**iTMO** 

ITMO

# Shuffle, duffle, muzzle, muff. Fista, wista, mista-cuff

overpressured but trained

The 2024 ICPC Northern Eurasia Finals
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# Mathematics (1)

## 1.1 Equations

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

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

$$ax + by = e$$

$$cx + dy = f$$

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

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

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

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

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

## 1.2 Recurrences

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

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

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

## 1.3 Trigonometry

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

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

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

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

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

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

## 1.4 Geometry

## 1.4.1 Triangles

Side lengths: a, b, c

Semiperimeter:  $p = \frac{a+b+c}{2}$ 

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

Circumradius:  $R = \frac{abc}{4A}$ 

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

Length of median (divides triangle into two equal-area triangles):  $m_a = \frac{1}{2}\sqrt{2b^2 + 2c^2 - a^2}$ 

Length of bisector (divides angles in two):

$$s_a = \sqrt{bc \left[ 1 - \left( \frac{a}{b+c} \right)^2 \right]}$$

Law of sines:  $\frac{\sin \alpha}{a} = \frac{\sin \beta}{b} = \frac{\sin \gamma}{c} = \frac{1}{2R}$ Law of cosines:  $a^2 = b^2 + c^2 - 2bc\cos \alpha$ 

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

## 1.4.2 Quadrilaterals

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

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

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

# 1.5 Derivatives/Integrals

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

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

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

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

Integration by parts:

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

## 1.6 Sums

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

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

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

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

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

## 1.7 Series

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

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

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

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

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

## 1.8 Probability theory

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

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

## 1.9 Markov chains

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

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

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

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

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

# Data structures (2)

#### StatisticTree.h

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

## HashMap.h

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

StatisticTree HashMap Segtree LazySegmentTree UnionFindRollback

#### Segtree.h

Description: tree for mx and += ops

Time:  $\mathcal{O}(\log N)$ d41d8c, 46 lines struct segtree { int n; vector<11> mx, upd; segtree() { } segtree(int n): n(n) { mx.resize(n \* 4);upd.resize(n \* 4); void add(int v, int v1, int vr, int 1, int r, 11 x) { **if** (vr <= 1 || vl >= r) return; if (1 <= v1 && vr <= r) { mx[v] += x;upd[v] += x;return; int vm = (v1 + vr) >> 1; add(v \* 2, v1, vm, 1, r, x); add(v \* 2 + 1, vm, vr, 1, r, x); mx[v] = max(mx[v \* 2], mx[v \* 2 + 1]) + upd[v];void add(int 1, int r, 11 x) { add(1, 0, n, 1, r, x); 11 getMax(int v, int vl, int vr, int l, int r) { **if** (vr <= 1 || v1 >= r) return INT64 MIN; if (1 <= v1 && vr <= r) return mx[v]; int vm = (v1 + vr) >> 1; return max(getMax(v \* 2, v1, vm, 1, r), getMax(v \* 2 + 1, vm, vr, 1, r)) + upd[v]; 11 getMax(int 1, int r) { return getMax(1, 0, n, 1, r);

## LazySegmentTree.h

};

ll getMax() {

return mx[1];

