1	Contest	1	8.5 3D				
2	Mathematics 2.1 Equations 2.2 Recurrences 2.3 Trigonometry 2.4 Geometry 2.5 Derivatives/Integrals 2.6 Sums 2.7 Series 2.8 Probability theory 2.9 Markov chains	1 1 2 2 2 2 2 2 2 2 2 3	9 Strings 19 10 Various 20 10.1 Intervals 20 10.2 Misc. algorithms 21 10.3 Dynamic programming 21 10.4 Debugging tricks 21 10.5 Optimization tricks 21 Contest (1)				
3	Data structures	3	instructions.txt 30 lines				
	Numerical 4.1 Polynomials and recurrences	5 5 5 6 7	Compilation: 1. mkdir WF 2. vi .bashrc 3. Add the line: export PATH="\$PATH:\$HOME/WF" 4. cd WF && vi cf.sh -> Write the compilation commands 5. mv cf.sh cf && chmod +x cf 6. Restart terminal				
5	Number theory 5.1 Modular arithmetic	8 8 8 9 9 9 9	<pre>Kate: 1. Theme: Settings->Configure Kate->Color Themes 2. Vim mode: Settings->Configure Kate->Editing ->Default input mode. Then Vi Input mode->Insert mode->jk = <esc> 3. Word wrap: Settings->Configure Kate-> Appearance->Turn off dynamic w.w. 4. Terminal: Make sure View->Tool Views->Show sidebars is on. Go to Settings->Configure Kate->Terminal and turn</esc></pre>				
6	Combinatorial 6.1 Permutations	9 10	5. Hotkey for terminal: Change Focus Terminal Panel to F4. Click "Reassign" when it says it collides with Show Terminal Panel. Fast Compile, Template, Debug:				
7	Graph 7.1 Fundamentals	10 10 11 12 12 13 13 14 15	1. cd WF && mkdir bits 2. Insert stdc++.h 3. Compile using the flags of cf.sh 4. cd and write template.cpp Windows: 1. Using cmd: echo %PATH%. Using Powershell: echo \$env:PATH 2. Add path using cmd: set PATH=%PATH%;C:\ Program Files\CodeBlocks\MinGW\bin It should be the directory where g++ is. 3. If we're using g++ of CodeBlocks, fsanitize won't be available:(4. Write cf.bat at some directory. Ensure that directory is in PATH.				
8	Geometry 8.1 Geometric primitives	15 15 16 17 18	cf.sh 6 lines #!/bin/bash prog_name=\$1				

```
g++ "${prog_name}.cpp" -o $prog_name -std=c
     ++17 -q -Wall -Wshadow -fsanitize=address,
     undefined &&
   "./$prog_name"
stdc++.h
                                           34 lines
#include <bits/stdc++.h>
using namespace std;
template <typename T> constexpr
void __print (const T &x);
template<typename T, typename V>
void __print(const pair<T, V> &x) {
  cerr << "{"; __print(x.first);</pre>
  cerr << ", "; __print(x.second); cerr << "}"
template <typename T> constexpr
void __print (const T &x) {
  if constexpr (is_arithmetic_v<T> ||
    is same v<T, const char*> || is same v<T,
        bool>
    || is_same_v<T, string>) cerr << x;</pre>
  else {
    int f = 0; cerr << '{';</pre>
    for (auto &i: x)
      cerr << (f++ ? ", " : ""), __print(i);
    cerr << "}";
void _print() { cerr << "]\n"; }</pre>
template <typename T, typename... V>
void _print(T t, V... v) {
  __print(t);
  if (sizeof...(v)) cerr << ", ";</pre>
  _print(v...);
#ifdef DeBuG
#define dbg(x...) cerr << "\t\e[93m"<< func
     <<":"<<__LINE__<<" [" << #x << "] = [";
     print(x); cerr << "\e[0m";
#endif
template.cpp
                                           19 lines
#include "bits/stdc++.h"
using namespace std;
#ifndef DeBuG
 #define dbg(...)
#endif
#define rep(i, a, b) for(int i = a; i < (b);
#define all(x) begin(x), end(x)
#define sz(x) (int)(x).size()
using 11 = long long; using vi = vector<int>;
using pii = pair<int,int>; using pll = pair<ll</pre>
     ,11>;
template<class T> using V = vector<T>;
int main() {
  ios_base::sync_with_stdio(false);
  cin.tie(0); cout.tie(0);
```

```
cf.bat
```

hash.sh

Hashes a file, ignoring all whitespace and comments. Use for

stress.sh

21 lines

#!/bin/bash

proq_A=\$1

 $\begin{tabular}{ll} \#\ prog_A\ and\ prog_B\ are\ the\ executables\ to\\ compare \end{tabular}$

```
prog_B=$2
generator=$3

inp_file="inp_${generator}.txt"
out_file1="outA_${generator}.txt"
out_file2="outB_${generator}.txt"
```

```
for ((i = 1; ; ++i)) do
    echo $i
    "./$generator" > $inp_file
    "./$prog_A" < $inp_file > $out_file1
    "./$prog_B" < $inp_file > $out_file2
    diff -w "${out_file1}" "${out_file2}" ||
         break
```

done

notify-send "bug found!!!!"

Mathematics (2)

2.1 Equations

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

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

$$ax + by = e$$

$$cx + dy = f$$

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

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

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

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

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

2.2 Recurrences

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

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

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

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

$$s_a = \sqrt{bc \left[1 - \left(\frac{a}{b+c} \right)^2 \right]}$$
Law of sines:
$$\frac{\sin \alpha}{a} = \frac{\sin \beta}{b} = \frac{\sin \gamma}{c} = \frac{1}{2R}$$
Law of cosines:
$$a^2 = b^2 + c^2 - 2bc \cos \alpha$$

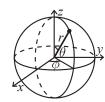
2.4.2 Quadrilaterals

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

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

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

2.4.3 Spherical coordinates



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

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

$$z = r \cos \theta \qquad \phi = a\tan(2y, x)$$

2.5 Derivatives/Integrals

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

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

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

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

$$\int \tan ax = -\frac{\ln|\cos ax|}{a}$$

$$\int x\sin ax = \frac{\sin ax - ax\cos ax}{a^2}$$

$$\int e^{-x^2} = \frac{\sqrt{\pi}}{2}\operatorname{erf}(x)$$

$$\int xe^{ax}dx = \frac{e^{ax}}{a^2}(ax-1)$$

Integration by parts:

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

2.6 Sums

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

$$\sum_{k=1}^{n} k = \frac{n(n+1)}{2}$$

$$\sum_{k=1}^{n} k^2 = \frac{n(2n+1)(n+1)}{6}$$

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

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

2.7 Series

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

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

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

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

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

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 σ is the standard deviation. If X is instead continuous it will have a probability density function $f_X(x)$ and the sums above will instead be integrals with $p_X(x)$ replaced by $f_X(x)$.

Expectation is linear:

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

For independent X and Y,

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

2.8.1 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

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

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

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

First success distribution

The number of trials needed to get the first success in independent yes/no experiments, each which yields success with probability p is $Fs(p), 0 \le p \le 1.$

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

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

Poisson distribution

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

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

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

2.8.2 Continuous distributions Uniform distribution

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

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

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

OrderStatisticTree HashMap SegmentTree segtree UnionFindRollback

Exponential distribution

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

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

Normal distribution

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

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

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

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

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.

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

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

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

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

Data structures (3)

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

782797, 16 lines

```
#include <bits/extc++.h>
using namespace __gnu_pbds;

template<class T>
using Tree = tree<T, null_type, less<T>,
        rb_tree_tag,
        tree_order_statistics_node_update>;

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

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) 1.7 lines

```
#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 = l1(4e18 * acos(0)) | 71;
    l1 operator()(1l x) const { return
        _builtin_bswap64(x*C); }
};
_gnu_pbds::gp_hash_table<11,int,chash> h({},{},{},{},{},{1<<16});</pre>
```

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

```
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 /=
     if (b % 2) ra = f(ra, s[b++]);
     if (e % 2) rb = f(s[--e], rb);
   return f(ra, rb);
};
segtree.cpp
                                    deb606, 92 lines
template<class S> struct segtree {
 int n; vector<S> t;
```

```
void init(int ) { n = ; t.assign(n+n-1, S
  void init(const vector<S>& v) {
    n = sz(v); t.assign(n + n - 1, S());
    build(0,0,n-1,v);
  } template <typename... T>
  void upd(int 1, int r, const T&... v) {
    assert(0 <= 1 && 1 <= r && r < n);
    upd(0, 0, n-1, 1, r, v...);
 S get(int 1, int r) {
    assert(0 <= 1 && 1 <= r && r < n);
   return get(0, 0, n-1, 1, r);
private:
  inline void push(int u, int b, int e) {
    if (t[u].lazy == 0) return;
    int mid = (b+e)>>1, rc = u+((mid-b+1)<<1);</pre>
   t[u+1].upd(b, mid, t[u].lazy);
   t[rc].upd(mid+1, e, t[u].lazy);
   t[u].lazy = 0;
  void build(int u, int b, int e, const vector
       <S>& V) {
    if (b == e) return void(t[u] = v[b]);
    int mid = (b+e)>>1, rc = u+((mid-b+1)<<1);</pre>
    build(u+1, b, mid, v); build(rc, mid+1, e,
          v);
    t[u] = t[u+1] + t[rc];
  } template<typename... T>
  void upd(int u, int b, int e, int l, int r,
      const T&... v) {
    if (1 <= b && e <= r) return t[u].upd(b, e</pre>
         , v...);
    push(u, b, e);
    int mid = (b+e)>>1, rc = u+((mid-b+1)<<1);</pre>
    if (1 <= mid) upd(u+1, b, mid, 1, r, v...)</pre>
    if (mid < r) upd(rc, mid+1, e, 1, r, v...)</pre>
   t[u] = t[u+1] + t[rc];
 S get(int u, int b, int e, int 1, int r) {
    if (1 <= b && e <= r) return t[u];</pre>
    push(u, b, e);
    S res; int mid = (b+e)>>1, rc = u+((mid-b
         +1) <<1);
```

