

University of Calgary

Codetoads

Rosa Shah, Charlie Zheng, Kingsley Zhong

1	Contest	<pre>typedef long long ll; typedef pair<int, int=""> pii; typedef vector<int> vi;</int></int,></pre>				
2	Data structures	typedef vector <vi> vvi;</vi>				
3	Graph 3.1 Fundamentals 3.2 Network flow 3.2 Network flow 3.3 Matching 3.4 DFS algorithms 3.5 Coloring 3.5 Heuristics 3.7 Trees 3.8 Math 3.8 Math	4 6 7 8 8 8 8	<pre>int main() { cin.tie(0)->sync_with_stdio(0); cin.exceptions(cin.failbit); // cout.precision(17); // freopen(".in", "r", stdin); // freopen(".out", "w", stdout); } bashre</pre>			
4	Mathematics 4.1 Geometry	10 11 11 11	alias c='g++ -Wall -Wconversion -Wfatal-errors -g -std=c++17 \ -fsanitize=undefined, address' hash.sh # Hashes a file, ignoring all whitespace and comments. Use for # verifying that code was correctly typed. # Usage: ./hash.sh < FILE (make executable first: chmod +x hash. sh) # cpp -dD -P -fpreprocessed tr -d '[:space:]' md5sum cut -c-cpp -dD -P -fpreprocessed tr -d '[:space:]' md5sum awk '{			
5	Numerical5.1 Polynomials and recurrences5.2 Optimization5.3 Matrices5.4 Fourier transforms	12 13	stresstest.sh # stresstest.sh			
	Number theory 6.1 Modular arithmetic 6.2 Primality 6.3 Divisibility 6.4 Fractions 6.5 Pythagorean Triples 6.6 Primes 6.7 Estimates 6.8 Mobius Function Combinatorial	15 15 15 16 16 16	<pre>g++ \$3 -o optimal_executable # running loop for n times (N files) for ((i=1; i<=n; ++i)) do python testcase.py ./brute_executable < testcase.txt > brute_out.txt ./optimal_executable < testcase.txt > optimal_out.txt if [[\$(diff brute_out.txt optimal_out.txt)]] then echo "\$(diff -Z brute_out.txt optimal_out.txt)" ></pre>			
	7.1 Permutations	16	done echo "Done" troubleshoot.txt 53 line			
	Geometry 8.1 Geometric primitives	18 19 20 20	Pre-submit: Write a few simple test cases if sample is not enough. Are time limits close? If so, generate max cases. Is the memory usage fine? Could anything overflow? Make sure to submit the right file. Wrong answer: Print your solution! Print debug output, as well. Are you clearing all data structures between test cases? Can your algorithm handle the whole range of input?			
	Strings Various 10.1 Intervals	23 23 24	Read the full problem statement again. Do you handle all corner cases correctly? Have you understood the problem correctly? Any uninitialized variables? Any overflows? Confusing N and M, i and j, etc.? Are you sure your algorithm works? What special cases have you not thought of? Are you sure the STL functions you use work as you think? Add some assertions, maybe resubmit. Create some testcases to run your algorithm on.			
<pre>Contest (1) template.cpp #include <bits stdc++.h=""> using namespace std;</bits></pre>		lines	Go through the algorithm for a simple case. Go through this list again. Explain your algorithm to a teammate. Ask the teammate to look at your code. Go for a small walk, e.g. to the toilet. Is your output format correct? (including whitespace) Rewrite your solution from the start or let a teammate do it.			
#de	<pre>efine rep(i, a, b) for (int i = a; i < (b); ++i) efine all(x) begin(x), end(x) efine sz(x) (int)(x).size()</pre>		Runtime error: Have you tested all corner cases locally? Any uninitialized variables?			

```
Are you reading or writing outside the range of any vector?
Any assertions that might fail?
Any possible division by 0? (mod 0 for example)
Any possible infinite recursion?
Invalidated pointers or iterators?
Are you using too much memory?
Debug with resubmits (e.g. remapped signals, see Various).
Time limit exceeded:
Do you have any possible infinite loops?
What is the complexity of your algorithm?
Are you copying a lot of unnecessary data? (References)
How big is the input and output? (consider scanf)
Avoid vector, map. (use arrays/unordered_map)
What do your teammates think about your algorithm?
\hbox{\tt Did you do anything unorthodox like string concatenation?}
Memory limit exceeded:
What is the max amount of memory your algorithm should need?
Are you clearing all data structures between test cases?
```

Data structures (2)

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

f2b2fb, 16 lines

HashMap.h

Description: Hash map with mostly the same API as unordered_map, but $\sim 3x$ faster. Uses 1.5x memory. Initial capacity must be a power of 2 (if_7powided).

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

SegmentTree.h

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

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

0f4bdb, 19 l

```
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);
}</pre>
```

LazySegmentTree.h

Description: Segment tree with ability to add or set values of large intervals, and compute max of intervals. Can be changed to other things. Use with a bump allocator for better performance, and SmallPtr or implicit indices to save memory.

```
Usage: Node* tr = new Node(v, 0, sz(v)); 

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

"../various/BumpAllocator.h"
```

34ecf5, 50 lines

```
const int inf = 1e9;
struct Node {
  Node *1 = 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) {</pre>
      int mid = lo + (hi - lo)/2;
      1 = new Node(v, lo, mid); r = new Node(v, mid, hi);
      val = max(1->val, r->val);
    else val = v[lo];
  int query(int L, int R) {
  if (R <= lo || hi <= L) return -inf;</pre>
    if (L <= lo && hi <= R) return val;
    push();
    return max(1->query(L, R), r->query(L, R));
  void set(int L, int R, int x) {
  if (R <= lo || hi <= L) return;</pre>
    if (L <= lo && hi <= R) mset = val = x, madd = 0;
      push(), l->set(L, R, x), r->set(L, R, x);
       val = max(1->val, r->val);
  void add(int L, int R, int x) {
    if (R <= lo || hi <= L) return;</pre>
    if (L <= lo && hi <= R) {
      if (mset != inf) mset += x;
      else madd += x;
      val += x;
    else {
      push(), 1->add(L, R, x), r->add(L, R, x);
       val = max(1->val, r->val);
  void push() {
    if (!1) {
      int mid = 10 + (hi - 10)/2;
      1 = new Node(lo, mid); r = new Node(mid, hi);
    if (mset != inf)
      1->set(lo,hi,mset), r->set(lo,hi,mset), mset = inf;
    else if (madd)
      1-add(lo,hi,madd), r-add(lo,hi,madd), madd = 0;
};
```

UnionFindRollback.h

Description: Disjoint-set data structure with undo. If undo is not needed, skip st, time() and rollback().

```
Usage: int t = uf.time(); ...; uf.rollback(t); Time: \mathcal{O}(\log(N))
```

```
struct RollbackUF {
  vi e; vector<pii> st;
  {\tt 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]); } // Path
       Compression
  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); // Optimization
    st.push_back({a, e[a]});
    st.push_back({b, e[b]});
    e[a] += e[b]; e[b] = a;
    return true;
};
```

Matrix.h

```
Description: Basic operations on square matrices.

Usage: Matrix<int, 3> A;

A.d = {{{{1,2,3}}, {{4,5,6}}, {{7,8,9}}}};

vector<int> vec = {1,2,3};

vec = (A^N) * vec;

template<class T, int N> struct Matrix {
    typedef Matrix M;
    array<array<T, N>, N> d{};
```

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

LineContainer.h

Description: Container where you can add lines of the form kx+m, and query maximum values at points x. 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; } }; struct LineContainer : multiset<Line, less<>>> { // (for doubles, use inf = 1/.0, div(a,b) = a/b) static const ll inf = LLONG_MAX; 11 div(11 a, 11 b) { // floored division
 return a / b - ((a ^ b) < 0 && a % b); }</pre> 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 1 = *lower_bound(x);

return 1.k * x + 1.m;

Description: A short self-balancing tree. It acts as a sequential container with log-time splits/joins, and is easy to augment with additional data. Time: $\mathcal{O}(\log N)$

9556fc, 55 lines struct Node { Node *1 = 0, *r = 0;int val, y, c = 1; Node(int val) : val(val), y(rand()) {} void recalc(); int cnt(Node* n) { return n ? n->c : 0; } void Node::recalc() { c = cnt(1) + cnt(r) + 1; } template<class F> void each(Node* n, F f) { if (n) { each(n->1, f); f(n->val); each(n->r, f); } pair<Node*, Node*> split(Node* n, int k) { if (!n) return {}; if $(cnt(n->1) >= k) { // "n=>val>= k" for lower_bound(k)}$ auto pa = split(n->1, k); n->1 = pa.second; n->recalc(); return {pa.first, n}; auto pa = split (n->r, k - cnt(n->1) - 1); // and just "k" n->r = pa.first; n->recalc();

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

FenwickTree.h

Description: Computes partial sums a[0] + a[1] + ... + a[pos - 1], and updates single elements a[i], taking the difference between the old and new value. **Time:** Both operations are $\mathcal{O}(\log N)$.

