; v < n; v++) {
        if(!paired[v]) {
            dist[v] = 0;
            que[fi++] = v;
   bool rep = false;
    while(st < fi) {</pre>
        int v = que[st++];
        for(auto e: q[v]) {
            int u = match[e];
            rep |= u == -1;
            if(u != -1 && dist[v] + 1 < dist[u]) {</pre>
                dist[u] = dist[v] + 1;
                 que[fi++] = u;
    return rep;
auto matching() {
    while(bfs()) {
        ptr.assign(n, 0);
        for (int v = 0; v < n; v++) {
            if(!paired[v]) {
                kuhn(v);
    vector<pair<int, int>> ans;
    for(int u = 0; u < m; u++) {
        if (match[u] != −1) {
            ans.emplace_back(match[u], u);
    return ans;
```

#### MinimumVertexCover.h

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

```
"DFSMatching.h"
                                                     da4196, 20 lines
vi cover(vector<vi>& q, int n, int m) {
  vi match (m, -1);
  int res = dfsMatching(g, match);
  vector<bool> lfound(n, true), seen(m);
  for (int it : match) if (it != -1) lfound[it] = false;
  vi q, cover;
  rep(i,0,n) if (lfound[i]) q.push_back(i);
  while (!q.empty()) {
   int i = q.back(); q.pop_back();
   lfound[i] = 1;
   for (int e : q[i]) if (!seen[e] && match[e] != -1) {
     seen[e] = true;
     q.push_back(match[e]);
  rep(i,0,n) if (!lfound[i]) cover.push_back(i);
  rep(i,0,m) if (seen[i]) cover.push_back(n+i);
  assert(sz(cover) == res);
```

```
return cover;
```

#### WeightedMatching.h

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

```
Time: \mathcal{O}(N^2M)
pair<int, vi> hungarian(const vector<vi> &a) {
 if (a.empty()) return {0, {}};
 int n = sz(a) + 1, m = sz(a[0]) + 1;
 vi u(n), v(m), p(m), ans(n - 1);
 rep(i,1,n) {
   p[0] = i;
    int j0 = 0; // add "dummy" worker 0
    vi dist(m, INT_MAX), pre(m, -1);
    vector<bool> done(m + 1);
    do { // dijkstra
     done[j0] = true;
      int i0 = p[j0], j1, delta = INT_MAX;
      rep(j,1,m) if (!done[j]) {
        auto cur = a[i0 - 1][j - 1] - u[i0] - v[j];
        if (cur < dist[j]) dist[j] = cur, pre[j] = j0;
       if (dist[j] < delta) delta = dist[j], j1 = j;</pre>
      rep(j,0,m) {
        if (done[j]) u[p[j]] += delta, v[j] -= delta;
        else dist[j] -= delta;
      j0 = j1;
    } while (p[j0]);
    while (j0) { // update alternating path
     int j1 = pre[j0];
     p[j0] = p[j1], j0 = j1;
 rep(j,1,m) if (p[j]) ans[p[j] - 1] = j - 1;
  return {-v[0], ans}; // min cost
```

#### GeneralMatching.h

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

```
cb1912, 40 lines
"../numerical/MatrixInverse-mod.h"
vector<pii> generalMatching(int N, vector<pii>& ed) {
 vector<vector<ll>> mat(N, vector<ll>(N)), A;
 for (pii pa : ed) {
   int a = pa.first, b = pa.second, r = rand() % mod;
   mat[a][b] = r, mat[b][a] = (mod - r) % mod;
 int r = matInv(A = mat), M = 2*N - r, fi, fj;
 assert (r % 2 == 0);
 if (M != N) do {
   mat.resize(M, vector<ll>(M));
   rep(i,0,N) {
     mat[i].resize(M);
     rep(j,N,M) {
       int r = rand() % mod;
        mat[i][j] = r, mat[j][i] = (mod - r) % mod;
 } while (matInv(A = mat) != M);
 vi has (M, 1); vector<pii> ret;
```

```
rep(it,0,M/2) {
  rep(i,0,M) if (has[i])
    rep(j,i+1,M) if (A[i][j] && mat[i][j]) {
      fi = i; fj = j; goto done;
  } assert(0); done:
  if (fj < N) ret.emplace_back(fi, fj);</pre>
  has[fi] = has[fj] = 0;
  rep(sw,0,2) {
    11 a = modpow(A[fi][fj], mod-2);
    rep(i,0,M) if (has[i] && A[i][fj]) {
      ll b = A[i][fj] * a % mod;
      rep(j, 0, M) A[i][j] = (A[i][j] - A[fi][j] * b) % mod;
    swap(fi,fj);
return ret;
```

```
DFS algorithms
SCC.h
Description: SCC.h
vector<bool> visited; // keeps track of which vertices are
     already visited
// runs depth first search starting at vertex v.
// each visited vertex is appended to the output vector when
     dfs leaves it.
void dfs(int v, vector<vector<int>> const& adi, vector<int> &
    output) {
    visited[v] = true;
    for (auto u : adj[v])
       if (!visited[u])
            dfs(u, adj, output);
    output.push_back(v);
// input: adj — adjacency list of G
// output: components — the strongy connected components in G
// output: adj_cond — adjacency list of G^SCC (by root
void strongly_connected_components(vector<vector<int>> const&
    adj,
                                  vector<vector<int>> &
                                      components,
                                  vector<vector<int>> &adj_cond
   int n = adj.size();
    components.clear(), adj_cond.clear();
    vector<int> order; // will be a sorted list of G's vertices
         by exit time
    visited.assign(n, false);
    // first series of depth first searches
    for (int i = 0; i < n; i++)
       if (!visited[i])
            dfs(i, adj, order);
    // create adjacency list of G^T
    vector<vector<int>> adj_rev(n);
    for (int v = 0; v < n; v++)</pre>
       for (int u : adj[v])
            adj_rev[u].push_back(v);
    visited.assign(n, false);
    reverse(order.begin(), order.end());
```

#### BiconnectedComponents 2sat EulerWalk EdgeColoring