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

```
Usage: Node* tr = new Node(v, 0, sz(v));
Time: \mathcal{O}(\log N).
"../various/BumpAllocator.h"
                                                         d41d8c, 50 lines
const int inf = 1e9;
struct Node {
  Node *1 = 0, *r = 0;
  int lo, hi, mset = inf, madd = 0, val = -inf;
  Node (int lo, int hi):lo(lo), hi(hi) {} // Large interval of -inf
  Node (vi& v, int lo, int hi) : lo(lo), hi(hi) {
    if (lo + 1 < hi) {
      int mid = lo + (hi - lo)/2;
      l = new Node(v, lo, mid); r = new Node(v, mid, hi);
      val = max(1->val, r->val);
    else val = v[lo];
  int query(int L, int R) {
    if (R <= lo || hi <= L) return -inf;</pre>
    if (L <= lo && hi <= R) return val;</pre>
    push();
    return max(l->query(L, R), r->query(L, R));
  void set(int L, int R, int x) {
    if (R <= lo || hi <= L) return;</pre>
    if (L <= lo && hi <= R) mset = val = x, madd = 0;</pre>
      push(), l\rightarrow set(L, R, x), r\rightarrow set(L, R, x);
      val = max(1->val, r->val);
  void add(int L, int R, int x) {
    if (R <= lo || hi <= L) return;</pre>
    if (L <= lo && hi <= R) {
      if (mset != inf) mset += x;
      else madd += x;
      val += x;
      push(), l\rightarrow add(L, R, x), r\rightarrow add(L, R, x);
      val = max(1->val, r->val);
  void push() {
      int mid = lo + (hi - lo)/2;
      l = new Node(lo, mid); r = new Node(mid, hi);
    if (mset != inf)
      1->set(lo,hi,mset), r->set(lo,hi,mset), mset = inf;
      1- add (lo, hi, madd), r- add (lo, hi, madd), madd = 0;
};
```

## UnionFindRollback.h

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

d41d8c, 21 lines

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

```
struct RollbackUF {
  vi e; vector<pii> st;
  RollbackUF(int n) : e(n, -1) {}
  int size(int x) { return -e[find(x)]; }
  int find(int x) { return e[x] < 0 ? x : find(e[x]); }
  int time() { return sz(st); }
  void rollback(int t) {
    for (int i = time(); i --> t;)
        e[st[i].first] = st[i].second;
```

```
ITMO(Ivchenko, Konovalov, Matveev)
    st.resize(t);
  bool join(int a, int b) {
    a = find(a), b = find(b);
    if (a == b) return false;
    if (e[a] > e[b]) swap(a, b);
    st.push back({a, e[a]});
    st.push_back({b, e[b]});
    e[a] += e[b]; e[b] = a;
    return true;
};
SubMatrix.h
Description: Calculate submatrix sums quickly, given upper-left and lower-
right corners (half-open).
Usage: SubMatrix<int> m(matrix);
m.sum(0, 0, 2, 2); // top left 4 elements
Time: O(N^2 + Q)
                                                      d41d8c, 13 lines
template<class T>
struct SubMatrix {
  vector<vector<T>> p;
  SubMatrix(vector<vector<T>>& v) {
    int R = sz(v), C = sz(v[0]);
```

```
p.assign(R+1, vector<T>(C+1));
    rep(r, 0, R) rep(c, 0, C)
     p[r+1][c+1] = v[r][c] + p[r][c+1] + p[r+1][c] - p[r][c];
  T sum(int u, int 1, int d, int r) {
    return p[d][r] - p[d][l] - p[u][r] + p[u][l];
};
```

#### Matrix.h

Description: Basic operations on square matrices.

```
Usage: Matrix<int, 3> A;
A.d = \{\{\{1,2,3\}\}, \{\{4,5,6\}\}, \{\{7,8,9\}\}\}\};
vector<int> vec = \{1,2,3\};
```

```
vec = (A^N) * vec;
                                                      d41d8c, 26 lines
template < class T, int N> struct Matrix {
  typedef Matrix M;
  array<array<T, N>, N> d{};
  M operator*(const M& m) const {
   Ma;
    rep(i,0,N) rep(j,0,N)
     rep(k,0,N) a.d[i][j] += d[i][k]*m.d[k][j];
    return a;
  vector<T> operator*(const vector<T>& vec) const {
   vector<T> ret(N);
    rep(i, 0, N) rep(j, 0, N) ret[i] += d[i][j] * vec[j];
    return ret;
  M operator^(ll p) const {
    assert (p >= 0);
   M a, b(*this);
    rep(i, 0, N) a.d[i][i] = 1;
    while (p) {
     if (p&1) a = a*b;
     b = b*b;
     p >>= 1;
    return a;
};
```