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

FenwickTree2d.h

Description: Computes sums a[i,j] for all i<I, j<J, and increases single elements a[i,j]. Requires that the elements to be updated are known in advance (call fakeUpdate() before init()). **Time:** $\mathcal{O}\left(\log^2 N\right)$. (Use persistent segment trees for $\mathcal{O}\left(\log N\right)$.)

```
"FenwickTree.h"
                                                             157f07, 22 lines
struct FT2 {
  vector<vi> ys; vector<FT> ft;
  FT2(int limx) : ys(limx) {}
  void fakeUpdate(int x, int y) {
    for (; x < sz(ys); x = x + 1) ys[x].push_back(y);
  void init() {
    for (vi& v : ys) sort(all(v)), ft.emplace_back(sz(v));
  int ind(int x, int v) {
    return (int) (lower_bound(all(ys[x]), y) - ys[x].begin()); }
  void update(int x, int y, 11 dif) {
  for (; x < sz(ys); x |= x + 1)</pre>
      ft[x].update(ind(x, y), dif);
  11 query(int x, int y) {
    11 \text{ sum} = 0;
    for (; x; x &= x - 1)
      sum += ft[x-1].query(ind(x-1, y));
    return sum:
 }
};
```

```
Description: Sparse Table. Returns min(V[a], V[a + 1], ... V[b - 1]) in con-
stant time.
```

Usage: RMQ rmq(values); rmg.query(inclusive, exclusive);

Time: $\mathcal{O}(|V|\log|V|+Q)$ 510c32, 16 lines

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

Graph (3)

3.1Fundamentals

Dijkstra.h

Description: SSSP Weighted, without negative cycles.

Time: $\mathcal{O}((E+V)\log V)$

1b221b, 18 lines

```
typedef pair<int, int> pii;
typedef vector<pii> vpi;
vector<vpi> adj;
int dist[N]; // N = number of nodes
priority_queue<pii, vector<pii>, greater<pii>> dja;
dja.emplace(0, START_NODE);
fill(dist, dist+N, INT_MAX);
dist[START_NODE] = 0;
while (!dja.empty()) {
  pii pt = dja.top(); dja.pop();
if (pt.first != dist[pt.second]) continue;
   for (pii ps : adj[pt.second]) {
     if (pt.first + ps.second < dist[ps.first]) {</pre>
       dist[ps.first] = pt.first + ps.second;
       dja.emplace(dist[ps.first], ps.first);
     }
  }
```

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: $\mathcal{O}\left(VE\right)$ 830a8f, 23 lines const ll inf = LLONG_MAX; struct Ed { int a, b, w, s() { return a < b ? a : -a; }}; struct Node { ll dist = inf; int prev = -1; }; void bellmanFord(vector<Node>& nodes, vector<Ed>& eds, int s) { nodes[s].dist = 0;sort(all(eds), [](Ed a, Ed b) { return a.s() < b.s(); });</pre> int lim = sz(nodes) / 2 + 2; // /3+100 with shuffled vertices rep(i,0,lim) for (Ed ed : eds) { Node cur = nodes[ed.a], &dest = nodes[ed.b]; if (abs(cur.dist) == inf) continue; 11 d = cur.dist + ed.w; if (d < dest.dist) {</pre> dest.prev = ed.a; dest.dist = (i < lim-1 ? d : -inf);rep(i,0,lim) for (Ed e : eds) { if (nodes[e.a].dist == -inf) nodes[e.b].dist = -inf;

FloydWarshall.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 if no path, or -inf if the path goes through a negativeweight cycle Time: $\mathcal{O}(N^3)$

```
const 11 inf = 1LL << 62;
```

```
void floydWarshall(vector<vector<11>>& m) {
```

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

TopoSort.h

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

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.

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

ArticulationBridge.h

Description: Find vertex/edge that would disconnect the graph.

```
{\bf Usage:} // N is number of vertices
cnt = 0;
adj.assign(N, VI()); // fill adj
dfs_num.assign(N, -1);
dfs_low.resize(N); // initialization not necessary
for (int n = 0; n < N; n++) dfs(n, n, -1);
Time: \mathcal{O}\left(V+E\right)
```

```
52f4d0, 29 lines
vvi adj;
vi dfs_low, dfs_num;
int cnt;
void dfs(int i, int r, int p) { // (cur, root, parent)
  if (dfs_num[i] != -1) return;
  dfs_low[i] = dfs_num[i] = cnt++;
  for (int j : adj[i])
  if (j != p) { // change cond if parallel edges
  if (dfs_num[j] == -1) {
      dfs(j, r, i);
      if (dfs_low[j] >= dfs_num[i]) ap++;
      if (dfs_low[j] > dfs_num[i]) {
        // (i,j) is a bridge
// each **UNORDERED** pair
        // will occur exactly once
      dfs_low[i] = min(dfs_low[i], dfs_low[j]);
      dfs_low[i] = min(dfs_low[i], dfs_num[j]);
  if (ap >= 2) {
      i is an articulation point
      each vertex will only occur once
}
```

3.2 Network flow

PushRelabel.h

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

```
Time: \mathcal{O}\left(V^2\sqrt{E}\right)
                                                               0ae1d4, 48 lines
struct PushRelabel {
  struct Edge {
    int dest, back;
    11 f, c;
  vector<vector<Edge>> g;
  vector<ll> ec;
  vector<Edge*> cur;
  vector<vi> hs; vi H;
  PushRelabel(int n) : g(n), ec(n), cur(n), hs(2*n), H(n) {}
```

531245, 12 lines

```
void addEdge(int s, int t, ll cap, ll rcap=0) {
  if (s == t) return;
  g[s].push_back({t, sz(g[t]), 0, cap});
  g[t].push_back({s, sz(g[s])-1, 0, rcap});
void addFlow(Edge& e, ll f) {
  Edge &back = g[e.dest][e.back];
  if (!ec[e.dest] && f) hs[H[e.dest]].push_back(e.dest);
  e.f += f; e.c -= f; ec[e.dest] += f;
  back.f = f; back.c += f; ec[back.dest] == f;
ll calc(int s, int t) {
  int v = sz(g); H[s] = v; ec[t] = 1;
  vi co(2*v); co[0] = v-1;
rep(i,0,v) cur[i] = g[i].data();
  for (Edge& e : g[s]) addFlow(e, e.c);
  for (int hi = 0;;) {
    while (hs[hi].empty()) if (!hi--) return -ec[s];
    int u = hs[hi].back(); hs[hi].pop_back(); while (ec[u] > 0) // discharge u
       if (\operatorname{cur}[u] == g[u].\operatorname{data}() + \operatorname{sz}(g[u])) {
         H[u] = 1e9;
         for (Edge& e : g[u]) if (e.c && H[u] > H[e.dest]+1)
           H[u] = H[e.dest]+1, cur[u] = &e;
         if (++co[H[u]], !--co[hi] && hi < v)
rep(i,0,v) if (hi < H[i] && H[i] < v)</pre>
                -co[H[i]], H[i] = v + 1;
         hi = H[u];
       } else if (cur[u] \rightarrow c \&\& H[u] == H[cur[u] \rightarrow dest]+1)
         addFlow(*cur[u], min(ec[u], cur[u]->c));
       else ++cur[u];
  }
bool leftOfMinCut(int a) { return H[a] >= sz(g); }
```

MinCostMaxFlow.h

Description: Min-cost max-flow. cap[i][j] != cap[j][i] is allowed; double edges are not. If costs can be negative, call setpi before maxflow, but note that negative cost cycles are not supported. To obtain the actual flow, look at positive values only.

