1 Contest	1	7.3 Network flow	.vimrc 6 lines		
2 Mathematics	1	7.4 Matching			
	1	7.5 DFS algorithms	" Select region and then type : Hash to hash your selection.		
2.1 Equations		7.6 Heuristics	1 3 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1		
2.2 Recurrences	- 1	7.7 Trees	ca Hash w !cpp -dD -P -fpreprocessed \ tr -d '[:space:]' \ \ md5sum \ cut -c-6		
2.3 Trigonometry		o G			
2.4 Geometry		8 Geometry 18	11dS11,S11		
2.5 Derivatives/Integrals		8.1 Geometric primitives	# Hashes a file, ignoring all whitespace and comments. Use for		
2.6 Sums		8.2 Circles			
2.7 Series		8.3 Polygons			
2.8 Probability theory	3	8.4 Misc. Point Set Problems			
2.9 Markov chains	3	8.5 3D	Pre-submit:		
			Write a few simple test cases, if sample is not enough.		
3 Data structures	3	9 Strings 22	Are time limits close? If so, generate max cases. Is the memory usage fine?		
3.1 Matrix	3	9.1 KMP	Could anything overflow?		
3.2 Treap	4	9.2 Zfunc			
3.3 RMQ	4	9.3 Manacher	Wrong answer:		
3.4 WaveletTree		9.4 MinRotation	Print your solution! Print debug output, as well.		
3.5 KDTree		9.5 SuffixArray	Are you clearing all datastructures between test cases? Can your algorithm handle the whole range of input?		
3.6 PersistentSegTree		9.6 SuffixTree	Read the full problem statement again.		
5.0 Telaboration grant in the second	0	9.7 Hashing	Do you handle all corner cases correctly? Have you understood the problem correctly?		
4 Numerical	6	9.8 AhoCorasick			
4.1 GoldenSectionSearch	6		Any overflows?		
4.2 Polynomial		10 Various 25	Confusing N and M, i and j, etc.? Are you sure your algorithm works?		
4.3 LinearRecurrence		10.1 Intervals	What special cases have you not thought of?		
		10.2 Misc. algorithms	I are voll sure the STL tunctions voll use work as voll think?		
e e e e e e e e e e e e e e e e e e e		10.3 Dynamic programming	Create some testcases to run your algorithm on.		
4.5 Matrix		10.4 Debugging tricks			
4.6 SolveLinear		10.5 Optimization tricks	Explain your algorithm to a team mate.		
4.7 Fourier transforms	8	10.0 Optimization tricks 2-	Ask the team mate to look at your code.		
₩ MT 1 (1		$\underline{\text{Contest}}$ (1)	Go for a small walk, e.g. to the toilet. Is your output format correct? (including whitespace)		
5 Number theory	9	$\frac{\text{Comecs}}{\text{Composite}}$	Rewrite your solution from the start or let a team mate do it.		
5.1 Modular arithmetic		template.cpp 15 line	Runtime error:		
5.2 Primality		#include <bits stdc++.h=""></bits>	Have you tested all corner cases locally?		
5.3 Divisibility	- 1	using namespace std;	Any uninitialized variables? Are you reading or writing outside the range of any vector?		
5.4 Fractions	- 1	#define rep(i, a, b) for(int i = a; i < (b); ++i)	Any assertions that might fail?		
5.5 Pythagorean Triples		#define trav(a, x) for(auto& a : x)	Any possible division by 0? (mod 0 for example) Any possible infinite recursion?		
5.6 Primes	10	#define all(x) begin(x), end(x) #define sz(x) (int)(x).size()	Invalidated pointers or iterators?		
5.7 Estimates	11	typedef long long 11;	Are you using too much memory?		
		<pre>typedef pair<int, int=""> pii; typedef vector<int> vi;</int></int,></pre>	Debug with resubmits (e.g. remapped signals, see Various).		
6 Combinatorial	11	typeder vector(int) vi,	Time limit exceeded:		
6.1 Permutations		int main() {	Do you have any possible infinite loops? What is the complexity of your algorithm?		
6.2 Partitions and subsets	11	<pre>cin.sync_with_stdio(0); cin.tie(0); cin.exceptions(cin.failbit);</pre>	Are you copying a lot of unnecessary data? (References)		
6.3 General purpose numbers	11	}	How big is the input and output? (consider scanf) Avoid vector, map. (use arrays/unordered_map)		
• •		11	What do your team mates think about your algorithm?		
7 Graph	12	.bashrc 3 line	Memory limit exceeded:		
7.1 Fundamentals	12	alias c='g++ -Wall -Wconversion -Wfatal-errors -g -std=c++14 \ -fsanitize=undefined,address'	What is the max amount of memory your algorithm should need?		
7.2 Euler walk	12	exmodmap -e 'clear lock' -e 'keycode 66=less greater' $\#caps = \bigcirc$	Are you clearing all datastructures between test cases?		

Mathematics (2)

2.1 Equations

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

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

$$ax + by = e$$

$$cx + dy = f$$

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

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

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.

Recurrences

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

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

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

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: $a^2 = b^2 + c^2 - 2bc \cos \alpha$

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

2.4.2 Quadrilaterals

With side lengths a, b, c, d, diagonals e, f, diagonals angle 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



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

Derivatives/Integrals

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

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

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

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

Integration by parts:

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

2.6Sums

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

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

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

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

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

2.7Series

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

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

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

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

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

2.8 Probability theory

Let X be a discrete random variable with probability $p_X(x)$ of assuming the value x. It will then have an expected value (mean) $\mu = \mathbb{E}(X) = \sum_x x p_X(x)$ and variance $\sigma^2 = V(X) = \mathbb{E}(X^2) - (\mathbb{E}(X))^2 = \sum_x (x - \mathbb{E}(X))^2 p_X(x)$ where 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 $Bin(n, p), n = 1, 2, ..., 0 \le p \le 1$.

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

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

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

First success distribution

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

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

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

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 $\text{Po}(\lambda), \ \lambda = t\kappa$.

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

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

2.8.2 Continuous distributions

Uniform distribution

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

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

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

Exponential distribution

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

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

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

Normal distribution

Most real random values with mean 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.

3

c43c7d, 26 lines

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 **A** and **G**, such that all states in **A** are absorbing $(p_{ii} = 1)$, and all states in **G** leads to an absorbing state in **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)

3.1 Matrix

p >>= 1;

return a;

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;
```

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;

Treap FenwickTree2d RMQ WaveletTree KDTree

Description: Computes sums a[i,j] for all i < I, j < J, and increases single ele-

ments a[i,j]. Requires that the elements to be updated are known in advance

Treap

Treap.h

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

```
Time: \mathcal{O}(\log N)
                                                      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};
    auto pa = split(n->r, k - cnt(n->1) - 1); // and just "k"
    n->r = pa.first;
   n->recalc();
    return {n, pa.second};
Node* merge(Node* 1, Node* r) {
  if (!1) return r;
  if (!r) return 1;
  if (1->y > r->y) {
   1->r = merge(1->r, r);
   1->recalc();
    return 1;
  } else {
    r->1 = merge(1, r->1);
    r->recalc();
    return r;
Node* ins(Node* t, Node* n, int pos) {
  auto pa = split(t, pos);
  return merge(merge(pa.first, n), pa.second);
// Example application: move the range [l, r) to index k
void move(Node*& t, int 1, int r, int k) {
 Node *a, *b, *c;
  tie(a,b) = split(t, 1); tie(b,c) = split(b, r - 1);
 if (k \le 1) t = merge(ins(a, b, k), c);
```

else t = merge(a, ins(c, b, k - r));

```
(call fakeUpdate() before init()).
Time: \mathcal{O}(\log^2 N). (Use persistent segment trees for \mathcal{O}(\log N).)
"FenwickTree.h"
struct FT2 {
 vector<vi> ys; vector<FT> ft;
 FT2(int limx) : ys(limx) {}
 void fakeUpdate(int x, int y) {
    for (; x < sz(ys); x = x + 1) ys[x].push_back(y);
   trav(v, vs) sort(all(v)), ft.emplace back(sz(v));
 int ind(int x, int y) {
    return (int) (lower_bound(all(ys[x]), y) - ys[x].begin()); }
 void update(int x, int y, ll dif) {
    for (; x < sz(vs); 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;
};
```

3.3 RMQ

FenwickTree2d.h

RMQ.h

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

Time: $\mathcal{O}\left(|V|\log|V|+Q\right)$

1f8996, 17 lines

```
template < class T>
struct RMO {
 vector<vector<T>> jmp;
 RMQ(const vector<T>& V) {
   int N = sz(V), on = 1, depth = 1;
   while (on < N) on \star= 2, depth++;
    jmp.assign(depth, V);
   rep(i, 0, depth-1) rep(i, 0, N)
     jmp[i+1][j] = min(jmp[i][j],
      jmp[i][min(N - 1, j + (1 << i))]);
 T query(int a, int b) {
   assert (a < b); // or return inf if a == b
   int dep = 31 - __builtin_clz(b - a);
   return min(jmp[dep][a], jmp[dep][b - (1 << dep)]);</pre>
};
```

3.4 WaveletTree

Wavelet Tree.h

1cd45c, 55 lines

```
template <class T>
struct wavelet{
 struct node
   vector<int> b;
   T lo, hi, md;
 vector<node> t;
```

```
void build(const vector<T> &c, T *A, T *B, int v, int i, int
    t[v].b.resize(B-A+1);
   t[v].lo = c[i], t[v].hi = c[j], t[v].md = c[(i+j)/2];
    for(int i = 0; A+i != B; ++i)
     t[v].b[i+1] = t[v].b[i] + (A[i] \le t[v].md);
    if(i == j) return;
    T *p = stable\_partition(A,B,[=](int x){return x <= t[v].md}
   build(c, A, p, v <<1, i, (i+j)/2);
   build(c,p,B,v << 1 | 1, (i+j)/2+1, j);
 void init(T *A, int n) {
   vector<T> c(A,A+n);
   sort(c.begin(),c.end());
   c.erase(unique(c.begin(),c.end()),c.end());
    int N = c.size();
   t.resize(N<<2);
   build(c, A, A+n, 1, 0, N-1);
 wavelet(){}
 wavelet(T *A, int n) {init(A, n);}
 //kth smallest element in [l. r]
 T kth(int 1, int r, int k, int v = 1) {
   if(t[v].lo == t[v].hi) return t[v].lo;
    int lb = t[v].b[l-1], rb = t[v].b[r], il = rb-lb;
    return (k < i1) ? kth(lb+1,rb,k,v<<1) : kth(l-lb,r-rb,k-i1,
         v << 1 | 1):
  //number of elements in [l,r] \leq to a
 int leg(int 1, int r, T a, int v = 1) {
   if(a < t[v].lo) return 0;</pre>
   if(t[v].hi <= a) return r-1+1;</pre>
   int lb = t[v].b[l-1], rb = t[v].b[r];
    return leq(lb+1, rb, a, v<<1) + leq(l-lb, r-rb, a, v<<1|1);</pre>
  //number of elements in [l, r] equal to a
 int count(int 1, int r, T a, int v = 1) {
   if(a < t[v].lo || a > t[v].hi) return 0;
   if(t[v].lo == t[v].hi) return r-l+1;
   int lb = t[v].b[1-1], rb = t[v].b[r];
   if(a <= t[v].md) return count(lb+1,rb,a,v<<1);</pre>
   return count (1-1b, r-rb, a, v<<1|1);
};
```

KDTree

KDTree.h

