

PDPM Indian Institute of Information Technology Design & Manufacturing, Jabalpur

# BLE

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# Contest (1)

```
template.cpp
                                         31 lines
#include <bits/stdc++.h>
using namespace std;
#define rep(i, a, b) for (int i = a; i < (b)
   ; ++i)
#define all(x) begin(x), end(x)
#define sz(x) (int)(x).size()
typedef long long 11;
typedef pair<int, int> pii;
typedef vector<int> vi;
// pbds
#include <ext/pb_ds/assoc_container.hpp>
#include <ext/pb_ds/tree_policy.hpp>
using namespace __gnu_pbds;
template <typename T>
using ordset =
    tree<T, null_type, less<T>, rb_tree_tag,
         tree order statistics node update>;
             // order\_of\_key, find\_by\_order
// GNU hash table
gp hash table<int, int> table;
mt19937 rng(chrono::steady clock::now().
   time_since_epoch().count());
// hash for pairs
struct chash {
  int operator()(pii x) const { return x.
     first * 31 + x.second; }
gp_hash_table<pii, int, chash> table;
int main() {
  cin.tie(0)->sync with stdio(0);
  cin.exceptions(cin.failbit);
.bashrc
alias c='q++ -Wall -Wconversion -pedantic -
   Wfatal-errors -q -std=c++17 \
  -fsanitize=undefined,address -D ONPC'
```

```
stress.sh
#!/usr/bin/env bash
for ((testNum=0;testNum<$4;testNum++))</pre>
    echo $testNum
    ./$3 > input
    ./$2 < input > outSlow
    ./$1 < input > outWrong
    if !(cmp -s "outWrong" "outSlow")
    then
        echo "Error found!"
        echo "Input:"
        cat input
        echo "Wrong Output:"
        cat outWrong
        echo "Slow Output:"
        cat outSlow
        exit
    fi
done
echo Passed $4 tests
troubleshoot.txt
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
Are you clearing all data structures between
    test cases?
Can your algorithm handle the whole range of
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.
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.
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
Avoid vector, map. (use arrays/unordered_map
What do your teammates think about your
   algorithm?
Memory limit exceeded:
```

What is the max amount of memory your algorithm should need?

Are you clearing all data structures between test cases?

# Mathematics (2)

# 2.1 Equations

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

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

$$ax + by = e$$

$$cx + dy = f$$

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

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

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

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

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

# 2.2 Recurrences

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

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

Non-distinct roots r become polynomial factors, e.g.  $a_n = (d_1n + d_2)r^n$ .

# 2.3 Geometry

# 2.3.1 Quadrilaterals

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

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

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

2.4 Sums

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

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

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

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

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

# 2.5 Series

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

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

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

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

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

# 2.6 Probability theory

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

Expectation is linear:

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

For independent X and Y,

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

# Data structures (3)

SegmentTree.h

**Description:**  $O(\log N)Status: stress - tested$ 

91fcdf, 34 line

```
struct segtree {
  int len, maxxn = 1e18;
  vector<int> minn;
  void init(int n) {
    len = 1;
    while (n > len)
      len \star= 2:
    minn.assign(2 * len, maxxn);
  void set(int i, int v, int x, int lx, int
     rx) {
    if (rx - 1x == 1)
      minn[x] = v;
      int m = (lx + rx) / 2;
      if (i < m)
        set(i, v, 2 * x + 1, 1x, m);
        set(i, v, 2 * x + 2, m, rx);
      minn[x] = min(minn[2 * x + 1], minn[2
         * x + 21);
  void set(int i, int v) { set(i, v, 0, 0,
     len); }
  int range(int 1, int r, int x, int lx, int
      rx) {
    if (1 >= rx or r <= lx)
      return maxxn;
    if (1 \le 1x \text{ and } rx \le r)
      return minn[x];
    int m = (lx + rx) / 2;
    auto a = range(1, r, 2 * x + 1, 1x, m);
    auto b = range(1, r, 2 * x + 2, m, rx);
    return min(a, b);
  int range(int 1, int r) { return range(1,
     r + 1, 0, 0, len); }
};
```

```
Usage: Node* tr = new Node(v, 0, sz(v));
Time: \mathcal{O}(\log N).
"../various/BumpAllocator.h"
                                        4f2aa3, 58 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) {
      int mid = lo + (hi - lo) / 2;
      l = new Node(v, lo, mid);
      r = new Node(v, mid, hi);
      val = max(l->val, r->val);
    } else
      val = v[lo];
  int query(int L, int R) {
    if (R <= lo || hi <= L)
      return -inf;
    if (L <= lo && hi <= R)
      return val;
    push();
    return max(l->query(L, R), r->query(L, R)
       ));
  void set(int L, int R, int x) {
    if (R <= lo || hi <= L)
      return;
    if (L <= lo && hi <= R)
      mset = val = x, madd = 0;
    else {
      push(), l\rightarrow set(L, R, x), r\rightarrow set(L, R, x)
         x);
      val = max(l->val, r->val);
    }
  void add(int L, int R, int x) {
    if (R <= lo || hi <= L)
      return;
    if (L <= lo && hi <= R) {
      if (mset != inf)
```

```
mset += x;
      else
        madd += x;
      val += x;
    } else {
      push(), l->add(L, R, x), r->add(L, R,
         x);
      val = max(l->val, r->val);
  void push() {
    if (!1) {
      int mid = lo + (hi - lo) / 2;
      1 = new Node(lo, mid);
      r = new Node (mid, hi);
    if (mset != inf)
      l->set(lo, hi, mset), r->set(lo, hi,
         mset), mset = inf;
    else if (madd)
      l->add(lo, hi, madd), r->add(lo, hi,
         madd), madd = 0;
 }
};
UnionFind.h
Description: Disjoint-set data structure.
Time: \mathcal{O}(\alpha(N))
                                       32c6e8, 14 lines
struct UF {
  vi e;
  UF (int n) : e(n, -1) {}
 bool sameSet(int a, int b) { return find(a
     == find(b);
  int size(int x) { return -e[find(x)]; }
  int find(int x) { return e[x] < 0 ? x : e[</pre>
     x] = find(e[x]);
 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);
```

e[a] += e[b]; e[b] = a;

return true;

};

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

ab6430, 30 lin

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

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

UnionFindRollback.h

struct RollbackUF {

->r, f); }

needed, skip st, time() and rollback().

