

International Institute Of Information Technology - Hyderabad

tesla_protocol

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		Ü	template.cpp 27 lines		
5	Number theory	9	<pre>#include <bits stdc++.h=""></bits></pre>	se	
_	5.1 Modular arithmetic	9	<pre>using namespace std;</pre>	if	
	5.2 Primality	9	#define REP(i, a, b) for(int i = (a); i < (b); ++i)		
	5.3 Divisibility	10	#define ALL(x) begin(x), end(x)	el	
	5.4 Fractions	10	<pre>#define SZ(x) ((int)(x).size()) #define SET(a,v) memset((a), (v), sizeof(a))</pre>	``	
	5.5 Pythagorean Triples	10	#define PB push_back	er	
	5.6 Primes	10	#define EB emplace_back #define MP make_pair	if	
	5.7 Estimates	10	#define fir make_pail #define fst first	er	
	5.8 Mobius Function	10	#define snd second	se	
	5.6 Mobius Function	10	<pre>using ll = long long; using dbl = double;</pre>	fo	
6	Graph	10	using PII = pair <int, int="">;</int,>		
U	6.1 Fundamentals	11	<pre>using VI = vector<int>;</int></pre>	er	
	6.2 Network flow	11	<pre>int main() {</pre>	$ _{\rm C}$	
	6.3 Matching	12	#ifdef LOCAL_EXEC		
	6.4 DFS algorithms	13	<pre>// freopen("sample.in", "r", stdin); // freopen("sample.out" "v" stdout);</pre>		
	6.5 Coloring	14	<pre>// freopen("sample.out", "w", stdout); #else</pre>		
	6.6 Heuristics	14	<pre>cin.tie(0)->sync_with_stdio(0);</pre>		
	6.7 Trees	15	<pre>cin.exceptions(cin.failbit); #endif</pre>		
	6.8 Math	16	return 0;		
	2.20011	10	}		
7	Strings	16		$ \mathbf{S} $	
•	~~		extras.cpp 32 lines		
8	Geometry	18	#define trace() {f(#VA_ARGS,VA_ARGS); }		
	8.1 Geometric primitives	18	<pre>template<typename arg=""> voidf(const char* name, Arg&& arg) { cerr << name << " = " << arg << endl;</typename></pre>		
	8.2 Circles	19	}		
	8.3 Polygons	19	template <typename arg1,="" args="" typename=""> void f(const char* names, Arg1&& arg1, Args&& args) {</typename>		
	8.4 Misc. Point Set Problems	20	const char* comma = strchr(names + 1, ',');		
	8.5 3D	21	cerr.write(names, comma - names) << " = " << arg1<<" ";		
			f(comma+1, args);		
9	Mathematics	22			
	9.1 Equations	22	<pre>#include <ext assoc_container.hpp="" pb_ds=""> #include <ext pb_ds="" tree_policy_hpp=""></ext></ext></pre>		
	9.2 Recurrences	22	<pre>#include <ext pb_ds="" tree_policy.hpp=""> using namespace gnu pbds; // OST: find by order, order of key</ext></pre>		
	9.3 Trigonometry	22	<pre>typedef tree<int, less<="" null_type,="">, rb_tree_tag,</int,></pre>		
	9.4 Geometry	22	<pre>tree_order_statistics_node_update> ordered_set; // #pragma GCC optimize("Ofast")</pre>		
	9.5 Derivatives/Integrals	22	// optimize("unroll-loops")		
	•		•		

'MakeLists.txt

```
18 lines
nake minimum required(VERSION 3.15)
roject(contest)
et(CMAKE CXX STANDARD 17)
et(CMAKE_CXX_FLAGS, "-std=c++17 -Wall -pedantic -Wconversion \
-Wshadow -Wfloat-equal")
 (DEBUG EXEC) # set in 'Debug'
set(CMAKE CXX FLAGS "${CMAKE CXX FLAGS} -00 -g -fsanitize=
     address, undefined")
set(CMAKE CXX FLAGS "${CMAKE CXX FLAGS} -02")
ndif()
 (LOCAL EXEC) # set in 'Debug' and 'Run'
set(CMAKE CXX FLAGS "${CMAKE CXX FLAGS} -DLOCAL EXEC")
et(CMAKE RUNTIME OUTPUT DIRECTORY ..) # folder with source
et(srcs a b c d e f q h i j k l m) # adjust names as necessary
reach(F IN LISTS srcs)
add executable(${F}.out ${F}.cpp)
ndforeach()
```

CLion setup:

- File \rightarrow New Project \rightarrow C++ executable
- File \rightarrow Settings \rightarrow Keymap \rightarrow Sublime Text
- Settings \rightarrow Build... \rightarrow CMake
- Add profiles: Debug, Run, OJ; use default toolchain.

Sublime Keybindings

- Add Selection for Next Occurrence: Ctrl-D
- Move Line Up: Ctrl-Shift-UP
- Move Line Down: Ctrl-Shift-DOWN
- Delete Line: Ctrl-Shift-K
- Duplicate Line or Selection: Ctrl-Shift-D
- Start New Line: Ctrl+Enter
- Start New Line Before Current: Ctrl+Shift+Enter
- Move Caret to Matchcing Brace: Ctrl-M
- Terminal: Ctrl-T

.vimrc

set cin aw ai is ts=2 sw=2 tm=50 nu noeb bg=dark ru cul sy on \mid im jk <esc> " Select region and then type :Hash to hash your selection. " Useful for verifying that there aren't mistypes. ca Hash w !cpp -dD -P -fpreprocessed \| tr -d '[:space:]' \ \| md5sum \| cut -c-6

hash.sh

Hashes a file, ignoring all whitespace and comments. Use for # verifying that code was correctly typed. cpp -dD -P -fpreprocessed | tr -d '[:space:]'| md5sum |cut -c-6

Combinatorial (2)

2.1 Permutations

2.1.1 Factorial

			-		8	-	10	
							3628800	
						16		
n!	4.0e7	′ 4.8e	8 6.2e	9 8.7e	10 1.3e	12 2.1e	l3 3.6e14	
n	20	25	30	40	50 10	00 15	0 171	
n!	2e18	2e25	3e32	$8e47 \ 3$	e64 9e	157 6e2	62 >dbl_ma	X

IntPerm.h

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

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

261010, 6 lines

2.1.2 Cycles

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

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

2.1.3 Derangements

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

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

2.1.4 Burnside's lemma

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

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

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

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

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

2.2 Partitions and subsets

2.2.1 Partition function

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

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

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

2.2.2 Lucas' Theorem

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

2.2.3 Binomials

multinomial.h

return c;

Description: Computes $\binom{k_1 + \dots + k_n}{k_1, k_2, \dots, k_n} = \frac{(\sum k_i)!}{k_1! k_2! \dots k_n!}$. 037a49, 6 lines limultinomial(VI& v) { ll c = 1, m = v.empty() ? 1 : v[0]; REP(i,1,5Z(v)) REP(j,0,v[i]) c = c * ++m / (i+1);

.3 General purpose numbers

2.3.1 Bernoulli numbers

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

Sums of powers:

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

Euler-Maclaurin formula for infinite sums:

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

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

2.3.2 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 $c(n,2) = 0, 0, 1, 3, 11, 50, 274, 1764, 13068, 109584, \dots$

2.3.3 Eulerian numbers

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

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

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

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

2.3.4 Stirling numbers of the second kind

Partitions of n distinct elements into exactly k groups.

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

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

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

2.3.5 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,

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

2.3.6 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 : $(n-2)!/((d_1-1)!\cdots(d_n-1)!)$

2.3.7 Catalan numbers

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

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

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

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

```
2.4 DP Optimizations
1D-1D.h
Description: Applicable if dp_i = min_{i>i}(dp_i + cost(i, j) \text{ s.t. } opt_i \leq opt_j
when i < j (which holds if quadrangle)
Time: FIXME
                                                       d76777, 27 lines
#define until first
#define opt second
11 dp[100000];
ll cost(int i, int j) {
  return dp[j] /* + cost to jump from i to j*/;
void solve(int n) {
  dp[n] = 0;
  vector<PII> v;
  v.EB(n - 1, n);
  for (int i = n - 1, ipos = 0; i >= 1; i--) {
    while (ipos+1 < SZ(v) && i <= v[ipos+1].until) ipos++;
    dp[i] = cost(i, v[ipos].opt);
    while (v.back().until < i</pre>
      && cost(v.back().until, i)
         <= cost(v.back().until, v.back().opt))
      { v.pop back(); }
    int l = 1, r = min(i - 1, v.back().until);
    while (l <= r) {
      int mid = (l + r)/2;
      if (cost(mid, i) <= cost(mid, v.back().opt)) {</pre>
        l = mid + 1;
      } else { r = mid - 1; }
    if (l - 1 >= 1) v.EB(l - 1, i);
Divide-and-Conquer.h
Description: Works when dp_{k,i} = min_{j < i}(dp_{k-1,j} + cost(j,i)) and
opt_k(i) \leq opt_k(i+1). (This holds when quadrangle)
call solve(i, 1, n, 1, n) for i = 2 to n after finding dp_1
Time: FIXME
                                                       1db0cf, 15 lines
ll dp[100][100]; // set correctly
ll cost(int i, int j); // cost to go from i to j, 1-indexed.
void solve(int i, int l, int r, int optl, int optr) {
  const ll inf = 1e18; // set correctly
  if (l > r || optl > optr) return;
  int mid = (l + r)/2; pair<ll, int> best = \{inf, -1\};
  for (int j = optl; j <= min(mid, optr); j++) {</pre>
    pair<ll,int> cand(dp[i - 1][j] + cost(j, mid), j);
    if (best.second == -1) best = cand;
    else best = min(best, cand):
  dp[i][mid] = best.first;
  solve(i, l, mid - 1, optl, best.second);
  solve(i, mid + 1, r, best.second, optr);
Li-Chao.h
Description: FIXME
Time: FIXME
                                                       b35345, 19 lines
int n, tree[1 << 16]; // set limits correctly</pre>
int eval(int f, int x); // evaluate f(x)
```

// Add func f to set which crosses every other func atmost once

```
void update(int f, int root = 1, int s = 1, int e = n) {
 int mid = (s + e)/2:
  bool lless = eval(f, s) < eval(tree[root], s);</pre>
  bool mless = eval(f, mid) < eval(tree[root], mid);</pre>
  if (mless) swap(f, tree[root]);
  if (s == e) return:
  if (lless != mless) update(f, 2*root, s, mid);
  else update(f, 2*root + 1, mid + 1, e);
// Evaluate mininum at x over all functions.
Il query(int x, int root = 1, int s = 1, int e = n) {
  int mid = (s + e)/2; ll ans = eval(tree[root], x);
  if (s == e) return ans;
  if (x <= mid) return min(query(x, 2*root, s, mid), ans);</pre>
  return min(query(x, 2*root + 1, mid + 1, e), ans);
Dynamic-CHT.h
Description: FIXME
Time: FIXME
                                                     f03adc, 55 lines
const dbl INF = 1e16:
struct HullDynamic {
#define CLREF const Line&
  struct Line {
    int a; ll b, val=0; dbl xLeft = -INF; bool type=0;
    Line(int a = 0, ll b = 0): a(a), b(b) {}
    ll eval(int x) const{ return a * 1ll * x + b; }
    bool operator< (CLREF l2) const {</pre>
      return l2.type ? (xLeft > l2.val) : (a < l2.a);
  using Iter = set<Line>::iterator;
  bool parallel(CLREF l1, CLREF l2) { return l1.a == l2.a; }
  dbl meetX(CLREF l1, CLREF l2) {
    return parallel(l1, l2) ? INF :
      (l2.b-l1.b) / (dbl(l1.a-l2.a));
  set<Line> hull;
  bool hasPrev(Iter it) { return it != hull.begin(); }
  bool hasNext(Iter it) {
    return it != hull.end() && next(it) != hull.end();
  bool bad(CLREF l1, CLREF l2, CLREF l3){
    return meetX(l1,l3) <= meetX(l1,l2);</pre>
  bool bad(Iter it) {
    return hasPrev(it) && hasNext(it)
        && (bad(*next(it), *it, *prev(it)));
  Iter upd left border(Iter it) {
    if(!hasNext(it)) return it;
    dbl val = meetX(*it, *next(it));
    Line buf(*it); it = hull.erase(it);
    buf.xLeft = val; return hull.insert(it, buf);
  void insert line(int a, ll b) {
    Line l3 = Line(a, b); auto it = hull.lower bound(l3);
    if (it != hull.end() && parallel(*it , 13) {
      if (it->b <= b) return;</pre>
      it = hull.erase(it):
    it = hull.insert(it, l3);
    if (bad(it)) { hull.erase(it); return; }
    while (hasPrev(it) && bad(prev(it))) hull.erase(prev(it));
    while (hasNext(it) && bad(next(it))) hull.erase(next(it));
    it = upd left border(it);
    if (hasPrev(it)) upd left border(prev(it));
    if (hasNext(it)) upd left border(next(it));
```

```
ll eval(int x) {
    Line q; q.val = x; q.type = 1;
    auto best = hull.lower bound(q);
    return (best == hull.end()) ? INF : best->eval(x);
};
```

Data structures (3)