```
if (r \le mid) res = get (u+1, b, mid, l, r)
    else if (mid < 1) res = get(rc, mid+1, e,</pre>
    else res = get(u+1, b, mid, l, r) + get(rc
        , mid+1, e, 1, r);
    t[u] = t[u+1] + t[rc]; return res;
};
/* Segment Tree
Inspiration: tourist, atcoder library
(1) Declaration:
  Create a node class (sample below).
  node class must have the following:
  * A constructor (to create empty nodes and
      also to make inplace nodes).
  * + operator: returns a node which contains
      the merged information of two nodes.
  * upd(b, e, ...): updates this node
      representing the range [b, e] using
      information from ...
  Now, segtree<node> T; declares the tree.
  You can use T. init(100) to create an empty
      tree of 100 nodes in [0, 100) range.
  You can also make a vector<node> v; Then put
        values in the vector v and make the
       tree using
   v by, T. init(v); This works in linear time
        and is faster than updating each
        individually.
(2) Usage:
  (2.1) init(int siz) or init(vector):
    Described above
  (2.2) \ upd(l, r, ...v):
    Update the range [l, r] with the
         information in ...
    Make sure the number of elements and the
         order of them you put here is the
         exact same
    as you declared in your node.upd()
         function.
struct node {
 11 sum;
  ll lazv;
  node(11 _a = 0, 11 _b = 0) : sum(_a), lazy(
  node operator+(const node &obj) {
   return {sum + obj.sum, 0};
  void upd(int b, int e, ll x) {
   sum += (e - b + 1) * x;
    lazv += x:
};
```

UnionFindRollback.h

Description: Disjoint-set data structure with undo. If

SubMatrix Matrix LineContainer Treap FenwickTree FenwickTree2d RMQ

```
undo is not needed, skip st, time() and rollback().
Usage: int t = uf.time(); ...; uf.rollback(t);
Time: \mathcal{O}(\log(N))
                                     de4ad0, 21 lines
struct RollbackUF {
  vi e; vector<pii> st;
  RollbackUF(int n) : e(n, -1) {}
  int size(int x) { return -e[find(x)]; }
  int find(int x) { return e[x] < 0 ? x : find
       (e[x]); }
  int time() { return sz(st); }
  void rollback(int t) {
    for (int i = time(); i --> t;)
     e[st[i].first] = st[i].second;
    st.resize(t);
  bool join(int a, int b) {
    a = find(a), b = find(b);
    if (a == b) return false;
    if (e[a] > e[b]) swap(a, b);
    st.push back({a, e[a]});
    st.push_back({b, e[b]});
    e[a] += e[b]; e[b] = a;
    return true;
};
SubMatrix.h
Description: Calculate submatrix sums quickly, given
upper-left and lower-right 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][
         1];
};
Matrix.h
Description: Basic operations on square matrices.
Usage: Matrix<int, 3> A;
A.d = \{\{\{1,2,3\}\}, \{\{4,5,6\}\}, \{\{7,8,9\}\}\}\};
vector < int > vec = \{1, 2, 3\};
vec = (A^N) * vec;
                                      c43c7d, 26 lines
template < class T, int N> struct Matrix {
  typedef Matrix M;
  array<array<T, N>, N> d{};
  M operator*(const M& m) const {
    rep(i,0,N) rep(j,0,N)
      rep(k, 0, N) \ a.d[i][j] += d[i][k] *m.d[k][j]
           ];
    return a;
```

```
vector<T> operator*(const vector<T>& vec)
       const {
    vector<T> ret(N);
    rep(i, 0, N) rep(j, 0, N) ret[i] += d[i][j] *
         vec[j];
    return ret:
  M operator^(ll p) const {
    assert (p >= 0);
    M a, b(*this);
    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. Use-
ful 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</pre>
        k < o.k; }
 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->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;
  \textbf{void} \ \text{add(ll } k \textbf{, 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(!emptv());
    auto 1 = *lower_bound(x);
    return 1.k * x + 1.m;
};
```

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

```
Time: \mathcal{O}(\log N)
                                     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) >= k) { // "n->val >= k" for
      lower\_bound(k)
    auto pa = split(n->1, k);
   n->1 = pa.second;
    n->recalc();
    return {pa.first, n};
  } else {
    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 <= 1) t = merge(ins(a, b, k), c);</pre>
 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 value.

Time: Both operations are $\mathcal{O}(\log N)$. e62fac, 22 lines

```
struct FT {
 vector<ll> s;
 FT(int n) : s(n) {}
 void update(int pos, 11 dif) { // a[pos] \neq =
   for (; pos < sz(s); pos |= pos + 1) s[pos]
         += dif;
 11 query (int pos) { // sum of values in [0,
   11 \text{ res} = 0;
   for (; pos > 0; pos &= pos - 1) res += s[
         pos-11;
   return res;
 int lower_bound(ll sum) {// min pos st sum
       of [0, pos] >= sum
    // Returns n if no sum is >= sum, or -1 if
          empty sum is.
   if (sum <= 0) return -1;</pre>
   int pos = 0;
   for (int pw = 1 << 25; pw; pw >>= 1) {
      if (pos + pw \leq sz(s) && s[pos + pw-1] \leq
        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"
                                     157f07, 22 lines
struct FT2 {
  vector<vi> ys; vector<FT> ft;
  FT2(int limx) : ys(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 y) {
    return (int) (lower_bound(all(ys[x]), y) -
         vs[x].begin()); }
  void update(int x, int y, ll dif) {
    for (; x < sz(ys); x | = x + 1)
      ft[x].update(ind(x, y), dif);
 11 query(int x, int y) {
    11 \text{ sum} = 0;
    for (; x; x &= x - 1)
      sum += ft[x-1].query(ind(x-1, y));
    return sum;
};
```

RMQ.h

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

```
Usage: RMO rmg(values);
rmg.querv(inclusive, exclusive);
Time: \mathcal{O}(|V|\log|V|+Q)
                                     510c32, 16 lines
template<class T>
struct RMO {
  vector<vector<T>> jmp;
  RMO(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])
              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 <<</pre>
          dep)]);
};
```

MoQueries.h

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

```
Time: \mathcal{O}(N\sqrt{Q})
```

```
a12ef4, 49 lines
void add(int ind, int end) { ... } // add a[
     ind \mid (end = 0 \text{ 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 > q.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]1 / 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; \} \setminus
                  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[
        il;
   return val;
 void diff() {
   rep(i,1,sz(a)) a[i-1] = i*a[i];
   a.pop_back();
 void divroot(double x0) {
   double b = a.back(), c; a.back() = 0;
   for(int i=sz(a)-1; i--;) c = a[i], a[i] =
        a[i+1]*x0+b, b=c;
   a.pop_back();
};
```

PolyRoots.h

```
Description: Finds the real roots to a polynomial.
Usage: polyRoots(\{\{2, -3, 1\}\}, -1e9, 1e9\}) // solve
x^2-3x+2 = 0
Time: \mathcal{O}\left(n^2\log(1/\epsilon)\right)
                                         b00bfe, 23 lines
"Polynomial.h"
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 1 = 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 \le 0) ^ sign) 1 = m;
      else h = m;
    ret.push_back((1 + h) / 2);
return ret;
```

PolyInterpolate.h

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

```
typedef vector<double> vd;
vd interpolate(vd x, vd y, int n) {
 vd res(n), temp(n);
 rep(k, 0, n-1) rep(i, k+1, n)
   v[i] = (v[i] - v[k]) / (x[i] - x[k]);
 double last = 0; temp[0] = 1;
 rep(k,0,n) rep(i,0,n) {
   res[i] += v[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 brute-forcing the first terms. Should work on any field, but numerical stability for floats is not guaranteed. Output will have size $\leq n$. Usage: berlekampMassey({0, 1, 1, 3, 5, 11}) // {1, 2}

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

```
"../number-theory/ModPow.h"
vector<ll> berlekampMassev(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) %
   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...>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)
```

```
f4e444, 26 lines
```

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

4.2 Optimization

GoldenSectionSearch.h

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

```
Usage:
                    double func(double x) { return
4+x+.3*x*x; }
double xmin = gss(-1000, 1000, func);
Time: \mathcal{O}(\log((b-a)/\epsilon))
                                          31d45b, 14 lines
```

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

```
} else {
    a = x1; x1 = x2; f1 = f2;
    x2 = a + r*(b-a); f2 = f(x2);
}
return a;
}
```

HillClimbing.h

Description: Poor man's optimization for unimodal functions.

8eeeaf, 14 lines

```
typedef array<double, 2> P;

template<class F> pair<double, P> hillClimb(P
    start, F f) {
    pair<double, P> cur(f(start), start);
    for (double jmp = 1e9; jmp > 1e-20; jmp /=
        2) {
        rep(j,0,100) rep(dx,-1,2) rep(dy,-1,2) {
            P p = cur.second;
            p[0] += dx*jmp;
            p[1] += dy*jmp;
            cur = min(cur, make_pair(f(p), p));
        }
    return cur;
```

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;
}
```

IntegrateAdaptive.h

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

```
Usage: double sphereVolume = quad(-1, 1, [](double x) { return quad(-1, 1, [&](double y) { return quad(-1, 1, [&](double z) { return x*x + y*y + z*z < 1; \});});}_{2dd79, 15 lines}
```

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

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

Simplex.h

int m, n;

vi N, B;

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

```
mance, 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 _{aa8530, 68 \text{ lines}}^{aa8530, 68 \text{ lines}}
```

```
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(j,0,m) {
```

if (D[i][s] <= eps) continue;</pre>

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

4.3 Matrices

Determinant.h

Description: Calculates determinant of a matrix. Destroys the matrix.

```
Time: \mathcal{O}\left(N^3\right) bd5cec, 15 lines
```

IntDeterminant.h

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

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

SolveLinear.h

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

```
e: O (n m) 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 SolveLinear, 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:; }
```

SolveLinearBinarv.h

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

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

fa2d7a, 34 lines

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

MatrixInverse.h

Description: Invert matrix A. Returns rank; result is stored in A unless singular (rank < n). Can easily be extended to prime moduli; for prime powers, repeatedly set $A^{-1} = A^{-1}(2I - AA^{-1}) \pmod{p^k}$ where A^{-1} starts as the inverse of A 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
  rep(i, 0, n) tmp[i][i] = 1, col[i] = i;
  rep(i,0,n) {
   int r = i, c = i;
   rep(j,i,n) rep(k,i,n)
     if (fabs(A[j][k]) > fabs(A[r][c]))
        r = j, c = k;
   if (fabs(A[r][c]) < 1e-12) return i;</pre>
```

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

Tridiagonal.h

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

$$\begin{pmatrix} b_0 \\ b_1 \\ b_2 \\ b_3 \\ \vdots \\ b_{n-1} \end{pmatrix}$$

$$= \begin{pmatrix} d_0 & p_0 & 0 & 0 & \cdots & 0 \\ q_0 & d_1 & p_1 & 0 & \cdots & 0 \\ 0 & q_1 & d_2 & p_2 & \cdots & 0 \\ \vdots & \vdots & \ddots & \ddots & \ddots & \vdots \\ 0 & 0 & \cdots & q_{n-3} & d_{n-2} & p_{n-2} \\ 0 & 0 & \cdots & 0 & q_{n-2} & d_{n-1} \end{pmatrix}$$

$$= \begin{pmatrix} x_0 \\ x_1 \\ x_2 \\ x_3 \\ \vdots \\ x_{n-1} \end{pmatrix} .$$

This is useful for solving problems on the type

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

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

$$\{a_i\}$$
 = tridiagonal($\{1, -1, -1, ..., -1, 1\}, \{0, c_1, c_2, ..., c_n\}, \{b_1, b_2, ..., b_n, 0\}, \{a_0, d_1, d_2, ..., d_n, d_n\}$

Fails if the solution is not unique.

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

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

8f9fa8, 26 lines typedef double T;

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

4.4 Fourier transforms

FastFourierTransform.h

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