```
<br/>
<br/>
dits/stdc++.h>
                                                       ff695b, 182 lines
// constructs from n points in O(n lq^2 n) time
// handles nearest-neighbor query in O(lq n) if points are well
      distributed
// O(\log n) average nearest neighbor, O(n) worst in
     pathological case
using namespace std;
// number type for coordinates, and its maximum value
typedef long long ntype;
```

const ntype sentry = numeric_limits<ntype>::max();

```
// point structure for 2D-tree, can be extended to 3D
struct point {
   ntype x, y;
    point(ntype xx = 0, ntype yy = 0) : x(xx), y(yy) {}
bool operator == (const point &a, const point &b) {
    return a.x == b.x && a.y == b.y;
// sorts points on x-coordinate
bool on x(const point &a, const point &b) {
    return a.x < b.x;</pre>
// sorts points on y-coordinate
bool on_y(const point &a, const point &b) {
    return a.y < b.y;</pre>
// squared distance between points
ntype pdist2 (const point &a, const point &b) {
    ntype dx = a.x-b.x, dy = a.y-b.y;
    return dx*dx + dy*dy;
// bounding box for a set of points
struct bbox {
    ntype x0, x1, y0, y1;
   bbox(): x0(sentry), x1(-sentry), y0(sentry), y1(-sentry)
        { }
    // computes bounding box from a bunch of points
    void compute(const vector<point> &v) {
        for (int i = 0; i < v.size(); ++i) {</pre>
           x0 = min(x0, v[i].x); x1 = max(x1, v[i].x);
            y0 = min(y0, v[i].y); y1 = max(y1, v[i].y);
    // squared distance between a point and this bbox, 0 if
         inside
   ntype distance (const point &p) {
       if (p.x < x0) {
           if (p.y < y0)
                                return pdist2(point(x0, y0), p)
            else if (p.y > y1) return pdist2(point(x0, y1), p)
            else
                                return pdist2(point(x0, p.y), p
                );
        else if (p.x > x1) {
                                return pdist2(point(x1, y0), p)
           if (p.y < y0)
            else if (p.y > y1) return pdist2(point(x1, y1), p)
            else
                                return pdist2(point(x1, p.y), p
                );
        else {
            if (p.y < y0)
                                return pdist2(point(p.x, y0), p
            else if (p.y > y1) return pdist2(point(p.x, y1), p
                );
            else
                                return 0;
```

```
// stores a single node of the kd-tree, either internal or leaf
struct kdnode {
                    // true if this is a leaf node (has one
   bool leaf;
        point)
    point pt;
                    // the single point of this is a leaf
    bbox bound;
                    // bounding box for set of points in
         children
    kdnode *first, *second; // two children of this kd-node
    kdnode() : leaf(false), first(0), second(0) {}
   ~kdnode() { if (first) delete first; if (second) delete
         second; }
    // intersect a point with this node (returns squared
    ntype intersect (const point &p) {
        return bound.distance(p);
    // recursively builds a kd-tree from a given cloud of
         points
    void construct(vector<point> &vp) {
        // compute bounding box for points at this node
        bound.compute(vp);
        // if we're down to one point, then we're a leaf node
       if (vp.size() == 1) {
            leaf = true;
            pt = vp[0];
       else
            // split on x if the bbox is wider than high (not
                 best heuristic...)
            if (bound.x1-bound.x0 >= bound.y1-bound.y0)
                sort(vp.begin(), vp.end(), on_x);
            // otherwise split on y-coordinate
            else
                sort(vp.begin(), vp.end(), on_y);
            // divide by taking half the array for each child
            // (not best performance if many duplicates in the
                 middle)
            int half = vp.size()/2;
            vector<point> vl(vp.begin(), vp.begin()+half);
            vector<point> vr(vp.begin()+half, vp.end());
            first = new kdnode(); first->construct(v1);
            second = new kdnode(); second->construct(vr);
// simple kd-tree class to hold the tree and handle queries
struct kdtree {
   kdnode *root:
    // constructs a kd-tree from a points (copied here, as it
         sorts them)
    kdtree(const vector<point> &vp) {
       vector<point> v(vp.begin(), vp.end());
       root = new kdnode();
       root->construct(v);
    ~kdtree() { delete root; }
    // recursive search method returns squared distance to
```

nearest point

```
ntype search (kdnode *node, const point &p) {
        if (node->leaf) {
            // commented special case tells a point not to find
               if (p = node > pt) return sentry;
               else
                return pdist2(p, node->pt);
        ntype bfirst = node->first->intersect(p);
        ntype bsecond = node->second->intersect(p);
        // choose the side with the closest bounding box to
             search first
        // (note that the other side is also searched if needed
        if (bfirst < bsecond) {</pre>
            ntype best = search(node->first, p);
            if (bsecond < best)</pre>
                best = min(best, search(node->second, p));
            return best;
            ntype best = search(node->second, p);
            if (bfirst < best)</pre>
                best = min(best, search(node->first, p));
            return best:
    // squared distance to the nearest
    ntype nearest(const point &p) {
        return search(root, p);
};
// some basic test code here
int main() {
    // generate some random points for a kd-tree
    vector<point> vp;
    for (int i = 0; i < 100000; ++i)</pre>
        vp.push_back(point(rand()%100000, rand()%100000));
    kdtree tree(vp);
    // query some points
    for (int i = 0; i < 10; ++i) {
        point g(rand()%100000, rand()%100000);
        cout << "Closest squared distance to (" << q.x << ", "
            << q.y << ")"
             << " is " << tree.nearest(q) << endl;
    return 0;
```



3.6 PersistentSegTree

```
PersistentSegTree.h
<iostream>, <vector>
                                                     44b362, 82 lines
using namespace std;
/* persistent segment tree w/ sum query */
struct per seq {
 vector<int> lc,rc,lx,rx,val,roots = {0};
  //CHANGE ME
  inline static int combine(int a, int b) {
   return a+b;
  //build PST on indices i ... j of array a
  int build(int i, int j, int* a) {
   int v = lc.size();
   lc.push_back(-1);
    rc.push back (-1);
    lx.push_back(i);
    rx.push back(j);
    val.push_back(a[i]);
    if(i != j) {
     int 1 = build(i, (i+j)/2, a), r = build((i+j)/2+1, j, a);
     lc[v] = 1;
     rc[v] = r;
     val[v] = combine(val[1], val[r]);
    return v;
  int q(int v, int i, int j) {
    if(j < lx[v] \mid | rx[v] < i)
     return 0;
   if(i <= lx[v] && rx[v] <= j)</pre>
     return val[v];
    return combine(q(lc[v],i,j),q(rc[v],i,j));
  int u(int v, int i, int a) {
   if(i < lx[v] || rx[v] < i)
     return v;
    int w = lc.size();
    lc.push_back(lc[v]);
    rc.push_back(rc[v]);
    lx.push_back(lx[v]);
    rx.push_back(rx[v]);
    //CHANGE ME
    val.push_back(val[v]+a);
    if(lx[v] != rx[v]) {
     int 1 = u(lc[v],i,a), r = u(rc[v],i,a);
     lc[w] = 1;
     rc[w] = r;
     val[w] = combine(val[1], val[r]);
    return w;
  //sum from i to j after t updates
  int query(int i, int j, int t = -1) {
   if(t == -1) t = roots.size()-1;
    return q(roots[t],i,j);
```

//add a to position i at time t

void update(int i, int a, int t) {

roots.push_back(u(roots[t],i,a));

```
};
/* USAGE */
int main() {
 int a[5] = {1,1,1,1,1};
 per_seq p;
 p.build(0,4,a);
 p.update(1,1);
 p.update(3,2);
 cout << p.query(2,2,5) << "\n";
 cout << p.query(0,2,5) << "\n";
 cout << p.query(1,0,3) << "\n"
 return 0;
```

Numerical (4)

GoldenSectionSearch

GoldenSectionSearch.h

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

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

x2 = a + r*(b-a); f2 = f(x2);

```
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 maximum
     b = x2; x2 = x1; f2 = f1;
      x1 = b - r*(b-a); f1 = f(x1);
      a = x1; x1 = x2; f1 = f2;
```

4.2 Polynomial

return a;

```
Polynomial.h
                                                     c9b7b0, 17 lines
struct Poly {
 vector<double> a;
 double operator() (double x) const {
   double val = 0;
   for(int i = sz(a); i--;) (val *= x) += a[i];
   return val:
 void diff() {
   rep(i, 1, sz(a)) a[i-1] = i*a[i];
   a.pop_back();
 void divroot (double x0) {
   double b = a.back(), c; a.back() = 0;
   for(int i=sz(a)-1; i--;) c = a[i], a[i] = a[i+1]*x0+b, b=c;
   a.pop_back();
```

```
};
PolyRoots.h
Description: Finds the real roots to a polynomial.
Usage: poly_roots(\{\{2, -3, 1\}\}, -1e9, 1e9\}) // solve x^2-3x+2=0
Time: \mathcal{O}\left(n^2\log(1/\epsilon)\right)
"Polynomial.h"
vector<double> poly_roots(Poly p, double xmin, double xmax)
 if (sz(p.a) == 2) { return {-p.a[0]/p.a[1]}; }
  vector<double> ret;
 Poly der = p_i
 der.diff();
  auto dr = poly_roots(der, xmin, xmax);
  dr.push_back(xmin-1);
  dr.push_back(xmax+1);
  sort (all (dr));
  rep(i, 0, sz(dr) - 1) {
    double l = dr[i], h = dr[i+1];
    bool sign = p(1) > 0;
    if (sign ^{\circ} (p(h) > 0)) {
      rep(it,0,60) { // while (h-l>1e-8)
        double m = (1 + h) / 2, f = p(m);
        if ((f \le 0) \hat{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: $\mathcal{O}\left(n^2\right)$