OrderStatisticTree.h

Description: A set (not multiset!) with support for finding the n'th element, and finding the index of an element. To get a map, change null_type. Time: $\mathcal{O}(\log N)$ 782797, 16 lines

```
#include <bits/extc++.h>
template<class T>
using Tree = tree<T, null_type, less<T>, rb_tree_tag,
    tree order statistics node update>;
void example() {
 Tree<int> t, t2; t.insert(8);
  auto it = t.insert(10).first;
  assert(it == t.lower bound(9)):
  assert(t.order_of_key(10) == 1);
  assert(t.order_of_key(11) == 2);
  assert(*t.find\ by\ order(0) == 8);
 t.join(t2); // assuming T < T2 or T > T2, merge t2 into t
```

HashMap.h

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

```
#include <bits/extc++.h>
// To use most bits rather than just the lowest ones:
struct chash { // large odd number for C
 const uint64 t C = 11(4e18 * acos(0)) | 71;
 ll operator()(ll x) const { return builtin bswap64(x*C); }
gnu pbds::gp hash table<ll,int,chash> h({},{},{},{},{1<<16});</pre>
```

SegmentTree.h

Description: Zero-indexed max-tree. Bounds are inclusive to the left and exclusive to the right. Can be changed by modifying T, f and unit. Time: $\mathcal{O}(\log N)$

```
0f4bdb, 19 lines
struct Tree {
  typedef int T;
 static constexpr T unit = INT MIN;
 T f(T a, T b) { return max(a, b); } // (any associative fn)
 vector<T> s: int n:
 Tree(int n = 0, T def = unit) : s(2*n, def), n(n) {}
  void update(int pos, T val) {
   for (s[pos += n] = val; pos /= 2;)
      s[pos] = f(s[pos * 2], s[pos * 2 + 1]);
 T query(int b, int e) { // query [b, e)
   T ra = unit, rb = unit;
    for (b += n, e += n; b < e; b /= 2, e /= 2) {
      if (b \% 2) ra = f(ra, s[b++]);
      if (e \% 2) rb = f(s[--e], rb);
    return f(ra, rb);
```

```
};
LazySegmentTree.h
Description: Segment tree with ability to add or set values of large inter-
vals, and compute max of intervals. Can be changed to other things. Use
with a bump allocator for better performance, and SmallPtr or implicit in-
dices to save memory.
Usage: Node* tr = new Node(v, 0, SZ(v));
Time: \mathcal{O}(\log N).
"../various/BumpAllocator.h"
                                                        f49770, 50 lines
const int inf = 1e9:
struct Node {
 Node *l = 0. *r = 0:
  int lo, hi, mset = inf, madd = 0, val = -inf;
  Node(int lo,int hi):lo(lo),hi(hi){} // Large interval of -inf
  Node(VI& v, int lo, int hi) : lo(lo), hi(hi) {
    if (lo + 1 < hi)
      int mid = lo + (hi - lo)/2;
      l = new Node(v, lo, mid); r = new Node(v, mid, hi);
      val = max(l->val, r->val);
    else val = v[lo];
  int query(int L, int R) {
    if (R <= lo || hi <= L) return -inf;</pre>
    if (L \le lo \&\& hi \le R) return val:
    push();
    return max(l->query(L, R), r->query(L, R));
  void set(int L, int R, int x) {
    if (R <= lo || hi <= L) return:</pre>
    if (L \le lo \&\& hi \le R) mset = val = x, madd = 0;
      push(), l->set(L, R, x), r->set(L, R, x);
      val = max(l->val, r->val);
  void add(int L, int R, int x) {
    if (R <= lo || hi <= L) return;</pre>
    if (L \le lo \&\& hi \le R)
      if (mset != inf) mset += x;
      else madd += x;
      val += x;
    else |
      push(), l->add(L, R, x), r->add(L, R, x);
      val = max(l->val, r->val);
  void push() {
    if (!1) {
      int mid = lo + (hi - lo)/2;
      l = new Node(lo, mid); r = new Node(mid, hi);
    if (mset != inf)
     l->set(lo,hi,mset), r->set(lo,hi,mset), mset = inf;
    else if (madd)
     l->add(lo,hi,madd), r->add(lo,hi,madd), madd = 0;
};
UnionFindRollback.h
skip st, time() and rollback().
Usage: int t = uf.time(); ...; uf.rollback(t);
Time: \mathcal{O}(\log(N))
```

```
Description: Disjoint-set data structure with undo. If undo is not needed,
```

f79afc, 21 lines

```
struct RollbackUF {
```

```
VI e; vector<PII> st;
  RollbackUF(int n) : e(n, -1) {}
  int size(int x) { return -e[find(x)]; }
  int find(int x) { return e[x] < 0 ? x : find(e[x]); }</pre>
  int time() { return SZ(st); }
  void rollback(int t) {
    for (int i = time(); i --> t;)
      e[st[i].first] = st[i].second;
    st.resize(t);
  bool join(int a, int b) {
    a = find(a), b = find(b);
    if (a == b) return false;
    if (e[a] > e[b]) swap(a, b);
    st.push back({a, e[a]});
    st.push back({b, e[b]});
    e[a] += e[b]; e[b] = a;
    return true;
};
SubMatrix.h
Description: Calculate submatrix sums quickly, given upper-left and lower-
right corners (half-open).
Usage: SubMatrix<int> m(matrix);
m.sum(0, 0, 2, 2); // top left 4 elements Time: \mathcal{O}\left(N^2+Q\right)
                                                        fe231e, 13 lines
template<class T>
struct SubMatrix {
  vector<vector<T>> p;
  SubMatrix(vector<vector<T>>& v) {
    int R = SZ(v), C = SZ(v[0]);
    p.assign(R+1, vector<T>(C+1));
    REP(r, 0, R) REP(c, 0, C)
      p[r+1][c+1] = v[r][c] + p[r][c+1] + p[r+1][c] - p[r][c];
  T sum(int u, int l, int d, int r) {
    return p[d][r] - p[d][l] - p[u][r] + p[u][l];
};
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;
                                                       d8e33f, 26 lines
template<class T, int N> struct Matrix {
  typedef Matrix M:
  array<array<T, N>, N> d{};
  M operator*(const M& m) const {
    REP(i,0,N) REP(i,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;

```
p >>= 1:
    return a;
};
```

LineContainer.h

Description: Container where you can add lines of the form kx+m, and query maximum values at points x. Useful for dynamic programming ("convex hull trick"). Time: $\mathcal{O}(\log N)$

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

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 *l = 0, *r = 0;
 int val, v, 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(l) + cnt(r) + 1; }
template<class F> void each(Node* n, F f) {
 if (n) { each(n->l, f); f(n->val); each(n->r, f); }
pair<Node*, Node*> split(Node* n, int k) {
 if (!n) return {};
 if (cnt(n->l) >= k) { // "n->val >= k" for lower bound(k)}
    auto pa = split(n->l, k);
   n->l = pa.second;
   n->recalc();
```

```
return {pa.first, n};
  } else {
    auto pa = split(n->r, k - cnt(n->l) - 1); // and just "k"
   n->r = pa.first:
   n->recalc();
    return {n, pa.second};
Node* merge(Node* l. Node* r) {
  if (!l) return r;
  if (!r) return l;
  if (l->y > r->y) {
   l->r = merge(l->r, r);
   l->recalc();
    return l;
  } else {
    r->l = merge(l, r->l);
    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 l, int r, int k) {
 Node *a, *b, *c;
  tie(a,b) = split(t, l); tie(b,c) = split(b, r - l);
  if (k \le l) t = merge(ins(a, b, k), c);
  else t = merge(a, ins(c, b, k - r));
FenwickTree.h
Description: Computes partial sums a[0] + a[1] + ... + a[pos - 1], and
updates single elements a[i], taking the difference between the old and new
Time: Both operations are \mathcal{O}(\log N).
```

```
b06af0, 22 lines
struct FT
  vector<ll> s:
  FT(int n) : s(n) {}
  void update(int pos, ll dif) { // a[pos] += dif
    for (; pos < SZ(s); pos |= pos + 1) s[pos] += dif;
  ll query(int pos) { // sum of values in [0, pos)
   11 \text{ res} = 0;
   for (; pos > 0; pos &= pos - 1) res += s[pos-1];
    return res:
  int lower bound(ll sum) {// min pos st sum of [0, pos] >= sum
    // Returns n if no sum is >= sum, or -1 if empty sum is.
   if (sum <= 0) return -1;
   int pos = 0:
    for (int pw = 1 << 25; pw; pw >>= 1) {
     if (pos + pw \leq SZ(s) && s[pos + pw-1] \leq sum)
       pos += pw, sum -= s[pos-1];
    return pos;
```

FenwickTree2d.h

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

```
Time: \mathcal{O}(\log^2 N). (Use persistent segment trees for \mathcal{O}(\log N).)
"FenwickTree.h"
                                                        d5c1b7, 22 lines
  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);
    for (VI& v : ys) sort(ALL(v)), ft.emplace back(SZ(v));
  int ind(int x, int y) {
    return (int)(lower bound(ALL(ys[x]), y) - ys[x].begin()); }
  void update(int x, int y, ll dif) {
    for (; x < SZ(ys); x = x + 1)
      ft[x].update(ind(x, y), dif);
  ll query(int x, int y) {
    ll sum = 0;
    for (; x; x \&= x - 1)
      sum += ft[x-1].query(ind(x-1, y));
    return sum:
};
RMQ.h
Description: Range Minimum Queries on an array. Returns min(V[a], V[a
+ 1], ... V[b - 1]) in constant time.
Usage: RMQ rmq(values);
rmq.query(inclusive, exclusive);
Time: \mathcal{O}(|V|\log|V|+Q)
                                                        9a1bbf, 16 lines
template<class T>
struct RMQ {
  vector<vector<T>> imp:
  RMO(const vector<T>& V) : jmp(1, V)
    for (int pw = 1, k = 1; pw * 2 \le SZ(V); pw *= 2, ++k) {
      jmp.emplace_back(SZ(V) - pw * 2 + 1);
      REP(j,0,SZ(jmp[k]))
        jmp[k][j] = min(jmp[k - 1][j], jmp[k - 1][j + pw]);
 T query(int a, int b) {
    assert(a < b); // or return inf if a == b
    int dep = 31 - builtin clz(b - a);
    return min(jmp[dep][a], jmp[dep][b - (1 << dep)]);</pre>
};
MoQueries.h
Description: Answer interval or tree path queries by finding an approxi-
mate TSP through the queries, and moving from one query to the next by
adding/removing points at the ends. If values are on tree edges, change step
to add/remove the edge (a, c) and remove the initial add call (but keep in).
Time: \mathcal{O}(N\sqrt{Q})
void add(int ind, int end) { ... } // add a[ind] (end = 0 or 1)
void del(int ind, int end) { ... } // remove a[ind]
int calc() { ... } // compute current answer
VI mo(vector<PII> 0) {
  int L = 0, R = 0, blk = 350; // \sim N/sqrt(Q)
 VI s(SZ(Q)), res = s;
#define K(x) PII(x.first/blk, x.second ^ -(x.first/blk & 1))
  iota(ALL(s), 0);
                                                                       Time: \mathcal{O}\left(n^2\log(1/\epsilon)\right)
  sort(ALL(s), [\&](int s, int t) \{ return K(Q[s]) < K(Q[t]); \});
                                                                       "Polynomial.h"
  for (int qi : s) {
                                                                       vector<double> polyRoots(Poly p, double xmin, double xmax) {
    PII q = Q[qi];
    while (L > q.first) add(--L, 0);
                                                                         if (SZ(p.a) == 2) { return {-p.a[0]/p.a[1]}; }
```

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

Numerical (4)

4.1 Polynomials and recurrences

```
Polynomial.h
                                                     2ad8d0, 17 lines
struct Polv {
  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: polyRoots(\{\{2,-3,1\}\},-1e9,1e9) // solve x^2-3x+2=0
```

```
vector<double> ret;
  Poly der = p;
  der.diff();
  auto dr = polyRoots(der, xmin, xmax);
  dr.push back(xmin-1);
  dr.push back(xmax+1);
  sort(ALL(dr));
  REP(i,0,SZ(dr)-1) {
    double l = dr[i], h = dr[i+1];
    bool sign = p(l) > 0;
    if (sign ^{\circ} (p(h) > 0)) {
      REP(it,0,60) { // while (h - l > 1e-8)
        double m = (l + h) / 2, f = p(m);
        if ((f \le 0) \land sign) l = m;
        else h = m;
      ret.push back((l + h) / 2);
  return ret;
PolyInterpolate.h
Description: Given n points (x[i], y[i]), computes an n-1-degree polynomial
p that passes through them: p(x) = a[0] * x^0 + ... + a[n-1] * x^{n-1}. For
numerical precision, pick x[k] = c * \cos(k/(n-1)*\pi), k = 0 \dots n-1.
Time: \mathcal{O}(n^2)
typedef vector<double> vd;
vd interpolate(vd x, vd y, int n) {
  vd res(n), temp(n);
  REP(k,0,n-1) REP(i,k+1,n)
   y[i] = (y[i] - y[k]) / (x[i] - x[k]);
  double last = 0; temp[0] = 1;
  REP(k,0,n) REP(i,0,n) {
    res[i] += v[k] * temp[i];
    swap(last, temp[i]);
    temp[i] -= last * x[k];
  return res:
BerlekampMassev.h
Description: Recovers any n-order linear recurrence relation from the first
2n terms of the recurrence. Useful for guessing linear recurrences after brute-
forcing the first terms. Should work on any field, but numerical stability for
floats is not guaranteed. Output will have size \leq n.
Usage: berlekampMassey(\{0, 1, 1, 3, 5, 11\}) // \{1, 2\}
Time: \mathcal{O}(N^2)
"../number-theory/ModPow.h"
                                                        6120d6, 20 lines
vector<ll> berlekampMassey(vector<ll> s) {
  int n = SZ(s), L = 0, m = 0;
  vector<ll> C(n), B(n), T;
  C[0] = B[0] = 1;
 ll b = 1;
  REP(i,0,n) \{ ++m;
   ll d = s[i] \% mod;
    REP(j,1,L+1) d = (d + C[j] * s[i - j]) % mod;
    if (!d) continue;
   T = C; 11 coef = d * modpow(b, mod-2) % mod;
    REP(j,m,n) C[j] = (C[j] - coef * B[j - m]) % mod;
   if (2 * L > i) continue;
   L = i + 1 - L; B = T; b = d; m = 0;
```