## LineContainer.h

**Description:** Container where you can add lines of the form kx+m, and query maximum values at points x. Useful for dynamic programming ("convex hull trick").

```
Time: \mathcal{O}(\log N)
                                                      d41d8c, 30 lines
struct Line {
 mutable 11 k, m, p;
 bool operator<(const Line& o) const { return k < o.k; }</pre>
 bool operator<(11 x) const { return p < x; }</pre>
struct LineContainer : multiset<Line, less<>>> {
 // (for doubles, use inf = 1/.0, div(a,b) = a/b)
 static const 11 inf = LLONG MAX;
 ll div(ll a, ll b) { // floored division
    return a / b - ((a ^ b) < 0 && a % b); }
 bool isect(iterator x, iterator y) {
   if (v == 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 = y;
    while (isect(v, z)) z = erase(z);
   if (x != begin() \&\& isect(--x, y)) isect(x, y = erase(y));
    while ((y = x) != begin() && (--x)->p >= y->p)
      isect(x, erase(v));
 ll query(ll x) {
   assert(!empty());
   auto 1 = *lower_bound(x);
   return 1.k * x + 1.m;
};
```

#### RationalLineContainer.h Description: CHT with rationals Time: $\mathcal{O}(\log N)$

d41d8c, 63 lines

```
struct CHT {
    struct Frac {
        11 a. b:
        bool operator <(const Frac &other) const {</pre>
            return (__int128_t)a * other.b < (__int128_t)other.
                 a * b:
        bool operator >(const Frac &other) const {
            return other < *this;
    };
    struct Line {
        11 k, b;
        11 get(ll x) const {
            return k * x + b;
        Frac intersect (const Line &other) const {
            Frac res = {other.b - b, k - other.k};
            if (res.b < 0) {
                res.a \star = -1;
                res.b \star = -1;
            return res;
```

```
vector<Frac> pt;
    vector<Line> lines;
    void addLine(11 k, 11 b) { // k[i] descrease
        Line l = \{k, b\};
        while (!pt.empty()) {
            const Line &back = lines.back();
            if (back.k == k) {
                if (b >= back.b)
                    return;
            } else if (l.intersect(back) > pt.back()) {
                break;
            pt.pop_back();
            lines.pop_back();
        if (!lines.empty() && lines.back().k == k) {
            if (b < lines.back().b)</pre>
                lines.back() = 1;
            return;
        if (!lines.empty())
            pt.push_back(l.intersect(lines.back()));
        lines.push_back(1);
    ll getMin(ll x) {
        if (lines.empty())
            return INT64 MAX;
        int pos = lower_bound(all(pt), Frac{x, 1}) - pt.begin()
        return lines[pos].get(x);
};
```

#### FenwickTree2d.h

};