```
vector<int> roots(n, 0); // gives the root vertex of a
    vertex's SCC
// second series of depth first searches
for (auto v : order)
   if (!visited[v]) {
        std::vector<int> component;
       dfs(v, adj_rev, component);
        components.push_back(component);
       int root = *min_element(begin(component), end(
            component));
       for (auto u : component)
           roots[u] = root;
// add edges to condensation graph
adj_cond.assign(n, {});
for (int v = 0; v < n; v++)
   for (auto u : adj[v])
       if (roots[v] != roots[u])
           adj_cond[roots[v]].push_back(roots[u]);
```

#### BiconnectedComponents.h

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

```
Usage: int eid = 0; ed.resize(N);
for each edge (a,b) {
ed[a].emplace_back(b, eid);
ed[b].emplace_back(a, eid++); }
bicomps([&](const vi& edgelist) {...});
```

c6b7c7, 32 lines

```
Time: \mathcal{O}\left(E+V\right)
vi num, st;
vector<vector<pii>> ed;
int Time;
template < class F>
int dfs(int at, int par, F& f) {
  int me = num[at] = ++Time, top = me;
  for (auto [y, e] : ed[at]) if (e != par) {
    if (num[y]) {
      top = min(top, num[y]);
      if (num[y] < me)
        st.push back(e);
    } else {
      int si = sz(st);
      int up = dfs(y, e, f);
      top = min(top, up);
      if (up == me) {
        st.push_back(e);
        f(vi(st.begin() + si, st.end()));
        st.resize(si);
      else if (up < me) st.push_back(e);</pre>
      else { /* e is a bridge */ }
  return top;
template<class F>
void bicomps(F f) {
  num.assign(sz(ed), 0);
  rep(i, 0, sz(ed)) if (!num[i]) dfs(i, -1, f);
```

```
2sat.h
Description: 2sat.h
                                                     3524b8, 68 lines
class TwoSAT {
private:
    std::vector<std::vector<int>> adj, adj t;
    std::vector<bool> used;
    std::vector<int> order, comp;
    std::vector<bool> assignment;
    void dfs1(int v) {
        used[v] = true;
        for (int u : adj[v]) {
            if (!used[u])
                dfs1(u);
        order.push_back(v);
    void dfs2(int v, int cl) {
        comp[v] = cl;
        for (int u : adj t[v]) {
            if (comp[u] == -1)
                dfs2(u, cl);
    TwoSAT(int size) : n(size), adj(2 * n), adj_t(2 * n), used
         (2 * n), comp(2 * n), assignment(n) {}
    bool solve() {
        order.clear();
        used.assign(2 * n, false);
        for (int i = 0; i < 2 * n; ++i) {
            if (!used[i])
                dfs1(i);
        comp.assign(2 * n, -1);
        for (int i = 0, j = 0; i < 2 * n; ++i) {</pre>
            int v = order[2 * n - i - 1];
            if (comp[v] == -1)
                dfs2(v, j++);
        assignment.assign(n, false);
        for (int i = 0; i < 2 * n; i += 2) {
            if (comp[i] == comp[i + 1])
                return false;
            assignment[i / 2] = comp[i] > comp[i + 1];
        return true;
    void add_disjunction(int a, bool na, int b, bool nb) {
        // na and nb signify whether a and b are to be negated
        a = 2 * a ^ na;
        b = 2 * b ^ nb;
        int neg_a = a ^ 1;
        int neq_b = b ^ 1;
        adj[neg_a].push_back(b);
        adj[neq_b].push_back(a);
        adj_t[b].push_back(neg_a);
        adj_t[a].push_back(neg_b);
    std::vector<bool> get assignment() {
```

```
return assignment;
};
```

#### EulerWalk.h

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

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

#### 7.5 Coloring

#### EdgeColoring.h

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

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

e210e2, 31 lines

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

#### 7.6 Heuristics

#### MaximalCliques.h

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

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

b0d5b1, 12 lines

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

#### MaximumClique.h

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

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

f7c0bc, 49 lines

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

```
vi maxClique() { init(V), expand(V); return qmax; }
 Maxclique(vb conn) : e(conn), C(sz(e)+1), S(sz(C)), old(S) {
   rep(i,0,sz(e)) V.push_back({i});
};
```

#### MaximumIndependentSet.h

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

#### 7.7Trees

#### BinaryLifting.h

Description: BinaryLifting.h

```
class Binary lift{
    public:
        int n, l, timer;
        vector<vector<int>> adj;
        vector<vector<int>> up;
        vector<vector<int>> min v;
        vector<int> depth;
        vector<int> tin;
        vector<int> tout;
        Binary lift(int n) {
            this -> n = n;
            this->1 = \log_2(n) + 1;
            adj.resize(n);
            up.resize(n, vector<int>(1, -1));
            min v.resize(n, vector<int>(1, inf));
            depth.resize(n);
            tin.resize(n);
            tout.resize(n);
            timer = 0;
        void set_min_v(vi& a) {
            fr(i,0,n){
                 \min_{v[i][0]} = a[i];
        void add_edge(int u, int v) {
            adj[u].push_back(v);
            adj[v].push_back(u);
        void dfs(int u, int p, vi& a, int d=0) {
            up[u][0] = p;
             // if(p!=-1) min_v[u][0] = min(min_v[u][0], a[p]);
            depth[u] = d;
            tin[u] = timer++;
            for(int i=1;i<1;i++) {</pre>
                 if (up[u][i-1] != -1) {
                     up[u][i] = up[up[u][i-1]][i-1];
                     min_v[u][i] = min(min_v[u][i-1], min_v[up[u][i-1]])
                          ][i-1]][i-1]);
             // cout << "u: "<< u<< \n ';
            // // print min_v
             // for(int i=0; i< l; i++){}
                    cout << min_v / u / / i / << ' ';
             // cout << \n';
```