C.resize(L + 1); C.erase(C.begin());

for (ll& x : C) x = (mod - x) % mod;

```
return C;
LinearRecurrence.h
Description: Generates the k'th term of an n-order linear recurrence
S[i] = \sum_{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)
                                                        f72958, 26 lines
typedef vector<ll> Poly;
ll linearRec(Poly S, Poly tr, ll k) {
 int n = SZ(tr);
  auto combine = [&](Poly a, Poly b) {
    Poly res(n * 2 + 1);
    REP(i,0,n+1) REP(j,0,n+1)
      res[i + j] = (res[i + j] + a[i] * b[j]) % mod;
    for (int i = 2 * n; i > n; --i) REP(j,0,n)
      res[i - 1 - j] = (res[i - 1 - j] + res[i] * tr[j]) % mod;
    res.resize(n + 1);
    return res:
  Poly pol(n + 1), e(pol);
  pol[0] = e[1] = 1;
  for (++k; k; k /= 2) {
    if (k % 2) pol = combine(pol, e);
    e = combine(e, e):
  ll res = 0:
  REP(i,0,n) res = (res + pol[i + 1] * S[i]) % mod;
  return res:
       Optimization
4.2
GoldenSectionSearch.h
Description: Finds the argument minimizing the function f in the inter-
val [a, b] assuming f is unimodal on the interval, i.e. has only one local
minimum. The maximum error in the result is eps. Works equally well for
maximization with a small change in the code. See TernarySearch.h in the
Various chapter for a discrete version.
Usage: double func(double x) { return 4+x+.3*x*x; }
double xmin = qss(-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);
    } else {
      a = x1; x1 = x2; f1 = f2;
      x2 = a + r*(b-a); f2 = f(x2);
  return a;
Description: Poor man's optimization for unimodal functions. fff601, 14 lines
typedef array<double, 2> P;
template < class F > pair < double, P > hillClimb(P start, F f) {
```

```
pair<double, P> cur(f(start), start);
  for (double jmp = 1e9; jmp > 1e-20; jmp /= 2) {
    REP(j,0,100) REP(dx,-1,2) REP(dy,-1,2) {
      P p = cur.second;
      p[0] += dx*jmp;
      p[1] += dy*jmp;
      cur = min(cur, make pair(f(p), p));
  return cur:
Integrate.h
Description: Simple integration of a function over an interval using Simp-
son's rule. The error should be proportional to h^4, although in practice you
will want to verify that the result is stable to desired precision when epsilon
changes.
                                                          f6be40, 7 lines
template<class F>
double quad(double a, double b, F f, const int n = 1000) {
  double h = (b - a) / 2 / n, v = f(a) + f(b);
  REP(i.1.n*2)
   v += f(a + i*h) * (i\&1 ? 4 : 2);
  return v * h / 3;
IntegrateAdaptive.h
Description: Fast integration using an adaptive Simpson's rule.
Usage: double sphereVolume = quad(-1, 1, [](double x)  {
return quad(-1, 1, [&](double y)
return quad(-1, 1, [&](double z)
return x*x + y*y + z*z < 1; ); ); ); });
typedef double d;
#define S(a,b) (f(a) + 4*f((a+b) / 2) + f(b)) * (b-a) / 6
template <class F>
d rec(F& f, d a, d b, d eps, d S) {
 d c = (a + b) / 2;
  d S1 = S(a, c), S2 = S(c, b), T = S1 + S2;
  if (abs(T - S) \le 15 * eps | | b - a < 1e-10)
    return T + (T - S) / 15;
  return rec(f, a, c, eps / 2, S1) + rec(f, c, b, eps / 2, S2):
template<class F>
d quad(d a, d b, F f, d eps = 1e-8) {
 return rec(f, a, b, eps, S(a, b));
Simplex.h
Description: Solves a general linear maximization problem: maximize c^T x
subject to Ax \leq b, x \geq 0. Returns -inf if there is no solution, inf if there
are arbitrarily good solutions, or the maximum value of c^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 \text{ val} = LPSolver(A, b, c).solve(x);
Time: \mathcal{O}(NM * \#pivots), where a pivot may be e.g. an edge relaxation.
\mathcal{O}(2^n) in the general case.
typedef double T; // long double, Rational, double + mod<P>...
typedef vector<T> vd;
typedef vector<vd> vvd;
const T eps = 1e-8, inf = 1/.0;
#define MP make pair
```

#define ltj(X) if $(s == -1 \mid | MP(X[j],N[j]) < MP(X[s],N[s])) s=j$

REP(j,i+1,n) if (fabs(a[j][i]) > fabs(a[b][i])) b = j;

if (v != 0) REP(k,i+1,n) a[j][k] -= v * a[i][k];

Description: Calculates determinant using modular arithmetics. Modulos

if (i != b) swap(a[i], a[b]), res *= -1;

double v = a[j][i] / a[i][i];

can also be removed to get a pure-integer version.

REP(i,0,n) {

return res;

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

IntDeterminant.h

const ll mod = 12345;

int b = i:

res *= a[i][i];

REP(j,i+1,n)

if (res == 0) return 0:

```
struct LPSolver {
  int m. n:
  VI N, B;
  vvd D:
  LPSolver(const vvd& A, const vd& b, const vd& c) :
    m(SZ(b)), n(SZ(c)), N(n+1), B(m), D(m+2, vd(n+2)) {
     REP(i,0,m) REP(j,0,n) D[i][j] = A[i][j];
      REP(i,0,m) \ \{ B[i] = n+i; D[i][n] = -1; D[i][n+1] = b[i]; \}
     REP(j,0,n) \{ N[j] = j; D[m][j] = -c[j]; \}
     N[n] = -1; D[m+1][n] = 1;
  void pivot(int r, int s) {
   T *a = D[r].data(), inv = 1 / a[s];
    REP(i,0,m+2) if (i != r && abs(D[i][s]) > eps) {
     T *b = D[i].data(), inv2 = b[s] * inv;
      REP(j,0,n+2) b[j] -= a[j] * inv2;
     b[s] = a[s] * inv2;
    REP(j,0,n+2) if (j != s) D[r][j] *= inv;
    REP(i,0,m+2) if (i != r) D[i][s] *= -inv;
   D[r][s] = inv;
    swap(B[r], N[s]);
  bool simplex(int phase) {
    int x = m + phase - 1;
    for (;;) {
      int s = -1:
      REP(j,0,n+1) if (N[j] != -phase) ltj(D[x]);
      if (D[x][s] >= -eps) return true;
      int r = -1;
      REP(i,0,m) {
       if (D[i][s] <= eps) continue;</pre>
       if (r == -1 || 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;
};
      Matrices
Determinant.h
Description: Calculates determinant of a matrix. Destroys the matrix.
Time: \mathcal{O}(N^3)
                                                     d14002, 15 lines
```

double det(vector<vector<double>>& a) {

int n = SZ(a); double res = 1;

```
ll det(vector<vector<ll>>& a) {
 int n = SZ(a); ll ans = 1;
 REP(i,0,n) {
    REP(j,i+1,n)
      while (a[j][i] != 0) { // gcd step
        ll t = a[i][i] / a[i][i];
        if (t) REP(k,i,n)
          a[i][k] = (a[i][k] - a[j][k] * t) % mod;
        swap(a[i], a[i]);
        ans *= -1;
    ans = ans * a[i][i] % mod;
   if (!ans) return 0;
 return (ans + mod) % mod;
SolveLinear.h
Description: Solves A * x = b. If there are multiple solutions, an arbitrary
one is returned. Returns rank, or -1 if no solutions. Data in A and b is lost.
Time: \mathcal{O}\left(n^2m\right)
                                                       d6dca7, 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[i][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)</pre>
               SolveLinear2.h
              Description: To get all uniquely determined values of x back from Solve-
              Linear, make the following changes:
               "SolveLinear.h"
                                                                        5cad07, 7 lines
669167, 18 lines
              REP(j,0,n) if (j != i) // instead of REP(j,i+1,n)
              // ... then at the end:
              x.assign(m, undefined);
              REP(i,0,rank)
                REP(j,rank,m) if (fabs(A[i][j]) > eps) goto fail;
                x[col[i]] = b[i] / A[i][i];
               fail:; }
              SolveLinearBinarv.h
              Description: Solves Ax = b over \mathbb{F}_2. If there are multiple solutions, one is
              returned arbitrarily. Returns rank, or -1 if no solutions. Destroys A and b.
              Time: \mathcal{O}\left(n^2m\right)
                                                                       c9c00b, 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;
                   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)</pre>
               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: $\mathcal{O}\left(n^3\right)$

```
43464f, 35 lines
int matInv(vector<vector<double>>& A) {
  int n = SZ(A); VI col(n);
  vector<vector<double>> tmp(n, vector<double>(n));
  REP(i.0.n) tmp[i][i] = 1. col[i] = i:
  REP(i,0,n) {
    int r = i, c = i;
   REP(j,i,n) REP(k,i,n)
     if (fabs(A[j][k]) > fabs(A[r][c]))
        r = j, c = k;
    if (fabs(A[r][c]) < 1e-12) return i;</pre>
    A[i].swap(A[r]); tmp[i].swap(tmp[r]);
   REP(i.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[i][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[i][i]:
   REP(k,0,n) tmp[j][k] -= v*tmp[i][k];
  REP(i,0,n) REP(i,0,n) A[col[i]][col[i]] = tmp[i][i];
  return n:
```

Tridiagonal.h

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

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

This is useful for solving problems on the type

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

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

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

Fails if the solution is not unique.

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

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

a74eda, 26 lines

```
typedef double T;
vector<T> tridiagonal(vector<T> diag, const vector<T>& super,
   const vector<T>& sub, vector<T> b) {
  int n = SZ(b); VI tr(n);
```

```
REP(i,0,n-1) {
  if (abs(diag[i]) < 1e-9 * abs(super[i])) { // diag[i] == 0}
    b[i+1] -= b[i] * diag[i+1] / super[i]:
    if (i+2 < n) b[i+2] -= b[i] * sub[i+1] / super[i];</pre>
    diag[i+1] = sub[i]; tr[++i] = 1;
  } else {
    diag[i+1] -= super[i]*sub[i]/diag[i];
    b[i+1] -= b[i]*sub[i]/diag[i];
for (int i = n; i--;) {
  if (tr[i]) {
    swap(b[i], b[i-1]);
    diag[i-1] = diag[i];
    b[i] /= super[i-1];
  } else {
    b[i] /= diag[i];
    if (i) b[i-1] -= b[i]*super[i-1];
return b;
```

4.4 Fourier transforms

FastFourierTransform.h

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

```
a333b0, 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 *= 2) {
   R.resize(n); rt.resize(n);
    auto x = polar(1.0L, acos(-1.0L) / k);
    REP(i,k,2*k) rt[i] = R[i] = i&1 ? R[i/2] * x : R[i/2];
 VI rev(n):
  REP(i,0,n) rev[i] = (rev[i / 2] | (i \& 1) << L) / 2;
  REP(i,0,n) if (i < rev[i]) swap(a[i], a[rev[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 + i] += 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 << L;</pre>
 vector<C > in(\overline{n}), out(\overline{n}):
  copy(ALL(a), begin(in));
  REP(i, 0, SZ(b)) in[i].imag(b[i]);
 fft(in):
 for (C\& x : in) x *= x;
 REP(i,0,n) out[i] = in[-i & (n - 1)] - coni(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<ll> vl:
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 = ll(real(outl[i])+.5), cv = ll(imag(outs[i])+.5);
   ll bv = ll(imag(outl[i])+.5) + ll(real(outs[i])+.5);
    res[i] = ((av % M * cut + bv) % M * cut + cv) % M;
  return res:
```

NumberTheoreticTransform.h

<< B;

int inv = modpow(n, mod - 2);

vl L(a), R(b), out(n);

L.resize(n), R.resize(n);

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

```
Time: \mathcal{O}(N \log N)
"../number-theory/ModPow.h"
const ll 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> vl;
void ntt(vl &a) {
 int n = SZ(a), L = 31 - builtin clz(n);
  static vl rt(2, 1);
  for (static int k = 2, s = 2; k < n; k *= 2, s++) {
    rt.resize(n);
    ||z|| = \{1, modpow(root, mod >> s)\}:
    REP(i,k,2*k) rt[i] = rt[i / 2] * z[i & 1] % mod:
  VI rev(n):
  REP(i,0,n) rev[i] = (rev[i / 2] | (i \& 1) << L) / 2;
  REP(i,0,n) if (i < rev[i]) swap(a[i], a[rev[i]]);
  for (int k = 1: k < n: k *= 2)
    for (int i = 0; i < n; i += 2 * k) REP(j,0,k) {
      ll z = rt[j + k] * a[i + j + k] % mod, &ai = a[i + j];
      a[i + j + k] = ai - z + (z > ai ? mod : 0);
      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
```