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

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

```
"FenwickTree.h"
                                                      d41d8c, 22 lines
struct FT2 {
 vector<vi> ys; vector<FT> ft;
 FT2(int limx) : vs(limx) {}
  void fakeUpdate(int x, int y) {
    for (; x < sz(ys); x |= x + 1) ys[x].push_back(y);</pre>
 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)
                                                       d41d8c, 65 lines
     auto fun = [\mathcal{E}](int \ i, \ int \ j) \ \{ \ return \ min(i, \ j); \ \};
     SparseTable < int, decltype(fun) > st(a, fun);
     SparseTable < int > st(a, [\&](int i, int j) { return min(i, j)}
template <typename T, class F = function<T(const T&, const T&)
class SparseTable {
public:
  int n:
  vector<vector<T>> mat:
  F func;
  SparseTable(const vector<T>& a, const F& f) : func(f) {
    n = static cast<int>(a.size());
    int max log = 32 - builtin clz(n);
    mat.resize(max_log);
    mat[0] = a;
    for (int j = 1; j < max_log; j++) {</pre>
     mat[j].resize(n - (1 << j) + 1);
     for (int i = 0; i <= n - (1 << j); i++) {
        mat[j][i] = func(mat[j-1][i], mat[j-1][i+(1 << (j-1)[i]))
              - 1))1):
  T get(int from, int to) const {
    assert (0 <= from && from <= to && to <= n - 1);
    int lg = 32 - __builtin_clz(to - from + 1) - 1;
    return func(mat[lg][from], mat[lg][to - (1 << lg) + 1]);</pre>
};
template <typename T, typename Func>
class DisjointSparseTable {
public:
  int _n;
  vector<vector<T>> _matrix;
  Func _func;
  DisjointSparseTable(const vector<T>& a, const Func& func) :
       _n(static_cast<int>(a.size())), _func(func) {
    _matrix.push_back(a);
    for (int layer = 1; (1 << layer) < _n; ++layer) {</pre>
      _matrix.emplace_back(_n);
      for (int mid = 1 << layer; mid < _n; mid += 1 << (layer +</pre>
            1)) {
        _{matrix[layer][mid - 1]} = a[mid - 1];
        for (int j = mid - 2; j >= mid - (1 << layer); --j) {</pre>
          _matrix[layer][j] = _func(a[j], _matrix[layer][j +
               11);
        _matrix[layer][mid] = a[mid];
        for (int j = mid + 1; j < min(_n, mid + (1 << layer));</pre>
          _matrix[layer][j] = _func(_matrix[layer][j - 1], a[j
```

```
RMQ MoQueries Polynomial
 T Query(int 1, int r) const {
    assert(0 <= 1 && 1 < r && r <= _n);
    if (r - 1 == 1) {
      return _matrix[0][1];
    int layer = 31 - __builtin_clz(1 ^ (r - 1));
    return _func(_matrix[layer][l], _matrix[layer][r - 1]);
};
MoQueries.h
Description: Answer interval or tree path queries by finding an approxi-
mate TSP through the queries, and moving from one query to the next by
adding/removing points at the ends. If values are on tree edges, change step
to add/remove the edge (a, c) and remove the initial add call (but keep in).
Time: \mathcal{O}(N\sqrt{Q})
void add(int ind, int end) { ... } // add a[ind] (end = 0 or 1)
void del(int ind, int end) { ... } // remove a[ind]
int calc() { ... } // compute current answer
vi mo(vector<pii> 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 = O[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>> 0, vector<vi>& ed, int root=0) {
 int N = sz(ed), pos[2] = {}, blk = 350; // \sim N/sqrt(Q)
 vi s(sz(Q)), res = s, I(N), L(N), R(N), in(N), par(N);
  add(0, 0), in[0] = 1;
  auto dfs = [&](int x, int p, int dep, auto& f) -> void {
    par[x] = p;
    L[x] = N;
    if (dep) I[x] = N++;
    for (int y : ed[x]) if (y != p) f(y, x, !dep, f);
    if (!dep) I[x] = N++;
    R[x] = N;
 dfs(root, -1, 0, dfs);
#define K(x) pii(I[x[0]] / blk, I[x[1]] ^ -(I[x[0]] / blk & 1))
```

sort(all(s), [&](int s, int t){ return K(Q[s]) < K(Q[t]); });

**else** { add(c, end); in[c] = 1; } a = c; }

**#define** step(c) { **if** (in[c]) { del(a, end); in[a] = 0; } \

iota(all(s), 0);

return res:

for (int qi : s) rep(end, 0, 2) {

I[i++] = b, b = par[b];

if (end) res[qi] = calc();

while (i--) step(I[i]);

while (a != b) step(par[a]);

int &a = pos[end], b = Q[qi][end], i = 0;