```
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];
                                    vi rev(n);
                                    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[</pre>
                                    for (int k = 1; k < n; k *= 2)
\{b_1,b_2,\ldots,b_n,0\},\{a_0,d_1,d_2,\ldots,d_n,a_{n+1}\} for (int i = 0; i < n; i += 2 * k) rep(j)
                                        Cz = rt[j+k] * a[i+j+k]; // (25\% faster)
                                               if hand-rolled)
```

a[i + j + k] = a[i + j] - z;

vd conv(const vd& a, const vd& b) {

a[i + j] += z;

```
if (a.empty() || b.empty()) return {};
vd res(sz(a) + sz(b) - 1);
int L = 32 - \underline{\quad} builtin_clz(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(
fft(out);
rep(i, 0, sz(res)) res[i] = imag(out[i]) / (4)
return res;
```

FastFourierTransformMod.h

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

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

```
"FastFourierTransform.h"
```

```
typedef vector<11> v1;
template<int M> vl convMod(const vl &a, const
    vl &b) {
  if (a.empty() || b.empty()) return {};
  vl res(sz(a) + sz(b) - 1);
  int B=32-__builtin_clz(sz(res)), n=1<<B, cut</pre>
       =int(sgrt(M));
  vector<C> L(n), R(n), outs(n), outl(n);
  rep(i,0,sz(a)) L[i] = C((int)a[i] / cut, (
       int)a[i] % cut);
  rep(i, 0, sz(b)) R[i] = C((int)b[i] / cut, (
       int)b[i] % cut);
  fft(L), fft(R);
  rep(i,0,n) {
    int j = -i \& (n - 1);
    outl[j] = (L[i] + conj(L[j])) * R[i] /
         (2.0 * n);
    outs[j] = (L[i] - conj(L[j])) * R[i] /
          (2.0 * n) / 1i;
  fft (outl), fft (outs);
  rep(i,0,sz(res)) {
    11 \text{ av} = 11(\text{real}(\text{outl}[i]) + .5), \text{ cv} = 11(\text{imag})
         (outs[i])+.5);
    11 \text{ bv} = 11(\text{imag}(\text{outl}[i]) + .5) + 11(\text{real}(
         outs[i])+.5);
    res[i] = ((av % M * cut + bv) % M * cut +
         cv) % M;
 return res:
```

NumberTheoreticTransform.h

Description: ntt(a) computes $\hat{f}(k) = \sum_x a[x]g^{xk}$ for all k, where $g = \operatorname{root}^{(mod-1)/N}$. N must be a power of 2. Useful for convolution modulo specific nice primes of the form 2^ab+1 , where the convolution result has size at most 2^a . For arbitrary modulo, see FFTMod. conv(a, b) = c, where $c[x] = \sum a[i]b[x-i]$. For manual convolution: NTT the inputs, multiply pointwise, divide by n, reverse(start+1, end), NTT back. Inputs must be in [0, mod).

```
Time: \mathcal{O}(N \log N)
"../number-theory/ModPow.h"
const 11 mod = (119 << 23) + 1, root = 62; //</pre>
    = 998244353
// For p < 2^30 there is also e.g. 5 << 25, 7
    << 26, 479 << 21
// and 483 << 21 (same root). The last two are
     > 10^9.
typedef vector<11> v1;
void ntt(vl &a) {
  int n = sz(a), L = 31 - __builtin_clz(n);
  static v1 rt(2, 1);
  for (static int k = 2, s = 2; k < n; k \neq 2,
       s++) {
   rt.resize(n);
   11 z[] = \{1, modpow(root, mod >> s)\};
    rep(i,k,2*k) rt[i] = rt[i / 2] * z[i & 1]
  vi rev(n);
  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
     11 z = rt[j + k] * a[i + j + k] % mod, &
          ai = a[i + j];
     a[i + j + k] = ai - z + (z > ai ? mod :
     ai += (ai + z >= mod ? z - mod : z);
vl conv(const vl &a, const vl &b) {
  if (a.emptv() || b.emptv()) return {};
  int s = sz(a) + sz(b) - 1, B = 32 -
       builtin clz(s), n = 1 \ll B;
  int inv = modpow(n, mod - 2);
  vl L(a), R(b), out(n);
  L.resize(n), R.resize(n);
  ntt(L), ntt(R);
  rep(i,0,n) out[-i & (n-1)] = (11)L[i] * R[
      il % mod * inv % mod;
  return {out.begin(), out.begin() + s};
```

FastSubsetTransform.h

Description: Transform to a basis with fast convolutions of the form $c[z] = \sum_{z=x \oplus y} a[x] \cdot b[y]$, where \oplus 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
```

Number theory (5)

5.1 Modular arithmetic

Modular Arithmetic.h

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

```
"euclid.h"
                                    35bfea, 18 lines
const 11 mod = 17; // change to something else
struct Mod {
 11 x;
 Mod(ll xx) : x(xx) \{ \}
  Mod operator+(Mod b) { return Mod((x + b.x)
      % mod); }
  Mod operator-(Mod b) { return Mod((x - b.x +
       mod) % mod); }
  Mod operator* (Mod b) { return Mod((x * b.x)
      % mod); }
  Mod operator/(Mod b) { return *this * invert
       (b); }
 Mod invert (Mod a) {
   ll x, y, g = euclid(a.x, mod, x, y);
   assert(q == 1); return Mod((x + mod) % mod
  Mod operator^(ll e) {
   if (!e) return Mod(1);
   Mod r = *this ^ (e / 2); r = r * r;
   return e&1 ? *this * r : r;
};
```

ModInverse.h

Description: Pre-computation of modular inverses. Assumes LIM \leq mod and that mod is a prime $_{66684f,\ 3\ lines}$

```
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

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

```
ll modpow(ll b, ll 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: $\mathcal{O}(\sqrt{m})$

ModSum.h

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

Time: $\log(m)$, with a large constant. 5c5bc5, 16 lines

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 by the lines

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

```
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"
ll sgrt(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)
 11 x = modpow(a, (s + 1) / 2, p);
 ll 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{ qs} = \text{modpow}(q, 1LL << (r - m - 1), p);
   q = qs * qs % p;
   x = x * qs % p;
```

5.2 Primality

b = b * q % p;

FastEratosthenes.h

 $\bf Description:$ Prime sieve for generating all primes smaller than LIM.

```
Time: LIM=1e9 \approx 1.5s 6b2912, 20 lines
```

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

MillerRabin Factor euclid CRT phiFunction ContinuedFractions FracBinarySearch

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)</pre>
       == 3;
  ull A[] = \{2, 325, 9375, 28178, 450775,
       9780504, 1795265022},
      s = \underline{\quad}builtin_ctzll(n-1), d = n >> s;
  for (ull a : A) { // ^ count trailing
       zeroes
    ull p = modpow(a%n, d, n), i = s;
    while (p != 1 && p != n - 1 && a % n && i
      p = modmul(p, p, n);
    if (p != n-1 && i != s) return 0;
  return 1;
```

Factor.h

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

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

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

Divisibility 5.3

euclid.h

Description: Finds two integers x and y, such that $ax + by = \gcd(a, b)$. If you just need gcd, use the built in $_gcd$ instead. If a and b are coprime, then x is the inverse of $a \pmod{b}$. 33ba8f, 5 lines

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

Description: Chinese Remainder Theorem.

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

```
"euclid.h"
                                     04d93a, 7 lines
11 crt(ll a, ll m, ll b, ll n) {
  if (n > m) swap(a, b), swap(m, n);
  ll x, y, g = euclid(m, n, x, y);
  assert((a - b) % g == 0); // else no
       solution
  x = (b - a) % n * x % n / g * m + a;
  return x < 0 ? x + m*n/q : x;
```

5.3.1 Bézout's identity

For $a \neq b \neq 0$, then d = 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 ϕ function is defined as $\phi(n) :=$ # of positive integers $\leq n$ that are coprime with n. $\phi(1)=1,\ p$ prime $\Rightarrow \phi(p^k)=(p-1)p^{k-1},\ m,n$ coprime $\Rightarrow \phi(mn) = \phi(m)\phi(n)$. If $n = p_1^{k_1} p_2^{k_2} ... p_r^{k_r}$ then $\phi(n) =$ $(p_1-1)p_1^{k_1-1}...(p_r-1)p_r^{k_r-1}.$ $\phi(n)=n\cdot\prod_{p\mid n}(1-1/p).$ $\sum_{d|n} \phi(d) = n, \sum_{1 \le k \le n, \gcd(k,n)=1} k = n\phi(n)/2, n > 1$

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]</pre>
    for (int j = i; j < LIM; j += i) phi[j] -=</pre>
          phi[j] / i;
```

5.4 Fractions

ContinuedFractions.h

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

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

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

```
typedef double d; // for N \sim 1e7; long double
    for N \sim 1e9
pair<ll, ll> approximate(d x, ll N) {
 11 LP = 0, LQ = 1, P = 1, Q = 0, inf =
      LLONG_MAX; d y = x;
  for (;;) {
```

```
ll lim = min(P ? (N-LP) / P : inf, O ? (N-LP)
     LO) / O : inf),
   a = (11) floor(v), 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)Q))?
    make_pair(NP, NQ) : make_pair(P, Q);
if (abs(y = 1/(y - (d)a)) > 3*N) {
  return {NP, NQ};
LP = P; P = NP;
LQ = Q; Q = NQ;
```

FracBinarySearch.h

Description: Given f and N, finds the smallest fraction $p/q \in [0,1]$ such that f(p/q) is true, and p,q < 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: $\mathcal{O}(\log(N))$ 27ab3e, 25 lines

```
struct Frac { ll p, q; };
```

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

5.5 Pythagorean Triples

The Pythagorean triples are uniquely generated

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

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

5.6 Primes

p = 962592769 is such that $2^{21} \mid p - 1$, which may be useful. For hashing use 970592641 (31-bit number), 31443539979727 (45-bit), 3006703054056749 (52-bit). There are 78498 primes less than 1000000.

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

5.7 Estimates

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

The number of divisors of n is at most around 100 for n < 5e4, 500 for n < 1e7, 2000 for n < 1e10, 200000 for n < 1e19.