### UnionFindRollback Treap FenwickTree FenwickTree2d

```
vi e; vector<pii> st;
  RollbackUF(int n) : e(n, -1) {}
  int size(int x) { return -e[find(x)]; }
  int find(int x) { return e[x] < 0 ? x :
      find(e[x]); }
  int time() { return sz(st); }
  void rollback(int t) {
    for (int i = time(); i --> t;)
      e[st[i].first] = st[i].second;
    st.resize(t);
  bool join(int a, int b) {
    a = find(a), b = find(b);
    if (a == b) return false;
    if (e[a] > e[b]) swap(a, b);
    st.push_back({a, e[a]});
    st.push_back({b, e[b]});
    e[a] += e[b]; e[b] = a;
    return true;
};
Treap.h
Description: A short self-balancing tree. It acts as a sequential con-
tainer with log-time splits/joins, and is easy to augment with additional
Time: \mathcal{O}(\log N)
                                         cad9a2, 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(l) + cnt(r) +
   1; }
template < class F > void each (Node* n, F f) {
  if (n) { each (n->1, f); f(n->val); each (n->1, f);
```

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

ca7d3d, 21 lines

Usage: int t = uf.time(); ...; uf.rollback(t);

```
pair<Node*, Node*> split(Node* n, int k) {
  if (!n) return {};
  if (cnt(n->1) >= k) \{ // "n-> val >= k" for
      lower\_bound(k)
    auto pa = split(n->1, k);
    n->1 = pa.second;
    n->recalc();
    return {pa.first, n};
  } else {
    auto pa = split (n->r, k - cnt(n->1) - 1)
       ; // and just "k"
    n->r = pa.first;
    n->recalc();
    return {n, pa.second};
Node* merge(Node* 1, Node* r) {
  if (!1) return r;
  if (!r) return 1;
 if (1->v > r->v) {
   1->r = merge(1->r, r);
    l->recalc();
    return 1;
  } else {
    r->1 = merge(1, r->1);
    r->recalc();
    return r;
Node* ins(Node* t, Node* n, int pos) {
  auto [l,r] = split(t, pos);
  return merge(merge(l, n), r);
// Example application: move the range [1, r
   ) to index k
void move(Node*& t, int 1, int r, int k) {
 Node *a, *b, *c;
  tie(a,b) = split(t, l); 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  $O(\log N)$ . Status: Stress-tested

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

# FenwickTree2d.h

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

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

```
for (; x < sz(ys); x = x + 1) ys[x].
        push_back(y);
  void init() {
    for (vi& v : ys) sort(all(v)), ft.
        emplace_back(sz(v));
  int ind(int x, int y) {
    return (int) (lower_bound(all(ys[x]), y)
        - ys[x].begin()); }
  void update(int x, int y, ll dif) {
    for (; x < sz(ys); x = x + 1)
      ft[x].update(ind(x, y), dif);
  11 query(int x, int y) {
    11 \text{ sum} = 0;
    for (; x; x &= x - 1)
      sum += ft[x-1].query(ind(x-1, y));
    return sum;
};
RMQ.h
Description: Range Minimum Queries on an array. Returns min(V[a],
V[a + 1], ... V[b - 1] in constant time.
Usage: RMQ rmq(values);
rmq.query(inclusive, exclusive);
Time: \mathcal{O}(|V|\log|V|+Q)
                                         ef1867, 16 lines
template<class T>
struct RMO {
  vector<vector<T>> jmp;
  RMQ(const vector<T>& V) : jmp(1, V) {
    for (int pw = 1, k = 1; pw * 2 <= sz(V);
         pw *= 2, ++k) {
      jmp.emplace_back(sz(V) - pw * 2 + 1);
      rep(j, 0, sz(jmp[k]))
         jmp[k][j] = min(jmp[k - 1][j], jmp[k]
             -1][\dot{1} + pw]);
    }
```

assert (a < b); // or return inf if a ==

return min(jmp[dep][a], jmp[dep][b - (1

int dep = 31 - \_\_builtin\_clz(b - a);

T query(int a, int b) {

<< dep)]);

```
MoQueries.h
Description: Answer interval or tree path queries by finding an ap-
proximate TSP through the queries, and moving from one query to the
next by adding/removing points at the ends. If values are on tree edges,
change step to add/remove the edge (a, c) and remove the initial add
call (but keep in).
Time: \mathcal{O}(N\sqrt{Q})
void add(int ind, int end) { ... } // add a/
    ind \mid (end = 0 \ or \ 1)
void del(int ind, int end) { ... } // remove
     a / ind /
int calc() { ... } // compute current answer
vi mo(vector<pii> Q) {
  int L = 0, R = 0, blk = 350; // \sim N/sqrt(Q)
  vi s(sz(Q)), res = s;
#define K(x) pii(x.first/blk, x.second ^ -(x
    .first/blk & 1))
  iota(all(s), 0);
  sort(all(s), [&](int s, int t) { return K(Q
      [s]) < K(Q[t]); );
  for (int qi : s) {
    pii q = Q[qi];
    while (L > q.first) add(--L, 0);
    while (R < q.second) add(R++, 1);</pre>
    while (L < q.first) del(L++, 0);
    while (R > q.second) del(--R, 1);
    res[qi] = calc();
  return res;
vi moTree(vector<array<int, 2>> Q, vector<vi</pre>
   >& ed, int root=0){
  int N = sz(ed), pos[2] = {}, blk = 350; //
       \sim N/sqrt(Q)
  vi s(sz(Q)), res = s, I(N), L(N), R(N), in
      (N), par(N);
  add(0, 0), in[0] = 1;
  auto dfs = [\&](int x, int p, int dep, auto
      & f) -> void {
    par[x] = p;
    L[x] = N;
    if (dep) I[x] = N++;
```

```
for (int y : ed[x]) if (y != p) f(y, x,
       !dep, f);
    if (!dep) I[x] = N++;
   R[x] = N;
 };
 dfs(root, -1, 0, dfs);
#define K(x) pii(I[x[0]] / blk, I[x[1]] ^ -(
   I[x[0]] / blk & 1))
 iota(all(s), 0);
  sort(all(s), [&](int s, int t) { return K(Q
     [s]) < K(Q[t]); \});
 for (int qi : s) rep(end, 0, 2) {
    int &a = pos[end], b = Q[qi][end], i =
#define step(c) { if (in[c]) { del(a, end);
   in[a] = 0;  \
                  else { add(c, end); in[c]
                     = 1;  } a = c;  }
   while (!(L[b] \le L[a] \&\& R[a] \le R[b]))
      I[i++] = b, b = par[b];
    while (a != b) step(par[a]);
    while (i--) step(I[i]);
    if (end) res[gi] = calc();
 return res;
```

# Numerical (4)

# 4.1 Polynomials and recurrences

Polynomial.h

```
a7945b, 17 lines
```

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

void diff() {
    rep(i,1,sz(a)) a[i-1] = i*a[i];
    a.pop_back();
```

```
void divroot(double x0) {
    double b = a.back(), c; a.back() = 0;
    for (int i=sz(a)-1; i--;) c = a[i], a[i]
        = a[i+1]*x0+b, b=c;
    a.pop_back();
};
PolyRoots.h
Description: Finds the real roots to a polynomial.
Usage: polyRoots(\{\{2, -3, 1\}\}, -1e9, 1e9\}) // solve x^2-3x+2
Time: \mathcal{O}\left(n^2\log(1/\epsilon)\right)
"Polynomial.h"
                                          9f92bf, 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  $(\mathbf{x}[\mathbf{i}], \mathbf{y}[\mathbf{i}])$ , computes an n-1-degree polynomial p that passes through them:  $p(x) = a[0] * x^0 + ... + a[n-1] * x^{n-1}$ . For numerical precision, pick  $x[k] = c * \cos(k/(n-1) * \pi), k = 0 \dots n-1$ . **Time:**  $\mathcal{O}(n^2)$ 

3714c0, 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  $\leq n$ . Usage: berlekampMassey( $\{0, 1, 1, 3, 5, 11\}$ ) //  $\{1, 2\}$  Time:  $\mathcal{O}(N^2)$ 