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

LinearRecurrence

LinearRecurrence.h

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

f4e444, 26 lines

```
typedef vector<11> Poly;
11 linearRec(Poly S, Poly tr, 11 k) {
 int n = sz(tr);
  auto combine = [&] (Poly a, Poly b) {
    Polv res(n \star 2 + 1);
```

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

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

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

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

4.4 HillClimbing

HillClimbing.h

Description: Poor man's optimization for unimodal functions_{f40e55, 16 lines}

```
typedef array<double, 2> P;
double func(P p);

pair<double, P> hillClimb(P start) {
  pair<double, P> cur(func(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(func(p), p));
    }
  }
  return cur;
}
```

4.5 Matrix

Determinant.h

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

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

IntDeterminant.h

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

```
Time: \mathcal{O}(N^3)
                                                       3313dc, 18 lines
const 11 mod = 12345;
11 det(vector<vector<ll>>& a) {
 int n = sz(a); ll ans = 1;
 rep(i,0,n) {
    rep(j,i+1,n) {
      while (a[j][i] != 0) { // gcd step
        ll t = a[i][i] / a[j][i];
        if (t) rep(k,i,n)
          a[i][k] = (a[i][k] - a[j][k] * t) % mod;
        swap(a[i], a[j]);
        ans \star = -1;
    ans = ans * a[i][i] % mod;
    if (!ans) return 0;
  return (ans + mod) % mod;
```

Simplex.h

swap(B[r], N[s]);

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.

```
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
```

Time: $\mathcal{O}\left(NM*\#pivots\right)$, where a pivot may be e.g. an edge relaxation. $\mathcal{O}\left(2^{n}\right)$ in the general case.

```
typedef double T; // long double, Rational, double + mod<P>...
typedef vector<T> vd;
typedef vector<vd> vvd;
const T eps = 1e-8, inf = 1/.0;
#define MP make pair
#define ltj(X) if (s == -1 \mid \mid MP(X[j], N[j]) < MP(X[s], N[s])) s=j
struct LPSolver {
 int m, n;
 vi N, B;
 vvd D:
 LPSolver (const vvd& A, const vd& b, const vd& c) :
    m(sz(b)), n(sz(c)), N(n+1), B(m), D(m+2), vd(n+2)) {
      rep(i, 0, m) rep(j, 0, n) D[i][j] = A[i][j];
      rep(i,0,m) \{ B[i] = n+i; D[i][n] = -1; D[i][n+1] = b[i]; \}
     rep(j,0,n) { N[j] = j; D[m][j] = -c[j]; }
     N[n] = -1; D[m+1][n] = 1;
  void pivot(int r, int s) {
    T *a = D[r].data(), inv = 1 / a[s];
    rep(i, 0, m+2) if (i != r \&\& abs(D[i][s]) > eps) {
     T *b = D[i].data(), inv2 = b[s] * inv;
      rep(j, 0, n+2) b[j] -= a[j] * inv2;
     b[s] = a[s] * inv2;
    rep(j, 0, n+2) if (j != s) D[r][j] *= inv;
    rep(i, 0, m+2) if (i != r) D[i][s] *= -inv;
    D[r][s] = inv;
```

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

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

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

SolveLinear

SolveLinear.h

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

44c9ab, 38 lines

```
typedef vector<double> vd;
const double eps = 1e-12;
int solveLinear(vector<vd>& A, vd& b, vd& x) {
  int n = sz(A), m = sz(x), rank = 0, br, bc;
 if (n) assert(sz(A[0]) == m);
  vi col(m); iota(all(col), 0);
  rep(i,0,n) {
   double v, bv = 0;
   rep(r,i,n) rep(c,i,m)
     if ((v = fabs(A[r][c])) > bv)
       br = r, bc = c, bv = v;
    if (bv <= eps) {
      rep(j, i, n) if (fabs(b[j]) > eps) return -1;
     break;
    swap(A[i], A[br]);
    swap(b[i], b[br]);
    swap(col[i], col[bc]);
    rep(j,0,n) swap(A[j][i], A[j][bc]);
   bv = 1/A[i][i];
    rep(j,i+1,n) {
     double fac = A[j][i] * bv;
     b[j] = fac * b[i];
     rep(k,i+1,m) A[j][k] = fac*A[i][k];
    rank++;
  x.assign(m, 0);
  for (int i = rank; i--;) {
   b[i] /= A[i][i];
   x[col[i]] = b[i];
   rep(j, 0, i) b[j] -= A[j][i] * b[i];
  return rank; // (multiple solutions if rank < m)
```

SolveLinear2.h

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

```
08e495, 7 lines
"SolveLinear.h"
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: $\mathcal{O}(n^2m)$ 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()) break;</pre>
    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)
```

Fourier transforms

FastFourierTransform.h

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

```
Time: \mathcal{O}(N \log N) with N = |A| + |B| \ (\sim 1s \text{ for } N = 2^{22})
                                                                                      aca9fa, 35 lines
```

```
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, M_PII / k); // M_PI, lower-case L
   rep(i,k,2*k) rt[i] = R[i] = i&1 ? R[i/2] * x : R[i/2];
 rep(i, 0, n) rev[i] = (rev[i / 2] | (i & 1) << L) / 2;
 rep(i,0,n) if (i < rev[i]) swap(a[i], a[rev[i]]);
 for (int k = 1; k < n; k *= 2)
```

```
for (int i = 0; i < n; i += 2 * k) rep(j, 0, k) {
     Cz = rt[j+k] * a[i+j+k]; // (25\% faster if hand-rolled)
      a[i + j + k] = a[i + j] - z;
      a[i + j] += z;
vd conv(const vd& a, const vd& b) {
 if (a.empty() || b.empty()) return {};
 vd res(sz(a) + sz(b) - 1);
  int L = 32 - builtin clz(sz(res)), n = 1 \ll L;
  vector<C> in(n), out(n);
  copy(all(a), begin(in));
  rep(i,0,sz(b)) in[i].imag(b[i]);
  fft(in);
  trav(x, in) x *= x;
  rep(i, 0, n) out[i] = in[-i & (n - 1)] - conj(in[i]);
  fft (out);
 rep(i, 0, sz(res)) res[i] = imag(out[i]) / (4 * n);
 return res;
```

FastFourierTransformMod.h

Description: Higher precision FFT, can be used for convolutions modulo arbitrary integers as long as $N \log_2 N \cdot \text{mod} < 8.6 \cdot 10^{14}$ (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=int(sqrt(M));</pre>
  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 av = 11(real(out1[i])+.5), cv = 11(imag(outs[i])+.5);
    11 \text{ bv} = 11(\text{imag}(\text{outl}[i]) + .5) + 11(\text{real}(\text{outs}[i]) + .5);
    res[i] = ((av % M * cut + bv) % M * cut + cv) % M;
  return res;
```

NumberTheoreticTransform.h

Description: Can be used for convolutions modulo specific nice primes of the form $2^a b + 1$, where the convolution result has size at most $2^{\bar{a}}$. Inputs must be in [0, mod).

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

```
"../number-theory/ModPow.h"
const 11 mod = (119 << 23) + 1, root = 62; // = 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<ll> v1;
void ntt(vl& a, vl& rt, vl& rev, int n) {
  rep(i,0,n) if (i < rev[i]) swap(a[i], a[rev[i]]);
 for (int k = 1; k < n; k *= 2)
```

11 z = rt[j + k] * a[i + j + k] % mod, &ai = a[i + j];

a[i + j + k] = (z > ai ? ai - z + mod : ai - z);

for (int i = 0; i < n; i += 2 * k) rep(j,0,k) {

```
ai += (ai + z >= mod ? z - mod : z);
}

vl conv(const vl& a, const vl& b) {
   if (a.empty() || b.empty())
      return {};
   int s = sz(a)+sz(b)-1, B = 32 - __builtin_clz(s), n = 1 << B;
   vl L(a), R(b), out(n), rt(n, 1), rev(n);
   L.resize(n), R.resize(n);
   rep(i,0,n) rev[i] = (rev[i / 2] | (i & 1) << B) / 2;
   ll curL = mod / 2, inv = modpow(n, mod - 2);
   for (int k = 2; k < n; k *= 2) {
      ll z[] = {1, modpow(root, curL /= 2)};
      rep(i,k,2*k) rt[i] = rt[i / 2] * z[i & 1] % mod;
   }
   ntt(L, rt, rev, n); ntt(R, rt, rev, n);
   rep(i,0,n) out[-i & (n-1)] = L[i] * R[i] % mod * inv % mod;
   ntt(out, rt, rev, n);
   return {out.begin(), out.begin() + s};
}</pre>
```

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

3de473, 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.

```
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. 6f684f, 3 lines

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

ModPow.h

b83e45, 8 lines

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

11 modpow(11 b, 11 e) {
    11 ans = 1;
    for (; e; b = b * b % mod, e /= 2)
        if (e & 1) ans = ans * b % mod;
    return ans;
}
```

ModLog.h

Description: Returns the smallest $x \ge 0$ s.t. $a^x = b \pmod{m}$. a and m must be coprime.

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

49d606, 10 lines

```
11 modLog(11 a, 11 b, 11 m) {
   assert(_gcd(a, m) == 1);
   11 n = (11) sqrt(m) + 1, e = 1, x = 1, res = LLONG_MAX;
   unordered_map<11, 11> f;
   rep(i,0,n) e = e * a % m;
   rep(i,0,n) x = x * e % m, f.emplace(x, i + 1);
   rep(i,0,n) if (f.count(b = b * a % m))
      res = min(res, f[b] * n - i - 1);
   return res;
}
```

ModSum.h

Description: Sums of mod'ed arithmetic progressions.

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

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

5c5bc5, 16 lines

```
typedef unsigned long long ull;
ull sumsq(ull to) { return to / 2 * ((to-1) | 1); }

ull divsum(ull to, ull c, ull k, ull m) {
    ull res = k / m * sumsq(to) + c / m * to;
    k %= m; c %= m;
    if (!k) return res;
    ull to2 = (to * k + c) / m;
    return res + (to - 1) * to2 - divsum(to2, m-1 - c, m, k);
}

ll modsum(ull to, ll c, ll k, ll m) {
    c = ((c % m) + m) % m;
    k = ((k % m) + m) % m;
    return to * c + k * sumsq(to) - m * divsum(to, c, k, m);
}
```

ModMulLL.h

```
Description: Calculate a \cdot b \mod c (or a^b \mod c) for 0 \le a, b < c < 2^{63}. Time: \mathcal{O}(1) for mod_mul, \mathcal{O}(\log b) for mod_pow 88c37a, 12 lines
```

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

ModSqrt.h

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

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

```
"ModPow.h"
                                                        19a793, 24 lines
ll sqrt(ll a, ll p) {
 a %= p; if (a < 0) a += p;
 if (a == 0) return 0;
 assert (modpow(a, (p-1)/2, p) == 1); // else no solution
 if (p % 4 == 3) return modpow(a, (p+1)/4, p);
  // a^{(n+3)/8} \text{ or } 2^{(n+3)/8} * 2^{(n-1)/4} \text{ works if } p \% 8 == 5
 11 s = p - 1, n = 2;
  int r = 0, m;
  while (s % 2 == 0)
   ++r, s /= 2;
  while (modpow(n, (p-1) / 2, p) != p-1) ++n;
  11 x = modpow(a, (s + 1) / 2, p);
  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{ gs} = \text{modpow}(g, 1LL << (r - m - 1), p);
    q = qs * qs % p;
    x = x * gs % p;
    b = b * q % p;
```

5.2 Primality

eratosthenes.h

Description: Prime sieve for generating all primes up to a certain limit. isprime[i] is true iff i is a prime.

Time: $\lim_{n\to\infty} 100'000'000 \approx 0.8 \text{ s.}$ Runs 30% faster if only odd indices are stored.