5.8 Mobius Function

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

Mobius Inversion:

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

Other useful formulas/forms:

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

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

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

Combinatorial (6)

6.1 Permutations

6.1.1 Factorial

n						9	
n!	1 2 6	24 1	20 720	5040	40320	362880	3628800
n	11	12	13	14	15	16	17
n!	4.0e7	′ 4.8e	8.6.2e9	8.7e	10 1.3e	12 2.1e1	3.6e1
n	20	25	30	40	50 1	00 150) 17
n!	2e18	2e25	3e32 8	3e47 3	e64 9e	157 6e26	52 > DBL

IntPerm.h

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

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

int permToInt(vi& v) { **int** use = 0, i = 0, r = 0; for(int x:v) r = r * ++i +__builtin_popcount(use & -(1 << x)), 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.

$$D(n) = (n-1)(D(n-1) + D(n-2)) = nD(n-1) + (n-1)(D(n-1) + D(n-1)) = nD(n-1) + (n-1)(D(n-1) + D(n-1)(D(n-1)) = nD(n-1) + (n-1)(D(n-1) + D(n-1)(D(n-1)) = nD(n-1) + (n-1)(D(n-1) + D(n-1)(D(n-1)) = nD(n-1) + (n-1)(D(n-1)(D(n-1)) = nD(n-1)(D(n-1)(D(n-1)) = nD(n-1)(D(n-1)(D(n-1)) = nD(n-1)(D(n-1)(D(n-1)(D(n-1)) = nD(n-1)(D(n-1)(D(n-1)(D(n-1)(D(n-1))) = nD(n-1)(D(n-1)(D(n-1)(D(n-1)(D(n-1)(D(n-1)(D(n-1)(D(n-1)(D(n-1)(D(n-1)(D(n-1)(D(n-1)(D(n-1)(D(n-1)(D(n-1)(D(n-1)(D(n-1)(D(n-1)(D(n-1)(D(n$$

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(q.x = x).

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

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

6.2 Partitions and subsets

6.2.1 Partition function

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

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

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

a0a312, 6 line

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

General purpose numbers

6.3.1 Bernoulli numbers

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

Sums of powers:

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

Euler-Maclaurin formula for infinite

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

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

6.3.2Stirling numbers of the first

Number of permutations on n items with k

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

$$p(0) = 1, \ p(n) = \sum_{k \in \mathbb{Z} \setminus \{0\}} (-1)^{k+1} p(n - k(3k - 1)/2) \begin{cases} c(8, k) = \\ 8, 0, 5040, 13068, 13132, 6769, 1960, 322, 28, 1 \\ c(n, 2) = \\ 0, 0, 1, 3, 11, 50, 274, 1764, 13068, 109584.... \end{cases}$$

6.3.3 Eulerian numbers

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

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

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

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

Stirling numbers of the second

Partitions of n distinct elements into exactly kgroups.

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

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

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

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

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

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

cycles.
$$C_n = \frac{1}{n+1} \binom{2n}{n} = \binom{2n}{n} - \binom{2n}{n+1} = \frac{(2n)!}{(n+1)!n!}$$
 FloydWarshall.h Description: Calculates all-pairs shortest path in a direction of the position of the positio

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

 $C_n = 1, 1, 2, 5, 14, 42, 132, 429, 1430, 4862, 16796, 5 \\ 8786 \text{ in } 11 \text{ inf = 1LL } << 62; \\ 12 \text{ inf = 1LL } << 62; \\ 13 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 14 \text{ inf = 1LL } << 62; \\ 1$

• sub-diagonal monotone paths in an $n \times n$ grid.

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

Graph (7)

7.1 Fundamentals

Ed>& eds, int s) {

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 ?</pre>

a : -a; }}; struct Node { ll dist = inf; int prev = -1; }; void bellmanFord(vector<Node>& nodes, vector<

nodes[s].dist = 0;sort(all(eds), [](Ed a, Ed b) { return a.s() < b.s(); }); 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;

put is an distance matrix m, where $m[i][j] = \inf if i$ and j are not adjacent. As output, m[i][j] is set to the shortest distance between i and j, inf if no path, or -inf if the path goes through a negative-weight cycle. Time: $\mathcal{O}(N^3)$

void floydWarshall (vector<vector<ll>>& m) { int n = sz(m);rep(i, 0, n) m[i][i] = min(m[i][i], OLL);

TopoSort PushRelabel MinCostMaxFlow EdmondsKarp MinCut GlobalMinCut

```
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], -
   m[i][j] = min(m[i][j], newDist);
rep(k, 0, n) if (m[k][k] < 0) rep(i, 0, n) rep(j)
  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|)
                                     66a137, 14 lines
vi topoSort(const vector<vi>& gr) {
  vi indeg(sz(gr)), ret;
  for (auto& li : gr) for (int x : li) indeg[x
  queue<int> q; // use priority_queue for
       lexic. largest ans.
  rep(i, 0, sz(gr)) if (indeg[i] == 0) q.push(i)
  while (!q.empty()) {
    int i = q.front(); // top() for priority
    ret.push_back(i);
    q.pop();
    for (int x : gr[i])
      if (--indeg[x] == 0) q.push(x);
  return ret;
```

7.2 Network flow

PushRelabel.h

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

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

0ae1d4, 48 lines

```
struct PushRelabel {
  struct Edge {
   int dest, back;
   11 f, c;
  };
  vector<vector<Edge>> g;
  vector<ll> 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;
  ll 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 : g[s]) addFlow(e, e.c);
    for (int hi = 0;;) {
      while (hs[hi].empty()) if (!hi--) return
            -ec[s];
      int u = hs[hi].back(); hs[hi].pop_back()
      while (ec[u] > 0) // discharge u
        if (cur[u] == q[u].data() + sz(q[u]))
          H[u] = 1e9;
          for (Edge& e : g[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] <
              --co[H[i]], H[i] = v + 1;
          hi = H[u];
        } 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
};
```

MinCostMaxFlow.h

Description: Min-cost max-flow. cap[i][j] != cap[j][i] is allowed; double edges are not. 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: Approximately \mathcal{O}(E^2)
```

```
fe85cc, 81 lines
#include <bits/extc++.h>
const 11 INF = numeric limits<11>::max() / 4;
typedef vector<11> VL;
struct MCMF {
 int N;
 vector<vi> ed, red;
 vector<VL> cap, flow, cost;
  vi seen;
  VL dist, pi;
  vector<pii> par;
  MCMF (int N) :
    N(N), ed(N), red(N), cap(N, VL(N)), flow(
        cap), cost (cap),
    seen(N), dist(N), pi(N), par(N) {}
  void addEdge(int from, int to, ll cap, ll
    this->cap[from][to] = cap;
```

```
this->cost[from][to] = cost;
  ed[from].push_back(to);
 red[to].push back(from);
void path(int s) {
 fill(all(seen), 0);
 fill(all(dist), INF);
 dist[s] = 0; ll di;
  __qnu_pbds::priority_queue<pair<11, int>>
 vector<decltype(q)::point_iterator> its(N)
 q.push({0, s});
  auto relax = [&](int i, ll cap, ll cost,
      int dir) {
   11 val = di - pi[i] + cost;
   if (cap && val < dist[i]) {
      dist[i] = val;
      par[i] = {s, dir};
      if (its[i] == q.end()) its[i] = q.push
           ({-dist[i], i});
      else q.modify(its[i], {-dist[i], i});
 };
 while (!q.empty()) {
   s = q.top().second; q.pop();
   seen[s] = 1; di = dist[s] + pi[s];
   for (int i : ed[s]) if (!seen[i])
     relax(i, cap[s][i] - flow[s][i], cost[
          s][i], 1);
   for (int i : red[s]) if (!seen[i])
      relax(i, flow[i][s], -cost[i][s], 0);
 rep(i,0,N) pi[i] = min(pi[i] + dist[i],
pair<11, 11> maxflow(int s, int t) {
 11 totflow = 0, totcost = 0;
 while (path(s), seen[t]) {
   11 fl = INF;
   for (int p,r,x = t; tie(p,r) = par[x], x
         ! = s; x = p)
      fl = min(fl, r ? cap[p][x] - flow[p][x]
          ] : flow[x][p]);
   totflow += fl;
    for (int p,r,x = t; tie(p,r) = par[x], x
         ! = s; x = p)
     if (r) flow[p][x] += fl;
      else flow[x][p] -= fl;
 rep(i,0,N) rep(j,0,N) totcost += cost[i][j]
      ] * flow[i][j];
 return {totflow, totcost};
// 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 (int to : ed[i]) if (cap[i][to])
```

```
if ((v = pi[i] + cost[i][to]) < pi[</pre>
               to])
            pi[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 482fe0. 35 lines

```
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 >
          par[e.first] = x;
          q[ptr++] = e.first;
          if (e.first == sink) goto out;
    return flow;
out:
    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].</pre>
           erase(y);
      graph[y][p] += inc;
```

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

GlobalMinCut.h

Description: Find a global minimum cut in an undirected graph, as represented by an adjacency matrix. Time: $\mathcal{O}(V^3)$

```
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

```
"PushRelabel.h"
```

```
0418b3, 13 lines
typedef array<11, 3> Edge;
vector<Edge> gomoryHu(int N, vector<Edge> ed)
  vector<Edge> tree;
 vi par(N);
  rep(i,1,N) {
   PushRelabel D(N); // Dinic also works
   for (Edge t : ed) D.addEdge(t[0], t[1], t
        [2], t[2]);
   tree.push_back({i, par[i], D.calc(i, par[i
        ])});
   rep(j,i+1,N)
     if (par[j] == par[i] && D.leftOfMinCut(j
          )) par[j] = i;
 return tree;
```

7.3 Matching

hopcroftKarp.h

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

```
Usage: vi btoa(m, -1); hopcroftKarp(g, btoa);
Time: \mathcal{O}\left(\sqrt{V}E\right)
                                                  f612e4, 42 lines
```

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

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

DFSMatching.h

Description: Simple 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); dfsMatching(q, btoa);
Time: \mathcal{O}(VE)
```

```
522b98, 22 lines
bool find(int j, vector<vi>& g, vi& btoa, vi&
 if (btoa[i] == -1) return 1;
  vis[j] = 1; int di = btoa[j];
  for (int e : q[di])
   if (!vis[e] && find(e, g, btoa, vis)) {
     btoa[e] = di;
      return 1;
  return 0;
int dfsMatching(vector<vi>& q, vi& btoa) {
  vi vis:
  rep(i, 0, sz(q)) {
   vis.assign(sz(btoa), 0);
    for (int j : g[i])
     if (find(j, g, btoa, vis)) {
        btoa[j] = i;
        break:
  return sz(btoa) - (int)count(all(btoa), -1);
```