```
for(int v: adj[u]){
        if(v != p){
            dfs(v, u,a,d+1);
    tout[u] = timer;
int lift(int u, int k){
    for (int i=1-1; i>=0; i--) {
        if(k >= (1<<i)) {
            u = up[u][i];
            k -= (1 << i);
    return u;
int lca(int u, int v) {
    if(depth[u] < depth[v]) {</pre>
        swap(u,v);
    u = lift(u, depth[u]-depth[v]);
    if(u == v){
        return u;
    for (int i=1-1; i>=0; i--) {
        if (depth[u] < (1 << i))
            continue;
        if(up[u][i] != up[v][i]){
            u = up[u][i];
            v = up[v][i];
    return up[u][0];
int get_kth_node_on_path(int u, int v, int k) {
    int lca = this->lca(u,v);
    int dist = this->depth[u] + this->depth[v] - 2*this
         ->depth[lca];
    if(k > dist){
        return -1;
    if(k == 0){
        return u;
    if(k == dist) {
        return v;
    if(this->depth[u] - this->depth[lca] >= k){
        return this->lift(u, k);
    return this->lift(v, dist-k);
int get_min_on_path(int u, int v) {
    int lca = this->lca(u,v);
    int ans = inf;
    for (int i=1-1; i>=0; i--) {
        if(this->depth[u] - (1<<i) >= this->depth[lca])
            ans = min(ans, this->min_v[u][i]);
            u = this->up[u][i];
```

```
for(int i=1-1; i>=0; i--) {
        if(this->depth[v] - (1<<i) >= this->depth[lca])
            ans = min(ans, this->min_v[v][i]);
            v = this->up[v][i];
        }
    ans = min(ans, this->min_v[u][0]);
    ans = min(ans, this->min_v[v][0]);
    return ans:
// returns the node number, we move from u to v
int first node less equal k on path (int u, int v, int k
     , vi& a) {
    if(a[u] <= k) return u;</pre>
    int lca = this->lca(u, v);
    // cout << u << '' << v << '' << lca << ' \n';
    for (int i=1-1; i>=0; i--) {
        if(this->depth[u] - (1<<i) >= this->depth[lca])
            if(this->min_v[u][i] <= k) continue;</pre>
            u = this->up[u][i];
            // cout<<u<' '<<tc<' '<<lca<<' '<<i<' '<<
                  this \rightarrow min_v[u]/i/<\langle n';
    // cout << v << '' << lca << 'n';
    int j = -1;
    if(u!=lca) return u;
    if(a[u] <= k) return u;</pre>
    for (int i=1-1; i>=0; i--) {
        if(this->depth[v] - (1<<i) >= this->depth[lca])
             int height = this->depth[v] - this->depth[
                 lca] - (1 << i);
            int node = this->lift(v, height);
            if(this->min v[node][i] <= k) v = node;</pre>
            else lca = up[v][i];
            break;
    // cout << u << '' << v << '' << lca << ^n ':
    for(int i=j;i>0;i--) {
        if(this->depth[v] - (1<<i) >= this->depth[lca])
            int node = this->up[v][i-1];
            if(this->min v[node][i-1] <= k) v = node;</pre>
             else lca = node;
    // assert(v=lca);
    return v:
int get_dist(int u, int v){
    return this->depth[u] + this->depth[v] - 2*this->
         depth[this->lca(u,v)];
```

### ${\bf CompressTree.h}$

};

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

Time:  $\mathcal{O}\left(|S|\log|S|\right)$ 

"LCA.h" 9775a0, 21 lines

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

#### HLD h

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

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

```
"../data-structures/LazySegmentTree.h"
                                                    03139d, 46 lines
template <bool VALS EDGES> struct HLD {
 int N, tim = 0;
 vector<vi> adj;
 vi par, siz, rt, pos;
 Node *tree:
 HLD(vector<vi> adj_)
   : N(sz(adj_)), adj(adj_), par(N, -1), siz(N, 1),
     rt(N),pos(N),tree(new Node(0, N)) { dfsSz(0); dfsHld(0); }
 void dfsSz(int v) {
   if (par[v] != -1) adj[v].erase(find(all(adj[v]), par[v]));
   for (int& u : adj[v]) {
     par[u] = v;
     dfsSz(u);
     siz[v] += siz[u];
     if (siz[u] > siz[adj[v][0]]) swap(u, adj[v][0]);
 void dfsHld(int v) {
   pos[v] = tim++;
   for (int u : adj[v]) {
     rt[u] = (u == adj[v][0] ? rt[v] : u);
     dfsHld(u);
 template <class B> void process(int u, int v, B op) {
   for (; rt[u] != rt[v]; v = par[rt[v]]) {
     if (pos[rt[u]] > pos[rt[v]]) swap(u, v);
     op(pos[rt[v]], pos[v] + 1);
   if (pos[u] > pos[v]) swap(u, v);
   op(pos[u] + VALS_EDGES, pos[v] + 1);
 void modifyPath(int u, int v, int val) {
   process(u, v, [&](int 1, int r) { tree->add(1, r, val); });
```

int queryPath(int u, int v) { // Modify depending on problem

```
int res = -1e9;
process(u, v, [&](int 1, int r) {
    res = max(res, tree->query(1, r));
});
return res;
}
int querySubtree(int v) { // modifySubtree is similar
    return tree->query(pos[v] + VALS_EDGES, pos[v] + siz[v]);
};
```

#### LinkCutTree.h

**Description:** Represents a forest of unrooted trees. You can add and remove edges (as long as the result is still a forest), and check whether two nodes are in the same tree.

**Time:** All operations take amortized  $\mathcal{O}(\log N)$ .

0fb462 90 lin