```
"../number-theory/ModPow.h"
vector<ll> berlekampMassey(vector<ll> s) {
 int n = sz(s), L = 0, m = 0;
 vector<ll> C(n), B(n), T;
 C[0] = B[0] = 1;
 11 b = 1;
 rep(i, 0, n) \{ ++m;
   ll d = s[i] % mod;
    rep(j,1,L+1) d = (d + C[j] * s[i - j]) %
        mod;
    if (!d) continue;
    T = C; 11 \text{ coef} = d * \text{modpow(b, mod-2)} %
       mod:
    rep(j,m,n) C[j] = (C[j] - coef * B[j - m]
       1) % 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 (l1& x : C) x = (mod - x) % mod;
```

# LinearRecurrence.h

return C;

**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 \star 2 + 1);
    rep(i, 0, n+1) rep(j, 0, n+1)
      res[i + j] = (res[i + j] + a[i] * b[j]
          1) % mod;
    for (int i = 2 * n; i > n; --i) rep(j,0,
      res[i - 1 - j] = (res[i - 1 - j] + res
          [i] * tr[i]) % mod;
    res.resize(n + 1);
    return res;
  };
  Poly pol(n + 1), e(pol);
  pol[0] = e[1] = 1;
  for (++k; k; k /= 2) {
    if (k % 2) pol = combine(pol, e);
    e = combine(e, e);
  11 \text{ res} = 0;
  rep(i, 0, n) res = (res + pol[i + 1] * S[i])
       % mod;
  return res;
```

# 4.2 Matrices

Determinant.h

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

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

```
res \star= a[i][i];
  if (res == 0) return 0;
  rep(j, i+1, n) {
    double v = a[j][i] / a[i][i];
    if (v != 0) rep(k, i+1, n) a[j][k] -= v
        * a[i][k];
return res;
```

### IntDeterminant.h

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

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

4940fc, 18 lines

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

# SolveLinear.h

**Description:** Solves A \* x = b. If there are multiple solutions, an arbitrary one is returned. Returns rank, or -1 if no solutions. Data in Aand b is lost.

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

d11d96, 38 lines

```
typedef vector<double> vd;
const double eps = 1e-12;
int solveLinear(vector<vd>& A, vd& b, vd& x)
 int n = sz(A), m = sz(x), rank = 0, br, bc
```

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

# SolveLinear2.h

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

```
rep(j,0,n) if (j != i) // instead of rep(j,i
   +1.n)
// ... then at the end:
x.assign(m, undefined);
rep(i, 0, rank) {
  rep(j, rank, m) if (fabs(A[i][j]) > eps)
     goto fail;
```

```
x[col[i]] = b[i] / A[i][i];
fail:; }
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].anv())
        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[i] ^= b[i];
      A[i] ^= A[i];
    rank++;
  x = bs();
  for (int i = rank; i--;) {
    if (!b[i]) continue;
    x[col[i]] = 1;
    rep(j, 0, i) b[j] ^= A[j][i];
```

return rank; // (multiple solutions if

rank < m)

#### MatrixInverse.h

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

```
685ba1, 35 lines
int matInv(vector<vector<double>>& A) {
  int n = sz(A); vi col(n);
  vector<vector<double>> tmp(n, vector<</pre>
     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[\dot{j}][k]) > fabs(A[r][c]))
        r = j, c = k;
    if (fabs(A[r][c]) < 1e-12) return i;</pre>
    A[i].swap(A[r]); tmp[i].swap(tmp[r]);
    rep(j,0,n)
      swap(A[j][i], A[j][c]), swap(tmp[j][i
         ], tmp[j][c]);
    swap(col[i], col[c]);
    double v = A[i][i];
    rep(j,i+1,n) {
      double f = A[j][i] / v;
      A[j][i] = 0;
      rep(k,i+1,n) A[j][k] -= f*A[i][k];
      rep(k,0,n) tmp[j][k] \rightarrow f*tmp[i][k];
    rep(j, i+1, n) A[i][j] /= v;
    rep(j,0,n) tmp[i][j] /= v;
    A[i][i] = 1;
  for (int i = n-1; i > 0; --i) rep(j, 0, i) {
    double v = A[i][i];
    rep(k,0,n) tmp[j][k] -= v*tmp[i][k];
  rep(i,0,n) rep(j,0,n) A[col[i]][col[j]] =
     tmp[i][j];
  return n;
```

# Fourier transforms

FastFourierTransform.h

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

```
Time: O(N \log N) with N = |A| + |B| (~1s for N = 2^{22})
```

```
typedef complex<double> C;
typedef vector<double> vd;
```

```
void fft(vector<C>& a) {
  int n = sz(a), L = 31 - builtin clz(n);
  static vector<complex<long double>> R(2,
  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</pre>
      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));

fft(in);

rep(i,0,sz(b)) in[i].imag(b[i]);

```
for (C& x : in) x \star = 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]) /
return res;
```

# Number theory (5)

#### Modular arithmetic 5.1

Modular Arithmetic.h

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

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

ModArithSimple.h

**Description:** Use this for simplicity in place of modular arithmeatic c.Status: Stress-tested

```
11 madd(ll a, ll b) { return (a + b) % mod;
ll msub(ll a, ll b) { return (((a - b) % mod
   ) + mod) % mod; }
ll mmul(ll a, ll b) { return ((a % mod) * (b
    % mod)) % mod; }
11 mpow(ll base, ll exp) {
  11 \text{ res} = 1;
  while (exp) {
    if (exp % 2 == 1) {
      res = (res * base) % mod;
    exp >>= 1;
    base = (base * base) % mod;
  return res;
11 minv(ll base) { return mpow(base, mod -
   2); }
ll mdiv(ll a, ll b) { return mmul(a, minv(b)
   ); }
```

# binaryexpinverse.h

```
const 11 mod = 1000000007; // faster if
   const
ll binpow(ll a, ll b) {
  ll ans = 1;
  while (b) {
    if (b & 1) {
      ans = (ans * a) % mod;
    a = (a * a) % mod;
    b = b >> 1;
  return (ans % mod);
ll modinverse(ll a, ll b) { return binpow(a,
    b - 2) % mod; }
```

### ModLog.h

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

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

```
// Description: Returns the smallest \$x > 0\$
    s.t. \$a^x = b \ pmod \ m\$, \ or
// * \$-1\$ if no such \$x\$ exists.
ll modLog(ll a, ll b, ll m) {
 ll n = (ll) sqrt(m) + 1, e = 1, f = 1, j =
     1;
  unordered_map<11, 11> A;
  while (j \le n \& \& (e = f = e * a % m) != b
     % m)
   A[e * b % m] = j++;
  if (e == b % m)
    return j;
 if (__gcd(m, e) == __gcd(m, b))
    rep(i, 2, n + 2) if (A.count(e = e * f %
        m)) return n * i - A[e];
  return -1:
```

### ModMuLL.h

Description: Calculate  $a \cdot b \mod c$  (or  $a^b \mod c$ ) for  $0 \leq a,b \leq c \leq 7.2 \cdot 10^{18}$ . Time: O(1) for modmul,  $O(\log b)$  for modpow Status: stress-tested, proven correct Details: $< a, b < c < 2^{52} \approx 4.5 \cdot 10^{15}$ .