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 indepen-

```
"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}\left(N^2M\right)
                                     1e0fe9, 31 lines
pair<int, vi> hungarian(const vector<vi> &a) {
 if (a.empty()) return {0, {}};
 int n = sz(a) + 1, m = sz(a[0]) + 1;
 vi u(n), v(m), p(m), ans(n - 1);
 rep(i,1,n) {
   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] -
        if (cur < dist[j]) dist[j] = cur, pre[</pre>
             j] = j0;
        if (dist[i] < delta) delta = dist[i],</pre>
             j1 = j;
      rep(j,0,m) {
        if (done[j]) u[p[j]] += delta, v[j] -=
              delta;
        else dist[j] -= delta;
      j0 = j1;
    } while (p[j0]);
    while (j0) { // update alternating path
```

```
int j1 = pre[j0];
   p[j0] = p[j1], j0 = j1;
rep(j,1,m) if (p[j]) ans[p[j] - 1] = j - 1;
return {-v[0], ans}; // min cost
```

GeneralMatching.h

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

```
"../numerical/MatrixInverse-mod.h"
                                    cb1912, 40 lines
vector<pii> generalMatching(int N, vector<pii
  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, fi
 assert (r % 2 == 0);
 if (M != N) do {
   mat.resize(M, vector<11>(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) {
```

7.4 DFS algorithms

swap(fi,fj);

return ret:

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

 $ll \ a = modpow(A[fi][fj], mod-2);$

ll b = A[i][fj] * a % mod;

j] * b) % mod;

rep(i,0,M) if (has[i] && A[i][fj]) {

rep(j, 0, M) A[i][j] = (A[i][j] - A[fi][

```
Usage: scc(graph, [&](vi& v) { ... }) visits
all components
in reverse topological order. comp[i] holds
the component
index of a node (a component only has edges to
components with
lower index). ncomps will contain the number
of components.
Time: \mathcal{O}\left(E+V\right)
                                     76b5c9, 24 lines
vi val, comp, z, cont;
int Time, ncomps;
template < class G, class F> int dfs (int j, G& q
  int low = val[j] = ++Time, x; z.push back(j)
  for (auto e : g[j]) if (comp[e] < 0)</pre>
   low = min(low, val[e] ?: dfs(e,q,f));
  if (low == val[i]) {
    do {
      x = z.back(); z.pop_back();
      comp[x] = ncomps;
     cont.push back(x);
    } while (x != j);
    f(cont); cont.clear();
   ncomps++;
  return val[j] = low;
template<class G, class F> void scc(G& g, F f)
  int n = sz(g);
  val.assign(n, 0); comp.assign(n, -1);
  Time = ncomps = 0;
  rep(i,0,n) if (comp[i] < 0) dfs(i, g, f);
```

BiconnectedComponents.h

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

```
Usage: int eid = 0; ed.resize(N);
for each edge (a,b) {
ed[a].emplace_back(b, eid);
ed[b].emplace_back(a, eid++); }
bicomps([&](const vi& edgelist) {...});
Time: \mathcal{O}\left(E+V\right)
                                      2965e5, 33 lines
vi num, st;
vector<vector<pii>>> ed;
int Time;
template < class F>
int dfs(int at, int par, F& f) {
  int me = num[at] = ++Time, e, y, top = me;
  for (auto pa : ed[at]) if (pa.second != par)
    tie(y, e) = pa;
    if (num[y]) {
      top = min(top, num[y]);
      if (num[y] < me)
        st.push_back(e);
    } else {
      int si = sz(st);
      int up = dfs(y, e, f);
```

```
top = min(top, up);
   if (up == me) {
        st.push_back(e);
        f(vi(st.begin() + si, st.end()));
        st.resize(si);
    }
   else if (up < me) st.push_back(e);
   else { /* e is a bridge */ }
   }
} return top;
}

template<class F>
void bicomps(F f) {
   num.assign(sz(ed), 0);
   rep(i,0,sz(ed)) if (!num[i]) dfs(i, -1, f);
}
```

2sat.h

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

Usage: TwoSat ts(number of boolean variables); ts.either(0, \sim 3); // Var 0 is true or var 3 is false ts.setValue(2); // Var 2 is true ts.atMostOne($\{0, \sim 1, 2\}$); // <= 1 of vars 0, \sim 1 and 2 are true ts.solve(); // Returns true iff it is solvable ts.values[0..N-1] holds the assigned values to the vars

Time: $\mathcal{O}\left(N+E\right)$, where N is the number of boolean variables, and E is the number of clauses. _{5f9706, 56 lines}

```
struct TwoSat {
 int N;
  vector<vi> gr;
  vi values; // 0 = false, 1 = true
 TwoSat(int n = 0) : N(n), qr(2*n) {}
  int addVar() { // (optional)
    gr.emplace_back();
    gr.emplace_back();
    return N++;
  void either(int f, int j) {
    f = \max(2*f, -1-2*f);
    j = \max(2*j, -1-2*j);
   gr[f].push_back(j^1);
    gr[j].push_back(f^1);
  void setValue(int x) { either(x, x); }
  void atMostOne(const vi& li) { // (optional)
    if (sz(li) <= 1) return;</pre>
    int cur = \simli[0];
    rep(i,2,sz(li)) {
     int next = addVar();
     either(cur, ~li[i]);
     either(cur, next);
     either(~li[i], next);
      cur = ~next;
    either(cur, ~li[1]);
```

```
vi val, comp, z; int time = 0;
 int dfs(int i) {
   int low = val[i] = ++time, x; z.push_back(
        i);
   for(int e : gr[i]) if (!comp[e])
     low = min(low, val[e] ?: dfs(e));
   if (low == val[i]) do {
     x = z.back(); z.pop_back();
     comp[x] = low;
     if (values[x>>1] == -1)
       values[x>>1] = x&1;
   } while (x != i);
   return val[i] = low;
 bool solve() {
   values.assign(N, -1);
   val.assign(2*N, 0); comp = val;
   rep(i,0,2*N) if (!comp[i]) dfs(i);
   rep(i,0,N) if (comp[2*i] == comp[2*i+1])
        return 0;
   return 1:
};
```

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) = gr[x][it++];
   if (!eu[e]) {
     D[x] --, D[y] ++;
     eu[e] = 1; s.push_back(y);
 for (int x : D) if (x < 0 | | sz(ret) !=
      nedges+1) return {};
 return {ret.rbegin(), ret.rend()};
```

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: O(NM)

```
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 ^
  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.

MaximumClique.h

MaximumIndependentSet BinaryLifting LCA CompressTree HLD LinkCutTree

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 free bet, 49 lines

```
typedef vector<bitset<200>> vb;
struct Maxclique {
  double limit=0.025, pk=0;
  struct Vertex { int i, d=0; };
  typedef vector<Vertex> vv;
  vv V;
  vector<vi> C;
  vi qmax, q, S, old;
  void init(vv& r) {
    for (auto& v : r) v.d = 0;
    for (auto& v : r) for (auto i : 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:
     g.push_back(R.back().i);
     vv T;
      for(auto v:R) if (e[R.back().i][v.i]) T.
          push_back({v.i});
      if (sz(T)) {
       if (S[lev]++ / ++pk < limit) init(T);</pre>
       int j = 0, mxk = 1, mnk = max(sz(gmax))
              - 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][
          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
  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 MinimumVertexCover.

Trees

BinaryLifting.h

Description: Calculate power of two jumps in a tree, to support fast upward jumps and LCAs. Assumes the root node points to itself.

```
Time: construction \mathcal{O}(N \log N), queries \mathcal{O}_{\text{bree}_{5}}(\log N)
vector<vi> treeJump(vi& P) {
  int on = 1, d = 1;
  while (on < sz(P)) on *= 2, d++;
  vector<vi> jmp(d, P);
  rep(i,1,d) rep(j,0,sz(P))
    jmp[i][j] = jmp[i-1][jmp[i-1][j]];
  return jmp;
int jmp(vector<vi>& tbl, int nod, int steps){
  rep(i, 0, sz(tbl))
    if(steps&(1<<i)) nod = tbl[i][nod];
  return nod;
int lca(vector<vi>& tbl, vi& depth, int a, int
  if (depth[a] < depth[b]) swap(a, b);</pre>
  a = jmp(tbl, a, depth[a] - depth[b]);
  if (a == b) return a;
  for (int i = sz(tbl); i--;) {
    int c = tbl[i][a], d = tbl[i][b];
    if (c != d) a = c, b = d;
  return tbl[0][a];
```

LCA.h

Description: Data structure for computing lowest common ancestors in a tree (with 0 as root). C should be an adjacency list of the tree, either directed or undirected. Time: $\mathcal{O}(N \log N + Q)$

```
"../data-structures/RMQ.h"
                                     0f62fb, 21 lines
struct LCA {
 int T = 0;
 vi time, path, ret;
  RMQ<int> rmq;
  LCA(vector<vi>& C) : time(sz(C)), rmg((dfs(C
      ,0,-1), ret)) {}
  void dfs(vector<vi>& C, int v, int par) {
   time[v] = T++;
    for (int y : C[v]) if (y != par) {
      path.push_back(v), ret.push_back(time[v
          ]);
      dfs(C, y, v);
 int lca(int a, int b) {
   if (a == b) return a;
   tie(a, b) = minmax(time[a], time[b]);
    return path[rmq.query(a, b)];
  //dist(a,b){return depth[a] + depth[b] - 2*
       depth[lca(a,b)];}
```

```
CompressTree.h
```

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

```
"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] <</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)
                                     6f34db, 46 lines
"../data-structures/LazySegmentTree.h"
template <bool VALS EDGES> struct HLD {
 int N, tim = 0;
 vector<vi> adj;
 vi par, siz, depth, rt, pos;
 Node *tree:
 HLD(vector<vi> adj_)
   : N(sz(adj_)), adj(adj_), par(N, -1), siz(
        N, 1), depth(N),
     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, depth[u] = depth[v] + 1;
     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
   for (; rt[u] != rt[v]; v = par[rt[v]]) {
     if (depth[rt[u]] > depth[rt[v]]) swap(u,
     op(pos[rt[v]], pos[v] + 1);
   if (depth[u] > depth[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

```
Time: All operations take amortized \mathcal{O}\left(\log N\right)_{5909e2, 90 \text{ lines}}
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],
          \star 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];
```

 $z \rightarrow c[h ^1] = b ? x : this;$

DirectedMST Point lineDistance SegmentDistance SegmentIntersection

```
y - > c[i ^1] = b ? this : x;
    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,
    assert(!connected(u, v));
   makeRoot(&node[u]);
    node[u].pp = &node[v];
  void cut (int u, int v) { // remove an edge (
   Node *x = &node[u], *top = &node[v];
   makeRoot(top); x->splay();
    assert(top == (x->pp ?: x->c[0]));
    if (x->pp) x->pp = 0;
     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/arbores-
cence 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;
 ll delta;
  void prop() {
   key.w += delta;
   if (1) 1->delta += delta;
   if (r) r->delta += delta;
  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)));
void pop(Node*& a) { a->prop(); a = merge(a->1
    , a->r); }
pair<ll, vi> dmst(int n, int r, vector<Edge>&
    q) {
  RollbackUF uf(n);
  vector<Node*> heap(n);
  for (Edge e : g) heap[e.b] = merge(heap[e.b
       1, 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) {
      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]
        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.) $_{47ec0a, 28 \text{ lines}}$