```
struct Node { // Splay tree. Root's pp contains tree's parent.
 Node *p = 0, *pp = 0, *c[2];
 bool flip = 0;
 Node() { c[0] = c[1] = 0; fix(); }
  void fix() {
   if (c[0]) c[0]->p = this;
    if (c[1]) c[1]->p = this;
    // (+ update sum of subtree elements etc. if wanted)
  void pushFlip() {
    if (!flip) return;
    flip = 0; swap(c[0], c[1]);
    if (c[0]) c[0]->flip ^= 1;
    if (c[1]) c[1]->flip ^= 1;
  int up() { return p ? p->c[1] == this : -1; }
  void rot(int i, int b) {
    int h = i ^ b;
    Node *x = c[i], *y = b == 2 ? x : x -> c[h], *z = b ? y : x;
    if ((y->p = p)) p->c[up()] = y;
    c[i] = z -> c[i ^ 1];
    if (b < 2) {
      x->c[h] = y->c[h ^ 1];
      y - > c[h ^ 1] = x;
    z \rightarrow c[i ^1] = this;
    fix(); x->fix(); y->fix();
    if (p) p->fix();
    swap(pp, y->pp);
  void splay() {
    for (pushFlip(); p; ) {
      if (p->p) p->p->pushFlip();
      p->pushFlip(); pushFlip();
      int c1 = up(), c2 = p->up();
      if (c2 == -1) p->rot(c1, 2);
      else p->p->rot(c2, c1 != c2);
 Node* first() {
    pushFlip();
    return c[0] ? c[0]->first() : (splay(), this);
};
struct LinkCut {
  vector<Node> node;
 LinkCut(int N) : node(N) {}
  void link(int u, int v) { // add an edge (u, v)
    assert(!connected(u, v));
    makeRoot(&node[u]);
```

```
node[u].pp = &node[v];
  void cut(int u, int v) { // remove an edge (u, v)
   Node *x = &node[u], *top = &node[v];
   makeRoot(top); x->splay();
   assert(top == (x->pp ?: x->c[0]));
   if (x->pp) x->pp = 0;
   else {
     x->c[0] = top->p = 0;
     x \rightarrow fix();
  bool connected (int u, int v) { // are u, v in the same tree?
   Node* nu = access(&node[u])->first();
   return nu == access(&node[v])->first();
  void makeRoot(Node* u) {
   access(u);
   u->splay();
   if(u->c[0]) {
     u - c[0] - p = 0;
     u->c[0]->flip ^= 1;
     u - c[0] - pp = u;
     u - > c[0] = 0;
     u->fix();
  Node* access(Node* u) {
   u->splay();
    while (Node* pp = u->pp) {
     pp->splay(); u->pp = 0;
     if (pp->c[1]) {
       pp - c[1] - p = 0; pp - c[1] - pp = pp; 
     pp - c[1] = u; pp - fix(); u = pp;
    return u;
};
```

#### DirectedMST.h

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

```
Time: \mathcal{O}\left(E\log V\right)
"../data-structures/UnionFindRollback.h"
                                                         39e620, 60 lines
struct Edge { int a, b; ll w; };
struct Node {
  Edge key;
  Node *1, *r;
  11 delta:
  void prop() {
    key.w += delta;
    if (1) 1->delta += delta;
   if (r) r->delta += delta;
    delta = 0:
  Edge top() { prop(); return key; }
Node *merge(Node *a, Node *b) {
  if (!a || !b) return a ?: b;
  a->prop(), b->prop();
  if (a->key.w > b->key.w) swap(a, b);
  swap(a->1, (a->r = merge(b, a->r)));
  return a;
void pop (Node \star \& a) { a->prop(); a = merge(a->1, a->r); }
pair<ll, vi> dmst(int n, int r, vector<Edge>& g) {
  RollbackUF uf(n);
  vector<Node*> heap(n);
```

```
for (Edge e : q) heap[e.b] = merge(heap[e.b], new Node{e});
11 \text{ res} = 0:
vi seen(n, -1), path(n), par(n);
seen[r] = r;
vector<Edge> Q(n), in(n, \{-1,-1\}), comp;
deque<tuple<int, int, vector<Edge>>> cycs;
rep(s,0,n) {
  int u = s, qi = 0, w;
  while (seen[u] < 0) {</pre>
    if (!heap[u]) return {-1,{}};
    Edge e = heap[u]->top();
    heap[u]->delta -= e.w, pop(heap[u]);
    Q[qi] = e, path[qi++] = u, seen[u] = s;
    res += e.w, u = uf.find(e.a);
    if (seen[u] == s) {
      Node* cyc = 0;
      int end = qi, time = uf.time();
      do cyc = merge(cyc, heap[w = path[--qi]]);
      while (uf.join(u, w));
      u = uf.find(u), heap[u] = cyc, seen[u] = -1;
      cycs.push_front({u, time, {&Q[qi], &Q[end]}});
  rep(i, 0, qi) in[uf.find(Q[i].b)] = Q[i];
for (auto& [u,t,comp] : cycs) { // restore sol (optional)
 uf.rollback(t);
  Edge inEdge = in[u];
  for (auto& e : comp) in[uf.find(e.b)] = e;
  in[uf.find(inEdge.b)] = inEdge;
rep(i,0,n) par[i] = in[i].a;
return {res, par};
```

#### 7.8 Math

#### 7.8.1 Number of Spanning Trees

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

#### 7.8.2 Erdős–Gallai theorem

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

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

## Geometry (8)

### 8.1 Geometric primitives

#### Point.h

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

47ec0a, 28 lines

```
template <class T> int sgn(T x) { return (x > 0) - (x < 0); } template <class T>
```

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

#### lineDistance.h

#### Description:

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



```
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"</pre>
```

5c88f4, 6 lines

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

### SegmentIntersection.h

#### Description:

If a unique intersection point between the line segments going from s1 to e1 and from s2 to e2 exists then it is returned. If no intersection point exists an empty vector is returned. If infinitely many exist a vector with 2 elements is returned, containing the endpoints of the common line segment. The wrong position will be returned if P is Point<|1> and the intersection point does not have integer coordinates. Products of three coordinates are used in intermediate steps so watch out for overflow if using int or long long.