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

fe85cc, 81 lines

```
#include <bits/extc++.h>
const 11 INF = numeric limits<11>::max() / 4;
typedef vector<11> VL;
struct MCMF {
 int N:
  vector<vi> ed, red;
 vector<VL> cap, flow, cost;
  vi seen:
 VL dist, pi;
 vector<pii> par;
 MCMF (int N) :
    N\left(N\right), 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; 11 di;
     _gnu_pbds::priority_queue<pair<11, int>> q;
    vector<decltype(q)::point_iterator> its(N);
    q.push({0, s});
    auto relax = [&](int i, ll cap, ll cost, int dir) {
      11 val = di - pi[i] + cost;
      if (cap && val < dist[i]) {
        dist[i] = val;
        par[i] = {s, dir};
        if (its[i] == q.end()) its[i] = q.push({-dist[i], i});
        else q.modify(its[i], {-dist[i], i});
    };
    while (!q.empty()) {
      s = q.top().second; q.pop();
```

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

EdmondsKarp.h

Description: Flow algorithm with guaranteed complexity $O(VE^2)$. To get edge flow values, compare capacities before and after, and take the positive 482fe0, 35 lines

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

MinCut.h

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

GlobalMinCut.h

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

8b0e19, 21 lines

```
pair<int, vi> globalMinCut(vector<vi> mat) {
```

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

3.3 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(g, btoa);
Time: \mathcal{O}\left(\sqrt{V}E\right)
                                                          f612e4, 42 lines
bool dfs(int a, int L, vector<vi>& g, vi& btoa, vi& A, vi& B) {
  if (A[a] != L) return 0;
  A[a] = -1;
  for (int b : g[a]) if (B[b] == L + 1) {
    B[b] = 0;
    if (btoa[b] == -1 \mid \mid dfs(btoa[b], L + 1, g, btoa, A, B))
      return btoa[b] = a, 1;
  return 0;
int hopcroftKarp(vector<vi>& g, vi& btoa) {
  int res = 0;
  vi A(g.size()), B(btoa.size()), cur, next;
    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 : g[a]) {
        if (btoa[b] == -1) {
          B[b] = lay;
          islast = 1;
        else if (btoa[b] != a && !B[b]) {
          B[b] = lay;
          next.push_back(btoa[b]);
      if (islast) break;
      if (next.empty()) return res;
      for (int a : next) A[a] = lay;
      cur.swap(next);
```

DFSMatching.h

rep(a,0,sz(g))

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

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

res += dfs(a, 0, g, btoa, A, B);

```
Time: \mathcal{O}(VE) 522b98, 22 lines bool find(int j, vector<vi>& g, vi& btoa, vi& vis) { if (btoa[j] == -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>& g, vi& btoa) {
  vi vis;
  rep(i,0,sz(g)) {
    vis.assign(sz(btoa), 0);
    for (int j : g[i])
       if (find(j, g, btoa, vis)) {
        btoa[j] = i;
        break;
     }
  }
  return sz(btoa) - (int)count(all(btoa), -1);
}
```

${\bf Minimum Vertex Cover.h}$

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>& g, 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 : g[i]) if (!seen[e] && match[e] != -1) {
   seen[e] = true;
       q.push_back(match[e]);
  rep(i,0,n) if (!lfound[i]) cover.push_back(i);
  rep(i,0,m) if (seen[i]) cover.push_back(n+i);
  assert(sz(cover) == res);
  return cover;
```

WeightedMatching.h

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 $\cos[N][M]$, where $\cos[i][j] = \cos f$ for L[i] to be matched with R[j] and returns (min cost, match), where L[i] is matched with R[match[i]]. Negate costs for max cost. Requires $N \leq M$. **Time:** $\mathcal{O}\left(N^2M\right)$

```
1e0fe9, 31 lines
pair<int, vi> hungarian(const vector<vi> &a) {
  if (a.empty()) return {0, {}};
int n = sz(a) + 1, m = sz(a[0]) + 1;
  vi u(n), v(m), p(m), ans(n-1);
  rep(i,1,n) {
    p[0] = i;
    int j0 = 0; // add "dummy" worker 0
vi dist(m, INT_MAX), pre(m, -1);
vector<bool> done(m + 1);
    do { // dijkstra
      done[j0] = true;
int i0 = p[j0], j1, delta = INT_MAX;
rep(j,1,m) if (!done[j]) {
         auto cur = a[i0 - 1][j - 1] - u[i0] - v[j];
          if (cur < dist[j]) dist[j] = cur, pre[j] = j0;</pre>
         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: $\mathcal{O}\left(N^3\right)$

```
"../numerical/MatrixInverse-mod.h"

vector<pii> generalMatching(int N, vector<pii>& ed) {
 vector<vector<ll>> mat(N, vector<ll>< (N)), A;
 for (pii pa : ed) {
   int a = pa.first, b = pa.second, r = rand() % mod;
   mat[a][b] = r, mat[b][a] = (mod - r) % mod;
}</pre>
```

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

3.4 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: $sc(graph, [\&](vi\&v) \{ \dots \})$ 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}\left(E+V\right)
                                                             76b5c9, 24 lines
vi val, comp, z, cont;
int Time, ncomps;
template<class G, class F> int dfs(int j, G& 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);
```

${\bf Biconnected Components.h}$

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

```
Usage: int eid = 0; ed.resize(N); for each edge (a,b) { ed[a].emplace.back(b, eid); ed[b].emplace.back(a, eid++); } bicomps([&] (const vi& edgelist) \{...\}); Time: \mathcal{O}(E+V)
```

vi num, st;
vector<vector<pii>>> ed;
int Time;
template<class F>
int dfs(int at, int par, F& f) {
 int me = num[at] = ++Time, e, y, top = me;
 for (auto pa : ed[at]) if (pa.second != par) {

```
tie(y, e) = pa;
    if (num[v]) {
      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 */ }
 return top;
}
template<class F>
void bicomps(F f) {
  num.assign(sz(ed), 0);
  rep(i,0,sz(ed)) if (!num[i]) dfs(i, -1, f);
```

2sat h

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

```
Usage: TwoSat ts(number of boolean variables); ts.either(0, \sim3); // Var 0 is true or var 3 is false ts.setValue(2); // Var 2 is true ts.atMostOne(\{0, \sim 1, 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.

```
struct TwoSat {
 int N;
  vector<vi> gr;
  vi values; // 0 = false, 1 = true
  TwoSat(int n = 0) : N(n), qr(2*n) {}
  int addVar() { // (optional)
    gr.emplace_back();
    gr.emplace_back();
    return N++;
 void either(int f, int j) {
   f = max(2*f, -1-2*f);

j = max(2*j, -1-2*j);
    gr[f].push_back(j^1);
    gr[j].push_back(f^1);
 void setValue(int x) { either(x, x); }
  void atMostOne(const vi& li) { // (optional)
    if (sz(li) <= 1) return;</pre>
    int cur = ~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 : gr[i]) if (!comp[e])
      low = min(low, val[e] ?: dfs(e));
    if (low == val[i]) do {
      x = z.back(); z.pop_back();
      comp[x] = low;
      if (values[x>>1] == -1)
        values[x>>1] = x&1;
    } while (x != i);
    return val[i] = low;
 bool solve() {
```

values.assign(N, -1);

2965e5, 33 lines

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

3.5 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: $\mathcal{O}\left(NM\right)$

e210e2, 31 lines

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

3.6 Heuristics

MaximalCliques.h

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

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

b0d5b1, 12 lines

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

MaximumClique.h

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

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

f7c0bc, 49 lines

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

MaximumIndependentSet.h

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

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

bfce85, 25 lines

```
vector<vi> treeJump(vi& P){
  int on = 1, d = 1;
  while (on < sz(P)) on *= 2, d++;
  vector<vi> jmp(d, P);
  rep(i,1,d) rep(j,0,sz(P))
    jmp[i][j] = jmp[i-1][jmp[i-1][j]];
  return jmp;
int jmp(vector<vi>& tbl, int nod, int steps){
  rep(i,0,sz(tbl))
    if(steps&(1<<i)) nod = tbl[i][nod];</pre>
  return nod;
int lca(vector<vi>& tbl, vi& depth, int a, int b) {
 if (depth[a] < depth[b]) swap(a, b);</pre>
  a = jmp(tbl, a, depth[a] - depth[b]);
  if (a == b) return a;
  for (int i = sz(tbl); i--;) {
    int c = tbl[i][a], d = tbl[i][b];
    if (c != d) a = c, b = d;
```

```
return tbl[0][a];
}
```

LCA.h

Description: Data structure for computing lowest common ancestors in a tree (with 0 as root). C should be an adjacency list of the tree, either directed or undirected.

Time: $O(N \log N + Q)$

```
"../data-structures/RMQ.h"
                                                              0f62fb, 21 lines
struct LCA {
  int T = 0;
  vi time, path, ret;
  RMQ<int> rmq;
  \label{eq:local_local_local_local} \mbox{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[rmq.query(a, b)];
  //dist(a,b) {return depth[a] + depth[b] - 2*depth[lca(a,b)];}
```

CompressTree.h

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

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

```
"LCA.h"
                                                            9775a0, 21 lines
typedef vector<pair<int, int>> vpi;
vpi compressTree(LCA& lca, const vi& subset) {
  static vi rev; rev.resize(sz(lca.time));
  vi li = subset, &T = lca.time;
  auto cmp = [&](int a, int b) { return T[a] < T[b]; };</pre>
  sort(all(li), cmp);
  int m = sz(li)-1;
  rep(i,0,m) {
    int a = li[i], b = li[i+1];
    li.push_back(lca.lca(a, b));
  sort(all(li), cmp);
  li.erase(unique(all(li)), li.end());
rep(i,0,sz(li)) rev[li[i]] = i;
  vpi ret = {pii(0, li[0])};
  rep(i, 0, sz(li)-1) {
    int a = li[i], b = li[i+1];
    ret.emplace_back(rev[lca.lca(a, b)], b);
  return ret;
```

HLD.h

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

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

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