```
template <class T> int sqn(T x) { return (x >
    0) - (x < 0);
template<class T>
struct Point {
 typedef Point P;
 explicit Point (T x=0, T y=0) : x(x), y(y) {}
 bool operator<(P p) const { return tie(x,y)</pre>
      < tie(p.x,p.y); }
 bool operator==(P p) const { return tie(x,y)
      ==tie(p.x,p.y); }
 P operator+(P p) const { return P(x+p.x, y+p
 P operator-(P p) const { return P(x-p.x, y-p
 P operator*(T d) const { return P(x*d, y*d);
 P operator/(T d) const { return P(x/d, y/d);
 T dot(P p) const { return x*p.x + y*p.y; }
 T cross(P p) const { return x*p.y - y*p.x; }
 T cross(P a, P b) const { return (a-*this).
      cross(b-*this); }
  T dist2() const { return x*x + y*y; }
 double dist() const { return sqrt((double)
      dist2()); }
```

```
// angle to x-axis in interval [-pi, pi]
double angle() const { return atan2(y, x); }
P unit() const { return *this/dist(); } //
    makes dist()=1
P perp() const { return P(-y, x); } //
    rotates +90 degrees
P normal() const { return perp().unit(); }
// returns point rotated 'a' radians ccw
    around the origin
P rotate(double a) const {
    return P(x*cos(a)-y*sin(a),x*sin(a)+y*cos(a)); }
friend ostream& operator<<(ostream& os, P p)
    {
    return os << "(" << p.x << "," << p.y << "
)"; }
};</pre>
```

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. doublor 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 .dis on the result of the cross product.

```
p

s
"Point.h" f6bf6b, 4 lines

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

SegmentDistance.h

Description:

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

SegmentIntersection.h

Description:

If a unique intersection point between the line segments gdingescription: Returns where p is as seen from s towards from s1 to e1 and from s2 to e2 exists then it is returned. 1/0/-1 ⇔ left/on line/right. If the optional argument If no intersection point exists an empty vector is returned by is given 0 is returned if p is within distance eps from If infinitely many exist a vector with 2 elements is returned line. P is supposed to be Point<T> where T is e.g. containing the endpoints of the common line segment. Throuble or long long. It uses products in intermediate wrong position will be returned if P is Point<11> and the isteps so watch out for overflow if using int or long long. tersection point does not have integer coordinates. Productisage: bool left = sideOf(p1,p2,q) ==1; of three coordinates are used in intermediate steps so watchoint.h" out for overflow if using int or long long.

vector<P> inter =



Usage:

```
segInter(s1,e1,s2,e2);
if (sz(inter) == 1)
cout << "segments intersect at " << inter[0]</pre>
<< endl;
"Point.h", "OnSegment.h"
template < class P > vector < P > segInter (P a, P b,
     P c, P d) {
  auto oa = c.cross(d, a), ob = c.cross(d, b),
       oc = a.cross(b, c), od = a.cross(b, d);
  // Checks if intersection is single non-
       endpoint point.
  if (sgn(oa) * sgn(ob) < 0 && sgn(oc) * sgn(
      od) < 0
   return { (a * ob - b * oa) / (ob - oa) };
  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 streed Point double P; and s2,e2 exists {1, point} is returned. If no intersection point linear Transformation (g exists $\{0, (0,0)\}$ is returned and if infinitely many exists $\{-1, -1, -1\}$ (0,0)} is returned. The wrong position will be returned if P is Point<||> and the intersection point does not have integer p dp = p1-p0, dq = q1-q0, num(dp.cross(dq), coordinates. Products of three coordinates are used in intermediate steps so watch out for overflow if using int or ll.



```
Usage: auto res = lineInter(s1,e1,s2,e2);
if (res.first == 1)
cout << "intersection point at " <<
res.second << endl;
"Point.h"
                                      a01f81, 8 lines
```

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

sideOf.h

```
template<class P>
int sideOf(P s, P e, P p) { return sgn(s.cross
     (e, p)); }
template<class P>
int sideOf(const P& s, const P& e, const P& p,
     double eps) {
  auto a = (e-s).cross(p-s);
  double l = (e-s).dist()*eps;
  return (a > 1) - (a < -1);
```

OnSegment.h

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

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

linearTransformation.h

Description:

Apply the linear transformation (translation, rotation and if (a.t180() < r) r.t--; spalin, which takes line p0-p1 to line q0-q1 to point r. res

```
03a306, 6 lines
linearTransformation(const P& p0, const P&
  const P& q0, const P& q1, const P& r) {
     dp.dot(dq));
return q0 + P((r-p0).cross(num), (r-p0).dot(
```

num))/dp.dist2();

Angle.h

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

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

```
struct Angle {
 int x, y;
 int t;
 Angle(int x, int y, int t=0) : x(x), y(y), t
```

```
Angle operator-(Angle b) const { return {x-b
       .x, y-b.y, t}; }
  int half() const {
   assert(x || y);
    return y < 0 || (y == 0 && x < 0);
  Angle t90() const { return {-y, x, t + (half
       () && x >= 0)}; }
  Angle t180() const { return \{-x, -y, t +
      half()}; }
 Angle t360() const { return {x, y, t + 1}; }
bool operator<(Angle a, Angle b) {</pre>
  // add a. dist2() and b. dist2() to also
      compare distances
  return make_tuple(a.t, a.half(), a.y * (11)b
         make_tuple(b.t, b.half(), a.x * (11)b
              .y);
// Given two points, this calculates the
     smallest angle between
// them, i.e., the angle that covers the
     defined line segment.
pair<Angle, Angle> segmentAngles(Angle a,
    Angle b) {
  if (b < a) swap(a, b);
  return (b < a.t180() ?
          make_pair(a, b) : make_pair(b, a.
              t360()));
Angle operator+(Angle a, Angle b) { // point a
     + vector b
  Angle r(a.x + b.x, a.y + b.y, a.t);
  return r.t180() < a ? r.t360() : r;
Angle angleDiff(Angle a, Angle b) \{ // angle b \}
     - angle a
  int tu = b.t - a.t; a.t = b.t;
  return {a.x*b.x + a.y*b.y, a.x*b.y - a.y*b.x
      , tu - (b < a);
```

8.2 Circles

CircleIntersection.h

Description: Computes the pair of points at which two circles intersect. Returns false in case of no intersection. 84d6d3, 11 lines

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

CircleTangents.h

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

```
"Point.h"
                                    b0153d, 13 lines
template<class P>
vector<pair<P, P>> tangents(P c1, double r1, P
     c2, double r2) {
  P d = c2 - c1;
  double dr = r1 - r2, d2 = d.dist2(), h2 = d2
       - dr * dr:
  if (d2 == 0 || h2 < 0) return {};</pre>
  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, g) atan2(p.cross(g), p.dot(g))
double circlePoly(P c, double r, vector<P> ps)
 auto tri = [&](P p, P q) {
   auto r2 = r * r / 2;
   Pd = 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, g) * r2;</pre>
   auto s = max(0., -a-sqrt(det)), t = min
         (1., -a+sgrt(det));
   if (t < 0 \mid | 1 \le s) return arg(p, q) * r2
   P u = p + d * s, v = p + d * t;
   return arg(p,u) * r2 + u.cross(v)/2 + arg(
        v,q) * r2;
 auto sum = 0.0;
 rep(i, 0, sz(ps))
   sum += tri(ps[i] - c, ps[(i + 1) % sz(ps)]
         - c);
 return sum;
```

circumcircle.h

Description:

The circumcirle of a triangle is the circle intersecting all three vertices. ccRadius returns the radius of the circle gding through points A, B and C and ccCenter returns the center - ((a.y<p[i].y) - (a.y<q.y)) * a. of the same circle.



1caa3a, 9 lines

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

Minimum Enclosing Circle.h

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

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

```
"circumcircle.h"
                                    09dd0a, 17 lines
pair<P, double> mec(vector<P> ps) {
  shuffle(all(ps), mt19937(time(0)));
  P \circ = ps[0];
  double r = 0, EPS = 1 + 1e-8;
  rep(i, 0, sz(ps)) if ((o - ps[i]).dist() > r *
       EPS) {
    o = ps[i], r = 0;
    rep(j,0,i) if ((o - ps[j]).dist() > r *
         EPS) {
      o = (ps[i] + ps[j]) / 2;
      r = (o - ps[i]).dist();
      rep(k, 0, j) if ((o - ps[k]).dist() > r *
           EPS) {
        o = ccCenter(ps[i], ps[j], ps[k]);
        r = (o - ps[i]).dist();
   }
  return {o, r};
```

8.3 Polygons

InsidePolygon.h

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

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

"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) \le eps)
      return !strict;
      cross(p[i], q) > 0;
return cnt;
```

PolygonArea.h

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

```
f12300, 6 lines
template<class T>
T polygonArea2(vector<Point<T>>& v) {
 T = v.back().cross(v[0]);
  rep(i,0,sz(v)-1) a += v[i].cross(v[i+1]);
  return a;
```

PolygonCenter.h

"Point.h"

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

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

PolygonCut.h

Description:

Returns 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));
                                       f2b7d4, 13 lines
"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;</pre>
    if (side != (s.cross(e, prev) < 0))</pre>
      res.push_back(lineInter(s, e, cur, prev)
           .second);
    if (side)
      res.push_back(cur);
  return res;
```

ConvexHull.h

```
Returns a vector of the points of the convex hull in counter-return sqn(1[a].cross(1[b], p)) < r;
clockwise order. Points on the edge of the hull between two
other points are not considered part of the hull.
Time: O(n \log n)
"Point.h"
                                       310954, 13 lines
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 \geq s + 2 && h[t-2].cross(h[t
           -1], p) <= 0) t--;
```

HullDiameter.h

9706dc, 9 lines

h[t++] = p;

&& h[0] == h[1]);

Description: Returns the two points with max distance on a convex hull (ccw, no duplicate/collinear points). Time: $\mathcal{O}(n)$ "Point.h" c571b8, 12 lines

return {h.begin(), h.begin() + t - (t == 2

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

return res.second;

PointInsideHull.h

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

71446b, 14 lines

```
Time: \mathcal{O}(\log N)
"Point.h", "sideOf.h", "OnSegment.h"
```

```
typedef Point<11> P;
bool inHull(const vector<P>& 1, P p, bool
    strict = true) {
  int a = 1, b = sz(1) - 1, r = !strict;
 if (sz(1) < 3) return r && onSegment(1[0], 1</pre>
       .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;

```
LineHullIntersection.h
```