while (!(L[b] <= L[a] && R[a] <= R[b]))</pre>

# Numerical (3)

## 3.1 Polynomials and recurrences

Polynomial.h

**Description:** Polynomial operations

d41d8c, 290 lines

```
namespace Polynomial {
    template<typename base>
    vector<base> derivative(vector<base> a) {
        int n = a.size();
        for (int i = 0; i < n - 1; ++i) {
            a[i] = a[i + 1] * (i + 1);
        a.pop_back();
        return a;
    template<typename base>
    vector<base> integral(vector<base> a) {
        int n = a.size();
        a.push back(0);
        for (int i = n; i > 0; --i) {
            a[i] = a[i - 1] / i;
        a[0] = 0;
        return a;
    template<typename base>
    vector<base> add(vector<base> a, const vector<base> &b) {
        int n = a.size(), m = b.size();
        a.resize(max(n, m));
        for (int i = 0; i < max(n, m); ++i) {
            a[i] = (i \ge a.size() ? 0 : a[i]) + (i \ge b.size()
                 ? 0 : b[i]);
        return a;
    template<typename base>
    vector<base> sub(vector<base> a, const vector<base> &b) {
        int n = a.size(), m = b.size();
        a.resize(max(n, m));
        for (int i = 0; i < max(n, m); ++i) {</pre>
            a[i] = (i \ge a.size() ? 0 : a[i]) - (i \ge b.size()
                 ? 0 : b[i]);
        return a;
    namespace NTT {
        const int MOD = 998244353;
        const int q = 3;
        vector<int> R;
        void NTT(vector<Mint<MOD>>& a, int n, int on) {
            for (int i = 0; i < n; i++)</pre>
                if (i < R[i])
                    swap(a[i], a[R[i]]);
            Mint<MOD> wn, u, v;
            for (int i = 1, m = 2; i < n; i = m, m <<= 1) {
                wn = Mint < MOD > :: binpow(q, (MOD - 1) / m);
                if (on == -1)
                    wn = 1 / wn;
                for (int j = 0; j < n; j += m) {
                    Mint < MOD > w = 1;
```

```
for (int k = 0; k < i; k++, w *= wn) {
                    u = a[j + k], v = w * a[i + j + k];
                    a[j + k] = u + v;
                    a[i + j + k] = u - v;
           }
        if (on == -1) {
           Mint<MOD> k = Mint<MOD>(1) / Mint<MOD>(n);
            for (int i = 0; i < n; i++)</pre>
                a[i] = a[i] * k;
    template<typename base>
    vector<base>& A, vector<base>& B) {
        static_assert(std::is_same_v<base, Mint<MOD>>);
        assert (A.size() == B.size() && __builtin_popcount (A
             .size()) == 1);
        int n = A.size();
        int L = __builtin_ctz(n);
       if (R.size() != n) {
           R.assign(n, 0);
            for (int i = 0; i < n; i++)</pre>
                R[i] = (R[i >> 1] >> 1) | ((i & 1) << (L -
       NTT(A, n, 1);
       NTT(B, n, 1);
        for (int i = 0; i < n; i++)
           A[i] \star = B[i];
       NTT(A, n, -1);
        return A;
int get_lim(int n) {
   int res = 1;
    while (res < n) {
        res <<= 1;
    return res;
template<typename base>
vector<br/>base> mul(vector<br/>base> a, vector<br/>base> b, int size)
    int l = get_lim(a.size() + b.size());
   a.resize(1);
   b.resize(1);
   auto res = NTT::mul(a, b);
    res.resize(size);
    return res;
template<tvpename base>
vector<base> mul(vector<base> a, base scalar) {
    for (auto& val : a)
       val *= scalar:
    return a;
template<typename base>