```
for (int u : adj[v]) {
      rt[u] = (u == adj[v][0] ? rt[v] : u);
      dfsHld(u);
    }
  template <class B> void process(int u, int v, 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 1, int r) { tree->add(1, r, val); });
  int queryPath(int u, int v) { // Modify depending on problem
    int res = -1e9;
    process(u, v, [&](int 1, int r) {
        res = max(res, tree->query(1, r));
    });
    return res:
  int querySubtree (int v) { // modifySubtree is similar
    return tree->query(pos[v] + VALS_EDGES, pos[v] + siz[v]);
};
```

LinkCutTree.h

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

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

5909e2, 90 lines

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

DirectedMST.h

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

```
Time: \mathcal{O}\left(E\log V\right)
"../data-structures/UnionFindRollback.h"
                                                           39e620, 60 lines
struct Edge { int a, b; ll w; };
struct Node {
  Edge key;
  Node *1, *r;
  ll delta:
  void prop() {
    key.w += delta;
    if (1) 1->delta += delta;
    if (r) r->delta += delta;
    delta = 0:
  Edge top() { prop(); return key; }
Node *merge(Node *a, Node *b) {
  if (!a || !b) return a ?: b;
  a->prop(), b->prop();
  if (a->key.w > b->key.w) swap(a, b);
  swap(a->1, (a->r = merge(b, a->r)));
void pop(Node*\& a) { a->prop(); a = merge(a->1, a->r); }
pair<ll, vi> dmst(int n, int r, vector<Edge>& g) {
  RollbackUF uf(n);
  vector<Node*> heap(n);
  for (Edge e : g) heap[e.b] = merge(heap[e.b], new Node{e});
  11 \text{ res} = 0;
  vi seen(n, -1), path(n), par(n);
  seen[r] = r;
  vector<Edge> Q(n), in(n, \{-1,-1\}), comp;
  deque<tuple<int, int, vector<Edge>>> cycs;
  rep(s,0,n) {
    int u = s, qi = 0, w;
    while (seen[u] < 0) {
      if (!heap[u]) return {-1,{}};
      Edge e = heap[u] -> top();
      heap[u]->delta -= e.w, pop(heap[u]);
Q[qi] = e, path[qi++] = u, seen[u] = s;
      res += e.w, u = uf.find(e.a);
      if (seen[u] == s) {
        Node* cyc = 0;
         int end = qi, time = uf.time();
        do cyc = merge(cyc, heap[w = path[--qi]]);
        while (uf.join(u, w));
        u = uf.find(u), heap[u] = cyc, seen[u] = -1;
        cycs.push_front({u, time, {&Q[qi], &Q[end]}});
    rep(i, 0, qi) in[uf.find(Q[i].b)] = Q[i];
```

```
for (auto& [u,t,comp] : cycs) { // restore sol (optional)
  uf.rollback(t);
  Edge inEdge = in[u];
  for (auto& e : comp) in[uf.find(e.b)] = e;
  in[uf.find(inEdge.b)] = inEdge;
rep(i,0,n) par[i] = in[i].a;
return {res, par};
```

3.8 Math

3.8.1 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).

3.8.2 Erdős–Gallai theorem

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

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

Mathematics (4)

4.1 Geometry

Triangles 4.1.1

Side lengths: a, b, c

Semiperimeter: $p = \frac{a+b+c}{2}$ Area: $A = \sqrt{p(p-a)(p-b)(p-c)}$

Circumradius: $R = \frac{abc}{dt}$

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

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

4.1.2 Quadrilaterals

With side lengths a, b, c, d, diagonals e, f, diagonals angle θ , area Aand 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 $A = \sqrt{(p-a)(p-b)(p-c)(p-d)}$.

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

Euler/Fermat little's theorem

For any prime p and integer a, $a^p \equiv a \pmod{p}$. If a is not divisible by p, then $a^{p-1} \equiv 1 \pmod{p}$ and a^{p-2} is the modular inverse of a modulo p. More generally, for any coprime n and a, $a^{\varphi(n)} \equiv 1$ (mod n). $\varphi(n)$ is Euler's totient function, the number positive integers up to a given integer n that are relatively prime to n. If gcd(m, n) = 1, then $\varphi(mn) = \varphi(m)\varphi(n)$ (multiplicative property). For all n and m, and $e \ge \log_2(m)$, it holds that n^e $\pmod{m} \equiv n^{\varphi(m)+e \mod{\varphi(m)}} \pmod{m}$. Starting from $n=1, \varphi(n)$ values: 1, 1, 2, 2, 4, 2, 6, 4, 6.

Great-circle distance

Let ϕ_1, λ_1 and ϕ_2, λ_2 be the geographical latitude and longitude of two points 1 and 2, and R be sphere radius. This Haversine Formula provides higher numerical precision.

$$a = \sin^2\left(\frac{\phi_2 - \phi_1}{2}\right) + \cos(\phi_1)\cos(\phi_2)\sin^2\left(\frac{\lambda_2 - \lambda_1}{2}\right)$$

Great-circle distance = $2 \times R \times \text{atan2} (\sqrt{a}, \sqrt{1-a})$

4.5 Lagrange multiplier

To maximize/minimize f(x,y) subject to g(x,y)=0, you may be able to set partial derivatives of \mathcal{L} to zero, where $\mathcal{L}(x, y, \lambda) = f(x, y) - \lambda \cdot g(x, y)$

Game theory

Applicable to impartial games under the normal play

- Grundy number: Represents Nim pile size. $g(x) = \min (n \ge 0 : n \ne g(y) \,\forall y \in f(x))$
- Remoteness: Moves left if winner forces a win as soon as possible and loser tries to lose as slowly as possible. r(x) = 0if terminal, $1 + \text{least even } r(k), k \in f(x)$ if such exists, otherwise $1 + \text{greatest odd } r(k), k \in f(x)$.
- Suspense function: Moves left if winner tries to play as long as possible and loser tries to lose as soon as possible. s(x) = 0 if terminal, $1 + \text{greatest even } r(k), k \in f(x)$ if such exists, otherwise $1 + \text{least odd } r(k), k \in f(x)$.

Losing conditions:

• Ordinary sum of games: The player whose turn it is must choose one of the games and make a move in it. A player who is not able to move in all the games loses. $g_1 \oplus g_2 \oplus \cdots \oplus g_n = 0.$

- Union of games: The player whose turn it is must choose at least one of these games and make one move in every chosen one. A player who is not able to move loses. $\forall i \ g_i = 0$.
- Selective compound: The player whose turn it is must choose at least one of these subgames, but he cannot choose all of them and then make one move in every chosen one. A player who is not able to move loses. $g_1 = g_2 = ... = g_n$.
- Conjunctive compound: The player whose turn it is must make a move in every subgame. A player who is not able to move loses. $\min(r_1, r_2, \cdots, r_n)$ is even.
- Continued conjunctive compound: The player whose turn it is must make a move in every subgame he can and the game ends and a player loses only if he cannot move anywhere. $\max(s_1, s_2, \cdots, s_n)$ is even.

Numerical (5)

5.1Polynomials and recurrences

Polynomial.h

c9b7b0, 17 lines

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

PolyRoots.h

Description: Finds the real roots to a polynomial.

Usage: polyRoots($\{\{2,-3,1\}\},-1e9,1e9$) // solve $x^2-3x+2=0$ Time: $\mathcal{O}\left(n^2\log(1/\epsilon)\right)$

```
b00bfe, 23 lines
vector<double> polyRoots(Poly p, double xmin, double xmax) {
  if (sz(p.a) == 2) { return {-p.a[0]/p.a[1]}; }
  vector<double> ret;
  Poly der = p;
  der.diff();
  auto dr = polyRoots(der, xmin, xmax);
 dr.push_back(xmin-1);
  dr.push_back(xmax+1);
  sort(all(dr));
  rep(i, 0, sz(dr)-1) {
    double l = dr[i], h = dr[i+1];
    bool sign = p(1) > 0;
if (sign ^ (p(h) > 0)) {
      rep(it, 0, 60) { // while (h - l > 1e-8) double m = (1 + h) / 2, f = p(m);
         if ((f \le 0) ^ sign) 1 = m;
        else h = m;
      ret.push_back((1 + h) / 2);
    }
 return ret;
```

PolyInterpolate.h

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

08bf48, 13 lines

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