```
5f1d88, 12 lines
```

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

# ModSart.h

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

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

```
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,
                                                                                                                                                                                                        // a^{(n+3)/8} \text{ or } 2^{(n+3)/8} * 2^{(n-1)/4}
                                                                                                                                                                                                                      works if p \% 8 == 5
                                                                                                                                                                                                       11 s = p - 1, n = 2;
                                                                                                                                                                                                        int r = 0, m;
                                                                                                                                                                                                       while (s % 2 == 0)
                                                                                                                                                                                                                ++r, s /= 2;
                                                                                                                                                                                                        while (modpow(n, (p-1) / 2, p) != p-1)
                                                                                                                                                                                                       11 x = modpow(a, (s + 1) / 2, p);
                                                                                                                                                                                                       ll b = modpow(a, s, p), g = modpow(n, s, p)
                                                                                                                                                                                                                     );
                                                                                                                                                                                                       for (;; r = m) {
                                                                                                                                                                                                               11 t = b;
                                                                                                                                                                                                                for (m = 0; m < r && t != 1; ++m)
                                                                                                                                                                                                                      t = t * t % p;
                                                                                                                                                                                                                if (m == 0) return x;
                                                                                                                                                                                                                ll qs = modpow(q, 1LL \ll (r - m - 1), p)
                                                                                                                                                                                                                q = qs * qs % p;
This runs\ 2x faster than the naive ({}_{int128_t})a*bdoc/mod mul-proof. tex. A near live version of {\tt \pm heproof}, for {\tt growthept} becode used a*b/(long double) M, is indoc/mod mul-proof. tex. A near live version of {\tt \pm heproof}, for {\tt growthept} becode used a*b/(long double) M, is indoc/mod mul-proof. tex. A near live version of {\tt \pm heproof}, for {\tt growthept} becode used a*b/(long double) M, is indoc/mod mul-proof. tex. A near live version of {\tt \pm heproof}, for {\tt growthept} becode used a*b/(long double) M, is indoc/mod mul-proof. tex. A near live version of {\tt \pm heproof}, for {\tt growthept} becode used a*b/(long double) M, is indoc/mod mul-proof. Text. A near live version of {\tt \pm heproof}, for {\tt growthept} becode used a*b/(long double) M, is indoc/mod mul-proof. Text. A near live version of {\tt \pm heproof}, for {\tt growthept} becode used a*b/(long double) M, is indoc/mod mul-proof. Text. A near live version of {\tt \pm heproof}, for {\tt growthept} becode used a*b/(long double) M, is indoc/mod mul-proof. Text. A near live version of {\tt \pm heproof}, for {\tt growthept} becode used a*b/(long double) M, is indoc/mod mul-proof. Text. A near live version of {\tt \pm heproof}, for {\tt growthept} becode used a*b/(long double) M, is indoc/mod mul-proof. Text. A near live version of {\tt \pm heproof}, for {\tt growthept} becode used a*b/(long double) M, is indoc/mod mul-proof. Text. A near live version of {\tt \pm heproof}, for {\tt growthept} becode used a*b/(long double) M, is indoc/mod mul-proof. Text. A near live version of {\tt \pm heproof}, for {\tt growthept} becode used a*b/(long double) M, is indoc/mod mul-proof. Text. A near live version of {\tt \pm heproof}, for {\tt growthept} becode used a*b/(long double) M, is indoc/mod mul-proof. Text. A near live version of {\tt \pm heproof}, for {\tt growthept} becode used a*b/(long double) M, is indoc/mod mul-proof. Text. A near live version of {\tt \pm heproof}, for {\tt \pm hepro
                                                                                                                                                                                                               b = b * q % p;
```

# Primality

FastEratosthenes.h

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

```
ccd47a, 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[i] = 1;
for (int L = 1; L \le R; L += S) {
 array<bool, S> block{};
 for (auto &[p, idx] : cp)
    for (int i=idx; i < S+L; idx = (i+=p))
        block[i-L] = 1;
 rep(i, 0, min(S, R - L))
    if (!block[i]) pr.push_back((L + i) *
       2 + 1);
for (int i : pr) isPrime[i] = 1;
return pr;
```

### MillerRabin.h

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

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

```
"ModMulLL.h"
                                      ef9b53, 12 lines
bool isPrime(ull n) {
  if (n < 2 | | n % 6 % 4 != 1) return (n |
     1) == 3;
  ull A[] = \{2, 325, 9375, 28178, 450775,
     9780504, 1795265022},
      s = builtin ctzll(n-1), d = n >> s;
  for (ull a : A) { // ^ count trailing
     zeroes
    ull p = modpow(a%n, d, n), i = s;
    while (p != 1 && p != n - 1 && a % n &&
       i --)
      p = modmul(p, p, n);
    if (p != n-1 && i != s) return 0;
  return 1;
```

# Factor.h

**Description:** Pollard-rho randomized factorization algorithm. Returns prime factors of a number, in arbitrary order (e.g.  $2299 \rightarrow \{11, 19, 11\}$ ). **Time:**  $\mathcal{O}(n^{1/4})$ , less for numbers with small factors.

```
"ModMulLL.h", "MillerRabin.h"
                                               d8f0f9, 18 lines
ull pollard(ull n) {
  ull x = 0, y = 0, t = 30, prd = 2, i = 1,
      q;
```

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

#### Divisibility 5.3

euclid.h

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

```
7e9eaf, 17 lines
ll euclid(ll a, ll b, ll &x, ll &y) {
  if (!b)
    return x = 1, y = 0, a;
  ll d = euclid(b, a % b, y, x);
  return v -= a / b * x, d;
ll generalized_mod_inverse(ll a, ll m) {
 11 x, v;
  int g = euclid(a, m, x, y);
 if (q != 1) {
    return -1;
  } else {
    x = (x % m + m) % m;
    return x;
```

**Description:** Chinese Remainder Theorem.

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

```
Time: \log(n)
```

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

# 5.3.1 Bézout's identity

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

$$ax + by = d$$

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

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

phiFunction.h

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

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

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

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

#### 5.4 Fractions

FracBinarySearch.h

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