```
const int MAX_PR = 5'000'000;
bitset<MAX_PR> isprime;
vi eratosthenes_sieve(int lim) {
  isprime.set(); isprime[0] = isprime[1] = 0;
  for (int i = 4; i < lim; i += 2) isprime[i] = 0;
  for (int i = 3; i*i < lim; i += 2) if (isprime[i])
    for (int j = i*i; j < lim; j += i*2) isprime[j] = 0;
  vi pr;
  rep(i,2,lim) if (isprime[i]) pr.push_back(i);
  return pr;
}
```

MillerRabin.h

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



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

Factor.h

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

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

```
"ModMulLL.h", "MillerRabin.h"
                                                     f5adaa, 18 lines
ull pollard(ull n) {
  auto f = [n](ull x) { return (mod_mul(x, x, n) + 1) % n; };
  if (!(n & 1)) return 2;
  for (ull i = 2;; i++) {
   ull x = i, y = f(x), p;
    while ((p = \_gcd(n + y - x, n)) == 1)
     x = f(x), y = f(f(y));
   if (p != n) return p;
vector<ull> factor(ull n) {
 if (n == 1) return {};
 if (isPrime(n)) return {n};
  ull x = pollard(n);
  auto 1 = factor(x), r = factor(n / x);
 1.insert(1.end(), all(r));
  return 1;
```

5.3 Divisibility

euclid.h

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

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

Euclid.java

```
Description: Finds \{x, y, d\} s.t. ax + by = d = gcd(a, b). <sub>6aba01, 11 lines</sub>
```

```
static BigInteger[] euclid(BigInteger a, BigInteger b) {
   BigInteger x = BigInteger.ONE, yy = x;
   BigInteger y = BigInteger.ZERO, xx = y;
   while (b.signum() != 0) {
      BigInteger q = a.divide(b), t = b;
      b = a.mod(b); a = t;
      t = xx; xx = x.subtract(q.multiply(xx)); x = t;
      t = yy; yy = y.subtract(q.multiply(yy)); y = t;
   }
   return new BigInteger[]{x, y, a};
}
```

CRT.h

5.3.1 Bézout's identity

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

$$ax + by = d$$

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

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

phiFunction.h

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

```
const int LIM = 5000000;
int phi[LIM];

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

5.4 Fractions

ContinuedFractions.h

Description: Given N and a real number $x \ge 0$, finds the closest rational approximation p/q with $p,q \le N$. It will obey $|p/q - x| \le 1/qN$. For consecutive convergents, $p_{k+1}q_k - q_{k+1}p_k = (-1)^k$. $(p_k/q_k$ alternates

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

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

```
if (a > b) {
    // If b > a/2, we have a semi—convergent that gives us a
    // better approximation; if b = a/2, we *may* have one.
    // Return {P, Q} here for a more canonical approximation.
    return (abs(x - (d)NP / (d)NQ) < abs(x - (d)P / (d)Q)) ?
    make_pair(NP, NQ) : make_pair(P, Q);
}
if (abs(y = 1/(y - (d)a)) > 3*N) {
    return {NP, NQ};
}
LP = P; P = NP;
LQ = Q; Q = NQ;
}
```

FracBinarySearch.h

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

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

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

5.5 Pythagorean Triples

The Pythagorean triples are uniquely generated by

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

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

5.6 Primes

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

IntPerm binomialModPrime multinomial

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

Estimates

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

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

Combinatorial (6)

Permutations

6.1.1 Factorial

						9		
n!	1 2 6	24 1	20 720	0 5040	40320	362880	3628800 17	
n!	4.0e7	4.8e	8 6.2e	9 8.7e	10 1.3e	12 2.1e1	3 3.6e14	
n	20	25	30	40	50 1	00 150	3 3.6e14) 171 32 >DBL_MA	
n!	2e18	2e25	3e32	8e47.3	664 9e	157 6e26	$52 > DBL_M$	AΧ

IntPerm.h

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

int permToInt(vi& v) { int use = 0, i = 0, r = 0; $trav(x, v) r = r * ++i + \underline{\quad builtin popcount(use & -(1 << x)),}$ // (note: minus. not \sim !) return r:

Cycles 6.1.2

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

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

Derangements

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

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

6.1.4 Burnside's lemma

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

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

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

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

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

Partitions and subsets

6.2.1 Partition function

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

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

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

6.2.2 Binomials

binomialModPrime.h

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

Time: $O(\log_n n)$

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

Il multinomial (vi& v) {
```

```
11 c = 1, m = v.empty() ? 1 : v[0];
rep(i, 1, sz(v)) rep(j, 0, v[i])
```

General purpose numbers

6.3.1 Bernoulli numbers

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

Sums of powers:

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

Euler-Maclaurin formula for infinite sums:

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

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

6.3.2 Stirling numbers of the first kind

Number of permutations on n items with k cycles.

$$c(n,k) = c(n-1,k-1) + (n-1)c(n-1,k), \ c(0,0) = 1$$

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

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

6.3.3 Eulerian numbers

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

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

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

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

6.3.4 Stirling numbers of the second kind

Partitions of n distinct elements into exactly k groups.

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

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

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, For p prime,

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

6.3.6 Labeled unrooted trees

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

6.3.7 Catalan numbers

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

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

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

- sub-diagonal monotone paths in an $n \times n$ grid.
- \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.
- \bullet permutations of [n] with no 3-term increasing subseq.

Graph (7)

7.1 Fundamentals

BellmanFord.h

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

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

FlovdWarshall.h

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

Description: Calculates all-pairs shortest path in a directed graph that might have negative edge weights. Input is an distance matrix m, where $m[i][j] = \inf$ 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.

```
const 11 inf = 1LL << 62;
void floydWarshall(vector<vector<1l>>& m) {
  int n = sz(m);
  rep(i,0,n) m[i][i] = min(m[i][i], 0LL);
  rep(k,0,n) rep(i,0,n) rep(j,0,n)
  if (m[i][k] != inf && m[k][j] != inf) {
    auto newDist = max(m[i][k] + m[k][j], -inf);
    m[i][j] = min(m[i][j], newDist);
  }
  rep(k,0,n) if (m[k][k] < 0) rep(i,0,n) rep(j,0,n)
  if (m[i][k] != inf && m[k][j] != inf) m[i][j] = -inf;
}</pre>
```

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.

```
return ret;
```

7.2 Euler walk

EulerWalk.h

Description: Eulerian undirected/directed path/cycle algorithm. 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, also put it->second in s (and then ret).

```
Time: \mathcal{O}(E) where E is the number of edges.
```

f8bd47, 27 lines

```
vector<pii> outs; // (dest, edge index)
 int nins = 0;
vi euler walk(vector<V>& nodes, int nedges, int src=0) {
 trav(n, nodes) c += abs(n.nins - sz(n.outs));
 if (c > 2) return {};
 vector<vector<pii>::iterator> its;
 trav(n, nodes)
   its.push_back(n.outs.begin());
 vector<bool> eu(nedges);
 vi ret, s = \{src\};
 while(!s.empty()) {
   int x = s.back();
   auto& it = its[x], end = nodes[x].outs.end();
   while(it != end && eu[it->second]) ++it;
   if(it == end) { ret.push_back(x); s.pop_back(); }
    else { s.push_back(it->first); eu[it->second] = true; }
 if(sz(ret) != nedges+1)
   ret.clear(); // No Eulerian cycles/paths.
  // else, non-cycle if ret.front() != ret.back()
 reverse(all(ret));
 return ret:
```

7.3 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)

typedef 11 Flow;

struct Edge {
  int dest, back;
  Flow f, c;
};

struct PushRelabel {
  vector<vector<Edge>> g;
  vector<Flow> ec;
  vector<Flow> ec;
  vector<vi> hs; vi H;
  PushRelabel(int n) : g(n), ec(n), cur(n), hs(2*n), H(n) {}

void add_edge(int s, int t, Flow cap, Flow rcap=0) {
  if (s == t) return;
  g[s].push_back({t, sz(g[t]), 0, cap});
```

 $q[t].push_back({s, sz(q[s])-1, 0, rcap});$

```
void add_flow(Edge& e, Flow f) {
    Edge &back = q[e.dest][e.back];
    if (!ec[e.dest] && f) hs[H[e.dest]].push_back(e.dest);
    e.f += f; e.c -= f; ec[e.dest] += f;
   back.f -= f; back.c += f; ec[back.dest] -= f;
  Flow maxflow(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] = q[i].data();
    trav(e, g[s]) add_flow(e, e.c);
    for (int hi = 0;;) {
      while (hs[hi].empty()) if (!hi--) return -ec[s];
      int u = hs[hi].back(); hs[hi].pop_back();
      while (ec[u] > 0) // discharge u
        if (cur[u] == g[u].data() + sz(g[u])) {
          H[u] = 1e9;
          trav(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] < v)
              --co[H[i]], H[i] = v + 1;
          hi = H[u];
        } else if (cur[u]->c && H[u] == H[cur[u]->dest]+1)
          add_flow(*cur[u], min(ec[u], cur[u]->c));
        else ++cur[u];
};
```

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 $\mathcal{O}(E^2)$

```
6915ce, 81 lines
#include <bits/extc++.h>
const 11 INF = numeric limits<11>::max() / 4;
typedef vector<ll> 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 cost) {
    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;
```