```
ШТН
  ntt(L), ntt(R);
  REP(i,0,n) out[-i & (n - 1)] = (ll)L[i] * R[i] % mod * inv %
  ntt(out);
  return {out.begin(), out.begin() + s};
FastSubsetTransform.h
Description: Transform to a basis with fast convolutions of the form
c[z]=\sum_{z=x\oplus y}a[x]\cdot b[y], where \oplus is one of AND, OR, XOR. The size of a must be a power of two.
Time: \mathcal{O}(N \log N)
                                                              b1dc68, 16 lines
```

```
void FST(VI& a, bool inv) {
 for (int n = SZ(a), step = 1; step < n; step *= 2) {
   for (int i = 0; i < n; i += 2 * step) REP(j,i,i+step) {
     int \&u = a[j], \&v = a[j + step]; tie(u, v) =
       inv ? PII(v - u, u) : PII(v, u + v); // AND
       inv ? PII(v, u - v) : PII(u + v, u); // OR
       PII(u + v, u - v);
 if (inv) for (int& x : a) x /= SZ(a); // XOR only
VI conv(VI a. VI b) {
 FST(a, 0); FST(b, 0);
 REP(i,0,SZ(a)) a[i] *= b[i];
 FST(a, 1); return a;
```

Number theory (5)

5.1 Modular arithmetic

Modular Arithmetic.h

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

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

ModInverse.h

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

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

```
ModPow.h
const ll mod = 1000000007; // faster if const
ll modpow(ll b, ll e) {
 ll ans = 1;
 for (; e; b = b * b % mod, e /= 2)
    if (e \& 1) ans = ans * b % mod:
  return ans:
ModLog.h
Description: Returns the smallest x > 0 s.t. a^x = b \pmod{m}, or -1 if no
such x exists. modLog(a,1,m) can be used to calculate the order of a.
```

```
Time: \mathcal{O}(\sqrt{m})
                                                             0ff368, 11 lines
Il modLog(ll a, ll b, ll m) {
  ll n = (ll) sqrt(m) + 1, e = 1, f = 1, j = 1;
  unordered map<ll, ll> A;
 while (j \le n \&\& (e = f = e * a % m) != b % m)
    A[e * b % m] = j++;
  if (e == b % m) return j;
 if (_gcd(m, e) == _gcd(m, b))
    REP(i,2,n+2) if (A.count(e = e * f % m))
      return n * i - A[e];
  return -1;
```

ModSum.h

Description: Sums of mod'ed arithmetic progressions.

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

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 to 2 = (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 * sumsa(to) - m * divsum(to, c, k, m):
```

ModMulLL.h

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

```
typedef unsigned long long ull;
ull modmul(ull a. ull b. ull M) {
 ll ret = a * b - M * ull(1.L / M * a * b);
 return ret + M * (ret < 0) - M * (ret >= (11)M);
ull modpow(ull b, ull e, ull mod) {
 ull ans = 1:
 for (; e; b = modmul(b, b, mod), e /= 2)
   if (e \& 1) ans = modmul(ans, b, mod);
 return ans:
```

```
ModSgrt.h
```

Description: Tonelli-Shanks algorithm for modular square roots. Finds x s.t. $x^2 = a \pmod{p}$ (-x gives the other solution). **Time:** $\mathcal{O}(\log^2 p)$ worst case, $\mathcal{O}(\log p)$ for most p

```
"ModPow.h"
                                                     19a793, 24 lines
ll sgrt(ll a, ll p) {
 a \% = p: if (a < 0) a += p:
 if (a == 0) return 0;
 assert(modpow(a, (p-1)/2, p) == 1); // else no solution
 if (p \% 4 == 3) return modpow(a, (p+1)/4, p);
  // a^{(n+3)/8} or 2^{(n+3)/8} * 2^{(n-1)/4} 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) {
    ll t = b;
    for (m = 0; m < r \&\& t != 1; ++m)
     t = t * t % p;
    if (m == 0) return x;
    ll qs = modpow(q, 1LL \ll (r - m - 1), p);
    q = qs * qs % p;
    x = x * gs % p;
    b = b * a % p:
```

5.2 Primality

FastEratosthenes.h

Description: Prime sieve for generating all primes smaller than LIM. Time: LIM=1e9 $\approx 1.5s$ bdcb93, 20 lines

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

MillerRabin.h

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

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

```
"ModMulLL.h"
                                                            60dcd1, 12 lines
bool isPrime(ull n) {
  if (n < 2 || n % 6 % 4 != 1) return (n | 1) == 3;
  ull A[] = \{2, 325, 9375, 28178, 450775, 9780504, 1795265022\},
       s = builtin ctzll(n-1), d = n >> s;
  for (ull a : A) { // ^ count trailing zeroes
  ull p = modpow(a%n, d, n), i = s;
```

```
while (p != 1 && p != n - 1 && a % n && i--)
   p = modmul(p, p, n);
 if (p != n-1 && i != s) return 0:
return 1;
```

Factor.h

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

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

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

Divisibility 5.3

euclid.h

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

33ba8f, 5 lines

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

CRT.h

Description: Chinese Remainder Theorem.

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

04d93a, 7 lines "euclid.h" ll crt(ll a, ll m, ll b, ll n) { **if** (n > m) swap(a, b), swap(m, n);

ll x, y, g = euclid(m, n, x, y); assert((a - b) % g == 0); // else no solution x = (b - a) % n * x % n / g * m + a;return x < 0 ? x + m*n/g : x;

5.3.1 Bézout's identity

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

ax + by = d

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

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

phiFunction.h

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

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

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

e4742a, 8 lines

```
const int LTM = 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 > 0, finds the closest rational approximation p/q with p, q < N. It will obey |p/q - x| < 1/qN.

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

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

```
typedef double d: // for N \sim 1e7: long double for N \sim 1e9
pair<ll, ll> approximate(d x, ll N) {
 II LP = 0, LQ = 1, P = 1, Q = 0, inf = LLONG MAX; d y = x;
 for (::) {
   ll lim = min(P ? (N-LP) / P : inf, Q ? (N-LQ) / Q : inf),
       a = (11)floor(y), b = min(a, lim),
       NP = b*P + LP \cdot NO = b*O + LO:
   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: $\mathcal{O}(\log(N))$ 27ab3e, 25 lines

```
struct Frac { ll p, q; };
template<class F>
Frac fracBS(F f, ll N) {
 bool dir = 1, A = 1, B = 1;
```

```
Frac lo{0}, 1, hi{1}, 1; // Set hi to 1/0 to search (0, N)
if (f(lo)) return lo:
assert(f(hi)):
while (A || B)
  ll adv = 0, step = 1; // move hi if dir, else lo
  for (int si = 0; step; (step *= 2) >>= si) {
    Frac mid{lo.p * adv + hi.p, lo.g * adv + hi.g};
    if (abs(mid.p) > N \mid \mid mid.q > N \mid \mid 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:
```

10

5.5 Pythagorean Triples

The Pythagorean triples are uniquely generated by

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

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

5.6 Primes

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

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

5.7 Estimates

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

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

Mobius Function

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

Mobius Inversion:

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

Other useful formulas/forms:

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

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

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

Graph (6)

6.1 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:** $\mathcal{O}\left(VE\right)$

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

FlovdWarshall.h

Description: Calculates all-pairs shortest path in a directed graph that might have negative edge weights. Input is an distance matrix m, where $m[i][j] = \inf$ if i and j are not adjacent. As output, m[i][j] is set to the shortest distance between i and j, \inf inf in opath, or -inf if the path goes through a negative-weight cycle.

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.

```
I topoSort(const vectorsvi>& gr) {
VI indeg(SZ(gr)), ret;
for (auto& li : gr) for (int x : li) indeg[x]++;
queue<int> q; // use priority_queue for lexic. largest ans.
REP(i,0,SZ(gr)) if (indeg[i] == 0) q.push(i);
while (!q.empty()) {
   int i = q.front(); // top() for priority queue
```

```
ret.push_back(i);
q.pop();
for (int x : gr[i])
   if (--indeg[x] == 0) q.push(x);
}
return ret;
```

6.2 Network flow

Dinic.h

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

```
struct Dinic {
 struct Edge {
   int to, rev;
    ll c, oc;
    Il flow() { return max(oc - c, OLL); } // if you need flows
  VI lvl, ptr, q;
  vector<vector<Edge>> adi;
 Dinic(int n) : lvl(n), ptr(n), q(n), adj(n) {}
  void addEdge(int a, int b, ll c, ll rcap = 0) {
   adj[a].push back({b, SZ(adj[b]), c, c});
    adj[b].push back({a, SZ(adj[a]) - 1, rcap, rcap});
 ll dfs(int v, int t, ll f) {
    if (v == t || !f) return f;
    for (int\& i = ptr[v]: i < SZ(adi[v]): i++)
      Edge& e = adj[v][i];
      if (lvl[e.to] == lvl[v] + 1)
       if (ll p = dfs(e.to, t, min(f, e.c))) {
         e.c -= p, adi[e.to][e.rev].c += p;
          return p;
    return 0;
 ll calc(int s, int t) {
    II flow = 0; q[0] = s;
    REP(L,0,31) do { // 'int L=30' maybe faster for random data
      lvl = ptr = VI(SZ(q));
     int qi = 0, qe = lvl[s] = 1;
      while (qi < qe && !lvl[t]) {</pre>
       int v = q[qi++];
       for (Edge e : adj[v])
         if (!lvl[e.to] && e.c >> (30 - L))
            q[qe++] = e.to, lvl[e.to] = lvl[v] + 1;
     while (ll p = dfs(s, t, LLONG MAX)) flow += p;
    } while (lvl[t]);
    return flow:
 bool leftOfMinCut(int a) { return lvl[a] != 0; }
```

MCMF-SPFA.h

Description: Multiedges and negative costs allowed. **Time:** Approximately $\mathcal{O}(V^2E^2)$

```
template <typename FLOW, typename COST> struct MCMF {
  const COST INFC = 1e9, EPSC = 0;
  const FLOW INFF = 1e9, EPSF = 0;
  struct Edge {
   int from, to;
   FLOW flow, cap;
  COST cost;
}
```

```
int nodes, src, dest, m = 0;
  vector<vector<int>> adi:
  vector<Edae> edaes:
  void add(int u, int v, FLOW cap, COST cost) {
    edges.EB(u, v, 0, cap, cost);
    adi[u].PB(m++);
    edges.EB(v, u, 0, 0, -cost);
    adj[v].PB(m++);
  vector<COST> dis;
  vector<bool> in0:
  VI par;
  pair<FLOW. COST> SPFA() {
    fill(ALL(dis), INFC);
    fill(ALL(inQ), false);
    queue<int> Q;
    Q.push(src), dis[src] = 0, inQ[src] = true;
    while (!Q.empty()) {
      int u = Q.front(); Q.pop();
      inQ[u] = false;
      for (int i : adj[u]) {
        auto &e = edges[i];
        if (e.cap - e.flow > EPSF
            && dis[e.to] - (dis[u] + e.cost) > EPSC) {
          dis[e.to] = dis[u] + e.cost;
          par[e.to] = i;
          if (!inQ[e.to]) { Q.push(e.to), inQ[e.to] = true; }
    if (dis[dest] + EPSC >= INFC) return {0, 0};
    FLOW aug = INFF;
    for (int u = dest; u != src; u = edges[par[u]].from) {
      aug = min(aug, edges[par[u]].cap - edges[par[u]].flow);
    for (int u = dest; u != src; u = edges[par[u]].from) {
      edges[par[u]].flow += aug;
      edges[par[u] ^ 1].flow -= aug;
    return {aug, aug * dis[dest]};
  MCMF(int n. int s. int t)
    : nodes(n), src(s), dest(t), adj(n), dis(n), inQ(n), par(n)
  pair<FLOW, COST> mincostmaxflow() {
    pair<FLOW, COST> ans(0, 0);
    while (true) {
      auto cur = SPFA():
      if (cur.first <= EPSF)</pre>
      ans.first += cur.first:
      ans.second += cur.second;
    return ans;
};
```

MinCostMaxFlow.h

19b593, 63 lines

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.

```
const ll INF = numeric_limits<ll>::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;
     qnu pbds::priority queue<pair<ll, int>> q;
   vector<decltype(q)::point iterator> its(N);
   q.push({0, s});
    auto relax = [&](int i, ll cap, ll cost, int dir) {
     ll 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 (!a.emptv()) {
     s = q.top().second; q.pop();
     seen[s] = 1; di = dist[s] + pi[s];
     for (int i : ed[s]) if (!seen[i])
       relax(i, cap[s][i] - flow[s][i], cost[s][i], 1);
     for (int i : red[s]) if (!seen[i])
       relax(i, flow[i][s], -cost[i][s], 0);
   REP(i,0,N) pi[i] = min(pi[i] + dist[i], INF);
  pair<ll, ll> maxflow(int s, int t) {
   ll totflow = 0, totcost = 0;
   while (path(s), seen[t]) {
     ll 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(i,0,N) totcost += cost[i][i] * flow[i][i];
   return {totflow, totcost};
  // If some costs can be negative, call this before maxflow:
  void setpi(int s) { // (otherwise, leave this out)
   fill(ALL(pi), INF); pi[s] = 0;
   int it = N, ch = 1; ll v;
   while (ch-- && it--)
```

```
REP(i,0,N) if (pi[i] != INF)
        for (int to : ed[i]) if (cap[i][to])
          if ((v = pi[i] + cost[i][to]) < pi[to])
             pi[to] = v, ch = 1;
    assert(it >= 0); // negative cost cycle
};
MinCut.h
Description: After running max-flow, the left side of a min-cut from s to t
is given by all vertices reachable from s, only traversing edges with positive
residual capacity.
GlobalMinCut.h
```

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

```
1d69cc, 21 lines
pair<int, VI> globalMinCut(vector<VI> mat) {
 pair<int, VI> best = {INT MAX, {}};
 int n = SZ(mat):
 vector<VI> co(n);
 REP(i,0,n) co[i] = \{i\};
 REP(ph,1,n) |
   VI w = mat[0]:
   size t s = 0. t = 0:
    REP(\overline{i}t,0,n-ph) { // O(V^2) -> O(E \log V) with prio. queue
     w[t] = INT MIN:
     s = t, t = max element(ALL(w)) - w.begin();
      REP(i,0,n) w[i] += mat[t][i];
    best = min(best, \{w[t] - mat[t][t], co[t]\});
    co[s].insert(co[s].end(), ALL(co[t]));
    REP(i,0,n) mat[s][i] += mat[t][i];
    REP(i,0,n) mat[i][s] = mat[s][i];
   mat[0][t] = INT MIN;
 return best;
```

GomoryHu.h

Description: Given a list of edges representing an undirected flow graph, returns edges of the Gomory-Hu tree. The max flow between any pair of vertices is given by minimum edge weight along the Gomory-Hu tree path.