```
Usage:
        fracBS([](Frac f) { return f.p>=3*f.q; }, 10);
// {1,3}
```

```
Time: \mathcal{O}(\log(N))
struct Frac { ll p, q; };
```

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

# 5.5 Pythagorean Triples

The Pythagorean triples are uniquely generated by

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

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

# 5.6 Primes

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

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

# 5.7 Estimates

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

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

# Combinatorial (6)

# 6.1 Permutations

# 6.1.1 Factorial

IntPerm.h

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

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

# **6.1.2** Cycles

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

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

# 6.1.3 Derangements

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

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

# 6.1.4 Burnside's lemma

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

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

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

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

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

# 6.2 Partitions and subsets

# 6.2.1 Partition function

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

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

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

Let n, m be non-negative integers and p a prime. Write  $n = n_k p^k + ... + n_1 p + n_0$  and  $m = m_k p^k + ... + m_1 p + m_0$ .

Then  $\binom{n}{k}$  Bindals (mod p).

multinomial.h

return c;

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# 6.3 General purpose numbers

# 6.3.1 Bernoulli numbers

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

Sums of powers:

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

Euler-Maclaurin formula for infinite sums:

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

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

# 6.3.2 Stirling numbers of the first kind

Number of permutations on n items with k cycles.

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

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

# 6.3.3 Eulerian numbers

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

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

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

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

# 6.3.4 Stirling numbers of the second kind

Partitions of n distinct elements into exactly k groups.

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

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

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

# 6.3.5 Bell numbers

Total number of partitions of n distinct elements. B(n) = 1, 1, 2, 5, 15, 52, 203, 877, 4140, 21147, .... For <math>p prime,

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

### 6.3.6 Labeled unrooted trees

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

### 6.3.7 Catalan numbers

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

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

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

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

# Graph (7)

# 7.1 Fundamentals

BellmanFord.h

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

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

```
const ll inf = LLONG_MAX;

struct Ed { int a, b, w, s() { return a < b

? a : -a; }};

struct Node { ll dist = inf; int prev = -1;

};
```

```
void bellmanFord(vector<Node>& nodes, vector
   <Ed>& eds, int s) {
  nodes[s].dist = 0;
  sort(all(eds), [](Ed a, Ed b) { return a.s
     () < b.s(); \});
  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;
    ll 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;
```

# 7.2 Network flow

MinCostMaxFlow.h

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

**Time:**  $\mathcal{O}(FE \log(V))$  where F is max flow.  $\mathcal{O}(VE)$  for setpi.

```
#include <bits/extc++.h>

const ll INF = numeric_limits<ll>::max() /
4;

struct MCMF {
    struct edge {
        int from, to, rev;
        ll cap, cost, flow;
    };
    int N;
    vector<vector<edge>> ed;
    vi seen;
    vector<eldge*> par;
```

```
MCMF(int N) : N(N), ed(N), seen(N), dist(N
   ), pi(N), par(N) {}
void addEdge(int from, int to, ll cap, ll
   cost) {
  if (from == to) return;
  ed[from].push_back(edge{ from, to, sz(ed[
     to]),cap,cost,0 });
  ed[to].push_back(edge{ to,from,sz(ed[
     from])-1,0,-cost,0 });
void path(int s) {
  fill(all(seen), 0);
  fill(all(dist), INF);
  dist[s] = 0; ll di;
  gnu pbds::priority queue<pair<ll, int</pre>
     >> q;
  vector<decltype(q)::point_iterator> its(
     N);
  q.push({ 0, s });
  while (!q.empty()) {
    s = q.top().second; q.pop();
    seen[s] = 1; di = dist[s] + pi[s];
    for (edge& e : ed[s]) if (!seen[e.to])
      ll val = di - pi[e.to] + e.cost;
      if (e.cap - e.flow > 0 && val < dist</pre>
         [e.to]) {
        dist[e.to] = val;
        par[e.to] = &e;
        if (its[e.to] == q.end())
          its[e.to] = q.push({ -dist[e.to
             1, e.to });
        else
          q.modify(its[e.to], { -dist[e.to
             1, e.to });
    }
  rep(i, 0, N) pi[i] = min(pi[i] + dist[i],
     INF);
```

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

# 7.3 Matching

# DFSMatching.h

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

f43bd9, 22 lines

```
Usage: vi btoa(m, -1); dfsMatching(q, btoa);
Time: \mathcal{O}(VE)
```

```
vi match (m, -1);
int res = dfsMatching(q, match);
vector<bool> lfound(n, true), seen(m);
for (int it : match) if (it != -1) lfound[
   itl = 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]);
```

# MinimumVertexCover.h

-1);

& vis) {

return 0;

vi vis;

**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.

return sz(btoa) - (int) count(all(btoa),

bool find(int j, vector<vi>& q, vi& btoa, vi

if (!vis[e] && find(e, g, btoa, vis)) {

int dfsMatching(vector<vi>& q, vi& btoa) {

if (find(j, g, btoa, vis)) {

**if** (btoa[j] == -1) **return** 1;

vis[j] = 1; int di = btoa[j];

vis.assign(sz(btoa), 0);

**for** (**int j** : **q**[i])

btoa[j] = i;

for (int e : q[di])

return 1;

rep(i, 0, sz(q)) {

break;

btoa[e] = di;

```
"DFSMatching.h"
vi cover(vector<vi>& q, int n, int m) {
```

```
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;
```

#### DFS algorithms 7.4

### 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 vand vice versa.

```
scc(graph, [&](vi& v) { ... }) visits all
Usage:
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)
```

```
fc214c, 24 lines
vi val, comp, z, cont;
int Time, ncomps;
template < class G, class F> int dfs (int j, G&
    q, F& f) {
  int low = val[j] = ++Time, x; z.push_back(
     j);
  for (auto e : q[\dot{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(q);
  val.assign(n, 0); comp.assign(n, -1);
  Time = ncomps = 0;
  rep(i, 0, n) if (comp[i] < 0) dfs(i, q, f);
```

```
FindingBridges.h
Description: Finds bridges in a graph. A bridge is an edge which when
Time: \mathcal{O}(N+M)
                                       bae75b, 37 lines
void IS BRIDGE(int v, int to); // some
   function to process the found bridge
                                 // number of
int n;
   nodes
                                 // adjacency
vector<vector<int>> adj;
   list of graph
vector<bool> visited;
vector<int> tin, low;
int timer;
void dfs(int v, int p = -1) {
  visited[v] = true;
  tin[v] = low[v] = timer++;
  bool parent_skipped = false;
  for (int to : adj[v]) {
    if (to == p && !parent_skipped) {
      parent_skipped = true;
      continue;
    if (visited[to]) {
      low[v] = min(low[v], tin[to]);
    } else {
      dfs(to, v);
      low[v] = min(low[v], low[to]);
      if (low[to] > tin[v])
        IS_BRIDGE(v, to);
void find_bridges() {
  timer = 0;
  visited.assign(n, false);
  tin.assign(n, -1);
  low.assign(n, -1);
  for (int i = 0; i < n; ++i) {
    if (!visited[i])
      dfs(i);
```

#### 2sat.h

Description: Calculates a valid assignment to boolean variables a, b, c,... to a 2-SAT problem, so that an expression of the type (a||b)&&(!a||c)&&(d||!b)&&... becomes true, or reports that it is unsatisfiable. Negated variables are represented by bit-inversions ( $\sim x$ ). Usage: TwoSat ts(number of boolean variables); ts.either(0,  $\sim$ 3); // Var 0 is true or var 3 is false ts.setValue(2); // Var 2 is true  $ts.atMostOne(0,\sim 1,2)$ ;  $// \le 1$  of vars 0,  $\sim 1$  and 2 are true ts.solve(); // Returns true iff it is solvable ts.values[0..N-1] holds the assigned values to the vars

**Time:**  $\mathcal{O}(N+E)$ , where N is the number of boolean variables, and E is the number of clauses. Status: stress-tested

14