__gnu_pbds::priority_queue<pair<11, int>> q;

```
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];
     trav(i, ed[s]) if (!seen[i])
       relax(i, cap[s][i] - flow[s][i], cost[s][i], 1);
     trav(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], INF);
 pair<11, 11> maxflow(int s, int t) {
   11 \text{ totflow} = 0, \text{ totcost} = 0;
    while (path(s), seen[t]) {
     11 fl = INF;
      for (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; 11 v;
   while (ch-- && it--)
     rep(i,0,N) if (pi[i] != INF)
       trav(to, ed[i]) if (cap[i][to])
         if ((v = pi[i] + cost[i][to]) < pi[to])</pre>
           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 979bb9, 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;
    fill(all(par), -1);
    par[source] = 0;
    int ptr = 1;
    q[0] = source;
```

```
rep(i,0,ptr) {
      int x = q[i];
      trav(e, graph[x]) {
        if (par[e.first] == -1 && e.second > 0) {
          par[e.first] = x;
          q[ptr++] = e.first;
          if (e.first == sink) goto out;
    return flow:
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].erase(y);</pre>
      graph[y][p] += inc;
```

MinCut.h

Description: After running max-flow, the left side of a min-cut from s to tis 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: $\mathcal{O}(V^3)$

03261f, 31 lines

```
pair<int, vi> GetMinCut(vector<vi>& weights) {
 int N = sz(weights);
  vi used(N), cut, best_cut;
  int best weight = -1;
  for (int phase = N-1; phase >= 0; phase--) {
    vi w = weights[0], added = used;
    int prev, k = 0;
    rep(i,0,phase){
     prev = k;
      k = -1;
      rep(j,1,N)
       if (!added[j] && (k == -1 || w[j] > w[k])) k = j;
      if (i == phase-1) {
        rep(j,0,N) weights[prev][j] += weights[k][j];
        rep(j,0,N) weights[j][prev] = weights[prev][j];
        used[k] = true;
        cut.push_back(k);
        if (best_weight == -1 || w[k] < best_weight) {</pre>
          best_cut = cut;
          best_weight = w[k];
     } else {
        rep(j,0,N)
         w[j] += weights[k][j];
        added[k] = true;
 return {best_weight, best_cut};
```



7.4 Matching

hopcroftKarp.h

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

Usage: vi btoa(m, -1); hopcroftKarp(g, btoa);

```
Time: \mathcal{O}\left(\sqrt{V}E\right)
bool dfs(int a, int L, vector<vi>& q, vi& btoa, vi& A, vi& B) {
  if (A[a] != L) return 0;
  A[a] = -1;
  trav(b, g[a]) if (B[b] == L + 1) {
   B[b] = 0;
    if (btoa[b] == -1 || dfs(btoa[b], L + 1, g, btoa, A, B))
      return btoa[b] = a, 1;
  return 0;
int hopcroftKarp(vector<vi>& q, vi& btoa) {
  int res = 0;
  vi A(g.size()), B(btoa.size()), cur, next;
  for (;;) {
    fill(all(A), 0);
    fill(all(B), 0);
    cur.clear();
    trav(a, btoa) if(a != -1) A[a] = -1;
    rep(a, 0, sz(q)) if(A[a] == 0) cur.push_back(a);
    for (int lav = 1;; lav++) {
     bool islast = 0;
      next.clear();
      trav(a, cur) trav(b, g[a]) {
        if (btoa[b] == -1) {
          B[b] = lav;
          islast = 1;
        else if (btoa[b] != a && !B[b]) {
          B[b] = lav;
          next.push_back(btoa[b]);
      if (islast) break;
      if (next.empty()) return res;
      trav(a, next) A[a] = lay;
      cur.swap(next);
    rep(a, 0, sz(q))
      res += dfs(a, 0, g, 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(g, btoa); **Time:** $\mathcal{O}(VE)$

```
bool find(int j, vector<vi>& g, vi& btoa, vi& vis) {
   if (btoa[j] == -1) return 1;
   vis[j] = 1; int di = btoa[j];
   trav(e, g[di])
   if (!vis[e] && find(e, g, btoa, vis)) {
     btoa[e] = di;
    return 1;
}
```

```
return 0;
}
int dfsMatching(vector<vi>& g, vi& btoa) {
  vi vis;
  rep(i,0,sz(g)) {
    vis.assign(sz(btoa), 0);
    trav(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 independent set.

```
"DFSMatching.h"
                                                     d0b3f2, 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);
 trav(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;
   trav(e, g[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

R[j] = i;

Description: Min cost bipartite matching. Negate costs for max cost. **Time:** $\mathcal{O}\left(N^3\right)$ 055ca9, 75 lines

```
typedef vector<double> vd;
bool zero(double x) { return fabs(x) < 1e-10; }</pre>
double MinCostMatching(const vector<vd>& cost, vi& L, vi& R) {
 int n = sz(cost), mated = 0;
 vd dist(n), u(n), v(n);
 vi dad(n), seen(n);
  rep(i,0,n) {
   u[i] = cost[i][0];
    rep(j,1,n) u[i] = min(u[i], cost[i][j]);
  rep(j,0,n) {
    v[j] = cost[0][j] - u[0];
    rep(i,1,n) \ v[j] = min(v[j], cost[i][j] - u[i]);
  L = R = vi(n, -1);
 rep(i,0,n) rep(j,0,n) {
    if (R[j] != -1) continue;
    if (zero(cost[i][j] - u[i] - v[j])) {
     L[i] = j;
```

```
mated++;
    break;
for (; mated < n; mated++) { // until solution is feasible</pre>
  int s = 0;
  while (L[s] !=-1) s++;
  fill(all(dad), -1);
  fill(all(seen), 0);
  rep(k,0,n)
    dist[k] = cost[s][k] - u[s] - v[k];
  int j = 0;
  for (;;) {
    \dot{1} = -1;
    rep(k,0,n){
      if (seen[k]) continue;
      if (j == -1 || dist[k] < dist[j]) j = k;</pre>
    seen[j] = 1;
    int i = R[i];
    if (i == -1) break;
    rep(k,0,n) {
      if (seen[k]) continue;
      auto new_dist = dist[j] + cost[i][k] - u[i] - v[k];
      if (dist[k] > new_dist) {
        dist[k] = new_dist;
        dad[k] = j;
  rep(k,0,n) {
    if (k == j || !seen[k]) continue;
    auto w = dist[k] - dist[j];
    v[k] += w, u[R[k]] -= w;
  u[s] += dist[j];
  while (dad[j] >= 0) {
   int d = dad[j];
    R[j] = R[d];
    L[R[j]] = j;
    j = d;
  R[j] = s;
  L[s] = j;
auto value = vd(1)[0];
rep(i,0,n) value += cost[i][L[i]];
return value:
```

GeneralMatching.h

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

SCC BiconnectedComponents 2sat MaximalCliques

```
if (M != N) do {
 mat.resize(M, vector<ll>(M));
  rep(i,0,N) {
   mat[i].resize(M);
   rep(j,N,M) {
     int r = rand() % mod;
      mat[i][j] = r, mat[j][i] = (mod - r) % mod;
} while (matInv(A = mat) != M);
vi has(M, 1); vector<pii> ret;
rep(it,0,M/2) {
  rep(i,0,M) if (has[i])
   rep(j,i+1,M) if (A[i][j] && mat[i][j]) {
      fi = i; fj = j; goto done;
  } assert(0); done:
  if (fj < N) ret.emplace_back(fi, fj);</pre>
  has[fi] = has[fj] = 0;
  rep(sw.0.2) {
   11 a = modpow(A[fi][fj], mod-2);
   rep(i,0,M) if (has[i] && A[i][fj]) {
     ll b = A[i][fj] * a % mod;
      rep(j, 0, M) A[i][j] = (A[i][j] - A[fi][j] * b) % mod;
   swap(fi,fj);
return ret;
```

DFS algorithms

SCC.h

Description: 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

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}(E+V)
                                                     bb2963, 24 lines
vi val, comp, z, cont;
int Time, ncomps;
template < class G, class F> int dfs (int j, G& g, F& f) {
  int low = val[j] = ++Time, x; z.push_back(j);
  trav(e,g[j]) if (comp[e] < 0)
   low = min(low, val[e] ?: dfs(e,g,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(q);
  val.assign(n, 0); comp.assign(n, -1);
 Time = ncomps = 0;
  rep(i,0,n) if (comp[i] < 0) dfs(i, q, 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}(E+V)
                                                      cca7e6, 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;
 trav(pa, ed[at]) if (pa.second != par) {
   tie(y, e) = pa;
    if (num[y]) {
      top = min(top, num[y]);
      if (num[v] < me)</pre>
        st.push_back(e);
    } else {
      int si = sz(st);
      int up = dfs(y, e, f);
      top = min(top, up);
      if (up == me) {
        st.push back(e);
        f(vi(st.begin() + si, st.end()));
        st.resize(si);
      else if (up < me) st.push back(e);</pre>
      else { /* e is a bridge */ }
  return top;
template<class F>
void bicomps (F f) {
 num.assign(sz(ed), 0);
  rep(i,0,sz(ed)) if (!num[i]) dfs(i, -1, f);
```

2sat.h

Description: 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, \sim3); // Var 0 is true or var 3 is false
ts.set_value(2); // Var 2 is true
ts.at_most_one(\{0, \sim 1, 2\}); // <= 1 of vars 0, \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: $\mathcal{O}(N+E)$, where N is the number of boolean variables, and E is the number of clauses. 0911c1, 56 lines

```
struct TwoSat {
 int N:
 vector<vi> gr;
 vi values; // 0 = false, 1 = true
 TwoSat(int n = 0) : N(n), gr(2*n) {}
 int add_var() { // (optional)
```

```
gr.emplace_back();
   gr.emplace_back();
    return N++;
 void either(int f, int i) {
   f = \max(2*f, -1-2*f);
   j = \max(2*j, -1-2*j);
   gr[f].push_back(j^1);
    gr[i].push back(f^1);
 void set_value(int x) { either(x, x); }
 void at_most_one(const vi& li) { // (optional)
   if (sz(li) <= 1) return;</pre>
    int cur = \simli[0];
    rep(i,2,sz(li)) {
     int next = add var();
     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);
   trav(e, qr[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;
};
```

7.6Heuristics

MaximalCliques.h

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

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

b0d5b1, 12 lines

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

d952a9, 93 lines



MaximumClique.h

struct Maxclique {

typedef vector<bitset<200>> vb;

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.

fbbef1, 49 lines

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

MaximumIndependentSet.h

};

C[k].push_back(v.i);

expand(T, lev + 1);

q.pop_back(), R.pop_back();

rep(i,0,sz(e)) V.push_back({i});

if (j > 0) T[j - 1].d = 0;

rep(k, mnk, mxk + 1) trav(i, C[k])

} else if (sz(q) > sz(qmax)) qmax = q;

vi maxClique() { init(V), expand(V); return qmax; }

T[j].i = i, T[j++].d = k;

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

Maxclique(vb conn) : e(conn), C(sz(e)+1), S(sz(C)), old(S) {

7.7 Trees

TreePower.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 $\mathcal{O}(\log N)$

bfce85, 25 lines

```
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];</pre>
 return nod:
int lca(vector<vi>& tbl, vi& depth, int a, int b) {
 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. Can also find the distance between two nodes.