Time: $\mathcal{O}(V)$ Flow Computations

```
b8cecf, 13 lines
"PushRelabel.h"
typedef array<ll, 3> Edge;
vector<Edge> gomorvHu(int N, vector<Edge> ed) {
 vector<Edge> tree;
 VI par(N);
 REP(i,1,N)
    PushRelabel D(N); // Dinic also works
    for (Edge t : ed) D.addEdge(t[0], t[1], t[2], t[2]);
    tree.push back({i, par[i], D.calc(i, par[i])});
     if (par[j] == par[i] && D.leftOfMinCut(j)) par[j] = i;
 return tree:
```

6.3 Matching

hopcroftKarp.h

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

```
Time: \mathcal{O}\left(\sqrt{V}E\right)
bool dfs(int a, int L, vector<VI>& g, VI& btoa, VI& A, VI& B) {
 if (A[a] != L) return 0:
  A[a] = -1;
  for (int b : q[a]) if (B[b] == L + 1) {
    B[b] = 0:
    if (btoa[b] == -1 \mid | dfs(btoa[b], L + 1, g, btoa, A, B))
      return btoa[b] = a, 1;
  return 0;
int hoperoftKarp(vector<VI>& q, VI& btoa) {
 int res = 0:
  VI A(g.size()), B(btoa.size()), cur, next;
  for (;;) {
    fill(ALL(A), 0);
    fill(ALL(B), 0);
    cur.clear();
    for (int a : btoa) if(a != -1) A[a] = -1;
    REP(a,0,SZ(g)) if (A[a] == 0) cur.push back(a);
    for (int lay = 1;; lay++) {
      bool islast = 0:
      next.clear();
      for (int a : cur) for (int b : q[a]) {
        if (btoa[b] == -1) {
          B[b] = lay;
          islast = 1:
        else if (btoa[b] != a && !B[b]) {
          B[b] = lay;
          next.push back(btoa[b]);
      if (islast) break;
      if (next.empty()) return res;
      for (int a : next) A[a] = 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(q, btoa);

```
Time: \mathcal{O}(VE)
                                                      d47d74, 22 lines
bool find(int j, vector<VI>& q, VI& btoa, VI& vis) {
 if (btoa[i] == -1) return 1;
  vis[j] = 1; int di = btoa[j];
  for (int e : g[di])
    if (!vis[e] && find(e, g, btoa, vis)) {
      btoa[e] = di;
      return 1;
  return 0;
int dfsMatching(vector<VI>& q, VI& btoa) {
  VI vis:
  REP(i, 0, SZ(g)) {
    vis.assign(SZ(btoa), 0);
    for (int j : q[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" VI cover(vector<VI>& q, int n, int m) { VI match(m, -1); int res = dfsMatching(g, match); vector<bool> lfound(n, true), seen(m); for (int it : match) if (it != -1) lfound[it] = false: VI q, cover; REP(i,0,n) if (lfound[i]) q.push back(i); while (!q.empty()) { int i = q.back(); q.pop back(); lfound[i] = 1;for (int e : q[i]) if (!seen[e] && match[e] != -1) { seen[e] = true; q.push back(match[e]); REP(i,0,n) if (!lfound[i]) cover.push back(i); REP(i,0,m) if (seen[i]) cover.push back(n+i); assert(SZ(cover) == res); return cover:

WeightedMatching.h

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

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

```
REP(j,1,m) if (p[j]) ans[p[j] - 1] = j - 1;
 return {-v[0], ans}; // min cost
GeneralMatching.h
Description: Matching for general graphs. Fails with probability N/mod.
Time: \mathcal{O}(N^3)
"../numerical/MatrixInverse-mod.h"
                                                      69fab6, 40 lines
vector<PII> generalMatching(int N. vector<PII>& ed) {
  vector<vector<ll>>> mat(N, vector<ll>(N)), A;
 for (PII pa : ed) {
   int a = pa.first, b = pa.second, r = rand() % mod;
    mat[a][b] = r, mat[b][a] = (mod - r) % mod;
 int r = matInv(A = mat), M = 2*N - r, fi, fi;
 assert(r % 2 == 0);
 if (M != N) do {
    mat.resize(M, vector<ll>(M));
    REP(i,0,N) {
      mat[i].resize(M);
      REP(j,N,M) {
        int r = rand() % mod;
        mat[i][j] = r, mat[j][i] = (mod - r) % mod;
 } while (matInv(A = mat) != M);
  VI has(M, 1); vector<PII> ret;
 REP(it,0,M/2) {
    REP(i,0,M) if (has[i])
      REP(j,i+1,M) if (A[i][j] && mat[i][j]) {
        fi = i; fj = j; goto done;
     assert(0); done:
    if (fj < N) ret.emplace back(fi, fj);</pre>
    has[fi] = has[fj] = 0;
    REP(sw.0.2) {
      11 a = modpow(A[fi][fj], mod-2);
      REP(i,0,M) if (has[i] && A[i][fj]) {
        ll b = A[i][fj] * a % mod;
        REP(j,0,M) A[i][j] = (A[i][j] - A[fi][j] * b) % mod;
      swap(fi,fj);
 return ret;
```

6.4 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 versa.

Usage: $scc(graph, [\&](VI\& v) \{ \dots \})$ 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: $\mathcal{O}(E+V)$

```
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);
   for (auto e : g[j]) if (comp[e] < 0)
   low = min(low, val[e] ?: dfs(e,g,f));
```

```
if (low == val[j]) {
    do {
        x = z.back(); z.pop_back();
        comp[x] = ncomps;
        cont.push_back(x);
    } while (x != j);
    f(cont); cont.clear();
    ncomps++;
}
return val[j] = low;
}
template<class G, class F> void scc(G& g, F f) {
    int n = SZ(g);
    val.assign(n, 0); comp.assign(n, -1);
    Time = ncomps = 0;
    REP(i,0,n) if (comp[i] < 0) dfs(i, g, f);
}</pre>
```

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) {...}); \\ \hline \textbf{Time: } \mathcal{O}(E+V) \\ \hline \textbf{VI num, st;} \\ vector<vector<\mathbf{PII}>> ed;
```

```
int Time:
template<class F>
int dfs(int at, int par, F& f) {
 int me = num[at] = ++Time, e, y, top = me;
  for (auto pa : ed[at]) if (pa.second != par) {
    tie(y, e) = pa;
    if (num[y]) {
      top = min(top, num[y]);
      if (num[y] < me)</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 *\overline{/} }
  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 unsatis-
fiable. 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.setValue(2): // Var 2 is true
ts.atMostOne(\{0,\sim1,2\}); // <= 1 of vars 0, \sim1 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.
                                                          def624, 56 lines
struct TwoSat {
  int N;
  vector<VI> qr;
  VI values; // 0 = false, 1 = true
```

```
TwoSat(int n = 0) : N(n), gr(2*n) {}
int addVar() { // (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[j].push back(f^1);
void setValue(int x) { either(x, x); }
void atMostOne(const VI& li) { // (optional)
 if (SZ(li) <= 1) return;</pre>
  int cur = \sim li[0];
  REP(i,2,SZ(li)) {
   int next = addVar();
   either(cur, ~li[i]);
   either(cur, next);
   either(~li[i], next);
   cur = ~next:
  either(cur, ~li[1]);
VI val, comp, z; int time = 0;
int dfs(int i)
  int low = val[i] = ++time, x; z.push back(i);
  for(int e : 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:
```

EulerWalk.h

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

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

6.5 Coloring

EdgeColoring.h

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

```
Time: \mathcal{O}(NM)
                                                     8618ee, 31 lines
VI edgeColoring(int N, vector<PII> eds) {
 VI cc(N + 1), ret(SZ(eds)), fan(N), free(N), loc;
 for (PII e : eds) ++cc[e.first], ++cc[e.second];
 int u, v, ncols = *max element(ALL(cc)) + 1;
 vector<VI> adj(N, VI(ncols, -1));
 for (PII e : eds) {
   tie(u, v) = e;
    fan[0] = v;
    loc.assign(ncols, 0);
    int at = u, end = u, d, c = free[u], ind = 0, i = 0;
    while (d = free[v], !loc[d] \&\& (v = adj[u][d]) != -1)
     loc[d] = ++ind, cc[ind] = d, fan[ind] = v;
    cc[loc[d]] = c;
    for (int cd = d; at != -1; cd ^= c ^ d, at = adj[at][cd])
     swap(adj[at][cd], adj[end = at][cd ^ c ^ d]);
    while (adj[fan[i]][d] != -1) {
     int left = fan[i], right = fan[++i], e = cc[i];
      adi[u][e] = left;
      adj[left][e] = u;
     adi[right][e] = -1;
      free[right] = e;
    adj[u][d] = fan[i];
    adj[fan[i]][d] = u;
    for (int y : {fan[0], u, end})
     for (int& z = free[y] = 0; adj[y][z] != -1; z++);
 REP(i,0,SZ(eds))
   for (tie(u, v) = eds[i]; adj[u][ret[i]] != v;) ++ret[i];
  return ret:
```

6.6 Heuristics

MaximalCliques.h

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

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

```
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 \& \sim 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;
```

MaximumClique.h

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

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

```
typedef vector<bitset<200>> vb;
struct Maxclique
 double limit=0.025, pk=0;
 struct Vertex { int i, d=0; };
```

```
typedef vector<Vertex> vv;
vb e;
vv V;
vector<VI> C;
VI qmax, q, S, old;
void init(vv& r) {
  for (auto\& v : r) v.d = 0;
  for (auto \& v : r) for (auto j : r) v.d += e[v.i][j.i];
  sort(ALL(r), [](auto a, auto b) { return a.d > b.d; });
  int mxD = r[0].d;
  REP(i,0,SZ(r)) r[i].d = min(i, mxD) + 1;
void expand(vv\& R, int lev = 1)
  S[lev] += S[lev - 1] - old[lev];
  old[lev] = S[lev - 1]:
  while (SZ(R)) {
    if (SZ(q) + R.back().d <= SZ(qmax)) return;</pre>
    g.push back(R.back().i):
    vv T;
    for(auto v:R) if (e[R.back().i][v.i]) T.push back({v.i});
    if (SZ(T)) {
      if (S[lev]++ / ++pk < limit) init(T);</pre>
      int j = 0, mxk = 1, mnk = max(SZ(gmax) - SZ(g) + 1, 1);
      C[1].clear(), C[2].clear();
      for (auto v : T) {
        int k = 1:
        auto f = [&](int i) { return e[v.i][i]; };
        while (any of(ALL(C[k]), f)) k++;
        if (k > mxk) mxk = k, C[mxk + 1].clear();
        if (k < mnk) T[j++].i = v.i;
        C[k].push back(v.i);
      if (j > 0) T[j - 1].d = 0;
      REP(k, mnk, mxk + 1) for (int i : C[k])
       T[i].i = i, T[i++].d = k;
      expand(T, lev + 1);
```

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

```
q.pop back(), R.pop back();
  VI maxClique() { init(V), expand(V); return gmax; }
  Maxclique(vb conn) : e(conn), C(SZ(e)+1), S(SZ(C)), old(S) {
    REP(i,0,SZ(e)) V.push back({i});
};
```

MaximumIndependentSet.h

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

6.7 Trees

BinaryLifting.h

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

Time: construction $\mathcal{O}(N \log N)$, queries $\mathcal{O}(\log N)$ c4e44c, 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 imp(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.