```
temp[i] -= last * x[k];
}
return res;
}
```

BerlekampMassey.h

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

 $\begin{tabular}{ll} \textbf{Usage:} & \texttt{berlekampMassey}(\{ \^0,\ 1,\ 1,\ 3,\ 5,\ 11 \})\ //\ \{1,\ 2\} \\ \textbf{Time:} & \mathcal{O}\left(N^2\right) \end{tabular}$

```
"../number-theory/ModPow.h"
                                                                               96548b, 20 lines
vector<11> berlekampMassey(vector<11> s) {
  int n = sz(s), L = 0, m = 0;
vector<11> C(n), B(n), T;
  C[0] = B[0] = 1;
  11 b = 1;
   rep(i,0,n) { ++m;
     11 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;
     \texttt{rep}(\texttt{j},\texttt{m},\texttt{n}) \ \texttt{C}[\texttt{j}] \ = \ (\texttt{C}[\texttt{j}] \ - \ \texttt{coef} \ \star \ \texttt{B}[\texttt{j} \ - \ \texttt{m}]) \ \% \ \texttt{mod};
     if (2 * L > i) continue;
L = i + 1 - L; B = T; b = d; m = 0;
  }
  C.resize(L + 1); C.erase(C.begin());
  for (11& x : C) x = (mod - x) % mod;
  return C;
```

LinearRecurrence.h

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

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

```
typedef vector<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);
  11 \text{ res} = 0;
  rep(i, 0, n) res = (res + pol[i + 1] * S[i]) % mod;
```

5.2 Optimization

GoldenSectionSearch.h

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

```
Usage: double func(double x) { return 4+x+.3*x*x; }
double xmin = gss(-1000,1000,func);
```

```
Time: \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;
```

HillClimbing.h

Description: Poor man's optimization for unimodal functions. 8eeeaf, 14 lines

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

template<class F> pair<double, P> hillClimb(P start, F f) {
  pair<double, P> cur(f(start), start);
  for (double jmp = le9; jmp > le-20; jmp /= 2) {
    rep(j,0,100) rep(dx,-1,2) rep(dy,-1,2) {
        P p = cur.second;
        p[0] += dx*jmp;
        p[1] += dy*jmp;
        cur = min(cur, make_pair(f(p), p));
    }
  }
  return cur;
}
```

Integrate.h

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

```
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; }); }); }); }); }
                                                                92dd79, 15 lines
typedef double d;
#define S(a,b) (f(a) + 4*f((a+b) / 2) + f(b)) * (b-a) / 6
template <class F>
d rec(F\& f, d a, d b, d eps, d S) {
  dc = (a + b) / 2;
  d S1 = S(a, c), S2 = S(c, b), T = S1 + S2;
if (abs(T - S) <= 15 * eps || b - a < 1e-10)
  return T + (T - S) / 15;</pre>
  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^Tx subject to $Ax \leq b, x \geq 0$. Returns -inf if there is no solution, inf if there are arbitrarily good solutions, or the maximum value of c^Tx otherwise. The input vector is set to an optimal x (or in the unbounded case, an arbitrary solution fulfilling the constraints). Numerical stability is not guaranteed. For better performance, define variables such that x=0 is viable.

```
Usage: vvd A = {{1,-1}, {-1,1}, {-1,-2}};

vd b = {{1,1,-4}, c = {-1,-1}, x;

T val = LPSolver(A, b, c).solve(x);
```

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

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

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

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

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

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

5.3 Matrices

Determinant.h

Description: Calculates determinant of a matrix. Destroys the matrix. **Time:** $\mathcal{O}\left(N^3\right)$ bd5cec, 15 lines

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

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

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

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

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

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

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

SolveLinear2.h

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

SolveLinearBinary.h

Description: Solves Ax = b over \mathbb{F}_2 . If there are multiple solutions, one is returned arbitrarily. Returns rank, or -1 if no solutions. Destroys A and b. **Time:** $\mathcal{O}\left(n^2m\right)$

```
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;
    if (br == n) {
      rep(j,i,n) if(b[j]) return -1;
      break;
    int bc = (int)A[br]._Find_next(i-1);
    swap(A[i], A[br]);
swap(b[i], b[br]);
    swap(col[i], col[bc]);
rep(j,0,n) if (A[j][i] != A[j][bc]) {
     A[j].flip(i); A[j].flip(bc);
    rep(j,i+1,n) if (A[j][i]) {
      b[j] ^= b[i];
      A[j] ^= A[i];
   rank++;
 x = bs();
  for (int i = rank; i--;) {
   if (!b[i]) continue;
    x[col[i]] = 1;
    rep(j,0,i) b[j] ^= A[j][i];
  return rank; // (multiple\ solutions\ if\ rank < m)
```

MatrixInverse.h

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

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

rep(i,0,n) {
  int r = i, c = i;
  rep(j,i,n) rep(k,i,n)
```

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

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

```
00ced6, 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 + j] += z;
vd conv(const vd& a, const vd& b) {
  if (a.empty() || b.empty()) return {};
  vd res(sz(a) + sz(b) - 1);
  int L = 32 - \underline{\quad}builtin_clz(sz(res)), n = 1 << L;
 vector < C > in(n), out(n);
  copy(all(a), begin(in));
  rep(i,0,sz(b)) in[i].imag(b[i]);
  fft(in);
  for (C& x : in) x *= x;
  rep(i,0,n) out[i] = in[-i & (n - 1)] - conj(in[i]);
  fft(out);
  rep(i,0,sz(res)) res[i] = imag(out[i]) / (4 * n);
  return res;
```

Number theory (6)

6.1 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.

```
Mod operator/(Mod b) { return *this * invert(b); }
Mod invert(Mod a) {
    11 x, y, g = euclid(a.x, mod, x, y);
    assert(g == 1); return Mod((x + mod) % mod);
}
Mod operator^(11 e) {
    if (!e) return Mod(1);
    Mod r = *this ^ (e / 2); r = r * r;
    return e&1 ? *this * r : r;
}
};
```

ModInverse.h

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

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

ModPow.h

b83e45 8 lines

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

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

ModLog.h

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

11 modLog(11 a, 11 b, 11 m) {
 11 n = (11) sqrt(m) + 1, e = 1, f = 1, j = 1;
 unordered_map<11, 11> A;
 while (j <= n && (e = f = e * a % m) != b % m)
 A[e * b % m] = j++;
 if (e == b % m) return j;
 if (__gcd(m, e) == __gcd(m, b))
 rep(i, 2, n+2) if (A.count(e = e * f % m))
 return n * i - A[e];
 return -1;

ModSum.h

Description: Sums of mod'ed arithmetic progressions.

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

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

5c5bc5, 16 lines

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

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

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

ModMulLL.h

Description: Calculate $a \cdot b \mod c$ (or $a^b \mod c$) for $0 \le a, b \le c \le 7.2 \cdot 10^{18}$. **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 >= (ll)M);
}
ull modpow(ull b, ull e, ull mod) {
    ull ans = 1;
    for (; e; b = modmul(b, b, mod), e /= 2)
        if (e & 1) ans = modmul(ans, b, mod);
    return ans;
}
```

ModSqrt.h

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

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

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

11 x = modpow(a, (s+1) / 2, p);
  11 b = modpow(a, s, p), g = modpow(n, s, p);
  for (;; r = m) {
    11 t = b;
    for (m = 0; m < r \&\& t != 1; ++m)
      t = t * t % p;
    if (m == 0) return x;
    11 \text{ gs} = \text{modpow}(g, 1LL << (r - m - 1), p);
    g = gs * gs % p;
    x = x * gs % p;
    b = b * g % p;
```

6.2 Primality

FastEratosthenes.h

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

Time: LIM=1e9 ≈ 1.5 s 6b2912, 20 lines

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

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 $a^b \mod c$.

Factor.h

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

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

```
if (isPrime(n)) return {n};
ull x = pollard(n);
auto 1 = factor(x), r = factor(n / x);
l.insert(l.end(), all(r));
return 1;
```

6.3 Divisibility

euclid.h

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

```
11 euclid(l1 a, l1 b, l1 &x, l1 &y) {
   if (!b) return x = 1, y = 0, a;
   l1 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 $0 \le x < \operatorname{lcm}(m,n)$. Assumes $mn < 2^{62}$. Time: $\log(n)$

6.3.1 Bézout's identity

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

$$ax + by = d$$

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

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

phiFunction.h

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

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

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

's little thin: p prime $\Rightarrow a^x \equiv 1 \pmod{p} \ \forall a$.