vector<base> mul(const vector<base> &a, const vector<base>
    return mul(a, b, a.size() + b.size() - 1);
template <typename base>
```

```
vector<base> plug_minus_x(vector<base> a) {
    for (int i = 1; i < a.size(); i += 2) {</pre>
        a[i] *= -1;
    return a;
template <typename base>
void plug_x_squared_inplace(vector<base>& a) {
    a.resize(a.size() * 2);
    for (int i = (int) a.size() * 2 - 1; i >= 0; --i) {
        if (i % 2 != 0) a[i] = 0;
        else a[i] = a[i / 2];
template <typename base>
vector<base> plug_x_squared(const vector<base>& a) {
    vector<base> res(a.size() * 2);
    for (int i = 0; i < a.size(); ++i) {</pre>
        res[i * 2] = a[i];
    return res;
template <typename base>
void only_even_inplace(vector<base>& a) {
    for (int i = 0; i < a.size(); i += 2) {</pre>
        a[i / 2] = a[i];
    a.resize((a.size() + 1) / 2);
template <typename base>
vector<base> only_even(const vector<base>& a) {
    vector<base> res((a.size() + 1) / 2);
    for (int i = 0; i < a.size(); i += 2) {</pre>
        res[i / 2] = a[i];
    return res;
// O(n*log(n))
template<typename base>
void inverse_inplace(vector<base> &a, int size) {
    assert(!a.empty() && a[0] != 0);
    if (size == 0) {
        a = \{0\};
        return;
    if (size == 1) {
        a = \{1/a[0]\};
        return;
    auto op = plug_minus_x(a);
    auto T = mul(a, op);
    only_even_inplace(T);
    inverse_inplace(T, (size + 1) / 2);
    plug_x_squared_inplace(T);
    a = mul(op, T, size);
template <typename base>
vector<br/>base> inverse(const vector<br/>base>& a, int size) {
    assert(size > 0 && a[0] != 0);
    vector<base> Q{1/a[0]};
    for (int sz = 2;; sz *= 2) {
        Q = mul(Q, sub(\{2\}, mul(a, Q, sz)), sz);
```

```
if (sz >= size)
            break;
    O.resize(size);
    return Q;
  // O(n*log(n)) too slow, big constant factor
  template<typename base>
  vector<br/>
base> inverse(const vector<br/>
base> &a, int size) {
      assert(!a.empty() \&\& a[0] != 0);
       if (size = 0) f
          return \{0\};
      if (size == 1)  {
          return \{1/a[0]\};
      auto op = plug\_minus\_x(a);
      auto T = mul(a, op);
      T = only_{-even}(T);
      T = inverse(T, (size + 1) / 2);
      T = plug\_x\_squared(T);
      auto res = mul(op, T, size);
      return res;
template<typename base>
vector<br/>base> divide(const vector<br/>base> &a, const vector<
     base> &b, int size) {
    return mul(a, inverse(b, size), size);
// O(n*log(n))
template<typename base>
vector<base> ln(const vector<base> &a, int size) {
    auto res = integral(divide(derivative(a), a, size));
    res.resize(size);
    return res;
// O(n*log(n))
template<typename base>
vector<base> exp(const vector<base> &a, int size) {
    assert(size > 0 && a[0] == 0);
    vector<base> Q{1};
    for (int sz = 2;; sz *= 2) {
        Q = \text{mul}(Q, \text{sub}(\text{add}(a, \{1\}), \text{ln}(Q, \text{sz})), \text{sz});
        if (sz >= size)
            break;
    O.resize(size);
    return 0;
// O(n*log(n))
template<typename base>
vector<base> pow(vector<base> a, ll p, int size) {
    int i = 0;
    while (i < a.size()) {
        if (a[i] != 0)
            break:
        ++i;
    if (i == a.size()) {
        auto res = vector<base>(size, 0);
        if (p == 0)
            res[0] = 1;
        return res;
```