```
Usage: LCA lca(undirGraph);
lca.query(firstNode, secondNode);
lca.distance(firstNode, secondNode);
Time: O(N log N + O)
```

```
Time: \mathcal{O}(N \log N + Q)
"../data-structures/RMQ.h"
                                                      aa0d4d, 37 lines
typedef vector<pii> vpi;
typedef vector<vpi> graph;
struct LCA {
 vi time;
 vector<ll> dist:
 RMQ<pii> rmq;
 LCA(graph& C): time(sz(C), -99), dist(sz(C)), rmg(dfs(C)) {}
 vpi dfs(graph& C) {
   vector<tuple<int, int, int, 11>> q(1);
    vpi ret;
   int T = 0, v, p, d; 11 di;
    while (!q.empty()) {
      tie(v, p, d, di) = q.back();
      q.pop_back();
      if (d) ret.emplace_back(d, p);
     time[v] = T++;
      dist[v] = di;
      trav(e, C[v]) if (e.first != p)
        q.emplace_back(e.first, v, d+1, di + e.second);
    return ret;
```

```
int query(int a, int b) {
   if (a == b) return a;
   a = time[a], b = time[b];
   return rmq.query(min(a, b), max(a, b)).second;
}
ll distance(int a, int b) {
   int lca = query(a, b);
   return dist[a] + dist[b] - 2 * dist[lca];
}
};
```

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"
                                                     dabd75, 20 lines
vpi compressTree(LCA& lca, const vi& subset) {
 static vi rev; rev.resize(sz(lca.dist));
 vi li = subset, &T = lca.time;
 auto cmp = [&](int a, int b) { return T[a] < T[b]; };</pre>
  sort(all(li), cmp);
 int m = sz(1i)-1;
 rep(i,0,m) {
   int a = li[i], b = li[i+1];
   li.push back(lca.query(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.query(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. The function of the HLD can be changed by modifying T, LOW and f. f is assumed to be associative and commutative.

```
Usage: HLD hld(G);
hld.update(index, value);
tie(value, lca) = hld.query(n1, n2);
"../data-structures/SegmentTree.h"
typedef vectorpii> vpi;

struct Node {
  int d, par, val, chain = -1, pos = -1;
};

struct Chain {
  int par, val;
  vector<int> nodes;
  Tree tree;
};

struct HLD {
  typedef int T;
  const T LOW = -(1<<29);</pre>
```

void f(T& a, T b) { a = max(a, b); }

vector<Node> V; vector<Chain> C;

```
HLD(vector<vpi>& g) : V(sz(g)) {
 dfs(0, -1, q, 0);
 trav(c, C) {
   c.tree = {sz(c.nodes), 0};
   for (int ni : c.nodes)
     c.tree.update(V[ni].pos, V[ni].val);
void update(int node, T val) {
 Node& n = V[node]; n.val = val;
 if (n.chain != -1) C[n.chain].tree.update(n.pos, val);
int pard(Node& nod) {
 if (nod.par == -1) return -1;
 return V[nod.chain == -1 ? nod.par : C[nod.chain].par].d;
// query all *edges* between n1, n2
pair<T, int> querv(int i1, int i2) {
 T ans = LOW;
 while(i1 != i2) {
   Node n1 = V[i1], n2 = V[i2];
   if (n1.chain != -1 && n1.chain == n2.chain) {
     int lo = n1.pos, hi = n2.pos;
     if (lo > hi) swap(lo, hi);
     f(ans, C[n1.chain].tree.query(lo, hi));
     i1 = i2 = C[n1.chain].nodes[hi];
   } else {
     if (pard(n1) < pard(n2))
       n1 = n2, swap(i1, i2);
     if (n1.chain == -1)
       f(ans, n1.val), i1 = n1.par;
     else (
       Chain& c = C[n1.chain];
       f(ans, n1.pos ? c.tree.query(n1.pos, sz(c.nodes))
                     : c.tree.s[1]);
       i1 = c.par;
 return make_pair(ans, i1);
// query all *nodes* between n1, n2
pair<T, int> query2(int i1, int i2) {
 pair<T, int> ans = query(i1, i2);
 f(ans.first, V[ans.second].val);
 return ans;
pii dfs(int at, int par, vector<vpi>& q, int d) {
 V[at].d = d; V[at].par = par;
 int sum = 1, ch, nod, sz;
 tuple<int,int,int> mx(-1,-1,-1);
 trav(e, g[at]){
   if (e.first == par) continue;
   tie(sz, ch) = dfs(e.first, at, g, d+1);
   V[e.first].val = e.second;
   sum += sz;
   mx = max(mx, make_tuple(sz, e.first, ch));
 tie(sz, nod, ch) = mx;
 if (2*sz < sum) return pii(sum, -1);</pre>
 if (ch == -1) { ch = sz(C); C.emplace_back(); }
 V[nod].pos = sz(C[ch].nodes);
 V[nod].chain = ch;
```

```
C[ch].par = at;
   C[ch].nodes.push_back(nod);
   return pii(sum, ch);
};
```

LinkCutTree.h

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

```
Time: All operations take amortized \mathcal{O}(\log N).
                                                     693483, 90 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 push_flip() {
   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 \hat{b};
   Node *x = c[i], *y = b == 2 ? x : x -> c[h], *z = b ? y : x;
   if ((y->p = p)) p->c[up()] = y;
   c[i] = z -> c[i ^ 1];
    if (b < 2) {
     x->c[h] = y->c[h^1];
     z - > c[h ^1] = b ? x : this;
   y - c[i ^1] = b ? this : x;
   fix(); x->fix(); y->fix();
   if (p) p->fix();
   swap(pp, y->pp);
 void splay() {
   for (push_flip(); p; ) {
     if (p->p) p->p->push_flip();
     p->push_flip(); push_flip();
     int c1 = up(), c2 = p->up();
     if (c2 == -1) p->rot (c1, 2);
     else p->p->rot(c2, c1 != c2);
 Node* first() {
   push flip();
   return c[0] ? c[0]->first() : (splay(), this);
} ;
struct LinkCut {
 vector<Node> node;
 LinkCut(int N) : node(N) {}
 void link(int u, int v) { // add \ an \ edge \ (u, \ v)
   assert(!connected(u, v));
   make root(&node[u]);
   node[u].pp = &node[v];
 void cut (int u, int v) { // remove an edge (u, v)
   Node *x = &node[u], *top = &node[v];
   make_root(top); x->splay();
```

```
assert(top == (x->pp ?: x->c[0]));
    if (x->pp) x->pp = 0;
    else {
      x->c[0] = top->p = 0;
      x->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 make_root(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 \rightarrow fix():
  Node* access(Node* u) {
    u->splay();
    while (Node* pp = u->pp) {
      pp \rightarrow splay(); u \rightarrow 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 11:
};
```

DirectedMST.h

Description: Edmonds' algorithm for finding the weight of the minimum spanning tree/arborescence of a directed graph, given a root node. If no MST exists, returns -1.

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

```
"../data-structures/UnionFind.h"
                                                      a69883, 48 lines
struct Edge { int a, b; ll w; };
struct Node {
  Edge kev:
  Node *1, *r;
  ll delta;
  void prop()
   kev.w += delta;
   if (1) 1->delta += delta;
    if (r) r->delta += delta;
    delta = 0:
  Edge top() { prop(); return key; }
Node *merge(Node *a, Node *b) {
  if (!a || !b) return a ?: b;
  a->prop(), b->prop();
  if (a->key.w > b->key.w) swap(a, b);
  swap (a->1, (a->r = merge(b, a->r)));
  return a:
void pop(Node*\& a) { a->prop(); a = merge(a->1, a->r); }
11 dmst(int n, int r, vector<Edge>& g) {
  UF uf(n):
  vector<Node*> heap(n);
  trav(e, g) heap[e.b] = merge(heap[e.b], new Node{e});
  11 \text{ res} = 0;
  vi seen(n, -1), path(n);
```



```
seen[r] = r;
rep(s,0,n) {
 int u = s, qi = 0, w;
 while (seen[u] < 0) {
   path[qi++] = u, seen[u] = s;
   if (!heap[u]) return -1;
   Edge e = heap[u]->top();
   heap[u]->delta -= e.w, pop(heap[u]);
    res += e.w, u = uf.find(e.a);
    if (seen[u] == s) {
     Node \star cyc = 0;
      do cyc = merge(cyc, heap[w = path[--qi]]);
      while (uf.join(u, w));
     u = uf.find(u);
     heap[u] = cyc, seen[u] = -1;
return res;
```

Geometry (8)

8.1 Geometric primitives

Point.h

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

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

lineDistance.h

Description:

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

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 line segment from point s to e.

Usage: Point $\langle double \rangle$ a, b(2,2), p(1,1); bool onSegment = segDist(a,b,p) < 1e-10;

"Point.h" 5c88f4, 6 lines typedef Point < double > P; double segDist(P& s, P& e, P& p) { if (s==e) return (p-s).dist(); **auto** d = (e-s).dist2(), t = min(d, max(.0, (p-s).dot(e-s)));**return** ((p-s)*d-(e-s)*t).dist()/d;

SegmentIntersection.h

Description:

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

```
if (sz(inter) == 1)
cout << "segments intersect at " << inter[0] << endl;</pre>
"Point.h", "OnSegment.h"
template < class P > vector < P > seqInter (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

if (res.first == 1)

Description:

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



3af81c. 9 lines

```
cout << "intersection point at " << res.second << endl;</pre>
"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, s1);
  return {1, (s1 * p + e1 * a) / d};
```

sideOf.h

res

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

```
Usage: bool left = sideOf(p1,p2,q)==1;
"Point.h"
template<class P>
```

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

0.0

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

linearTransformation.h Description:

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

"Point.h" 03a306, 6 lines

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

Angle.h

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

Usage: vector<Angle> $\hat{v} = \{w[0], w[0].t360() ...\};$ // 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 $_{0f0602, 35 \ lines}$

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

CircleTangents.h

"Point.h"

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

circumcircle.h

Description:

"Point.h"

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



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

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

MinimumEnclosingCircle.h

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

```
rep(i,0,sz(ps)) if ((o - ps[i]).dist() > r * EPS) {
  o = ps[i], r = 0;
  rep(j,0,i) if ((o - ps[j]).dist() > r * EPS) {
    o = (ps[i] + ps[j]) / 2;
    r = (o - ps[i]).dist();
  rep(k,0,j) if ((o - ps[k]).dist() > r * EPS) {
    o = ccCenter(ps[i], ps[j], ps[k]);
    r = (o - ps[i]).dist();
    }
  }
}
return {o, r};
}
```

8.3 Polygons

InsidePolygon.h

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

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

"Point.h", "OnSegment.h", "SegmentDistance.h" 2bf504, 11 lines

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

PolygonArea.h

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

"Point.h"

f12300, 6 lines

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

PolygonCenter.h

Description: Returns the center of mass for a polygon.