```
Time: \mathcal{O}(N \log N + Q)
"../data-structures/RMQ.h"
                                                         cbd116, 21 lines
struct LCA {
  int T = 0:
  VI time, path, ret;
  RMQ<int> rmq;
  LCA(vector < VI > \& C) : time(SZ(C)), rmq((dfs(C,0,-1), ret)) {}
  void dfs(vector<VI>& C, int v, int par) {
    time[v] = T++;
    for (int y : C[v]) if (y != par) {
      path.push back(v), ret.push back(time[v]);
      dfs(C, y, v);
  int lca(int a, int b) {
```

```
if (a == b) return a;
    tie(a, b) = minmax(time[a], time[b]);
    return path[rmg.guery(a, b)];
 //dist(a,b) {return depth[a] + depth[b] - 2*depth[lca(a,b)];}
CompressTree.h
Description: Given a rooted tree and a subset S of nodes, compute the
minimal subtree that contains all the nodes by adding all (at most |S|-1)
pairwise LCA's and compressing edges. Returns a list of (par, orig_index)
representing a tree rooted at 0. The root points to itself.
Time: \mathcal{O}\left(|S|\log|S|\right)
"LCA.h"
                                                         bcf310, 21 lines
typedef vector<pair<int, int>> vpi;
vpi compressTree(LCA& lca, const VI& subset) {
  static VI rev; rev.resize(SZ(lca.time));
  VI li = subset, &T = lca.time;
  auto cmp = [&](int a, int b) { return T[a] < T[b]; };</pre>
  sort(ALL(li), cmp);
  int m = SZ(li) - 1;
  REP(i,0,m) {
    int a = li[i], b = li[i+1];
    li.push back(lca.lca(a, b));
  sort(ALL(li), cmp);
  li.erase(unique(ALL(li)), li.end());
  REP(i,0,SZ(li)) rev[li[i]] = i;
  vpi ret = {PII(0, li[0])};
  REP(i,0,SZ(li)-1) {
    int a = li[i], b = li[i+1];
    ret.emplace back(rev[lca.lca(a, b)], b);
  return ret;
HLD.h
Description: Decomposes a tree into vertex disjoint heavy paths and light
edges such that the path from any leaf to the root contains at most log(n)
light edges. Code does additive modifications and max queries, but can
support commutative segtree modifications/queries on paths and subtrees.
Takes as input the full adjacency list. VALS_EDGES being true means that
values are stored in the edges, as opposed to the nodes. All values initialized
to the segtree default. Root must be 0.
Time: \mathcal{O}\left((\log N)^2\right)
"../data-structures/LazySegmentTree.h"
                                                         e7978a, 46 lines
template <bool VALS EDGES> struct HLD {
  int N, tim = 0;
  vector<VI> adi:
  VI par, siz, depth, rt, pos:
  Node *tree:
  HLD(vector<VI> adj )
    : N(SZ(adj_)), adj(adj_), par(N, -1), siz(N, 1), depth(N),
      rt(N),pos(N),tree(new Node(0, N)){ dfsSz(0); dfsHld(0); }
  void dfsSz(int v) {
    if (par[v] != -1) adj[v].erase(find(ALL(adj[v]), par[v]));
    for (int& u : adj[v]) {
      par[u] = v, depth[u] = depth[v] + 1;
      dfsSz(u);
      siz[v] += siz[u];
      if (siz[u] > siz[adj[v][0]]) swap(u, adj[v][0]);
```

void dfsHld(int v) {

for (int u : adj[v]) {

rt[u] = (u == adj[v][0] ? rt[v] : u);

pos[v] = tim++;

dfsHld(u);

```
template <class B> void process(int u. int v. B op) {
    for (; rt[u] != rt[v]; v = par[rt[v]])
      if (depth[rt[u]] > depth[rt[v]]) swap(u, v);
      op(pos[rt[v]], pos[v] + 1);
    if (depth[u] > depth[v]) swap(u, v);
    op(pos[u] + VALS EDGES, pos[v] + 1);
 void modifyPath(int u, int v, int val) {
    process(u, v, [&](int l, int r) { tree->add(l, r, val); });
 int queryPath(int u, int v) { // Modify depending on problem
    int res = -1e9;
    process(u, v, [&](int l, int r) {
        res = max(res, tree->query(l, r));
    return res;
 int querySubtree(int v) { // modifySubtree is similar
    return tree->query(pos[v] + VALS EDGES, pos[v] + siz[v]);
};
```

LinkCutTree.h

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

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

```
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] \rightarrow p = this;
    if (c[1]) c[1]->p = this;
    // (+ update sum of subtree elements etc. if wanted)
  void pushFlip() {
    if (!flip) return;
    flip = 0; swap(c[0], c[1]);
    if (c[0]) c[0]->flip ^= 1;
    if (c[1]) c[1]->flip ^= 1;
  int up() { return p ? p->c[1] == this : -1: }
  void rot(int i, int b) {
    int h = i ^ b;
    Node *x = c[i], *y = b == 2 ? x : x -> c[h], *z = b ? y : x;
    if ((y->p = p)) p->c[up()] = y;
    c[i] = z -> c[i ^ 1];
    if (b < 2) {
      x - c[h] = y - c[h ^ 1];
      z - c[h \land 1] = b ? x : this;
    y - > c[i \land 1] = b ? this : x;
    fix(); x->fix(); y->fix();
    if (p) p->fix();
    swap(pp, y->pp);
  void splay() {
    for (pushFlip(); p; ) {
      if (p->p) p->p->pushFlip();
      p->pushFlip(); pushFlip();
      int c1 = up(), c2 = p->up();
      if (c2 == -1) p->rot(c1, 2);
      else p->p->rot(c2, c1 != c2);
```

DirectedMST KMP Zfunc Manacher

void $pop(Node*\& a) \{ a->prop(); a = merge(a->l, a->r); \}$

for (Edge e : g) heap[e.b] = merge(heap[e.b], new Node{e});

pair<ll, VI> dmst(int n, int r, vector<Edge>& g) {

Edge top() { prop(); return key; }

if (a->key.w > b->key.w) swap(a, b);
swap(a->l, (a->r = merge(b, a->r)));

VI seen(n, -1), path(n), par(n);

vector<Edge> Q(n), in(n, $\{-1,-1\}$), comp;

if (!heap[u]) **return** {-1,{}};

res += e.w, u = uf.find(e.a);

while (uf.join(u, w));

in[uf.find(inEdge.b)] = inEdge;

Edge e = heap[u]->top();

deque<tuple<int, int, vector<Edge>>> cycs;

heap[u]->delta -= e.w, pop(heap[u]);

int end = qi, time = uf.time();

REP(i.0.ai) in[uf.find(0[i].b)] = 0[i]:

for (auto& e : comp) in[uf.find(e.b)] = e;

Q[qi] = e, path[qi++] = u, seen[u] = s;

do cyc = merge(cyc, heap[w = path[--qi]]);

u = uf.find(u), heap[u] = cyc, seen[u] = -1;

cycs.push front($\{u, time, \{\&Q[qi], \&Q[end]\}\}\);$

for (auto& [u,t,comp] : cycs) { // restore sol (optional)

Node *merge(Node *a, Node *b) {
 if (!a || !b) return a ?: b;
 a->prop(), b->prop();

return a;

RollbackUF uf(n):

ll res = 0:

seen[r] = r;

REP(s,0,n) {

vector<Node*> heap(n);

int u = s, qi = 0, w;

while (seen[u] < 0) {

if (seen[u] == s) {

Node* cyc = 0;

uf.rollback(t);

return {res, par};

Edge inEdge = in[u]:

REP(i,0,n) par[i] = in[i].a;

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

DirectedMST.h

delta = 0;

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

```
Time: \mathcal{O}\left(E\log V\right)
".../data-structures/UnionFindRollback.h" 719f12, 60 lines

struct Edge { int a, b; ll w; };

struct Node {
    Edge key;
    Node *l, *r;
    ll delta;
    void prop() {
        key.w += delta;
        if (l) l->delta += delta;
        if (r) r->delta += delta;
```

```
6.8 Math
```

6.8.1 Number of Spanning Trees

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

6.8.2 Erdős–Gallai theorem

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

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

Strings (7)

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: $\mathcal{O}(n)$

```
VI pi(const string& s) {
    VI p(SZ(s));
    REP(i,1,SZ(s)) {
        int g = p[i-1];
        while (g && s[i] != s[g]) g = p[g-1];
        p[i] = g + (s[i] == s[g]);
    }
    return p;
}

VI match(const string& s, const string& pat) {
    VI p = pi(pat + '\0' + s), res;
    REP(i,SZ(p)-SZ(s),SZ(p))
        if (p[i] == SZ(pat)) res.push_back(i - 2 * SZ(pat));
    return res;
}
```

Zfunc.h

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

```
VI Z(string S) {
VI z(SZ(S));
int l = -1, r = -1;
REP(i,1,SZ(S)) {
    z[i] = i >= r ? 0 : min(r - i, z[i - l]);
    while (i + z[i] < SZ(S) && S[i + z[i]] == S[z[i]])
    z[i]++;
    if (i + z[i] > r)
        l = i, r = i + z[i];
}
return z;
```

Manacher.h

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

```
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]);
    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++;</pre>
```

```
if (R>r) l=L, r=R;
  return p;
MinRotation.h
Description: Finds the lexicographically smallest rotation of a string.
Usage: rotate(v.begin(), v.begin()+minRotation(v), v.end());
Time: \mathcal{O}(N)
int minRotation(string s) {
  int a=0, N=SZ(s); s += s;
  REP(b,0,N) REP(k,0,N) {
    if (a+k == b \mid | s[a+k] < s[b+k]) \{b += max(0, k-1); break; \}
    if (s[a+k] > s[b+k]) { a = b; break; }
  return a;
```

SuffixArray.h

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: $\mathcal{O}(n \log n)$ 3e14f8, 23 lines

```
struct SuffixArray {
  VI sa. lcp:
  SuffixArray(string& s, int lim=256) { // or basic string<int>
    int n = SZ(s) + 1, k = 0, a, b;
    VI x(ALL(s)+1), y(n), ws(max(n, lim)), 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) a = 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)</pre>
     for (k \&\& k--, j = sa[rank[i] - 1];
          s[i + k] == s[i + k]; k++);
};
```

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).

struct SuffixTree {

int toi(char c) { return c - 'a'; } string a; // v = cur node, q = cur positionint t[N][ALPHA], l[N], r[N], p[N], s[N], v=0, q=0, m=2; void ukkadd(int i, int c) { suff:

```
Time: \mathcal{O}(26N)
                                                          5e69b7, 50 lines
  enum { N = 200010, ALPHA = 26 }; // N \sim 2*maxlen+10
    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; goto suff;
  SuffixTree(string a) : a(a) {
    fill(r,r+N,SZ(a));
    memset(s, 0, sizeof s);
    memset(t, -1, sizeof t);
    fill(t[1],t[1]+ALPHA,0);
    s[0] = 1; l[0] = l[1] = -1; r[0] = r[1] = p[0] = p[1] = 0;
    REP(i,0,SZ(a)) ukkadd(i, toi(a[i]));
  // example: find longest common substring (uses ALPHA = 28)
  int lcs(int node, int i1, int i2, int olen) {
    if (l[node] <= i1 && i1 < r[node]) return 1;</pre>
    if (l[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.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(0,A,B) H operator O(H o) { ull r = x; asm \
  (A "addq %rdx, %0\n adcq $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 qet() const { return x + !\sim x; }
  bool operator==(H o) const { return get() == o.get(); }
  bool operator<(H o) const { return get() < o.get(); }</pre>
static const H C = (ll)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};
  REP(i,length,SZ(str)) {
    ret.push back(h = h * C + str[i] - pw * str[i-length]);
 return ret;
H hashString(string& s){H h{}; for(char c:s) h=h*C+c;return h;}
```

AhoCorasick.h

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

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

```
struct AhoCorasick {
 enum {alpha = 26, first = 'A'}; // change this!
 struct Node {
   // (nmatches is optional)
   int back, next[alpha], start = -1, end = -1, nmatches = 0;
   Node(int v) { memset(next, v, sizeof(next)); }
 vector<Node> N;
 VI backp;
 void insert(string& s, int j) {
   assert(!s.empty());
   int n = 0:
   for (char c : s) {
     int& m = N[n].next[c - first];
     if (m == -1) { n = m = SZ(N); N.emplace back(-1); }
      else n = m;
   if (N[n].end == -1) N[n].start = i:
   backp.push back(N[n].end);
   N[n].end = j;
   N[n].nmatches++:
 AhoCorasick(vector<string>\& pat) : N(1, -1) {
    REP(i,0,SZ(pat)) insert(pat[i], i);
   N[0].back = SZ(N);
   N.emplace back(0);
    queue<int> q;
   for (q.push(0); !q.empty(); q.pop()) {
     int n = q.front(), prev = N[n].back;
     REP(i,0,alpha)
       int &ed = N[n].next[i], y = N[prev].next[i];
       if (ed == -1) ed = y;
       else {
         N[ed].back = y;
          (N[ed].end == -1 ? N[ed].end : backp[N[ed].start])
```

3af81c, 9 lines

```
= N[y].end;
          N[ed].nmatches += N[y].nmatches;
          q.push(ed);
  VI find(string word) {
    int n = 0;
    VI res: // ll count = 0:
    for (char c : word) {
     n = N[n].next[c - first];
      res.push back(N[n].end):
     // count += N[n].nmatches;
    return res;
  vector<VI> findAll(vector<string>& pat, string word) {
   VI r = find(word):
    vector<VI> res(SZ(word));
   REP(i,0,SZ(word)) {
     int ind = r[i];
      while (ind != -1) {
       res[i - SZ(pat[ind]) + 1].push back(ind);
       ind = backp[ind];
    return res;
};
```

Geometry (8)

8.1 Geometric primitives

Point.h

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

47ec0a. 28 lines

```
template <class T> int 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 sqrt((double)dist2()); }
  // angle to x-axis in interval [-pi, pi]
  double angle() const { return atan2(v, 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.