```
const int LIM = 5000000;
int phi[LIM];
void calculatePhi() {
  rep(i,0,LIM) phi[i] = i&1 ? i : i/2;
  for (int i = 3; i < LIM; i += 2) if(phi[i] == i)
    for (int j = i; j < LIM; j += i) phi[j] -= phi[j] / i;
}
```

6.4 Fractions

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

ContinuedFractions.h

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

For consecutive convergents, $p_{k+1}q_k - q_{k+1}p_k = (-1)^k$. $(p_k/q_k$ alternates 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.

```
typedef double d; // for N \sim 1e7; long double for N \sim 1e9 pair<11, 11> approximate (d x, 11 N) {
    11 LP = 0, LQ = 1, P = 1, Q = 0, inf = LLONG_MAX; d y = x; for (;;) {
    11 lim = min(P? (N-LP) / P: inf, Q? (N-LQ) / Q: inf), a = (11) floor(y), b = min(a, lim), NP = b*P + LP, NQ = b*Q + LQ; if (a > b) {
        // If b > a/2, we have a semi-convergent that gives us a // better approximation; if b = a/2, we *may* have one. // Return \{P, Q\} here for a more canonical approximation. return (abs(x - (d)NP / (d)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, 11 N) {
 bool dir = 1, A = 1, B = 1;
Frac lo{0, 1}, hi{1, 1}; // Set hi to 1/0 to search (0, N]
  if (f(lo)) return lo;
  assert(f(hi));
  while (A || B) {
    11 adv = 0, step = 1; // move hi if dir, else lo
    for (int si = 0; step; (step *= 2) >>= si) {
      Frac mid{lo.p * adv + hi.p, lo.q * adv + hi.q};
      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;
```

6.5 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.

6.6 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 $1\,000\,000.\ 10^9+7$ and 10^9+9 are both primes.

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 $\mathbb{Z}_2 \times \mathbb{Z}_{2^{a-2}}$.

6.7 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.

6.8 Mobius Function

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

Mobius Inversion:

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

Other useful formulas/forms:

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

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

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

Combinatorial (7)

7.1 Permutations

7.1.1 Factorial

					8		10	
n!	1 2 6	3 24 1	20 72	0 5040	40320	362880	3628800	-
						16		
n!	4.0e7	7 4.8e	8 6.2e	9 8.7e	10 1.3e	12 2.1el	l3 3.6e14	
n	20	25	30	40	50 1	00 15	13 3.6e14 0 171	
n!	2e18	2e25	3e32	8e47 3	Be64 9e	157 6e2	$62 > DBL_M$	IAX

IntPerm.h

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

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

7.1.2 Cycles

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

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

7.1.3 Derangements

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

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

7.1.4 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

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

7.2 Partitions and subsets

7.2.1 Partition function

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

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

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

7.2.2 Lucas' Theorem

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

7.2.3 Binomials

BinomialMod.h

Description: Binomal coefficients modulo a prime power.

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

```
"./number-theory/euclid.h", "./number-theory/ModInverse.h" 9a67a2, 4 livid fact, invfact; // Precompute this
ll binom(ll n, ll k, ll m) {
    return fact[n] * invfact[k] % m * invfact[n-k] % m;
}
```

multinomial.h

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

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

7.3 General purpose numbers

7.3.1 Bernoulli numbers

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

Sums of powers:

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

Euler-Maclaurin formula for infinite sums:

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

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

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

7.3.3 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. $\pi(j) > j$.

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

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

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

7.3.4 Stirling numbers of the second kind

Partitions of n distinct elements into exactly k groups.

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

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

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

7.3.5 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,

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

7.3.6 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)!)
```

7.3.7 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.
- permutations of [n] with no 3-term increasing subseq.

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

lineDistance.h

Description

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



f6bf6b, 4 lines

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

SegmentDistance.h

typedef Point < double > P;

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);

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

5c88f4, 6 lines

double segDist(P& s, P& e, P& p) {

```
if (s==e) return (p-s).dist();
auto d = (e-s).dist2(), t = min(d, max(.0, (p-s).dot(e-s)));
return ((p-s)*d-(e-s)*t).dist()/d;
```

SegmentIntersection.h

Description:

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



```
Usage: vector<P> inter = segInter(s1,e1,s2,e2);
if (sz(inter) == 1)
cout << "segments intersect at " << inter[0] << endl;
"Point.h", "OnSegment.h"
                                                   9d57f2, 13 lines
template<class P> vector<P> segInter(P a, P b, P c, P d) {
 Checks if intersection is single non-endpoint point.
  \inf (sgn(oa) * sgn(ob) < 0 && sgn(oc) * sgn(od) < 0)
   return {(a * ob - b * oa) / (ob - oa)};
  if (onSegment(c, d, a)) s.insert(a);
  if (onSegment(c, d, b)) s.insert(b);
 if (onSegment(a, b, c)) s.insert(c);
 if (onSegment(a, b, d)) s.insert(d);
 return {all(s)};
```

lineIntersection.h

Description:

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



```
Usage: auto res = lineInter(s1,e1,s2,e2);
if (res.first == 1)
cout << "intersection point at " << res.second << endl;</pre>
"Point.h"
                                                             a01f81, 8 lines
template<class P>
pair<int, P> lineInter(P s1, P e1, P s2, P e2) {
  auto d = (e1 - s1).cross(e2 - s2);
if (d == 0) // if parallel
    return \{-(s1.cross(e1, s2) == 0), P(0, 0)\};
  auto p = s2.cross(e1, e2), q = s2.cross(e2, s1);
  return {1, (s1 * p + e1 * q) / d};
```

sideOf.h

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

```
Usage: bool left = sideOf(p1,p2,q) ==1;
"Point.h"
                                                         3af81c, 9 lines
template<class P>
int sideOf(P s, P e, P p) { return sgn(s.cross(e, p)); }
int sideOf(const P& s, const P& e, const P& p, double eps) {
 auto a = (e-s).cross(p-s);
 double l = (e-s).dist()*eps;
  return (a > 1) - (a < -1);
```

OnSegment.h

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

```
c597e8, 3 lines
template<class P> bool onSegment(P s, P e, P p) {
 return p.cross(s, e) == 0 && (s - p).dot(e - p) <= 0;
```

linearTransformation.h Description:

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



"Point.h" typedef Point < double > P;

```
03a306, 6 lines
```

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

Angle.h

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

```
Usage: vector < Angle > v = \{w[0], w[0].t360() ...\}; // sorted
int j = 0; rep(i,0,n) { while (v[j] < v[i].t180()) ++j; }
// sweeps j such that (j-i) represents the number of positively
oriented triangles with vertices at 0 and i
                                                        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 * (11)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 mid = a + vec*p, per = vec.perp() * sqrt(fmax(0, h2) / d2);
  *out = {mid + per, mid - per};
 return true;
```

CircleTangents.h

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

```
b0153d, 13 lines
"Point.h"
template<class P>
vector<pair<P, P>> tangents(P c1, double r1, P c2, double r2) {
 P d = c2 - c1;
 double dr = r1 - r2, d2 = d.dist2(), h2 = d2 - dr * dr;
  if (d2 == 0 || h2 < 0) return {};</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;
```

f12300, 6 lines

}

CirclePolygonIntersection.h

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

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

```
"../../content/geometry/Point.h"
                                                         a1ee63, 19 lines
typedef Point<double> P;
#define arg(p, q) atan2(p.cross(q), p.dot(q))
double circlePoly(P c, double r, vector<P> ps) {
 auto tri = [\&] (P p, P q) {
    auto r2 = r * r / 2;
    P d = q - p;
    auto a = d.dot(p)/d.dist2(), b = (p.dist2()-r*r)/d.dist2();
    auto det = a * a - b;
    if (det <= 0) return arg(p, q) * r2;</pre>
    auto s = max(0., -a-sqrt(det)), t = min(1., -a+sqrt(det));
    if (t < 0 \mid \mid 1 \le s) return arg(p, q) * r2;
    Pu = p + d * s, v = p + d * t;
    return arg(p,u) * r2 + u.cross(v)/2 + arg(v,q) * r2;
  };
  auto sum = 0.0;
 rep(i, 0, sz(ps))
   sum += tri(ps[i] - c, ps[(i + 1) % sz(ps)] - c);
  return sum;
```

circumcircle.h

Description:

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.



MinimumEnclosingCircle.h

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

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

8.3 Polygons

InsidePolygon.h

return cnt;

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.

```
mediate steps so watch out for overflow.
Usage: vector<P> v = {P{4,4}, P{1,2}, P{2,1}};
bool in = inPolygon(v, P{3, 3}, false);
Time: 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;
      cnt ^= ((a.y<p[i].y) - (a.y<q.y)) * a.cross(p[i], q) > 0;
}
```

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!

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

PolygonCenter.h

Description: Returns the center of mass for a polygon.