```
Usage: vector<P> inter = segInter(s1,e1,s2,e2);
if (sz(inter) == 1)
cout << "segments intersect at " << inter[0] << endl;</pre>
"Point.h", "OnSegment.h"
                                                      9d57f2, 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 (sqn(oa) * sqn(ob) < 0 && sqn(oc) * sqn(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 \$1,e1\$ and \$2,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<|1> 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;

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

**return** {1, (s1 \* p + e1 \* q) / d};

#### sideOf.h

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

#### OnSegment.h

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

```
linearTransformation.h
Description:
```

Apply the linear transformation (translation, rotation and scaling) which takes line p0-p1 to line q0-q1 to point r.  $\label{eq:p0-p1} \begin{subarray}{l} \tt "Point.h" \\ \end{subarray}$ 



#### 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 000002, 35 \ lines
```

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

#### 8.2 Circles

#### CircleIntersection.h

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

```
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

"Point.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.

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

#### CirclePolygonIntersection.h

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

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

```
"../../content/geometry/Point.h"
                                                      a1ee63, 19 lines
typedef Point<double> P;
#define arg(p, q) atan2(p.cross(q), p.dot(q))
double circlePoly(P c, double r, vector<P> ps) {
 auto tri = [&] (P p, P q) {
    auto r2 = r * r / 2;
    P d = q - p;
    auto a = d.dot(p)/d.dist2(), b = (p.dist2()-r*r)/d.dist2();
    auto det = a * a - b;
    if (det <= 0) return arg(p, q) * r2;</pre>
    auto s = max(0., -a-sqrt(det)), t = min(1., -a+sqrt(det));
    if (t < 0 || 1 <= s) return arg(p, q) * r2;</pre>
    Pu = 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:

"Point.h"

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



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

```
P ccCenter(const P& A, const P& B, const P& C) {
  P b = C-A, c = B-A;
  return A + (b*c.dist2()-c*b.dist2()).perp()/b.cross(c)/2;
}
```

#### MinimumEnclosingCircle.h

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

### 8.3 Polygons

#### InsidePolygon.h

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

```
Usage: vector P = \{P\{4,4\}, P\{1,2\}, P\{2,1\}\}; bool in = inPolygon(v, P\{3,3\}, false);

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

```
"Point.h", "OnSegment.h", "SegmentDistance.h"
```

\_\_2bf504, 11 lines

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

#### Polygon Area, h

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

```
"Point.h" f12300, 6 lines
```

```
template < class T>
T polygonArea2(vector<Point<T>>& v) {
  T 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}\left(n\right)
```

"Point.h" 9706dc, 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;</pre>
```

#### PolygonCut.h

#### Description:

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

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



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

#### ConvexHull.h

#### Description:

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



```
Time: \mathcal{O}(n \log n)
template \langle class T \rangle int sgn(T x) \{ return (x > 0) - (x < 0); \}
template<class T>
struct Point {
 typedef Point P;
  Тх, у;
  explicit Point (T x=0, T y=0) : x(x), y(y) {}
 bool operator<(P p) const { return tie(x,y) < tie(p.x,p.y); }</pre>
  bool operator==(P p) const { return tie(x,y)==tie(p.x,p.y); }
  P operator-(P p) const { return P(x-p.x, y-p.y); }
  T cross(P a, P b) const { return (a-*this).cross(b-*this); }
  T cross(P p) const { return x*p.y - y*p.x; }
 friend ostream& operator<<(ostream& os, P p) {</pre>
    return os << "(" << p.x << "," << p.y << ")"; }
typedef Point<11> P;
vector<P> convexHull(vector<P> pts) {
 if (sz(pts) <= 1) return pts;</pre>
  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 \ge s + 2 \&\& h[t-2].cross(h[t-1], p) \le 0) t--;
```

#### HullDiameter.h

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

**return** {h.begin(), h.begin() + t - (t == 2 && h[0] == h[1])};

#### 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}\left(\log N\right) "Point.h", "sideOf.h", "OnSegment.h" 71446b, 14 lines
```

```
typedef Point<11> P;

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

#### LineHullIntersection.h

**return** {-1, -1};

**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"
#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 (10 + 1 < hi) {
    int m = (lo + hi) / 2;
    if (extr(m)) return m;
    int 1s = cmp(1o + 1, 1o), ms = cmp(m + 1, m);
    (ls < ms \mid | (ls == ms \&\& ls == cmp(lo, m)) ? hi : lo) = m;
 return lo;
#define cmpL(i) sqn(a.cross(poly[i], b))
template <class P>
array<int, 2> lineHull(P a, P b, vector<P>& poly) {
 int endA = extrVertex(poly, (a - b).perp());
  int endB = extrVertex(poly, (b - a).perp());
 if (cmpL(endA) < 0 \mid \mid cmpL(endB) > 0)
```

#### ClosestPair kdTree FastDelaunay

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

#### 8.4 Misc. Point Set Problems

#### ClosestPair.h

Description: Finds the closest pair of points.

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

```
"Point.h"
                                                      ac41a6, 17 lines
typedef Point<11> P;
pair<P, P> closest(vector<P> v) {
  assert (sz(v) > 1);
  sort(all(v), [](P a, P b) { return a.y < b.y; });</pre>
  pair<11, pair<P, P>> ret{LLONG_MAX, {P(), P()}};
  int i = 0;
  for (P p : v) {
   P d{1 + (ll)sqrt(ret.first), 0};
    while (v[j].y \le p.y - d.x) S.erase(v[j++]);
    auto lo = S.lower_bound(p - d), hi = S.upper_bound(p + d);
    for (; lo != hi; ++lo)
     ret = min(ret, \{(*lo - p).dist2(), \{*lo, p\}\});
    S.insert(p):
```

#### kdTree.h

return ret.second;

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

```
bac5b0, 63 lines
typedef long long T;
typedef Point<T> P;
const T INF = numeric_limits<T>::max();
bool on_x(const P& a, const P& b) { return a.x < b.x; }</pre>
bool on_y(const P& a, const P& b) { return a.y < b.y; }</pre>
struct Node 4
  P pt; // if this is a leaf, the single point in it
  T x0 = INF, x1 = -INF, y0 = INF, y1 = -INF; // bounds
  Node *first = 0, *second = 0;
  T distance (const P& p) { // min squared distance to a point
    T x = (p.x < x0 ? x0 : p.x > x1 ? x1 : p.x);
    T y = (p.y < y0 ? y0 : p.y > y1 ? y1 : p.y);
    return (P(x,y) - p).dist2();
  Node (vector<P>&& vp) : pt(vp[0]) {
    for (P p : vp) {
      x0 = min(x0, p.x); x1 = max(x1, p.x);
      y0 = min(y0, p.y); y1 = max(y1, p.y);
```