```
b7e39d, 58 lines
struct TwoSat {
  int N;
 vector<vi> qr;
 vi values; // 0 = false, 1 = true
  TwoSat(int n = 0) : N(n), gr(2 * n) {}
 int addVar() { // (optional)
    gr.emplace back();
    gr.emplace back();
    return N++;
 void either(int f, int 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;
    int cur = \simli[0];
    rep(i, 2, sz(li)) {
      int next = addVar();
      either(cur, ~li[i]);
      either(cur, next);
      either(~li[i], next);
      cur = ~next;
    either(cur, ~li[1]);
 vi val, comp, z;
  int time = 0;
 int dfs(int i) {
    int low = val[i] = ++time, x;
```

z.push\_back(i);

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

## EulerWalk.h

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

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

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

```
eu[e] = 1; s.push back(y);
  }}
for (int x : D) if (x < 0 \mid \mid sz(ret) !=
   nedges+1) return {};
return {ret.rbegin(), ret.rend()};
```

#### Coloring 7.5

EdgeColoring.h

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

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

```
fd39d8, 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
          ^ d1);
   while (adj[fan[i]][d] != −1) {
     int left = fan[i], right = fan[++i], e
          = cc[i];
      adj[u][e] = left;
      adi[left][e] = u;
     adj[right][e] = -1;
      free[right] = e;
   adj[u][d] = fan[i];
```

```
adi[fan[i]][d] = u;
  for (int y : {fan[0], u, end})
    for (int& z = free[y] = 0; adj[y][z]
       ! = -1; z++);
rep(i, 0, sz(eds))
  for (tie(u, v) = eds[i]; adj[u][ret[i]]
     != v;) ++ret[i];
return ret;
```

#### 7.6 Heuristics

MaximalCliques.h

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

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

a99bda, 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 \mid 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. 2752a0, 49 lines

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

```
vi gmax, 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 \le sz(qmax))
       return:
    q.push_back(R.back().i);
    for(auto v:R) if (e[R.back().i][v.i])
       T.push back({v.i});
    if (sz(T)) {
      if (S[lev]++ / ++pk < limit) init(T)</pre>
      int j = 0, mxk = 1, mnk = max(sz(
         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 (\dot{j} > 0) T[\dot{j} - 1].d = 0;
      rep(k, mnk, mxk + 1) for (int i : C[k
         1)
        T[j].i = i, T[j++].d = k;
      expand(T, lev + 1);
    } else if (sz(q) > sz(qmax)) qmax = q;
    q.pop_back(), R.pop_back();
}
```

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

#### 7.7Trees

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

e0881a, 21 lines

```
7fc172, 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];
 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];
```

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

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

```
"../data-structures/RMQ.h"
```

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

### 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"
```

```
template <bool VALS EDGES> struct HLD {
  int N, tim = 0;
  vector<vi> adj;
 vi par, siz, rt, pos;
 Node *tree;
 HLD(vector<vi> adj )
    : N(sz(adj_)), adj(adj_), par(N, -1),
       siz(N, 1),
      rt(N), pos(N), tree(new Node(0, N)) {
         dfsSz(0); dfsHld(0); }
 void dfsSz(int v) {
    for (int& u : adj[v]) {
      adj[u].erase(find(all(adj[u]), v));
```

```
par[u] = v;
     dfsSz(u);
     siz[v] += siz[u];
     if (siz[u] > siz[adj[v][0]]) swap(u,
         adj[v][0]);
 void dfsHld(int v) {
   pos[v] = tim++;
   for (int u : adj[v]) {
     rt[u] = (u == adj[v][0] ? rt[v] : u);
     dfsHld(u);
   }
 template <class B> void process(int u, int
      v, B op) {
   for (;; v = par[rt[v]]) {
     if (pos[u] > pos[v]) swap(u, v);
     if (rt[u] == rt[v]) break;
      op(pos[rt[v]], pos[v] + 1);
   op(pos[u] + VALS\_EDGES, pos[v] + 1);
 void modifyPath(int u, int v, int val) {
   process(u, v, [&] (int l, int r) { tree->
       add(l, r, val); });
 int queryPath(int u, int v) { // Modify
     depending on problem
   int res = -1e9;
   process(u, v, [&](int 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]);
};
```

### Directed MST.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" f3d299, 60 lines

```
struct Edge { int a, b; ll w; };
struct Node {
 Edge key;
 Node *1, *r;
 ll delta;
 void prop() {
   kev.w += delta;
    if (1) 1->delta += delta;
    if (r) r->delta += delta;
    delta = 0;
 Edge top() { prop(); return key; }
Node *merge(Node *a, Node *b) {
 if (!a || !b) return a ?: b;
  a->prop(), b->prop();
 if (a->key.w > b->key.w) swap(a, b);
  swap(a->1, (a->r = merge(b, a->r)));
  return a:
void pop(Node*& a) { a->prop(); a = merge(a
   ->1, a->r); }
pair<ll, vi> dmst(int n, int r, vector<Edge</pre>
   >& q) {
  RollbackUF uf(n);
 vector<Node*> heap(n);
  for (Edge e : q) heap[e.b] = merge(heap[e.
     b], new Node{e});
  11 \text{ res} = 0;
  vi seen(n, -1), path(n), par(n);
  seen[r] = r;
  vector<Edge> Q(n), in(n, \{-1,-1\}), comp;
  deque<tuple<int, int, vector<Edge>>> cycs;
  rep(s,0,n) {
    int u = s, qi = 0, w;
    while (seen[u] < 0) {
      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[--
         qill);
      while (uf.join(u, w));
      u = uf.find(u), heap[u] = cyc, seen[
         ul = -1;
      cycs.push_front({u, time, {&Q[qi], &
         O[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)] =
  in[uf.find(inEdge.b)] = inEdge;
rep(i, 0, n) par[i] = in[i].a;
return {res, par};
```

#### 7.8 Math

# 7.8.1 Number of Spanning Trees

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

# Geometry (8)

#### Geometric primitives 8.1

Point.h

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

```
template <class T> int sqn(T x) { return (x
   > 0) - (x < 0);
template<class T>
struct Point {
  typedef Point P;
```