## PolyInterpolate BerlekampMassey

```
a.erase(a.begin(), a.begin() + i);
        auto f = a[0];
        for (auto& x : a) x \neq f;
        a = \exp(\text{mul}(\ln(a, \text{size}), (\text{base})p), \text{size});
        for (int j = size - 1; j >= 0; --j) {
             if ((i > 0 \&\& p >= size) || j - p * i < 0)
                 a[j] = 0;
             else
                 a[j] = a[j - i * p];
             a[j] *= base::binpow(f, p);
        return a;
int32 t main() {
    ios_base::sync_with_stdio(false);
    cin.tie(nullptr);
    const int MOD = 998244353;
    int n; cin >> n;
    11 m; cin >> m;
    vector<Mint<MOD>> a(n);
    for (auto& x : a) cin >> x;
    auto res = Polynomial::pow(a, m, n);
    for (auto x : res) cout << x << " ";</pre>
PolvInterpolate.h
Description: Given n points (x[i], y[i]), computes an n-1-degree polynomial
```

p that passes through them:  $p(x) = a[0] * x^0 + ... + a[n-1] * x^{n-1}$ . For numerical precision, pick  $x[k] = c * \cos(k/(n-1) * \pi), k = 0 \dots n-1$ . Time:  $\mathcal{O}(n^2)$ d41d8c, 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;
```

## BerlekampMassev.h

**Description:** Recovers any n-order linear recurrence relation from the first 2n terms of the recurrence. Useful for guessing linear recurrences after bruteforcing the first terms. Should work on any field, but numerical stability for floats is not guaranteed. Output will have size  $\leq n$ . Usage: berlekampMassev({0, 1, 1, 3, 5, 11}) // {1, 2}

```
Time: \mathcal{O}(N^2) and \mathcal{O}(nlog^2(k))
                                                         d41d8c, 143 lines
constexpr int mod = 1e9 + 7;
template<int32 t MOD>
struct modint {
    int32_t value;
    modint() = default;
    modint(int32_t value_) : value(value_) {}
    inline modint<MOD> operator+(modint<MOD> other) const {
```

int32\_t c = this->value + other.value;

return modint<MOD>(c >= MOD ? c - MOD : c);

```
inline modint<MOD> operator-(modint<MOD> other) const {
        int32 t c = this->value - other.value;
        return modint<MOD>(c < 0 ? c + MOD : c);</pre>
    inline modint<MOD> operator*(modint<MOD> other) const {
        int32_t c = (int64_t) this->value * other.value % MOD;
        return modint<MOD>(c < 0 ? c + MOD : c);</pre>
    inline modint<MOD> &operator+= (modint<MOD> other) {
        this->value += other.value;
        if (this->value >= MOD) this->value -= MOD;
        return *this;
    inline modint<MOD> &operator-= (modint<MOD> other) {
        this->value -= other.value;
        if (this->value < 0) this->value += MOD;
        return *this;
    inline modint<MOD> &operator*=(modint<MOD> other) {
        this->value = (int64_t) this->value * other.value % MOD
        if (this->value < 0) this->value += MOD;
        return *this;
    inline modint<MOD> operator-() const { return modint<MOD>(
        this->value ? MOD - this->value : 0); }
    modint<MOD> pow(uint64_t k) const {
        modint < MOD > x = *this, y = 1;
        for (; k; k >>= 1) {
            if (k & 1) y *= x;
            x *= x;
        return y;
    modint<MOD> inv() const { return pow(MOD - 2); } // MOD
         must be a prime
    inline modint<MOD> operator/(modint<MOD> other) const {
         return *this * other.inv(); }
    inline modint<MOD> operator/=(modint<MOD> other) { return *
         this *= other.inv(); }
    inline bool operator== (modint<MOD> other) const { return
         value == other.value; }
    inline bool operator!=(modint<MOD> other) const { return
         value != other.value; }
    inline bool operator<(modint<MOD> other) const { return
        value < other.value; }</pre>
    inline bool operator>(modint<MOD> other) const { return
        value > other.value; }
template<int32 t MOD>
modint<MOD> operator*(int64_t value, modint<MOD> n) { return
    modint<MOD>(value) * n; }
template<int32_t MOD>
```