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

PolygonCut.h Description:

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





ConvexHull.h

${\bf Description:}$

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



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

HullDiameter.h

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

c571b8, 12 lines

```
typedef Point<1l> P;
array<P, 2> hullDiameter(vector<P> S) {
  int n = sz(S), j = n < 2 ? 0 : 1;
  pair<1l, 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

"Point.h", "sideOf.h", "OnSegment.h"

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

Time: $O(\log N)$

```
typedef Point<11> P;

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

```
if (sideOf(1[0], 1[a], p) >= r || sideOf(1[0], 1[b], p) <= -r)
    return false;
while (abs(a - b) > 1) {
    int c = (a + b) / 2;
    (sideOf(1[0], 1[c], p) > 0 ? b : a) = c;
}
return sgn(1[a].cross(1[b], p)) < r;
}</pre>
```

LineHullIntersection.h

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

```
Time: \mathcal{O}\left(N + Q \log n\right)
```

```
"Point.h"
                                                     758f22, 39 lines
typedef array<P, 2> Line;
#define cmp(i, j) sqn(dir.perp().cross(poly[(i)%n]-poly[(j)%n]))
#define extr(i) cmp(i + 1, i) >= 0 && cmp(i, i - 1 + n) < 0
int extrVertex(vector<P>& poly, P dir) {
 int n = sz(poly), lo = 0, hi = n;
 if (extr(0)) return 0;
 while (lo + 1 < hi) {
   int m = (lo + hi) / 2;
   if (extr(m)) return m;
    int ls = cmp(lo + 1, lo), ms = cmp(m + 1, m);
    (1s < ms \mid | (1s == ms \&\& 1s == cmp(1o, m)) ? hi : 1o) = m;
 return lo;
#define cmpL(i) sqn(line[0].cross(poly[i], line[1]))
array<int, 2> lineHull(Line line, vector<P> poly) {
 int endA = extrVertex(poly, (line[0] - line[1]).perp());
 int endB = extrVertex(poly, (line[1] - line[0]).perp());
 if (cmpL(endA) < 0 \mid | cmpL(endB) > 0)
   return {-1, -1};
 array<int, 2> res;
 rep(i,0,2) {
   int lo = endB, hi = endA, n = sz(poly);
   while ((lo + 1) % n != hi) {
     int m = ((lo + hi + (lo < hi ? 0 : n)) / 2) % n;</pre>
      (cmpL(m) == cmpL(endB) ? lo : hi) = m;
    res[i] = (lo + !cmpL(hi)) % n;
    swap (endA, endB);
 if (res[0] == res[1]) return {res[0], -1};
 if (!cmpL(res[0]) && !cmpL(res[1]))
   switch ((res[0] - res[1] + sz(poly) + 1) % sz(poly)) {
     case 0: return {res[0], res[0]};
      case 2: return {res[1], res[1]};
 return res;
```

8.4 Misc. Point Set Problems

ClosestPair.h

71446b, 14 lines

```
Description: Finds the closest pair of points. Time: \mathcal{O}(n \log n)
```

```
"Point.h" d31bbf, 17 lines
typedef Point<1l> 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;
  trav(p, v) {
    P d{1 + (ll)sqrt(ret.first), 0};
    while (v[j].y \le p.y - d.x) S.erase(v[j++]);
    auto lo = S.lower bound(p - d), hi = S.upper bound(p + d);
    for (; lo != hi; ++lo)
      ret = min(ret, {(*lo - p).dist2(), {*lo, p}});
    S.insert(p);
  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; }
bool on_y(const P& a, const P& b) { return a.y < b.y; }
struct Node {
 P pt; // if this is a leaf, the single point in it
  T x0 = INF, x1 = -INF, y0 = INF, y1 = -INF; // bounds
  Node *first = 0, *second = 0;
  T distance (const P& p) { // min squared distance to a point
    T x = (p.x < x0 ? x0 : p.x > x1 ? x1 : p.x);
    T y = (p.y < y0 ? y0 : p.y > y1 ? y1 : p.y);
    return (P(x,y) - p).dist2();
  Node (vector<P>&& vp) : pt(vp[0]) {
    for (P p : vp) {
      x0 = min(x0, p.x); x1 = max(x1, p.x);
      y0 = min(y0, p.y); y1 = max(y1, p.y);
    if (vp.size() > 1) {
      // split on x if the box is wider than high (not best
           heuristic...)
      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:

return make_pair((p - node->pt).dist2(), node->pt);

// if $(p == node \rightarrow pt)$ return $\{INF, P()\}$;

Node *f = node -> first, *s = node -> second;

FastDelaunay PolyhedronVolume Point3D 3dHull

```
T bfirst = f->distance(p), bsec = s->distance(p);
    if (bfirst > bsec) swap(bsec, bfirst), swap(f, s);
    // search closest side first, other side if needed
    auto best = search(f, p);
    if (bsec < best.first)</pre>
      best = min(best, search(s, p));
    return best;
  // find nearest point to a point, and its squared distance
  // (requires an arbitrary operator< for Point)
  pair<T, P> nearest (const P& p) {
    return search(root, p);
};
FastDelaunav.h
Description: Fast Delaunay triangulation. Each circumcircle contains none
of the input points. There must be no duplicate points. If all points are on a
line, no triangles will be returned. Should work for doubles as well, though
there may be precision issues in 'circ'. Returns triangles in order {t[0][0],
t[0][1], t[0][2], \bar{t}[1][0], \dots\}, all counter-clockwise.
Time: \mathcal{O}(n \log n)
"Point.h"
                                                        bf87ec, 88 lines
typedef Point<ll> P;
typedef struct Ouad* O;
typedef __int128_t ll1; // (can be ll if coords are < 2e4)
P arb(LLONG_MAX, LLONG_MAX); // not equal to any other point
struct Quad {
  bool mark; Q o, rot; P p;
  P F() { return r()->p; }
  O r() { return rot->rot; }
  Q prev() { return rot->o->rot; }
  0 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) \starC + p.cross(b,c) \starA + p.cross(c,a) \starB > 0;
Q makeEdge(P orig, P dest) {
  Q q[] = \{new Quad\{0,0,0,oriq\}, new Quad\{0,0,0,arb\},
            new Quad{0,0,0,dest}, new Quad{0,0,0,arb}};
    q[i] \rightarrow o = q[-i \& 3], q[i] \rightarrow rot = q[(i+1) \& 3];
  return *q;
void splice(Q a, Q b) {
  swap(a->o->rot->o, b->o->rot->o); swap(a->o, b->o);
Q connect(Q a, Q b) {
  Q = makeEdge(a->F(), b->p);
  splice(q, a->next());
  splice(q->r(), b);
  return q;
pair<Q,Q> rec(const vector<P>& s) {
  if (sz(s) <= 3) {
    Q = makeEdge(s[0], s[1]), b = makeEdge(s[1], s.back());
    if (sz(s) == 2) return { a, a->r() };
    splice(a->r(), b);
    auto side = s[0].cross(s[1], s[2]);
    Q c = side ? connect(b, a) : 0;
    return {side < 0 ? c->r() : a, side < 0 ? c : b->r() };
```

```
#define H(e) e->F(), e->p
#define valid(e) (e->F().cross(H(base)) > 0)
 Q A, B, ra, rb;
 int half = sz(s) / 2;
 tie(ra, A) = rec({all(s) - half});
 tie(B, rb) = rec({sz(s) - half + all(s)});
 while ((B->p.cross(H(A)) < 0 \&& (A = A->next())) | |
         (A->p.cross(H(B)) > 0 && (B = B->r()->o)));
 Q base = connect (B->r(), A);
 if (A->p == ra->p) ra = base->r();
 if (B->p == rb->p) rb = base;
#define DEL(e, init, dir) O e = init->dir; if (valid(e)) \
    while (circ(e->dir->F(), H(base), e->F())) { \
     Q t = e->dir; \
     splice(e, e->prev()); \
     splice(e->r(), e->r()->prev()); \
     e = t; \
 for (;;) {
   DEL(LC, base->r(), o); DEL(RC, base, prev());
    if (!valid(LC) && !valid(RC)) break;
    if (!valid(LC) || (valid(RC) && circ(H(RC), H(LC))))
     base = connect(RC, base->r());
     base = connect(base->r(), LC->r());
 return { ra, rb };
vector<P> triangulate(vector<P> pts) {
 sort(all(pts)); assert(unique(all(pts)) == pts.end());
 if (sz(pts) < 2) return {};
 Q e = rec(pts).first;
 vector<Q> q = \{e\};
 int qi = 0;
 while (e->o->F().cross(e->F(), e->p) < 0) e = e->o;
#define ADD { Q c = e; do { c->mark = 1; pts.push_back(c->p); \
 q.push_back(c->r()); c = c->next(); } while (c != e); }
 ADD; pts.clear();
 while (qi < sz(q)) if (!(e = q[qi++])->mark) ADD;
 return pts;
       3D
8.5
PolyhedronVolume.h
Description: Magic formula for the volume of a polyhedron. Faces should
point outwards.
                                                     1ec4d3, 6 lines
template<class V, class L>
double signed_poly_volume(const V& p, const L& trilist) {
 double v = 0;
```

```
trav(i, trilist) v += p[i.a].cross(p[i.b]).dot(p[i.c]);
return v / 6:
```

Point3D.h

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

```
template<class T> struct Point3D {
 typedef Point3D P;
 typedef const P& R;
 explicit Point3D(T x=0, T y=0, T z=0) : x(x), y(y), z(z) {}
```