Description: Line-convex polygon intersection. The polygon must be ccw and have no collinear points. line-Hull(line, poly) returns a pair describing the intersection of a line with the polygon: \bullet (-1,-1) if no collision, • (i,-1) if touching the corner i, • (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"
                                     7cf45b, 39 lines
#define cmp(i,j) sgn(dir.perp().cross(poly[(i)
    %n]-poly[(j)%n]))
#define extr(i) cmp(i + 1, i) >= 0 \&\& cmp(i, i)
     -1 + n) < 0
template <class P> int extrVertex(vector<P>&
    poly, P dir) {
 int n = sz(poly), lo = 0, hi = n;
 if (extr(0)) return 0;
  while (lo + 1 < hi) {
   int m = (lo + hi) / 2;
   if (extr(m)) return m;
    int 1s = cmp(1o + 1, 1o), ms = cmp(m + 1,
    (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>
```

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

array<int, 2> lineHull(P a, P b, vector<P>&

int endA = extrVertex(poly, (a - b).perp());

polv) {

swap(endA, endB);

return res;

```
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]};
```

if (res[0] == res[1]) return {res[0], -1};

ClosestPair kdTree FastDelaunay PolyhedronVolume Point3D

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);
  set<P> S;
  sort(all(v), [](P a, P b) { return a.y < b.y</pre>
  pair<ll, pair<P, P>> ret{LLONG MAX, {P(), P
       () } };
  int j = 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,}
    S.insert(p):
  return ret.second;
```

```
kdTree.h
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: 1
bool on_y (const P& a, const P& b) { return a.y
      < b.v; }
struct Node {
  P pt; // if this is a leaf, the single point
       in it
  T x0 = INF, x1 = -INF, y0 = INF, y1 = -INF;
       // bounds
  Node *first = 0, *second = 0;
  T distance (const P& p) { // min squared
       distance to a point
    T x = (p.x < x0 ? x0 : p.x > x1 ? x1 : p.x
    T y = (p.y < y0 ? y0 : p.y > y1 ? y1 : p.y
    return (P(x,y) - p).dist2();
  Node (vector<P>&& vp) : pt(vp[0]) {
    for (P p : vp) {
      x0 = min(x0, p.x); x1 = max(x1, p.x);
      y0 = min(y0, p.y); y1 = max(y1, p.y);
    if (vp.size() > 1) {
      // split on x if width >= height (not
           ideal...)
      sort(all(vp), x1 - x0 >= y1 - y0 ? on_x
```

: on_y);

```
// divide by taking half the array for
           each child (not
      // best performance with many duplicates
            in the middle)
      int half = sz(vp)/2;
      first = new Node({vp.begin(), vp.begin()
            + half});
      second = new Node({vp.begin() + half, vp
           .end()});
};
struct KDTree {
  Node* root;
  KDTree (const vector < P > & vp) : root (new Node (
       {all(vp)})) {}
  pair<T, P> search(Node *node, const P& p) {
    if (!node->first) {
      // uncomment if we should not find the
           point itself:
      // if (p = node \rightarrow pt) return \{INF, P()\};
      return make_pair((p - node->pt).dist2(),
            node->pt);
    Node *f = node->first, *s = node->second;
    T bfirst = f->distance(p), bsec = s->
         distance(p);
    if (bfirst > bsec) swap(bsec, bfirst),
         swap(f, s);
    // search closest side first, other side
         if needed
    auto best = search(f, p);
    if (bsec < best.first)</pre>
      best = min(best, search(s, p));
    return best;
  // find nearest point to a point, and its
       squared distance
  // (requires an arbitrary operator< for
       Point)
  pair<T, P> nearest (const P& p) {
    return search(root, p);
};
FastDelaunav.h
Description: Fast Delaunay triangulation. Each cir-
```

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

```
P& F() { return r()->p; }
  O& 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->rot, r->p = arb, r->o = i
      & 1 ? r : r->r();
  r->p = orig; r->F() = dest;
 return r;
void splice(Q a, Q b) {
 swap(a->o->rot->o, b->o->rot->o); swap(a->o,
       b->o);
Q connect (Q a, Q b) {
 O g = makeEdge(a->F(), b->p);
  splice(q, a->next());
  splice(q->r(), b);
  return q;
pair<Q,Q> rec(const vector<P>& s) {
 if (sz(s) <= 3) {
    Q = makeEdge(s[0], s[1]), b = makeEdge(s[0], s[1])
         [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)
 Q A, B, ra, rb;
 int half = sz(s) / 2;
 tie(ra, A) = rec({all(s) - half});
  tie(B, rb) = rec({sz(s) - half + all(s)});
  while ((B->p.cross(H(A)) < 0 && (A = A->next)
       ())) ||
         (A->p.cross(H(B)) > 0 && (B = B->r()
              ->0)));
  O base = connect(B->r(), A);
 if (A->p == ra->p) ra = base->r();
  if (B->p == rb->p) rb = base;
#define DEL(e, init, dir) O 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());
   6156
     base = connect(base->r(), LC->r());
 return { ra, rb };
vector<P> triangulate(vector<P> pts) {
 sort(all(pts)); assert(unique(all(pts)) ==
      pts.end());
 if (sz(pts) < 2) return {};
 Q e = rec(pts).first;
 vector<Q> q = \{e\};
 int \alpha i = 0:
  while (e->o->F().cross(e->F(), e->p) < 0) e
      = e->o;
#define ADD { Q c = e; do { c->mark = 1; pts.
    push_back(c->p); \
 g.push_back(c->r()); c = c->next(); } while
      (c != e); }
 ADD; pts.clear();
 while (qi < sz(q)) if (!(e = q[qi++])->mark)
 return pts;
```

8.5 3D

PolyhedronVolume.h

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

3058c3, 6 lines

Point3D.h

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

8058ae, 32 lines

```
template < class T > struct Point3D {
  typedef Point3D P:
  typedef const P& R;
  T x, y, z;
  explicit Point3D (T x=0, T y=0, T z=0) : x(x)
       , y(y), z(z) \{ \}
  bool operator<(R p) const {</pre>
    return tie(x, y, z) < tie(p.x, p.y, p.z);
  \textbf{bool operator} {==} \; (\texttt{R} \;\; \texttt{p}) \;\; \textbf{const} \;\; \{
    return tie(x, y, z) == tie(p.x, p.y, p.z);
  P operator+(R p) const { return P(x+p.x, y+p
       .y, z+p.z); }
  P operator-(R p) const { return P(x-p.x, y-p
       .y, z-p.z); }
  P operator*(T d) const { return P(x*d, y*d,
       z*d); }
  P operator/(T d) const { return P(x/d, y/d,
       z/d); }
```

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

3dHull.h

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

```
Time: \mathcal{O}\left(n^2\right)
"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 l) {
   P3 q = (A[j] - A[i]).cross((A[k] - A[i]));
   if (q.dot(A[1]) > q.dot(A[i]))
     q = q * -1;
    F f{q, i, j, k};
    E(a,b).ins(k); E(a,c).ins(j); E(b,c).ins(i)
        );
   FS.push_back(f);
  rep(i,0,4) rep(j,i+1,4) rep(k,j+1,4)
   mf(i, j, k, 6 - i - j - k);
  rep(i, 4, sz(A)) {
```

rep(j,0,sz(FS)) {

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

sphericalDistance.h

Description: Returns the shortest distance on the sphere with radius radius between the points with azimuthal angles (longitude) f1 (ϕ_1) and f2 (ϕ_2) from x axis and zenith angles (latitude) t1 (θ_1) and t2 (θ_2) from z axis (0 = north pole). All angles measured in radians. The algorithm starts by converting the spherical coordinates to cartesian coordinates so if that is what you have you can use only the two last rows. dx*radius is then the difference between the two points in the x direction and d*radius is the total distance between the points. 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 = sgrt(dx*dx + dy*dy + dz*dz);
  return radius *2 *asin(d/2);
```

Strings (9)

KMP.h

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[x] 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=0,1=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: $\mathcal{O}(N)$

```
d07a42, 8 lines
int minRotation(string s) {
  int a=0, N=sz(s); s += s;
  rep(b, 0, N) rep(k, 0, N) {
    if (a+k == b | | s[a+k] < s[b+k]) {b += max}
         (0, k-1); break;}
    if (s[a+k] > s[b+k]) { a = b; break; }
  return a;
```

SuffixArray.h

Description: Builds suffix array for a string. sa[i] is the starting index of the suffix which is i'th in the sorted suffix array. The returned vector is of size n + 1, and sa[0] = n. The 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.

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

38db9f, 23 lines

```
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)+1), y(n), ws(max(n, lim)),
    sa = lcp = v, iota(all(sa), 0);
    for (int j = 0, p = 0; p < n; j = max(1, j)
         * 2), lim = p) {
      p = j, iota(all(y), n - j);
      rep(i, 0, n) if (sa[i] >= j) y[p++] = sa[i]
          ] - j;
      fill(all(ws), 0);
      rep(i, 0, n) ws[x[i]] ++;
      rep(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;
      rep(i,1,n) = sa[i-1], b = sa[i], x[b]
        (y[a] == y[b] && y[a + j] == y[b + j])
              ? p - 1 : p++;
    rep(i,1,n) rank[sa[i]] = i;
    for (int i = 0, j; i < n - 1; lcp[rank[i</pre>
         ++]] = k)
      for (k \& \& k--, j = sa[rank[i] - 1];
         s[i + k] == s[j + k]; k++);
};
```

SuffixTree.h

 $\textbf{Description:} \ \ \textbf{Ukkonen's algorithm for online suffix tree}$ 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).

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], l[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]; qoto 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])
      l[v]=q; p[v]=m; t[p[m]][toi(a[l[m]])]=
          m;
      v=s[p[m]]; q=l[m];
      while (q<r[m]) { v=t[v][toi(a[q])]; q+=</pre>
          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)
  int lcs(int node, int i1, int i2, int olen)
    if (l[node] <= i1 && i1 < r[node]) return</pre>
    if (1[node] <= i2 && i2 < r[node]) return</pre>
    int mask = 0, len = node ? olen + (r[node]
         - 1[node]) : 0;
    rep(c, 0, ALPHA) if (t[node][c] != -1)
     mask |= 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 hash-

```
2d2a67, 44 lines
// Arithmetic mod 2^64-1. 2x slower than mod 2
     ^64 and more
// code, but works on evil test data (e.g.
    Thue-Morse, where
// ABBA... and BAAB... of length 2^10 hash the
     same mod 2^64).
// "typedef ull H;" instead if you think test
     data is random,
// or work mod 10^9+7 if the Birthday paradox
    is not a problem.
typedef uint64_t ull;
struct H {
  ull x; H(ull x=0) : x(x) {}
  H operator+(H \circ) { return x + \circ.x + (x + \circ.x
       < x); }
  H operator-(H o) { return *this + ~o.x; }
  H operator*(H o) { auto m = (__uint128_t)x *
```