SegmentDistance.h

Description:

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

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

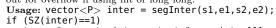
"Point.h" 5c88f4, 6 lines
typedef Point<double> P;

```
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<II> 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.



cout << "segments intersect at " << inter[0] << endl; "Point.h", "OnSegment.h" $$36\mathrm{c2d7},\ 13\ \mathrm{lir}$$

```
template<class P> vector<P> segInter(P a, P b, P c, P d) {
    auto oa = c.cross(d, a), ob = c.cross(d, b),
        oc = a.cross(b, c), od = a.cross(b, d);

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

lineIntersection.h

Description:

If a unique intersection point of the lines going through \$1,e1\$ and \$2,e2 exists $\{1, \text{ point}\}$ is returned. If no intersection point exists $\{0, (0,0)\}$ is returned and if infinitely many exists $\{-1, (0,0)\}$ is returned. The wrong position will be returned if P is Point<||scale="1">| Point<||scale="1



sideOf.h

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

```
Usage: bool left = sideOf(p1,p2,q)==1;
"Point.h"
```

```
template<class P>
int sideOf(P s, P e, P p) { return sgn(s.cross(e, p)); }

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

OnSegment.h

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

```
"Point.h" c597e8, 3 li
```

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

linearTransformation.h Description:

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

```
lation, rotation and \begin{array}{c} p_0 \\ q_0 \end{array} resolved to point r.
```

```
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

"Point.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<\hat{A}ngle> \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;
   int t;
   Angle(int x, int y, int t=0) : x(x), y(y), t(t) {}
   Angle operator-(Angle b) const { return {x-b.x, y-b.y, t}; }
   int half() const {
      assert(x || y);
   }
}
```

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

8.2Circles

CircleIntersection.h

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

```
84d6d3, 11 lines
typedef Point<double> P;
bool circleInter(P a,P b,double r1,double r2,pair<P, P>* out) {
 if (a == b) { assert(r1 != r2); return false; }
  P \text{ vec} = b - a;
  double d2 = vec.dist2(), sum = r1+r2, dif = r1-r2,
          p = (d2 + r1*r1 - r2*r2)/(d2*2), h2 = r1*r1 - p*p*d2;
  if (sum*sum < d2 || dif*dif > d2) return false;
  P \text{ mid} = a + \text{vec*p}, \text{ per} = \text{vec.perp}() * \text{sqrt}(\text{fmax}(0, h2) / d2);
  *out = {mid + per. mid - per}:
  return true;
```

CircleTangents.h

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

```
"Point.h"
                                                      b0153d, 13 lines
template<class P>
vector<pair<P, P>> tangents(P c1, double r1, P c2, double r2) {
 P d = c2 - c1:
  double dr = r1 - r2, d2 = d.dist2(), h2 = d2 - dr * dr;
  if (d2 == 0 || h2 < 0) return {};</pre>
  vector<pair<P, P>> out;
  for (double sign : {-1, 1}) {
   P v = (d * dr + d.perp() * sqrt(h2) * sign) / d2;
    out.push back(\{c1 + v * r1, c2 + v * r2\});
 if (h2 == 0) out.pop back();
  return out;
```

CirclePolygonIntersection.h

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

```
Time: \mathcal{O}(n)
                                                       f5c096, 19 lines
"../../content/geometry/Point.h"
typedef Point<double> P:
#define arg(p, q) atan2(p.cross(q), p.dot(q))
double circlePoly(P c, double r, vector<P> ps) {
 auto tri = [\&](Pp, Pq) {
    auto r2 = r * r / 2:
    Pd = q - p;
    auto a = d.dot(p)/d.dist2(), b = (p.dist2()-r*r)/d.dist2();
    auto det = a * a - b;
    if (det <= 0) return arg(p, q) * r2;</pre>
    auto s = max(0., -a-sqrt(det)), t = min(1., -a+sqrt(det));
    if (t < 0 || 1 <= s) return arg(p, q) * r2;</pre>
    P u = p + d * s, v = p + d * t;
    return arg(p,u) * r2 + u.cross(v)/2 + arg(v,q) * r2;
  auto sum = 0.0;
 REP(i, 0, SZ(ps))
   sum += tri(ps[i] - c, ps[(i + 1) % SZ(ps)] - c);
 return sum:
```

circumcircle.h

Description:

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



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

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

MinimumEnclosingCircle.h

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

```
"circumcircle.h"
                                                     69dd52, 17 lines
pair<P, double> mec(vector<P> ps) {
 shuffle(ALL(ps), mt19937(time(0)));
 P \circ = ps[0];
 double r = 0. EPS = 1 + 1e-8:
 REP(i,0,SZ(ps)) if ((o - ps[i]).dist() > r * EPS) {
   o = ps[i], r = 0;
    REP(j,0,i) if ((o - ps[j]).dist() > r * EPS) {
     o = (ps[i] + ps[j]) / 2;
     r = (o - ps[i]).dist();
     REP(k,0,j) if ((o - ps[k]).dist() > r * EPS) {
       o = ccCenter(ps[i], ps[j], ps[k]);
       r = (o - ps[i]).dist();
 return {o, r};
```

8.3 Polygons

```
InsidePolygon.h
```

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

```
bool in = inPolygon(v, P(3, 3), false);
Time: \mathcal{O}(n)
"Point.h", "OnSegment.h", "SegmentDistance.h"
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;</pre>
    cnt ^= ((a.y<p[i].y) - (a.y<q.y)) * a.cross(p[i], q) > 0;
  return cnt;
```

Usage: vector $P> v = \{P\{4,4\}, P\{1,2\}, P\{2,1\}\};$

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"

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

PolygonCenter.h

Description: Returns the center of mass for a polygon.

```
Time: \mathcal{O}\left(n\right)
"Point.h"
```

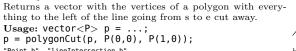
7d84e0, 9 lines

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

PolygonCut.h

Description:

thing to the left of the line going from s to e cut away.



```
"Point.h", "lineIntersection.h"
                                                       b09131, 13 lines
typedef Point<double> P:
vector<P> polygonCut(const vector<P>& poly, P s, P e) {
  vector<P> res:
  REP(i.0.SZ(polv)) {
    P cur = poly[i], prev = i ? poly[i-1] : poly.back();
    bool side = s.cross(e, cur) < 0;</pre>
    if (side != (s.cross(e, prev) < 0))</pre>
      res.push back(lineInter(s, e, cur, prev).second);
    if (side)
      res.push back(cur);
 return res;
```

ConvexHull.h

Description:

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



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

HullDiameter.h

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

PointInsideHull.h

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

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

```
"Point.h", "sideOf.h", "OnSegment.h" efb6da, 14 lines
```

typedef Point<ll> P;

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

LineHullIntersection.h

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

```
Time: \mathcal{O}(\log n)
"Point.h"
                                                     331463, 39 lines
#define cmp(i,j) sqn(dir.perp().cross(poly[(i)%n]-poly[(j)%n]))
#define extr(i) cmp(i + 1, i) >= 0 && cmp(i, i - 1 + n) < 0
template <class P> int extrVertex(vector<P>& poly, P dir) {
  int n = SZ(poly), lo = 0, hi = n;
  if (extr(0)) return 0;
  while (lo + 1 < hi) {
    int m = (lo + hi) / 2;
    if (extr(m)) return m;
    int ls = cmp(lo + 1, lo), ms = cmp(m + 1, m):
    (ls < ms \mid | (ls == ms \&\& ls == cmp(lo, m)) ? hi : lo) = m;
  return lo;
#define cmpL(i) sqn(a.cross(poly[i], b))
template <class P>
array<int, 2> lineHull(P a, P b, vector<P>& poly) {
  int endA = extrVertex(poly, (a - b).perp());
  int endB = extrVertex(poly, (b - a).perp());
  if (cmpL(endA) < 0 \mid | cmpL(endB) > 0)
    return {-1, -1};
  array<int, 2> res;
  REP(1.0.2) {
    int lo = endB. hi = endA. n = SZ(polv):
    while ((lo + 1) % n != hi) {
      int m = ((lo + hi + (lo < hi ? 0 : n)) / 2) % n;
      (cmpL(m) == cmpL(endB) ? lo : hi) = m:
    res[i] = (lo + !cmpL(hi)) % n:
    swap(endA, endB):
  if (res[0] == res[1]) return {res[0], -1}:
  if (!cmpL(res[0]) && !cmpL(res[1]))
    switch ((res[0] - res[1] + SZ(poly) + 1) % SZ(poly)) {
      case 0: return {res[0], res[0]};
      case 2: return {res[1], res[1]};
  return res:
```

8.4 Misc. Point Set Problems

```
ClosestPair.h
Description: Finds the closest pair of points.
Time: \mathcal{O}(n \log n)
"Point.h"
                                                       ac393c, 17 lines
typedef Point<ll> P;
pair<P. P> closest(vector<P> v) {
 assert(SZ(v) > 1);
 set<P> S:
 sort(ALL(v), [](P a, P b) \{ return a.y < b.y; \});
 pair<ll, pair<P, P>> ret{LLONG MAX, {P(), P()}};
 int j = 0:
 for (P p : v) {
    P d\{1 + (ll) sqrt(ret.first), 0\};
    while (v[i].v \le p.v - d.x) S.erase(v[i++]);
    auto lo = S.lower bound(p - d), hi = S.upper bound(p + d);
    for (; lo != hi; ++lo)
      ret = min(ret, {(*lo - p).dist2(), {*lo, p}});
    S.insert(p);
 return ret.second;
```

```
kdTree.h
Description: KD-tree (2d, can be extended to 3d)
"Point.h"
                                                     269f22, 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 <math>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 \times 0 = INF, \times 1 = -INF, y0 = INF, y1 = -INF; // bounds
  Node *first = 0, *second = 0;
 T distance(const P& p) { // min squared distance to a point
   T x = (p.x < x0 ? x0 : p.x > x1 ? x1 : p.x);
    T y = (p.y < y0 ? y0 : p.y > y1 ? y1 : p.y);
    return (P(x,y) - p).dist2();
  Node(vector<P>&& vp) : pt(vp[0]) {
    for (P p : vp) {
      x0 = min(x0, p.x); x1 = max(x1, p.x);
      y0 = min(y0, p.y); y1 = max(y1, p.y);
    if (vp.size() > 1) {
      // split on x if width >= height (not ideal...)
      sort(ALL(vp), x1 - x0 >= y1 - y0 ? on x : on y);
      // divide by taking half the array for each child (not
      // best performance with many duplicates in the middle)
      int half = SZ(vp)/2:
      first = new Node({vp.begin(), vp.begin() + half});
      second = new Node({vp.begin() + half, vp.end()});
};
struct KDTree {
 Node* root:
  KDTree(const vector<P>& vp) : root(new Node({ALL(vp)})) {}
  pair<T, P> search(Node *node, const P& p) {
    if (!node->first) {
      // uncomment if we should not find the point itself:
      // if (p == node -> pt) return \{INF, P()\};
      return make pair((p - node->pt).dist2(), node->pt);
    Node *f = node->first, *s = node->second;
    T bfirst = f->distance(p), bsec = s->distance(p):
    if (bfirst > bsec) swap(bsec, bfirst), swap(f, s);
    // search closest side first, other side if needed
    auto best = search(f, p);
    if (bsec < best.first)</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)</pre>
  pair<T, P> nearest(const P& p) {
    return search(root, p):
};
```

```
FastDelaunav.h
Description: Fast Delaunay triangulation. Each circumcircle contains none
of the input points. There must be no duplicate points. If all points are on a
line, no triangles will be returned. Should work for doubles as well, though
there may be precision issues in 'circ'. Returns triangles in order {t[0][0],
t[0][1], t[0][2], t[1][0], \dots\}, all counter-clockwise.
Time: \mathcal{O}(n \log n)
"Point.h"
                                                        aaea8f. 88 lines
typedef Point<ll> P;
typedef struct Ouad* 0:
typedef int128 t lll; // (can be ll if coords are < 2e4)
P arb(LLONG MAX, LLONG MAX); // not equal to any other point
struct Quad {
  Q rot, o; P p = arb; bool mark;
  P& F() { return r()->p; }
  Q& r() { return rot->rot; }
  Q prev() { return rot->o->rot; }
  Q next() { return r()->prev(); }
} *H;
bool circ(P p, P a, P b, P c) { // is p in the circumcircle?
  lll p2 = p.dist2(), A = a.dist2()-p2,
      B = b.dist2()-p2, C = c.dist2()-p2;
  return p.cross(a,b)*C + p.cross(b,c)*A + p.cross(c,a)*B > 0;
Q makeEdge(P orig, P dest) {
  Q r = H ? H : new Quad{new Quad{new Quad{new Quad{0}}}};
  H = r - > 0; r - > r() - > r() = r;
  REP(i,0,4) r = r -> rot, r -> p = arb, r -> o = i & 1 ? r : r -> r();
  r \rightarrow p = orig; r \rightarrow F() = dest;
  return r;
void splice(Q a, Q b) {
  swap(a->o->rot->o, b->o->rot->o); swap(a->o, b->o);
Q connect(Q a, Q b) {
  Q q = makeEdge(a->F(), b->p);
  splice(q, a->next());
  splice(q->r(), b);
  return q;
pair<0,0> rec(const vector<P>& s) {
  if (SZ(s) <= 3)
    Q a = makeEdge(s[0], s[1]), b = makeEdge(s[1], s.back());
    if (SZ(s) == 2) return { a, a->r() };
    splice(a->r(), b);
    auto side = s[0].cross(s[1], s[2]);
    Q c = side ? connect(b, a) : 0;
    return {side < 0 ? c->r() : a, side < 0 ? c : b->r() };
#define H(e) e->F(), e->p
#define valid(e) (e->F().cross(H(base)) > 0)
  0 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\rightarrow p.cross(H(A)) < 0 \&\& (A = A\rightarrow next())))
         (A->p.cross(H(B)) > 0 \&\& (B = B->r()->o)));
  Q base = connect(B->r(), A);
  if (A->p == ra->p) ra = base->r();
  if (B->p == rb->p) rb = base;
#define DEL(e, init, dir) Q e = init->dir; if (valid(e)) \
    while (circ(e->dir->F(), H(base), e->F())) { \
      0 t = e->dir; \
      splice(e, e->prev()); \
```