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

#### 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.



res

9e0fa5, 6 lines

template < class P > double lineDist (const P& a, const P& b,

# SegmentDistance.h

## Description:

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

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

"Point.h"

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

# SegmentIntersection.h

#### Description:



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"</pre>
2f7799, 13 lines

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

```
oc = a.cross(b, c), od = a.cross(b, d
          );

// Checks if intersection is single non-
          endpoint point.

if (sgn(oa) * sgn(ob) < 0 && sgn(oc) * sgn
          (od) < 0)
    return {(a * ob - b * oa) / (ob - oa)};

set<P> s;

if (onSegment(c, d, a)) s.insert(a);

if (onSegment(c, d, b)) s.insert(b);

if (onSegment(a, b, c)) s.insert(c);

if (onSegment(a, b, d)) s.insert(d);

return {all(s)};
```

### lineIntersection.h

#### Description:

"Point.h"

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<|| > and the intersection point does not have integer coordinates. Products of three coordinates are used in intermediate steps so watch out for overflow if using int or ll.

```
e2 r
```

c0ba2b, 9 lines

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

faee62, 8 lines

#### sideOf.h

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

```
Usage: bool left = sideOf(p1,p2,q)==1;
```

"Point.h"

template<class P>

lineDistance.h

int t;

b2beb9, 11 lines

```
int sideOf(P s, P e, P p) { return sqn(s.
    cross(e, p)); }
template<class P>
int sideOf(const P& s, const P& e, const P&
    p, double eps) {
  auto a = (e-s).cross(p-s);
  double l = (e-s).dist()*eps;
  return (a > 1) - (a < -1);
OnSegment.h
Description: Returns true iff p lies on the line segment from s to e. Use
(segDist(s,e,p) <=epsilon) instead when using Point <double>.
template < class P > bool on Segment (P s, P e, P
  return p.cross(s, e) == 0 \&\& (s - p).dot(e)
       - > (q - 1)
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"
                                             ade292, 6 lines
typedef Point<double> P;
P linearTransformation(const P& p0, const P&
     p1,
    const P& q0, const P& q1, const P& r) {
  P dp = p1-p0, dq = q1-q0, num(dp.cross(dq))
      , dp.dot(dq));
  return q0 + P((r-p0).cross(num), (r-p0).
      dot(num))/dp.dist2();
Angle.h
Description: A class for ordering angles (as represented by int points
and a number of rotations around the origin). Useful for rotational
sweeping. Sometimes also represents points or vectors.
Usage: vector<Angle> v = \{w[0], w[0].t360() ...\}; //
sorted
int j = 0; rep(i,0,n) { while (v[j] < v[i].t180()) ++j;
// sweeps j such that (j-i) represents the number of
positively oriented triangles with vertices at 0 and i
struct Angle {
  int x, y;
```

```
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 \mid | y);
    return v < 0 || (v == 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) <
         make tuple(b.t, b.half(), a.x \star (ll
            )b.v);
// Given two points, this calculates the
   smallest angle between
// them, i.e., the angle that covers the
   defined line segment.
pair<Angle, Angle> segmentAngles (Angle a,
   Angle b) {
  if (b < a) swap(a, b);
  return (b < a.t180() ?
         make pair(a, b) : make pair(b, a.
             t360()));
Angle operator+(Angle a, Angle b) { // point
    a + vector b
  Angle r(a.x + b.x, a.y + b.y, a.t);
  if (a.t180() < r) r.t--;
  return r.t180() < a ? r.t360() : r;
Angle angleDiff(Angle a, Angle b) { // angle
    b - angle a
  int tu = b.t - a.t; a.t = b.t;
  return {a.x*b.x + a.y*b.y, a.x*b.y - a.y*b
     .x, tu - (b < a);
```

# 8.2 Circles

### CircleIntersection.h

"Point.h"

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

```
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;
```

### 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.



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

# MinimumEnclosingCircle.h

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

```
Time: expected \mathcal{O}(n)
"circumcircle.h"
                                        3ed955, 17 lines
pair<P, double> mec(vector<P> ps) {
  shuffle(all(ps), mt19937(time(0)));
  P \circ = ps[0];
  double r = 0, EPS = 1 + 1e-8;
  rep(i, 0, sz(ps)) if ((o - ps[i]).dist() > r
       * EPS) {
    o = ps[i], r = 0;
    rep(j, 0, i) if ((o - ps[j]).dist() > r *
       EPS) {
      o = (ps[i] + ps[j]) / 2;
      r = (o - ps[i]).dist();
      rep(k, 0, j) if ((o - ps[k]).dist() > r
          * EPS) {
        o = ccCenter(ps[i], ps[j], ps[k]);
        r = (o - ps[i]).dist();
    }
  return {o, r};
```

# Polygons

# InsidePolygon.h

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

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

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

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

# PolygonArea.h

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

```
"Point.h"
                                                                           aa3c71, 6 lines
```

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

# PolygonCenter.h

**Description:** Returns the center of mass for a polygon.

"Point.h"

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

```
A += v[j].cross(v[i]);
return res / A / 3;
```

# PolygonCut.h

#### Description:

Returns a vector with the vertices of a polygon with evervthing 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"
                                                        e38095, 13 lines
typedef Point < double > P;
```

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

```
return res;
```

# ConvexHull.h

## Description:

dd5f93, 9 lines

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



5935f9, 13 lines

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

```
typedef Point<ll> P;
vector<P> convexHull(vector<P> pts) {
  if (sz(pts) <= 1) return pts;</pre>
  sort(all(pts));
  vector\langle P \rangle h(sz(pts)+1);
  int s = 0, t = 0;
  for (int it = 2; it--; s = --t, reverse(
     all(pts)))
    for (P p : pts) {
      while (t >= s + 2 \&\& h[t-2].cross(h[t
          -11, p) <= 0) t--;
      h[t++] = p;
  return {h.begin(), h.begin() + t - (t == 2
       && h[0] == h[1]);
```

#### HullDiameter.h

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

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

"Point.h"

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

```
pair<11, array<P, 2>> res({0, {S[0], S[0]}}
   });
rep(i,0,j)
  for (;; j = (j + 1) % n) {
    res = max(res, \{(S[i] - S[j]).dist2(),
        {S[i], S[j]}});
    if ((S[(j + 1) % n] - S[j]).cross(S[i
       + 1] - S[i]) >= 0)
      break;
return res.second;
```

3a707e, 16 lines

### PointInsideHull ClosestPair Point3D KMP

```
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"
                                                         5bef6c, 14 lines
typedef Point<ll> P;
```

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

# Misc. Point Set **Problems**

ClosestPair.h

**Description:** Finds the closest pair of points.

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

"Point.h" f85d3e, 17 lines typedef Point<ll> P; pair<P, P> closest(vector<P> v) { assert (sz(v) > 1); set < P > S; sort(all(v), [](P a, P b) { return a.y < b</pre> .v; }); pair<ll, pair<P, P>> ret{LLONG MAX, {P(), P()}}; int j = 0; **for** (P p : v) {  $P d\{1 + (ll) sqrt(ret.first), 0\};$ while  $(v[j].y \le p.y - d.x)$  S.erase $(v[j].y \le p.y - d.x)$ auto lo = S.lower\_bound(p - d), hi = S. upper\_bound(p + d);

```
for (; lo != hi; ++lo)
    ret = min(ret, {(*lo - p).dist2(), {*}}
       lo, p}});
 S.insert(p);
return ret.second;
```

#### 3D8.5

Point3D.h

or long long.