```
return H((ull)m) + (ull)(m >> 64); }
  ull get() const { return x + !~x; }
  bool operator==(H o) const { return get() ==
        o.get(); }
  bool operator<(H o) const { return get() < o</pre>
       .qet(); }
static const H C = (11)1e11+3; // (order \sim 3e9)
     ; random also ok)
struct HashInterval {
  vector<H> ha, pw;
  HashInterval(string& str) : ha(sz(str)+1),
      pw(ha) {
    pw[0] = 1;
    rep(i, 0, sz(str))
      ha[i+1] = ha[i] * C + str[i],
      pw[i+1] = pw[i] * C;
  H hashInterval(int a, int b) { // hash [a, b]
    return ha[b] - ha[a] * pw[b - a];
};
vector<H> getHashes(string& str, int length) {
 if (sz(str) < length) return {};</pre>
 H h = 0, pw = 1;
  rep(i,0,length)
  h = h * C + str[i], pw = pw * C;
  vector<H> ret = {h};
  rep(i,length,sz(str)) {
    ret.push\_back(h = h * C + str[i] - pw *
        str[i-length]);
  return ret;
H hashString(string& s) {H h{}; for(char c:s) h
    =h*C+c; return h; }
```

AhoCorasick.h

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

Time: construction takes $\mathcal{O}(26N)$, where N = sum oflength of patterns. find(x) is $\mathcal{O}(N)$, where N = length of x. findAll is $\mathcal{O}(NM)$. f35677, 66 lines

```
struct AhoCorasick {
  enum {alpha = 26, first = 'A'}; // change
      this!
  struct Node {
    // (nmatches is optional)
    int back, next[alpha], start = -1, end =
        -1, nmatches = 0;
    Node (int v) { memset (next, v, sizeof (next)
        ); }
  };
  vector<Node> N;
  vi backp;
```

```
void insert(string& s, int j) {
   assert(!s.empty());
   int n = 0;
   for (char c : s) {
     int& m = N[n].next[c - first];
     if (m == -1) { n = m = sz(N); N.
          emplace_back(-1); }
     else n = m;
   if (N[n].end == -1) N[n].start = i;
   backp.push_back(N[n].end);
   N[n].end = j;
   N[n].nmatches++;
 AhoCorasick(vector<string>& pat) : N(1, -1)
   rep(i,0,sz(pat)) insert(pat[i], i);
   N[0].back = sz(N);
   N.emplace_back(0);
   queue<int> q;
   for (q.push(0); !q.empty(); q.pop()) {
     int n = q.front(), prev = N[n].back;
     rep(i,0,alpha) {
       int &ed = N[n].next[i], y = N[prev].
            next[i];
       if (ed == -1) ed = y;
       else {
         N[ed].back = y;
          (N[ed].end == -1 ? N[ed].end : backp
              [N[ed].start])
           = N[y].end;
         N[ed].nmatches += N[y].nmatches;
         q.push(ed);
 vi find(string word) {
   int n = 0:
   vi res; // ll count = 0;
   for (char c : word) {
     n = N[n].next[c - first];
     res.push back(N[n].end);
     // count += 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 =
  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]; });
 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]));
      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.
              constantIntervals(0, sz(v), [&](int
x) {return v[x];}, [&] (int lo, int hi, T
val) {...};
Time: \mathcal{O}\left(k\log\frac{n}{k}\right)
```

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

LIS.h

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 \rightarrow i for longest non-
         decreasing subsequence
```

```
auto it = lower_bound(all(res), p{S[i], 0}
      );
 if (it == res.end()) res.emplace back(),
      it = res.end()-1;
  *it = {S[i], i};
 prev[i] = it == res.begin() ? 0 : (it-1) ->
int L = sz(res), cur = res.back().second;
while (L--) ans[L] = cur, cur = prev[cur];
return ans;
```

FastKnapsack.h

Description: Given N non-negative integer weights w and a non-negative target t, computes the maximum S <= t such that S is the sum of some subset of the weights. Time: $\mathcal{O}(N \max(w_i))$ 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)) {
   u = v;
   rep(x,0,m) v[x+w[i]] = max(v[x+w[i]], u[x
   for (x = 2*m; --x > m;) rep(j, max(0,u[x])
     v[x-w[j]] = max(v[x-w[j]], j);
 for (a = t; v[a+m-t] < 0; a--);
 return a;
```

10.3 Dynamic programming

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

DivideAndConquerDP.h

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

```
Time: \mathcal{O}((N + (hi - lo)) \log N)
                                      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); }
};
```

Debugging tricks

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

10.5 Optimization tricks

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

10.5.1 Bit hacks

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

10.5.2 **Pragmas**

• #pragma GCC optimize ("Ofast") will make GCC auto-vectorize loops and optimizes floating points better.

- #pragma GCC target ("avx2") can double performance of vectorized code, but causes crashes on old machines.
- #pragma GCC optimize ("trapv") kills the program on integer overflows (but is really

FastMod.h

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

```
typedef unsigned long long ull;
                                            struct FastMod {
                                             ull b, m;
                                             FastMod(ull b) : b(b), m(-1ULL / b) {}
signal(SIGSEGV, [](int) { _Exit(0); ^{ull}; ^{reduce(ull\ a)} { // a % b + (0 or b)
                                               réturn a - (ull) ((__uint128_t(m) * a) >>
                                                    64) * b;
                                            };
```

FastInput.h

Description: Read an integer from stdin. Usage requires your program to pipe in input from file. Usage: ./a.out < input.txt</pre>

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

```
7b3c70, 17 lines
inline char gc() { // like getchar()
  static char buf[1 << 16];</pre>
 static size_t bc, be;
 if (bc >= be) {
   buf[0] = 0, bc = 0;
   be = fread(buf, 1, sizeof(buf), stdin);
 return buf[bc++]; // returns 0 on EOF
int readInt() {
 int a, c;
 while ((a = qc()) < 40);
 if (a == '-') return -readInt();
 while ((c = qc()) >= 48) a = a * 10 + c -
      480;
 return a - 48;
```

BumpAllocator.h

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

```
// Either globally or in a single class:
static char buf[450 << 201;
void* operator new(size t s)
 static size t i = sizeof buf;
void operator delete(void*) {}
```

SmallPtr.h

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

"BumpAllocator.h" 2dd6c9, 10 lines

```
template<class T> struct ptr {
  unsigned ind;
  ptr(T*p = 0) : ind(p ? unsigned((char*)p -
      buf) : 0) {
    assert (ind < sizeof buf);
  T& operator*() const { return *(T*)(buf +
  T* operator->() const { return &**this; }
  T& operator[](int a) const { return (&**this
  explicit operator bool() const { return ind;
BumpAllocatorSTL.h
Description: BumpAllocator for STL containers.
Usage:
            vector<vector<int, small<int>>>
ed(N);
                                    bb66d4, 14 lines
char buf[450 << 20] alignas(16);</pre>
size t buf ind = sizeof buf;
template < class T > struct small {
  typedef T value_type;
  small() {}
  template < class U > small(const U&) {}
  T* allocate(size_t n) {
   buf_ind -= n * sizeof(T);
   buf_ind &= 0 - alignof(T);
   return (T*) (buf + buf_ind);
 void deallocate(T*, size_t) {}
};
SIMD.h
```

Description: Cheat sheet of SSE/AVX intrinsics, for doing arithmetic on several numbers at once. Can provide a constant factor improvement of about 4, orthogonal to loop unrolling. Operations follow the pattern "_mm(256)?_name_(si(128|256)|epi(8|16|32|64)|pd|ps)". Not all are described here; grep for _mm_ in /usr/lib/gcc/*/4.9/include/ for more. If AVX is unsupported, try 128-bit operations, "emmintrin.h" and #define __SSE_ and __MMX__ before including it. For aligned memory use _mm_malloc(size, 32) or int buf[N] alignas(32), but prefer loadu/stgrug, 43 lines

```
#pragma GCC target ("avx2") // or sse4.1
#include "immintrin.h"
typedef __m256i mi;
#define L(x) _mm256_loadu_si256((mi*)&(x))
// High-level/specific methods:
// load(u)?\_si256, store(u)?\_si256,
    setzero_si256, _mm_malloc
// blendv_{-}(epi8|ps|pd) (z?y:x), movemask_epi8
    (hibits of bytes)
// i32gather_epi32(addr, x, 4): map addr[]
    over 32-b parts of x
// sad_epu8: sum of absolute differences of u8
    , outputs 4xi64
// maddubs_epi16: dot product of unsigned i7's
    , outputs 16xi15
// madd_epi16: dot product of signed i16's,
    outputs 8xi32
// extractf128\_si256(, i) (256->128),
    cvtsi128_si32 (128->lo32)
```

```
// permute2f128\_si256(x,x,1) swaps 128-bit
// shuffle_epi32(x, 3*64+2*16+1*4+0) == x for
     each lane
// shuffle_epi8(x, y) takes a vector instead
     of an imm
// Methods that work with most data types (
     append\ e.g.\ \_epi32):
// set1, blend (i8?x:y), add, adds (sat.),
     mullo, sub, and/or,
// and not, abs, min, max, sign(1,x), cmp(gt | eq
     ), unpack(lo|hi)
int sumi32(mi m) { union {int v[8]; mi m;} u;
    u.m = m;
  int ret = 0; rep(i,0,8) ret += u.v[i];
      return ret; }
mi zero() { return _mm256_setzero_si256(); }
mi one() { return _mm256_set1_epi32(-1); }
bool all_zero(mi m) { return
    _mm256_testz_si256(m, m); }
bool all_one(mi m) { return _mm256_testc_si256
     (m, one()); }
11 example_filteredDotProduct(int n, short* a,
      short* b) {
  int i = 0; ll r = 0;
  mi zero = _mm256_setzero_si256(), acc = zero
  while (i + 16 <= n) {
    mi \ va = L(a[i]), \ vb = L(b[i]); \ i += 16;
    va = _mm256_and_si256(_mm256_cmpgt_epi16(
         vb, va), va);
    mi vp = _mm256_madd_epi16(va, vb);
    acc = _mm256_add_epi64(
         _mm256_unpacklo_epi32(vp, zero),
      _mm256_add_epi64(acc,
           mm256 unpackhi epi32(vp, zero)));
  union {11 v[4]; mi m;} u; u.m = acc; rep(i
      ,0,4) r += u.v[i];
  for (;i<n;++i) if (a[i] < b[i]) r += a[i]*b[</pre>
      il; // <- equiv
```

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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