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

#### FastDelaunav.h

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

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

```
"Point.h"
                                                     eefdf5, 88 lines
typedef Point<11> P;
typedef struct Ouad* O:
typedef __int128_t 111; // (can be ll if coords are < 2e4)
P arb(LLONG_MAX, LLONG_MAX); // not equal to any other point
struct Quad {
  Q rot, o; P p = arb; bool mark;
 P& F() { return r()->p; }
  Q& r() { return rot->rot; }
  O prev() { return rot->o->rot;
  Q next() { return r()->prev(); }
bool circ(P p, P a, P b, P c) { // is p in the circumcircle?
 111 p2 = p.dist2(), A = a.dist2()-p2,
      B = b.dist2()-p2, C = c.dist2()-p2;
  return p.cross(a,b)*C + p.cross(b,c)*A + p.cross(c,a)*B > 0;
Q makeEdge(P orig, P dest) {
  Q r = H ? H : new Quad{new Quad{new Quad{new Quad{0}}}};
```

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

#### 3D8.5

KTH

#### PolyhedronVolume.h

Description: Magic formula for the volume of a polyhedron. Faces should point outwards. 305<u>8c3, 6 lines</u>

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

#### Point3D.h

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

```
template<class T> struct Point3D {
  typedef Point3D P;
  typedef const P& R;
  T x, y, z;
  explicit Point3D(T x=0, T y=0, T z=0) : x(x), y(y), z(z) {}
  bool operator<(R p) const {</pre>
    return tie(x, y, z) < tie(p.x, p.y, p.z); }
  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 sgrt((double) dist2()); }
  //Azimuthal angle (longitude) to x-axis in interval [-pi, pi]
  double phi() const { return atan2(v, x); }
  //Zenith angle (latitude) to the z-axis in interval [0, pi]
  double theta() const { return atan2(sgrt(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;
};
```

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"
                                                                                  5b45fc, 49 lines
```

```
typedef Point3D<double> P3;
struct PR {
  void ins(int x) { (a == -1 ? a : b) = x; }
  void rem(int x) { (a == x ? a : b) = -1; }
  int cnt() { return (a != -1) + (b != -1); }
 int a, b;
struct F { P3 q; int a, b, c; };
vector<F> hull3d(const vector<P3>& A) {
  assert (sz(A) >= 4);
```

```
vector<vector<PR>>> E(sz(A), vector<PR>(sz(A), {-1, -1}));
#define E(x,y) E[f.x][f.y]
 vector<F> FS;
 auto mf = [\&] (int i, int j, int k, int 1) {
   P3 q = (A[j] - A[i]).cross((A[k] - A[i]));
   if (q.dot(A[1]) > q.dot(A[i]))
     q = q * -1;
   F f{q, i, j, k};
   E(a,b).ins(k); E(a,c).ins(j); E(b,c).ins(i);
   FS.push_back(f);
 rep(i, 0, 4) rep(j, i+1, 4) rep(k, j+1, 4)
   mf(i, j, k, 6 - i - j - k);
 rep(i,4,sz(A)) {
   rep(j,0,sz(FS)) {
     F f = FS[j];
     if(f.q.dot(A[i]) > f.q.dot(A[f.a])) {
       E(a,b).rem(f.c);
       E(a,c).rem(f.b);
       E(b,c).rem(f.a);
       swap(FS[j--], FS.back());
       FS.pop_back();
   int nw = sz(FS);
   rep(j,0,nw) {
     F f = FS[i];
#define C(a, b, c) if (E(a,b).cnt() != 2) mf(f.a, f.b, i, f.c);
     C(a, b, c); C(a, c, b); C(b, c, a);
 for (F& it : FS) if ((A[it.b] - A[it.a]).cross(
   A[it.c] - A[it.a]).dot(it.q) \ll 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. 611f07, 8 lines

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

### Strings (9)

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

```
Time: \mathcal{O}(n)
                                                            d4375c, 16 lines
vi pi(const string& s) {
 vi p(sz(s));
  rep(i,1,sz(s)) {
    int g = p[i-1];
```

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

#### Zfunc.h

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

```
ee09e2, 12 lines
vi Z(const string& S) {
 vi z(sz(S));
 int 1 = -1, r = -1;
 rep(i,1,sz(S)) {
   z[i] = i >= r ? 0 : min(r - i, z[i - 1]);
    while (i + z[i] < sz(S) \&\& S[i + z[i]] == S[z[i]])
     z[i]++;
    if (i + z[i] > r)
     1 = i, r = i + z[i];
 return z;
```

#### Manacher.h

**Description:** For each position in a string, computes p[0][i] = half length of longest even palindrome around pos i, p[1][i] = longest odd (half rounded down). Time:  $\mathcal{O}(N)$ 

```
e7ad79, 13 lines
array<vi, 2> manacher(const string& s) {
  int n = sz(s);
  array < vi, 2 > p = {vi(n+1), vi(n)};
  rep(z,0,2) for (int i=0, l=0, r=0; i < n; i++) {
    int t = r-i+!z;
    if (i<r) p[z][i] = min(t, p[z][l+t]);</pre>
    int L = i-p[z][i], R = i+p[z][i]-!z;
    while (L>=1 && R+1<n && s[L-1] == s[R+1])
      p[z][i]++, L--, R++;
    if (R>r) l=L, r=R;
 return p;
```