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

PolygonCut.h

Description:

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



f2b7d4, 13 lines

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

Usage: vector<P> p = ...;
```

```
p = polygonCut(p, P(0,0), P(1,0));
    "Point.h", "lineIntersection.h"

typedef Point<double> P;
vector<P> polygonCut(const vector<P>& poly, P s, leader of the polygonCut(const vector<P) res;</pre>
```

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

ConvexHull.h

Description:

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



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

HullDiameter.h

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

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

PointInsideHull.h

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

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

```
"Point.h", "sideOf.h", "OnSegment.h" 71446b, 14 lines
typedef Point<11> P;

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

${\bf Line Hull Intersection.h}$

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"
                                                           7cf45b, 39 lines
#define cmp(i,j) sgn(dir.perp().cross(poly[(i)%n]-poly[(j)%n]))
\#define \ extr(i) \ cmp(i + 1, i) >= 0 \&\& \ cmp(i, i - 1 + n) < 0
template <class P> int extrVertex(vector<P>& poly, P dir) {
  int n = sz(poly), lo = 0, hi = n;
  if (extr(0)) return 0;
  while (lo + 1 < hi) {
  int m = (lo + hi) / 2;
    if (extr(m)) return m;
    int 1s = cmp(1o + 1, 1o), ms = cmp(m + 1, m);
    (ls < ms \mid | (ls == ms \&\& ls == cmp(lo, m)) ? hi : lo) = m;
 return lo:
#define cmpL(i) sgn(a.cross(poly[i], b))
template <class P>
array<int, 2> lineHull(P a, P b, vector<P>& poly) {
 int endA = extrVertex(poly, (a - b).perp());
  int endB = extrVertex(poly, (b - a).perp());
  if (cmpL(endA) < 0 \mid \mid cmpL(endB) > 0)
    return {-1, -1};
  array<int, 2> res;
  rep(i,0,2) {
    int lo = endB, hi = endA, n = sz(poly);
    while ((lo + 1) % n != hi) {
      int m = ((lo + hi + (lo < hi ? 0 : n)) / 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: $\mathcal{O}(n \log n)$

```
struct Node
  P pt; // if this is a leaf, the single point in it
  T x0 = INF, x1 = -INF, y0 = INF, y1 = -INF; // bounds
  Node *first = 0, *second = 0;
  T distance(const P& p) { // min squared distance to a point
    T x = (p.x < x0 ? x0 : p.x > x1 ? x1 : p.x);

T y = (p.y < y0 ? y0 : p.y > y1 ? y1 : p.y);
    return (P(x,y) - p).dist2();
  Node(vector<P>&& vp) : pt(vp[0]) {
    for (P p : vp) {
      x0 = min(x0, p.x); x1 = max(x1, p.x);
      y0 = min(y0, p.y); y1 = max(y1, p.y);
    if (vp.size() > 1)
      // split on x if width >= height (not ideal...)
      sort(all(vp), x1 - x0 >= y1 - y0 ? on_x : on_y);

// divide by taking half the array for each child (not

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

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

Point3D.h

Description: Class to handle points in 3D space. T can be e.g. double or long long. \$8058ae, 32\$ lines

```
template<class T> struct Point3D {
  typedef Point3D P;
```

```
typedef const P& R;
  T x, v, 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}\left(n^2\right)
```

```
typedef Point3D<double> P3;
struct PR {
  void ins(int x) { (a == -1 ? a : b) = x; }
  void rem(int x) { (a == x ? a : b) = -1; }
  int cnt() { return (a != -1) + (b != -1); }
  int a, b;
struct F { P3 q; int a, b, c; };
vector<F> hull3d(const vector<P3>& A) {
  assert(sz(A) >= 4);
  \label{eq:vector} \text{vector} < \text{PR} >> \ \texttt{E} \left( \text{sz} \left( \texttt{A} \right) \,, \ \text{vector} < \text{PR} > \left( \text{sz} \left( \texttt{A} \right) \,, \ \left\{ -1 \,, \ -1 \right\} \right) \right) ;
#define E(x,y) E[f.x][f.y]
  vector<F> FS;
  auto mf = [&](int i, int j, int k, int l) {
  P3 q = (A[j] - A[i]).cross((A[k] - A[i]));
     if (q.dot(A[1]) > q.dot(A[i]))
       q = q * -1;
     F f{q, i, j, k};
     E(a,b).ins(k); E(a,c).ins(j); E(b,c).ins(i);
    FS.push_back(f);
  rep(i,0,4) rep(j,i+1,4) rep(k,j+1,4)
    mf(i, j, k, 6 - i - j - k);
  rep(i,4,sz(A)) {
     rep(j,0,sz(FS)) {
       F f = FS[j];
       if(f.q.dot(A[i]) > f.q.dot(A[f.a])) {
         E(a,b).rem(f.c);
          E(a,c).rem(f.b);
          E(b,c).rem(f.a);
          swap(FS[j--], FS.back());
          FS.pop_back();
     int nw = sz(FS);
     rep(j,0,nw) {
       F f = FS[j];
#define C(a, b, c) if (E(a,b).cnt() != 2) mf(f.a, f.b, i, f.c);
       C(a, b, c); C(a, c, b); C(b, c, a);
```

for (F& it : FS) if ((A[it.b] - A[it.a]).cross(
 A[it.c] - A[it.a]).dot(it.q) <= 0) swap(it.c, it.b);</pre>

sphericalDistance.h

return FS;

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);
}
```

Strings (9)

FYI: string.strstr() uses KMP but string.find() does not.

KMP.h

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

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.

```
vi pi(const string& s) {
vi pi(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

5b45fc, 49 lines

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(const 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 - 1]);
    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: 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=0,1=0,r=0; i < n; i++) {
    int t = r-i+!z;
    if (i<r) p[z][i] = min(t, p[z][i+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++;
    if (R>r) l=L, r=R;
  }
  return p;
}
```

MinRotation.h

```
int minRotation(string s) {
```

edce47, 23 lines

```
int a=0, N=sz(s); s += s;
rep(b,0,N) rep(k,0,N) {
  if (a+k == b \mid \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: $\mathcal{O}(n \log n)$ 38db9f, 23 lines

```
struct SuffixArray {
   vi sa, lcp;
   SuffixArray(string& s, int lim=256) { // or basic\_string < int > 
      int n = sz(s) + 1, k = 0, a, b;
vi x(all(s)+1), y(n), ws(max(n, lim)), rank(n);
       for (int j = 0, p = 0; p < n; j = max(1, j * 2), lim = p) {
   p = j, iota(all(y), n - j);
</pre>
           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];
          rep(1,1,11m) ws[1] +- ws[1 - 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)
  for (k && k--, j = sa[rank[i] - 1];
    s[i + k] == s[j + k]; k++);</pre>
};
```

SuffixTree.h

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

Time: $\mathcal{O}(26N)$ aae0b8, 50 lines struct SuffixTree { enum { N = 200010, ALPHA = 26 }; // $N \sim 2*maxlen+10$ enum { N = 200010, Alfrid - 20 }, // N = 200 int toi(char c) { return c - 'a'; } string a; // $v = cur \ node$, $q = cur \ position$ int t[N][ALPHA], 1[N], r[N], p[N], s[N], v=0, q=0, m=2; void ukkadd(int i, int c) { suff: $if (r[v] \le q) {$ 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[v]=q; p[v]=m; t[p[m]][toi(a[l[m]])]=m;v=s[p[m]]; q=l[m];while $(q \le r[m]) \{ v = t[v][toi(a[q])]; q + = r[v] - l[v]; \}$ if (q==r[m]) s[m]=v; else s[m]=m+2; q=r[v]-(q-r[m]); m+=2; goto suff;} SuffixTree(string a) : a(a) { fill(r,r+N,sz(a));memset(s, 0, sizeof s); memset(t, -1, sizeof t); fill(t[1],t[1]+ALPHA,0); s[0] = 1; 1[0] = 1[1] = -1; r[0] = r[1] = p[0] = p[1] = 0;rep(i,0,sz(a)) ukkadd(i, toi(a[i])); // example: find longest common substring (uses ALPHA = 28) int lcs(int node, int i1, int i2, int olen) { if (1[node] <= i1 && i1 < r[node]) return 1;</pre> if (1[node] <= i2 && i2 < r[node]) return 2;</pre> int mask = 0, len = node ? olen + (r[node] - 1[node]) : 0; rep(c, 0, ALPHA) if (t[node][c] != -1) mask |= 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:
};
```

HashKZ.h

Description: My own library, just use multiple hashes if you think test data is evil. See NT/Primes for prime list. Time: $\mathcal{O}(N)$

ce0e10, 21 lines class StringHash{ private: const 11 P=9973, M; vector<11> pow, hash; public: StringHash(const string& s, const ll &M): pow(s.size()+1), $hash(s.size()+1), M(M){$ const int n = s.size(); pow[0] = 1; hash[0] = 0;for(int i=1;i<=n;++i){ pow[i] = (pow[i-1] * P) % M;hash[i] = ((hash[i-1] * P) % M + s[i-1]) % M;} // S and F are 0 indexed and inclusive