```
splice(e->r(), e->r()->prev()); \
     e->0 = H: H = e: e = t: \
  for (;;) {
   DEL(LC, base->r(), o); DEL(RC, base, prev());
    if (!valid(LC) && !valid(RC)) break;
    if (!valid(LC) || (valid(RC) && circ(H(RC), H(LC))))
      base = connect(RC, base->r());
     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 {};
 0 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 \rightarrow mark = 1; pts.push back(c \rightarrow p); \
 q.push back(c->r()); c = c->next(); \} while (c = e); \}
 ADD; pts.clear();
 while (qi < SZ(q)) if (!(e = q[qi++]) -> mark) ADD;
 return pts;
8.5
      3D
PolyhedronVolume.h
Description: Magic formula for the volume of a polyhedron. Faces should
point outwards.
                                                      3058c3, 6 lines
template<class V, class L>
double signedPolvVolume(const V& p, const L& trilist) {
 double v = 0:
 for (auto i : trilist) v += p[i.a].cross(p[i.b]).dot(p[i.c]);
 return v / 6:
Point3D.h
Description: Class to handle points in 3D space. T can be e.g. double or
long long.
                                                     8058ae, 32 lines
template<class T> struct Point3D {
 typedef Point3D P;
 typedef const P& R;
 T x, y, z;
 explicit Point3D(T x=0, T y=0, T z=0) : x(x), y(y), z(z) {}
 bool operator<(R p) const {</pre>
    return tie(x, y, z) < tie(p.x, p.y, p.z); }
 bool operator==(R p) const {
    return tie(x, y, z) == tie(p.x, p.y, p.z); }
  P operator+(R p) const { return P(x+p.x, y+p.y, z+p.z); }
  P operator-(R p) const { return P(x-p.x, y-p.y, z-p.z); }
  P operator*(T d) const { return P(x*d, y*d, z*d); }
  P operator/(T d) const { return P(x/d, y/d, z/d); }
 T dot(R p) const { return x*p.x + y*p.y + z*p.z; }
 P cross(R p) const {
    return P(y*p.z - z*p.y, z*p.x - x*p.z, x*p.y - y*p.x);
 T dist2() const { return x*x + y*y + z*z; }
 double dist() const { return sqrt((double)dist2()); }
  //Azimuthal angle (longitude) to x-axis in interval [-pi, pi]
 double phi() const { return atan2(y, x); }
 //Zenith angle (latitude) to the z-axis in interval [0, pi]
  double theta() const { return atan2(sqrt(x*x+y*y),z); }
  P unit() const { return *this/(T)dist(); } //makes dist()=1
 //returns unit vector normal to *this and p
```

```
P normal(P p) const { return cross(p).unit(); }
  //returns point rotated 'angle' radians ccw around axis
  P rotate(double angle, P axis) const {
    double s = sin(angle), c = cos(angle); P u = axis.unit();
    return u*dot(u)*(1-c) + (*this)*c - cross(u)*s;
};
3dHull.h
Description: Computes all faces of the 3-dimension hull of a point set. *No
four points must be coplanar*, or else random results will be returned. All
faces will point outwards.
Time: \mathcal{O}(n^2)
"Point3D.h"
                                                      0754b0, 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); }
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[i] - A[i]).cross((A[k] - A[i]));
    if (q.dot(A[l]) > q.dot(A[i]))
      q = q * -1;
    F f{q, i, j, k};
    E(a,b).ins(k); E(a,c).ins(j); E(b,c).ins(i);
    FS.push back(f);
  REP(i,0,4) REP(j,i+1,4) REP(k,j+1,4)
    mf(i, j, k, 6 - i - j - k);
  REP(i,4,SZ(A)) {
    REP(j,0,SZ(FS)) {
      F f = FS[j];
      if(f.q.dot(A[i]) > f.q.dot(A[f.a])) {
        E(a,b).rem(f.c);
        E(a,c).rem(f.b);
        E(b,c).rem(f.a);
        swap(FS[i--], FS.back());
        FS.pop back();
    int nw = SZ(FS);
    REP(j,0,nw)
      F f = FS[j];
#define C(a, b, c) if (E(a,b).cnt() != 2) mf(f.a, f.b, i, f.c);
      C(a, b, c); C(a, c, b); C(b, c, a);
  for (F& it : FS) if ((A[it.b] - A[it.a]).cross(
    A[it.c] - A[it.a]).dot(it.q) \ll 0) swap(it.c, it.b);
  return FS;
```

sphericalDistance.h

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

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

Mathematics (9)

9.1 Equations

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

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

$$ax + by = e$$

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

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

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

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

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

9.2 Recurrences

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

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

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

9.3 Trigonometry

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

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

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

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

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

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

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

9.4 Geometry

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

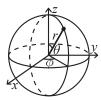
9.4.2 Quadrilaterals

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 $A = \sqrt{(p-a)(p-b)(p-c)(p-d)}$.

9.4.3 Spherical coordinates



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

9.5 Derivatives/Integrals

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

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

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

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

Integration by parts:

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

9.6 Sums

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

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

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

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

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

9.7Series

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

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

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

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

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

IntervalContainer IntervalCover

9.8 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,

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

9.8.1 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 $Po(\lambda)$, $\lambda = t\kappa$.

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

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

9.8.2 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 $\text{Exp}(\lambda)$, $\lambda > 0$.

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

Normal distribution

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

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

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

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

9.9 Markov chains

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

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

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

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

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

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.en\overline{d}() && it->first <= R) {
   R = max(R, it->second);
    before = it = is.erase(it);
  if (it != is.begin() && (--it)->second >= L) {
   L = min(L, it->first);
    R = max(R. it->second):
    is.erase(it);
 return is.insert(before, {L,R});
void removeInterval(set<PII>& is, int L, int R) {
 if (L == R) return;
 auto it = addInterval(is, L, R);
  auto r2 = it->second;
 if (it->first == L) is.erase(it):
  else (int&)it->second = L:
 if (R != r2) is.emplace(R, r2);
```

IntervalCover.h

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

Time: $\mathcal{O}(N \log N)$ 354a6a, 19 lines

```
template<class T>
VI cover(pair<T, T> G, vector<pair<T, T>> I) {
    VI S(SZ(I)), R;
    iota(ALL(S), 0);
    sort(ALL(S), [&](int a, int b) { return I[a] < I[b]; });
    T cur = G.first;
    int at = 0;
    while (cur < G.second) { // (A)
        pair<T, int> mx = make pair(cur, -1);
        while (at < SZ(I) && I[S[at]].first <= cur) {
            mx = max(mx, make_pair(I[S[at]].second, S[at]));
            at++;
    }
    if (mx.second == -1) return {};
    cur = mx.first;
        R.push_back(mx.second);
}</pre>
```

```
IIITH
  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){...});
Time: \mathcal{O}\left(k\log\frac{n}{h}\right)
template<class F, class G, class T>
void rec(int from, int to, F& f, G& q, int& i, T& p, T q) {
  if (p == q) return;
  if (from == to) {
    q(i, to, p);
    i = to; p = q;
  } else {
    int mid = (from + to) >> 1;
    rec(from, mid, f, g, i, p, f(mid));
    rec(mid+1, to, f, q, i, p, q);
template<class F, class G>
void constantIntervals(int from, int to, F f, G g) {
  if (to <= from) return;</pre>
  int i = from; auto p = f(i), q = f(to-1);
  rec(from, to-1, f, g, i, p, q);
  g(i, to, q);
10.2 Misc. algorithms
TernarySearch.h
Description: Find the smallest i in [a,b] that maximizes f(i), assuming
that f(a) < \ldots < f(i) \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))
template<class F>
int ternSearch(int a, int b, F f) {
  assert(a <= b);</pre>
  while (b - a >= 5) {
    int mid = (a + b) / 2;
    if (f(mid) < f(mid+1)) a = mid; // (A)
```

```
a9cf52, 11 lines
  else b = mid+1:
REP(i,a+1,b+1) if (f(a) < f(i)) a = i; // (B)
return a;
```

LIS.h

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

```
template<class I> VI lis(const vector<I>& S) {
 if (S.empty()) return {};
 VI prev(SZ(S));
 typedef pair<I, int> p;
 vector res;
  REP(i,0,SZ(S)) {
   // change 0 -> i for longest non-decreasing subsequence
   auto it = lower bound(ALL(res), p{S[i], 0});
   if (it == res.end()) res.emplace back(), it = res.end()-1;
    *it = \{S[i], i\};
   prev[i] = it == res.begin() ? 0 : (it-1)->second;
  int L = SZ(res), cur = res.back().second;
```

```
VI ans(L);
while (L--) ans[L] = cur, cur = prev[cur];
return ans:
```

Dynamic programming 10.3

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

```
struct DP { // Modify at will:
 int lo(int ind) { return 0; }
 int hi(int ind) { return ind; }
 ll f(int ind, int k) { return dp[ind][k]; }
 void store(int ind, int k, ll v) { res[ind] = PII(k, v); }
 void rec(int L, int R, int LO, int HI) {
   if (L >= R) return:
   int mid = (L + R) \gg 1;
   pair<ll, int> best(LLONG MAX, L0);
    REP(k, max(L0,lo(mid)), min(HI,hi(mid)))
     best = min(best, make pair(f(mid, k), k));
    store(mid, best.second, best.first);
    rec(L, mid, L0, best.second+1);
    rec(mid+1, R, best.second, HI);
 void solve(int L, int R) { rec(L, R, INT MIN, INT MAX); }
```

10.4 Debugging tricks

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

10.5 Optimization tricks

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

10.5.1 Bit hacks

- x & -x is the least bit in x.
- for (int x = m; x;) { --x &= m; ...} loops over all subset masks of **m** (except **m** itself).
- c = x&-x, r = x+c; $(((r^x) >> 2)/c) | r$ is the next number after X with the same number of bits set.

```
REP(b,0,K) REP(i,0,(1 << K))</li>
   if (i \& 1 << b) D[i] += D[i^(1 << b)];
 computes all sums of subsets.
```

10.5.2 Pragmas

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

FastMod.h

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

```
typedef unsigned long long ull;
struct FastMod {
 ull b, m;
  FastMod(ull b) : b(b), m(-1ULL / b) {}
  ull reduce(ull a) \{ // a \% b + (0 \text{ or } b) \}
    return a - (ull)(( uint128 t(m) * a) >> 64) * b;
};
```

FastInput.h

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

```
Usage: ./a.out < input.txt
```

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

7b3c70, 17 lines

```
inline char gc() { // like getchar()
  static char buf[1 << 16];</pre>
  static size t bc, be;
  if (bc >= be) {
    buf[0] = 0, bc = 0;
    be = fread(buf, 1, sizeof(buf), stdin);
  return buf[bc++]: // returns 0 on EOF
int readInt() {
  int a, c;
  while ((a = gc()) < 40);
  if (a == '-') return -readInt();
  while ((c = gc()) >= 48) a = a * 10 + c - 480;
  return a - 48:
```

BumpAllocator.h

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

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

```
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);

bb66d4, 14 lines

```
char buf[450 << 20] alignas(16);
size_t buf_ind = sizeof buf;

template<class T> struct small {
    typedef T value_type;
    small() {}
    template<class U> small(const U&) {}
    T* allocate(size_t n) {
        buf_ind -= n * sizeof(T);
        buf_ind &= 0 - alignof(T);
        return (T*)(buf + buf_ind);
    }
    void deallocate(T*, size_t) {}
};
```

SIMD.h

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

```
#pragma GCC target ("avx2") // or sse4.1
#include "immintrin.h"
typedef m256i mi;
#define \overline{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?v:x), movemask epi8 (hibits of bytes)
// i32gather_epi32(addr, x, 4): map addr[] over 32-b parts of x
// sad epu8: sum of absolute differences of u8, outputs 4xi64
// maddubs epi16: dot product of unsigned i7's, outputs 16xi15
// madd epi16: dot product of signed i16's, outputs 8xi32
// extractf128 si256(, i) (256->128), cvtsi128 si32 (128->lo32)
// permute2f12\overline{8} si256(x,x,1) swaps 128-bit lanes
// shuffle epi3\overline{2}(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,
// andnot, abs, min, max, sign(1,x), cmp(gt|eq), unpack(lo|hi)
int sumi32(mi m) { union {int v[8]; mi m;} u; u.m = m;
  int ret = 0; REP(i,0,8) ret += u.v[i]; return ret; }
mi zero() { return mm256 setzero si256(); }
```

Techniques (A)

techniques.txt

Recursion

159 lines

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

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

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

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