#### MinRotation.h

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

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

### SuffixArray.h

#### SuffixTree Hashing AhoCorasick

Time:  $\mathcal{O}(n \log n)$ struct SuffixArray { vi sa, lcp; SuffixArray(string& s, int lim=256) { // or basic\_string<int> **int** n = sz(s) + 1, k = 0, a, b; vi x(all(s)), v(n), ws(max(n, lim)); $x.push_back(0)$ , sa = lcp = y, iota(all(sa), 0); for (int j = 0, p = 0; p < n; j = max(1, j \* 2),  $lim = p) {$ p = j, iota(all(v), n - j); fr(i,0,n) **if** (sa[i] >= j) y[p++] = sa[i] - j;fill(all(ws), 0); fr(i,0,n) ws[x[i]]++;fr(i,1,lim) ws[i] += ws[i - 1];for (int i = n; i--;) sa[--ws[x[y[i]]]] = y[i]; swap(x, y), p = 1, x[sa[0]] = 0;fr(i,1,n) = sa[i-1], b = sa[i], x[b] =(y[a] == y[b] && y[a + j] == y[b + j]) ? p - 1 : p++;for (int i = 0, j; i < n - 1; lcp[x[i++]] = k) for (k & & k--, j = sa[x[i] - 1];s[i + k] == s[j + k]; k++);int lower\_bound(string& t,vector<int> &a,string &s){ **int** 1 = 1, r = sz(a);while(1<r){ **int** m = (1+r)/2;if(s.substr(a[m],min(sz(s)-a[m],sz(t)+1)) >= t) r = m;**else** 1 = m+1;} return 1:} int upper\_bound(string& t, vector<int> &a, string &s) { **int** 1 = 1, r = sz(a); while(1<r){

Description: Builds suffix array for a string. sa[i] is the starting index

of the suffix which is i'th in the sorted suffix array. The returned vector

is of size n+1, and sa[0] = n. The lcp array contains longest common

prefixes for neighbouring strings in the suffix array: lcp[i] = lcp(sa[i],

sa[i-1]), lcp[0] = 0. The input string must not contain any zero bytes.

### return 1:} SuffixTree.h

**int** m = (1+r)/2;

**else** 1 = m+1;}

Description: Ukkonen's algorithm for online suffix tree construction. Each node contains indices [l, r) into the string, and a list of child nodes. Suffixes are given by traversals of this tree, joining [l, r) substrings. The root is 0 (has l = -1, r = 0), non-existent children are -1. To get a complete tree, append a dummy symbol – otherwise it may contain an incomplete path (still useful for substring matching, though).

if(s.substr(a[m],min(sz(a)-a[m],sz(t))) > t) r = m;

Time:  $\mathcal{O}(26N)$ 

aae0b8, 50 lines

```
struct SuffixTree {
 enum { N = 200010, ALPHA = 26 }; // N \sim 2*maxlen+10
 int toi(char c) { return c - 'a'; }
  string a; //v = cur \ node, q = cur \ position
 int t[N][ALPHA],1[N],r[N],p[N],s[N],v=0,q=0,m=2;
 void ukkadd(int i, int c) { suff:
   if (r[v]<=q) {
     if (t[v][c]==-1) { t[v][c]=m; l[m]=i;
       p[m++]=v; v=s[v]; q=r[v]; goto suff; }
     v=t[v][c]; q=l[v];
   if (q==-1 || c==toi(a[q])) q++; else {
     l[m+1]=i; p[m+1]=m; l[m]=l[v]; r[m]=q;
     p[m]=p[v]; t[m][c]=m+1; t[m][toi(a[q])]=v;
     1[v]=q; p[v]=m; t[p[m]][toi(a[1[m]])]=m;
     v=s[p[m]]; q=l[m];
```

```
while (q < r[m]) \{ v = t[v][toi(a[q])]; q + = r[v] - l[v]; \}
      if (q==r[m]) s[m]=v; else s[m]=m+2;
      q=r[v]-(q-r[m]); m+=2; goto suff;
  SuffixTree(string a) : a(a) {
    fill(r,r+N,sz(a));
    memset(s, 0, sizeof s);
    memset(t, -1, sizeof t);
    fill(t[1],t[1]+ALPHA,0);
    s[0] = 1; 1[0] = 1[1] = -1; r[0] = r[1] = p[0] = p[1] = 0;
    rep(i,0,sz(a)) ukkadd(i, toi(a[i]));
  // example: find longest common substring (uses ALPHA = 28)
  pii best;
  int lcs(int node, int i1, int i2, int olen) {
    if (1[node] <= i1 && i1 < r[node]) return 1;</pre>
    if (1[node] <= i2 && i2 < r[node]) return 2;</pre>
    int mask = 0, len = node ? olen + (r[node] - 1[node]) : 0;
    rep(c, 0, ALPHA) if (t[node][c] != -1)
      mask \mid = lcs(t[node][c], i1, i2, len);
    if (mask == 3)
     best = max(best, {len, r[node] - len});
    return mask;
  static pii LCS(string s, string t) {
    SuffixTree st(s + (char) ('z' + 1) + t + (char) ('z' + 2));
    st.lcs(0, sz(s), sz(s) + 1 + sz(t), 0);
    return st.best;
};
Hashing.h
Description: Self-explanatory methods for string hashing.
// Arithmetic mod 2^64-1. 2x slower than mod 2^64 and more
// code, but works on evil test data (e.g. Thue-Morse, where
// ABBA... and BAAB... of length 2^10 hash the same mod 2^64).
// "typedef ull H;" instead if you think test data is random,
// or work mod 10^9+7 if the Birthday paradox is not a problem.
typedef uint64_t ull;
struct H {
  ull x; H(ull x=0) : x(x) {}
```

H operator-(H o) { return \*this + ~o.x; }

return H((ull)m) + (ull)(m >> 64); }

ull get() const { return x + !~x; }

ha[i+1] = ha[i] \* C + str[i],

return ha[b] - ha[a] \* pw[b - a];

H hashInterval (int a, int b) { // hash (a, b)

vector<H> getHashes(string& str, int length) {

pw[i+1] = pw[i] \* C;

struct HashInterval {

rep(i, 0, sz(str))

vector<H> ha, pw;

pw[0] = 1;

};