```
Description: Class to handle points in 3D space. T can be e.g. double
                                      14b717, 32 lines
template < class T > struct Point 3D {
 typedef Point3D P;
 typedef const P& R;
  T x, y, z;
  explicit Point3D(T x=0, T y=0, T z=0) : x(
     x), y(y), z(z) {}
 bool operator<(R p) const {</pre>
    return tie(x, y, z) < tie(p.x, p.y, p.z)</pre>
       ; }
 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 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;
```

# Strings (9)

### KMP.h

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

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

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

# Zfunc Manacher MinRotation SuffixArray SuffixTree

```
Zfunc.h
```

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

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

7a443b, 12 lines

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

### Manacher.h

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

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

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

# MinRotation.h

**Description:** Finds the lexicographically smallest rotation of a string. Usage: rotate(v.begin(), v.begin()+minRotation(v), v.end());

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

18bdba, 8 lines

};

```
int minRotation(string s) {
 int a=0, N=sz(s); s += s;
```

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

# SuffixArrav.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:  $\mathcal{O}(n \log n)$ 

077d68, 22 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)), y(n), ws(max(n, lim));
   x.push\_back(0), sa = lcp = y, iota(all(
       sa), 0);
   for (int j = 0, p = 0; p < n; j = max(1,
        j * 2), lim = p) {
     p = j, iota(all(y), n - j);
      rep(i, 0, n) if (sa[i] >= j) y[p++] = sa
         [i] - j;
      fill(all(ws), 0);
      rep(i, 0, n) ws[x[i]] ++;
      rep(i,1,lim) ws[i] += ws[i-1];
      for (int i = n; i--;) sa[--ws[x[y[i
        ]]]] = y[i];
      swap(x, y), p = 1, x[sa[0]] = 0;
     rep(i,1,n) = sa[i-1], b = sa[i], x
         [d]
        (y[a] == y[b] && y[a + j] == y[b + j]
           ]) ? p - 1 : p++;
   for (int i = 0, j; i < n - 1; lcp[x[i
       ++]] = k)
     for (k \& \& k--, j = sa[x[i] - 1];
          s[i + k] == s[j + k]; k++);
```

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

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

memset(t, -1, sizeof t);

fill(t[1],t[1]+ALPHA,0);

= p[0] = p[1] = 0;

s[0] = 1; 1[0] = 1[1] = -1; r[0] = r[1]

rep(i, 0, sz(a)) ukkadd(i, toi(a[i]));

H operator\*(H o) { auto m = ( uint128 t)x

```
// example: find longest common substring
      (uses ALPHA = 28)
  pii best;
  int lcs(int node, int i1, int i2, int olen
     ) {
    if (l[node] <= i1 && i1 < r[node])</pre>
       return 1;
    if (l[node] <= i2 && i2 < r[node])</pre>
       return 2;
    int mask = 0, len = node ? olen + (r[
       node] - l[node]) : 0;
    rep(c, 0, ALPHA) if (t[node][c] != -1)
      mask |= lcs(t[node][c], i1, i2, len);
    if (mask == 3)
      best = max(best, {len, r[node] - len})
    return mask;
  static pii LCS(string s, string t) {
    SuffixTree st(s + (char)('z' + 1) + t +
        (char) ('z' + 2));
    st.lcs(0, sz(s), sz(s) + 1 + sz(t), 0);
    return st.best;
};
Hashing.h
Description: Self-explanatory methods for string hashing.
// Arithmetic mod 2^64-1. 2x slower than mod
    2^64 and more
// code, but works on evil test data (e.g.
   Thue-Morse, where
// ABBA... and BAAB... of length 2^10 hash
   the same mod 2^64).
// "typedef ull H;" instead if you think
   test data is random,
// or work mod 10^9+7 if the Birthday
   paradox is not a problem.
typedef uint64 t ull;
struct H {
  ull x; H(ull x=0) : x(x) \{ \}
  H operator+(H \circ) { return x + \circ .x + (x + \circ .x)
     .x < x); }
  H operator-(H o) { return *this + ~o.x; }
```

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

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

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

```
bdcee5, 66 lines
```

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

};

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

# Various (10)

# 10.1 Intervals

IntervalContainer.h

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

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

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

### IntervalCover.h

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

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

8c7b09, 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[
   al < I[b]; });
T cur = G.first;
int at = 0;
while (cur < G.second) \{ // (A) \}
  pair<T, int> mx = make pair(cur, -1);
  while (at < sz(I) && I[S[at]].first <=</pre>
     cur) {
    mx = max(mx, make_pair(I[S[at]].second
       , S[at]));
    at++;
  if (mx.second == -1) return {};
  cur = mx.first;
  R.push back (mx.second);
return R;
```

### ConstantIntervals.h

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

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

```
a510b2, 19 lines
template < class F, class G, class T>
void rec(int from, int to, F& f, G& q, int&
   i, T& p, T q) {
  if (p == q) return;
  if (from == to) {
    q(i, to, p);
    i = to; p = q;
  } else {
    int mid = (from + to) >> 1;
    rec(from, mid, f, g, i, p, f(mid));
    rec(mid+1, to, f, q, i, p, q);
template < class F, class G>
void constantIntervals(int from, int to, F f
   , G q) {
  if (to <= from) return;</pre>
```

int i = from; auto p = f(i), q = f(to-1);

rec(from, to-1, f, q, i, p, q);

q(i, to, q);

# 10.2 Misc. algorithms

# TernarvSearch.h

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

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

fb4284, 11 lines

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

# LIS.h

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

```
template<class I> vi lis(const vector<I>& S)
 if (S.empty()) return {};
 vi prev(sz(S));
 typedef pair<I, int> p;
 vector res;
 rep(i, 0, sz(S)) {
    // change \theta \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 nonnegative target t, computes the maximum S <= t such that S is the sum of some subset of the weights.

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

} **;** 

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

# 10.3 Dynamic programming

### KnuthDP.h

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

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

# DivideAndConquerDP.h

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

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

bdae29, 18 lines

```
struct DP { // Modify at will:
 int lo(int ind) { return 0; }
```

```
int hi(int ind) { return ind; }
11 f(int ind, int k) { return dp[ind][k];
void store(int ind, int k, ll v) { res[ind
   ] = pii(k, v); 
void rec(int L, int R, int LO, int HI) {
  if (L >= R) return;
  int mid = (L + R) >> 1;
  pair<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); }
```