```
bool operator<(R p) const {</pre>
   return tie(x, y, z) < tie(p.x, p.y, p.z); }
 bool operator==(R p) const {
   return tie(x, y, z) == tie(p.x, p.y, p.z); }
 P operator+(R p) const { return P(x+p.x, y+p.y, z+p.z); }
 P operator-(R p) const { return P(x-p.x, y-p.y, z-p.z); }
 P operator*(T d) const { return P(x*d, y*d, z*d); }
 P operator/(T d) const { return P(x/d, y/d, z/d); }
 T dot(R p) const { return x*p.x + y*p.y + z*p.z; }
 P cross(R p) const {
   return P(y*p.z - z*p.y, z*p.x - x*p.z, x*p.y - y*p.x);
 T dist2() const { return x*x + y*y + z*z; }
 double dist() const { return 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(sgrt(x*x+y*y),z); }
 P unit() const { return *this/(T) dist(); } //makes dist()=1
  //returns unit vector normal to *this and p
 P normal(P p) const { return cross(p).unit(); }
  //returns point rotated 'angle' radians ccw around axis
 P rotate (double angle, P axis) const {
    double s = sin(angle), c = cos(angle); P u = axis.unit();
    return u*dot(u)*(1-c) + (*this)*c - cross(u)*s;
};
```

3dHull.h

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

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

```
"Point3D.h"
                                                     c172e9, 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[j];
      if(f.q.dot(A[i]) > f.q.dot(A[f.a])) {
        E(a,b).rem(f.c);
        E(a,c).rem(f.b);
```

E(b,c).rem(f.a);

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

sphericalDistance.h

Description: Returns the shortest distance on the sphere with radius radius between the points with azimuthal angles (longitude) f1 (ϕ_1) and f2 (ϕ_2) from x axis and zenith angles (latitude) t1 (θ_1) and t2 (θ_2) from z axis. All angles measured in radians. The algorithm starts by converting the spherical coordinates to cartesian coordinates so if that is what you have you can use only the two last rows. dx*radius is then the difference between the two points in the x direction and d*radius is the total distance between the points. 611f07, 8 lines

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

Strings (9)

9.1 KMP

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 q = p[i-1];
   while (g \&\& s[i] != s[g]) g = p[g-1];
   p[i] = g + (s[i] == s[g]);
  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;
```

9.2 Zfunc

Zfunc.h

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

Description: z[x] computes the length of the longest common prefix of s[i:]and s, except z[0] = 0. (abacaba -> 0010301)

3ae52<u>6</u>, 12 lines vi Z(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

Manacher.h

Description: For each position in a string, computes p[0][i] = half length of longest even palindrome around pos i, p[1][i] = longest odd (half rounded down).

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

9.4 MinRotation

MinRotation.h

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

```
int min rotation(string s) {
 int a=0, N=sz(s); s += s;
 rep(b,0,N) rep(i,0,N) {
   if (a+i == b || s[a+i] < s[b+i]) {b += max(0, i-1); break;}</pre>
   if (s[a+i] > s[b+i]) { a = b; break; }
 return a;
```

SuffixArray 9.5

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

```
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)), rank(n);
    sa = lcp = y, iota(all(sa), 0);
    for (int j = 0, p = 0; p < n; j = max(1, j * 2), lim = p) {
     p = j, iota(all(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++]] = k)
     for (k \&\& k--, j = sa[rank[i] - 1];
         s[i + k] == s[j + k]; k++);
};
```

SuffixTree

memset(s, 0, sizeof s);

SuffixTree.h

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

```
Time: \mathcal{O}(26N)
                                                     aae0b8, 50 lines
struct SuffixTree {
  enum { N = 200010, ALPHA = 26 }; //N \sim 2*maxlen+10
 int toi(char c) { return c - 'a'; }
 string a; //v = cur \ node, q = cur \ position
 int t[N][ALPHA], 1[N], r[N], p[N], s[N], v=0, q=0, m=2;
  void ukkadd(int i, int c) { suff:
    if (r[v]<=q) {
      if (t[v][c]==-1) { t[v][c]=m; l[m]=i;
        p[m++]=v; v=s[v]; q=r[v]; goto suff; }
      v=t[v][c]; q=l[v];
    if (q==-1 || c==toi(a[q])) q++; else {
     l[m+1]=i; p[m+1]=m; l[m]=l[v]; r[m]=q;
      p[m]=p[v]; t[m][c]=m+1; t[m][toi(a[q])]=v;
      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+=r[v]-l[v]; }</pre>
      if (q==r[m]) s[m]=v; else s[m]=m+2;
      q=r[v]-(q-r[m]); m+=2; qoto suff;
 SuffixTree(string a) : a(a) {
    fill(r,r+N,sz(a));
```

Hashing AhoCorasick IntervalContainer IntervalCover

```
memset(t, -1, sizeof t);
    fill(t[1],t[1]+ALPHA,0);
    s[0] = 1; 1[0] = 1[1] = -1; r[0] = r[1] = p[0] = p[1] = 0;
    rep(i,0,sz(a)) ukkadd(i, toi(a[i]));
  // example: find longest common substring (uses ALPHA = 28)
  pii best;
  int lcs(int node, int i1, int i2, int olen)
    if (l[node] <= i1 && i1 < r[node]) return 1;</pre>
    if (1[node] <= i2 && i2 < r[node]) return 2;</pre>
    int mask = 0, len = node ? olen + (r[node] - l[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

rep(i,length,sz(str)) {

Hashing.h

```
Description: Self-explanatory methods for string hashing.
```

```
// Arithmetic mod 2^64-1. 2x slower than mod 2^64 and more
// code, but works on evil test data (e.g. Thue-Morse, where
// ABBA... and BAAB... of length 2^10 hash the same mod 2^64).
// "typedef ull H;" instead if you think test data is random,
// or work mod 10^9+7 if the Birthday paradox is not a problem.
struct H {
  typedef uint64_t ull;
  ull x; H(ull x=0) : x(x) {}
#define OP(O,A,B) H operator O(H o) { ull r = x; asm \
  (A "addg %%rdx, %0\n adcg $0,%0" : "+a"(r) : B); return r; }
  OP(+,,"d"(o.x)) OP(*,"mul %1\n", "r"(o.x) : "rdx")
  H operator-(H o) { return *this + ~o.x; }
  ull get() const { return x + !~x; }
  bool operator==(H o) const { return get() == o.get(); }
  bool operator<(H o) const { return get() < o.get(); }</pre>
static const H C = (11)1e11+3; // (order ~ 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};
```

```
ret.push_back(h = h * C + str[i] - pw * str[i-length]);
 return ret;
H hashString(string& s) { H h{}; trav(c,s) h=h*C+c; return h; }
```

9.8 AhoCorasick

AhoCorasick.h

Description: Aho-Corasick tree is used for multiple pattern matching. Initialize the tree with create(patterns). 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. **Time:** create is $\mathcal{O}(26N)$ where N is the sum of length of patterns. find is $\mathcal{O}(M)$ where M is the length of the word. findAll is $\mathcal{O}(NM)$. $_{716ac4, 67 \text{ lines}}$

```
struct AhoCorasick {
 enum {alpha = 26, first = 'A'};
 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;
 vector<int> backp;
 void insert(string& s, int i) {
   assert(!s.emptv());
   int n = 0;
   trav(c, s) {
     int& m = N[n].next[c - first];
     if (m == -1) { n = m = sz(N); N.emplace_back(-1); }
   if (N[n].end == -1) N[n].start = j;
   backp.push back(N[n].end);
   N[n].end = j;
   N[n].nmatches++;
 AhoCorasick(vector<string>& pat) {
   N.emplace_back(-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;
   trav(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)
set<pii>::iterator addInterval(set<pii>& is, int L, int R) {
 if (L == R) return is.end();
  auto it = is.lower_bound({L, R}), before = it;
  while (it != is.end() && it->first <= R) {</pre>
    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: $\mathcal{O}(N \log N)$

```
template<class T>
```

```
vi cover(pair<T, T> G, vector<pair<T, T>> I) {
 vi S(sz(I)), R;
 iota(all(S), 0);
 sort(all(S), [&](int a, int b) { return I[a] < I[b]; });</pre>
 T cur = G.first;
 int at = 0;
```



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

ConstantIntervals.h

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

```
Usage: constantIntervals(0, sz(v), [&](int x){return v[x];},
[&] (int lo, int hi, T val) \{\ldots\});
Time: \mathcal{O}\left(k\log\frac{n}{k}\right)
                                                                    753a4c, 19 lines
```

```
template<class F, class G, class T>
void rec(int from, int to, F& f, G& g, int& i, T& p, T q) {
  if (p == q) return;
  if (from == to) {
   g(i, to, p);
    i = to; p = q;
    int mid = (from + to) >> 1;
    rec(from, mid, f, g, i, p, f(mid));
   rec(mid+1, to, f, g, i, p, q);
template < class F, class G>
void constantIntervals(int from, int to, F f, G g) {
 if (to <= from) return;</pre>
  int i = from; auto p = f(i), q = f(to-1);
  rec(from, to-1, f, q, 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, 13 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)
     a = mid;
    else
     b = mid+1;
  rep(i,a+1,b+1) if (f(a) < f(i)) a = i; // (B)
 return a:
```

```
Description: Compute indices for the longest increasing subsequence.
Time: \mathcal{O}(N \log N)
```

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

10.3 Dynamic programming

KnuthDP.h

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

DivideAndConquerDP.h

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

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

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

10.4 Debugging tricks

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

10.5 Optimization tricks

10.5.1 Bit hacks

5b77eb, 17 lines

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

10.5.2 Pragmas

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

FastMod.h

Description: Compute a%b about 4 times faster than usual, where b is constant but not known at compile time. Fails for b = 1. c977c5, 10 lines

```
typedef unsigned long long ull;
typedef __uint128_t L;
struct FastMod {
 ull b, m;
 FastMod(ull b) : b(b), m(ull((L(1) << 64) / b)) {}
 ull reduce(ull a) {
   ull q = (ull) ((L(m) * a) >> 64), r = a - q * b;
   return r >= b ? r - b : r;
};
```



```
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 some-
thing like 0.05us + 16 bytes per allocation.
```

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

BumpAllocatorSTL.h

Description: BumpAllocator for STL containers.

```
Usage: vector<vector<int, small<int>>> ed(N);
                                                           bb66d<u>4, 14 lines</u>
```

```
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) {}
```

Unrolling.h

520e76, 5 lines

```
#define F { . . . ; ++i; }
int i = from;
while (i&3 && i < to) F // for alignment, if needed
while (i + 4 <= to) { F F F F }
while (i < to) F
```

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/qcc/*/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/s-551b82, 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)
// i32 qather\_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 lanes
// 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(qt|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; 11 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 {ll 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[i]; // <- equiv
  return r;
```