H operator\*(H o) { auto m = (\_\_uint128\_t)x \* o.x;

```
AhoCorasick.h
                                                                    Time: construction takes \mathcal{O}(26N), where N = \text{sum of length of patterns}.
                                                                     struct AhoCorasick {
                                                                       enum {alpha = 26, first = 'A'}; // change this!
                                                                       struct Node {
                                                                         // (nmatches is optional)
                                                                         Node(int v) { memset(next, v, sizeof(next)); }
                                                                      vector<Node> N:
                                                                      vi backp;
                                                                      void insert(string& s, int j) {
                                                                         assert(!s.empty());
                                                                         int n = 0;
                                                                         for (char c : s) {
                                                                           int& m = N[n].next[c - first];
                                                                           else n = m;
                                                                         if (N[n].end == -1) N[n].start = j;
                                                                         backp.push_back(N[n].end);
                                                                         N[n].end = j;
  H operator+(H \circ) { return x + \circ.x + (x + \circ.x < x); }
                                                                         N[n].nmatches++;
                                                                       AhoCorasick(vector<string>& pat) : N(1, -1) {
                                                                         rep(i, 0, sz(pat)) insert(pat[i], i);
                                                                         N[0].back = sz(N);
 bool operator==(H o) const { return get() == o.get(); }
                                                                         N.emplace_back(0);
 bool operator<(H o) const { return get() < o.get(); }</pre>
                                                                         queue<int> q;
static const H C = (11)1e11+3; // (order \sim 3e9; random \ also \ ok)
                                                                         for (q.push(0); !q.empty(); q.pop()) {
                                                                           int n = q.front(), prev = N[n].back;
                                                                           rep(i,0,alpha) {
 HashInterval(string& str) : ha(sz(str)+1), pw(ha) {
                                                                             if (ed == -1) ed = y;
                                                                             else {
                                                                               N[ed].back = y;
                                                                                 = N[y].end;
                                                                               N[ed].nmatches += N[y].nmatches;
                                                                               q.push(ed);
                                                                      vi find(string word) {
```

```
if (sz(str) < length) return {};</pre>
H h = 0, pw = 1;
rep(i,0,length)
 h = h * C + str[i], pw = pw * C;
vector<H> ret = {h};
rep(i,length,sz(str)) {
```

ret.push\_back(h = h \* C + str[i] - pw \* str[i-length]);

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

return ret:

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

 $\operatorname{find}(x)$  is  $\mathcal{O}(N)$ , where  $N = \operatorname{length}$  of x. findAll is  $\mathcal{O}(NM)$ .

```
int back, next[alpha], start = -1, end = -1, nmatches = 0;
 if (m == -1) { n = m = sz(N); N.emplace_back(-1); }
   int &ed = N[n].next[i], y = N[prev].next[i];
     (N[ed].end == -1 ? N[ed].end : backp[N[ed].start])
```

```
int n = 0;
    vi res; // ll count = 0:
    for (char c : word) {
     n = N[n].next[c - first];
     res.push_back(N[n].end);
     // count += N[n].nmatches;
   return res:
  vector<vi> findAll(vector<string>& pat, string word) {
   vi r = find(word);
    vector<vi> res(sz(word));
    rep(i, 0, sz(word)) {
     int ind = r[i];
     while (ind !=-1) {
       res[i - sz(pat[ind]) + 1].push_back(ind);
       ind = backp[ind];
    return res;
};
```

### Various (10)

#### 10.1 Intervals

IntervalContainer.h

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

Time:  $\mathcal{O}(\log N)$ edce47, 23 lines

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

#### IntervalCover.h

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

9e9d8d, 19 lines

```
template < class T>
vi cover(pair<T, T> G, vector<pair<T, T>> I) {
 vi S(sz(I)), R;
  iota(all(S), 0);
```

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

#### ConstantIntervals.h

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

Usage: constantIntervals(0, sz(v), [&](int x){return v[x];}, [&] (int lo, int hi, T val) $\{\ldots\}$ ); Time:  $\mathcal{O}\left(k\log\frac{n}{k}\right)$ 

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

### 10.2 Misc. algorithms

#### TernarySearch.h

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

```
Usage: int ind = ternSearch(0, n-1, [&] (int i) {return a[i];});
Time: \mathcal{O}(\log(b-a))
                                                              9155b4, 11 lines
```

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

**Description:** Compute indices for the longest increasing subsequence. Time:  $\mathcal{O}(N \log N)$ 2932a0, 17 lines

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

#### FastKnapsack.h

753a4c, 19 lines

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))
                                                                               b20ccc, 16 lines
```

```
int knapsack(vi w, int t) {
 int a = 0, b = 0, x;
 while (b < sz(w) && a + w[b] <= t) a += w[b++];
 if (b == sz(w)) return a;
 int m = *max element(all(w));
 vi u, v(2*m, -1);
 v[a+m-t] = b;
 rep(i,b,sz(w)) {
   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:  $a[i][j] = \min_{i < k < j} (a[i][k] + a[i][k])$ a[k][j] + f(i,j), where the (minimal) optimal k increases with both i and j, one can solve intervals in increasing order of length, and search k = p[i][j] for a[i][j] only between p[i][j-1] and p[i+1][j]. This is known as Knuth DP. Sufficient criteria for this are if  $f(b,c) \leq f(a,d)$  and  $f(a,c) + f(b,d) \le f(a,d) + f(b,c)$  for all  $a \le b \le c \le d$ . Consider also: LineContainer (ch. Data structures), monotone queues, ternary search. Time:  $\mathcal{O}(N^2)$ 

#### DivideAndConquerDP.h

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

```
Time: \mathcal{O}((N + (hi - lo)) \log N)
                                                                            d38d2b, 18 lines
```

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

```
best = min(best, make_pair(f(mid, k), k));
store(mid, best.second, best.first);
rec(L, mid, LO, best.second+1);
rec(mid+1, R, best.second, HI);
}
void solve(int L, int R) { rec(L, R, INT_MIN, INT_MAX); }
};
```

# Techniques (A)

#### techniques.txt

Combinatorics

159 lines

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

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

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