Various (10)

};

return (ret % M + M) % M;

10.1 Intervals

IntervalContainer.h

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

set<pii>::iterator addInterval(set<pii>& is, int L, int R) {

if (L == R) return is.end(); auto it = is.lower_bound({L, R}), before = it; while (it != is.end() && it->first <= R) { 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);

IntervalCover.h

else (int&)it->second = L;

if (R != r2) is.emplace(R, r2);

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: O(N \log N)
                                                        9e9d8d, 19 lines
template<class T>
vi cover(pair<T, T> G, vector<pair<T, T>> I) {
 vi S(sz(I)), R;
  iota(all(S), 0);
  sort(all(S), [&](int a, int b) { return I[a] < I[b]; });</pre>
  T cur = G.first;
  int at = 0;
  while (cur < G.second) { // (A)
   pair<T, int> mx = make_pair(cur, -1);
    while (at < sz(I) && I[S[at]].first <= cur) {
     mx = max(mx, make_pair(I[S[at]].second, S[at]));
```

9155b4, 11 lines

b20ccc, 16 lines

```
if (mx.second == -1) return {};
  cur = mx.first;
  R.push back(mx.second);
return R:
```

ConstantIntervals.h

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

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

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

Misc. algorithms 10.2

MergeSort.h

Description: Merge sort with inversion count

Time: $O(N \log N)$

178c99, 25 lines

```
vi merge_sort(vi& arr, int s, int f, ll& inv){
    if(s == f){
        vi ret = {arr[s]};
        return ret;
    int mid = s + (f - s)/2;
    vi v1 = merge_sort(arr, s, mid, inv);
    vi v2 = merge_sort(arr, mid+1, f, inv);
    int n = f-s+1;
    11 s1 = v1.size(), s2 = v2.size();
    vi ret(n);
    11 p1=0, p2=0;
    for(int i=0;i<n;++i){</pre>
        if((p2 == s2 or v1[p1] \le v2[p2]) and p1 != s1){
            ret[i] = v1[p1];
            p1++;
        else{
            ret[i] = v2[p2];
            p2++;
            inv += s1 - p1;
    return ret;
```

BinarySearch.h

Description: Simple binary search

8ff8c7, 16 lines

```
int binarySearch(const vi& arr, int t) {
   int 1 = 0;
    int r = arr.size() - 1;
    int ans = -1;
    while (1 <= r) {
       int mid = 1 + (r - 1) / 2;
       if (arr[mid] >= t) {// lower bound, switch around
            conditions for upper bound
           ans = mid:
           r = mid - 1;
       else {
           1 = mid + 1:
    return ans;
```

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

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

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

FastKnapsack.h

Description: Given N non-negative integer weights w and a non-negative target t, computes the maximum S <= t such that S is the sum of some subset of Time: $O(N \max(w_i))$

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

10.3Dynamic programming

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

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

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: $\mathcal{O}((N + (hi - lo)) \log N)$

```
d38d2b, 18 lines
struct DP { // Modify at will:
 int lo(int ind) { return 0; }
  int hi(int ind) { return ind; }
  11 f(int ind, int k) { return dp[ind][k]; }
  void store(int ind, int k, ll v) { res[ind] = pii(k, v); }
  void rec(int L, int R, int LO, int HI) {
   if (L >= R) return;
    int mid = (L + R) >> 1;
```

ConstantInter-

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

Debugging tricks 10.4

- 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).

Optimization tricks

_builtin_ia32_ldmxcsr(40896); disables denormals (which make floats 20x slower near their minimum value).

10.5.1 GCC Builtins

For these builtins, you can append 1 or 11 to the function names to get the long or long long version.

- int __builtin_ffs(int x) Returns one plus the index of the least significant 1-bit of x. Returns 0 if x = 0.
- int __builtin_clz(unsigned int x) Returns the number of leading 0-bits in $x \neq 0$, starting at the most significant bit position. $|\log_2(n)| = 31 - \text{_builtin_clz}(n)$
- int __builtin_ctz(unsigned int x) Returns the number of trailing 0-bits in $x \neq 0$, starting at the most significant position.
- ullet int __builtin_clrsb(int x) Returns the number of leading redundant sign bits in x, i.e. the number of bits following the most significant bit that are identical to it. There are no special cases for 0 or other values.
- int __builtin_popcount(unsigned int x) Returns the number of 1-bits in x. (Slow on x86 without SSE4 flag)
- ullet int __builtin_parity(unsigned int x) Returns the parity of x, i.e. the number of 1-bits in x modulo 2.
- uintN_t __builtin_bswapN(uintN_t x) Returns x with the order of the bytes reversed. N = 16, 32, 64

10.5.2 Bit hacks

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

10.5.3 Pragmas

- #pragma GCC optimize ("Ofast") will make GCC auto-vectorize loops and optimizes floating points better.
- #pragma GCC target ("avx2") can double performance of vectorized code, but causes crashes on old machines.

• #pragma GCC optimize ("trapv") kills the program on integer overflows (but is really slow).

int128.h

Description: int128 manipulation

10e0c6, 18 lines

```
string printint128(__int128 a) { // prints as decimal
  if (!a) return "0";
  string s;
  while (a) {
   s = char(llabs(a % 10) + '0') + s;
    if (-10 < a && a < 0) s = '-' + s;
    a /= 10;
  return s:
__int128 parseint128(string s) { // parses decimal number
   _{int128 a = 0, sgn = 1;}
  for (char c : s) {
   if (c == '-') sgn *= -1; else a = a * 10 + sgn * (c - '0');
  return a;
```

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 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];
  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() {
  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.

745db2, 8 lines

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

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

```
"BumpAllocator.h"
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; }
```

Techniques (A)

techniques.txt

Hensel lifting

159 lines

```
Recursion
Divide and conquer
 Finding interesting points in N \log N
Algorithm analysis
 Master theorem
 Amortized time complexity
Greedy algorithm
  Scheduling
 Max contiguous subvector sum
  Invariants
 Huffman encoding
Graph theory
 Dynamic graphs (extra book-keeping)
Breadth first search
 Depth first search
  * Normal trees / DFS trees
  Dijkstra's algorithm
 MST: Prim's algorithm
  Bellman-Ford
  Konig's theorem and vertex cover
  Min-cost max flow
  Lovasz toggle
  Matrix tree theorem
  Maximal matching, general graphs
  Hopcroft-Karp
  Hall's marriage theorem
  Graphical sequences
  Floyd-Warshall
  Euler cycles
  Flow networks
  * Augmenting paths
  * Edmonds-Karp
  Bipartite matching
 Min. path cover
  Topological sorting
  Strongly connected components
  2-SAT
  Cut vertices, cut-edges and biconnected components
 Edge coloring
  * Trees
  Vertex coloring
  * Bipartite graphs (=> trees)
  * 3^n (special case of set cover)
 Diameter and centroid
  K'th shortest path
  Shortest cycle
Dynamic programming
  Knapsack
  Coin change
  Longest common subsequence
  Longest increasing subsequence
  Number of paths in a dag
  Shortest path in a dag
  Dynprog over intervals
 Dynprog over subsets
 Dynprog over probabilities
 Dynprog over trees
  3^n set cover
 Divide and conquer
  Knuth optimization
  Convex hull optimizations
  RMQ (sparse table a.k.a 2^k-jumps)
  Bitonic cycle
  Log partitioning (loop over most restricted)
Combinatorics
  Computation of binomial coefficients
  Pigeon-hole principle
  Inclusion/exclusion
  Catalan number
  Pick's theorem
Number theory
  Integer parts
  Divisibility
  Euclidean algorithm
 Modular arithmetic
  * Modular multiplication
  * Modular inverses
  * Modular exponentiation by squaring
  Chinese remainder theorem
  Fermat's little theorem
  Euler's theorem
  Phi function
  Frobenius number
  Quadratic reciprocity
  Pollard-Rho
  Miller-Rabin
```

```
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
  {\tt Root-finding\ with\ binary/ternary\ search}
  Golden section search
Matrices
  Gaussian elimination
  Exponentiation by squaring
Sorting
  Radix sort
Geometry
  Coordinates and vectors
  * Cross product
  * Scalar product
  Convex hull
  Polygon cut
  Closest pair
  Coordinate-compression
  Ouadtrees
  KD-trees
  All segment-segment intersection
Sweeping
  Discretization (convert to events and sweep)
  Angle sweeping
  Line sweeping
  Discrete second derivatives
  Longest common substring
  Palindrome subsequences
  Knuth-Morris-Pratt
  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\star
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
```