```
modint<MOD> operator*(int32_t value, modint<MOD> n) { return
     modint<MOD>(value % MOD) * n; }
template<int32 t MOD>
istream &operator>>(istream &in, modint<MOD> &n) { return in >>
      n.value: }
template<int32 t MOD>
ostream &operator<<(ostream &out, modint<MOD> n) { return out
     << n.value; }
using mint = modint<mod>;
vector<mint> BerlekampMassey(vector<mint> S) {
    int n = (int) S.size(), L = 0, m = 0;
    vector<mint> C(n), B(n), T;
    C[0] = B[0] = 1;
    mint b = 1;
    for (int i = 0; i < n; i++) {</pre>
        ++m:
        mint d = S[i];
        for (int j = 1; j <= L; j++) d += C[j] * S[i - j];</pre>
        if (d == 0) continue;
        T = C;
        mint coef = d * b.inv();
        for (int j = m; j < n; j++) C[j] -= coef * B[j - m];</pre>
        if (2 * L > i) continue;
        L = i + 1 - L;
        B = T;
        b = d;
        m = 0;
    C.resize(L + 1);
    C.erase(C.begin());
    for (auto &x: C) x *= -1;
    return C;
vector<mint> combine(int n, vector<mint> &a, vector<mint> &b,
     vector<mint> &tr) {
    vector<mint> res(n * 2 + 1, 0);
    for (int i = 0; i < n + 1; i++) {</pre>
        for (int j = 0; j < n + 1; j++) res[i + j] += a[i] * b[
    for (int i = 2 * n; i > n; --i) {
        for (int j = 0; j < n; j++) res[i - 1 - j] += res[i] *
             tr[j];
    res.resize(n + 1);
    return res;
};
// transition \rightarrow for(i = 0; i < x; i++) f[n] += tr[i] * f[n-i]
     -1
// S contains initial values, k is 0 indexed
mint LinearRecurrence(vector<mint> &S, vector<mint> &tr, long
     long k) {
    int n = S.size();
    assert(n == (int) tr.size());
    if (n == 0) return 0;
    if (k < n) return S[k];</pre>
    vector<mint> pol(n + 1), e(pol);
    pol[0] = e[1] = 1;
    for (++k; k; k /= 2) {
        if (k % 2) pol = combine(n, pol, e, tr);
        e = combine(n, e, e, tr);
    mint res = 0;
```

```
for (int i = 0; i < n; i++) res += pol[i + 1] * S[i];</pre>
    return res;
int32_t main() {
    vector<mint> a{1, 1, 2, 3, 5, 8}; // precalc for small
         values
    int n = 10;
    auto tr = BerlekampMassey(a);
    a.resize(tr.size());
    cout << LinearRecurrence(a, tr, n);</pre>
```

## 3.2 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  $\star = -1$ ; res  $\star$ = a[i][i]; if (res == 0) return 0;  $rep(j, i+1, n) {$ **double** v = a[i][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:  $\mathcal{O}(N^3)$ 

d41d8c, 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 \star = -1;
   ans = ans * a[i][i] % mod;
   if (!ans) return 0;
  return (ans + mod) % mod;
```

## SolveLinear.h

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

d41d8c, 38 lines

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

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

#### SolveLinear2.h

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

```
"SolveLinear.h"
                                                       d41d8c, 7 lines
rep(j,0,n) if (j!= i) // instead of rep(j, i+1,n)
// ... then at the end:
x.assign(m, undefined);
rep(i,0,rank) {
 rep(j,rank,m) if (fabs(A[i][j]) > eps) goto fail;
 x[col[i]] = b[i] / A[i][i];
fail:: }
```

#### SolveLinearBinarv.h

**Description:** Solves Ax = b over  $\mathbb{F}_2$ . If there are multiple solutions, one is returned arbitrarily. Returns rank, or -1 if no solutions. Destroys A and b. Time:  $\mathcal{O}\left(n^2m\right)$ d41d8c, 34 lines

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

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

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

```
d41d8c, 35 lines
int matInv(vector<vector<double>>& A) {
 int n = sz(A); vi col(n);
 vector<vector<double>> tmp(n, vector<double>(n));
  rep(i, 0, n) tmp[i][i] = 1, col[i] = i;
  rep(i,0,n) {
    int r = i, c = i;
    rep(j,i,n) rep(k,i,n)
      if (fabs(A[j][k]) > fabs(A[r][c]))
        r = j, c = k;
    if (fabs(A[r][c]) < 1e-12) return i;</pre>
    A[i].swap(A[r]); tmp[i].swap(tmp[r]);
      swap(A[j][i], A[j][c]), swap(tmp[j][i], tmp[j][c]);
    swap(col[i], col[c]);
    double v = A[i][i];
    rep(j,i+1,n) {
      double f = A[j][i] / v;
      A[j][i] = 0;
      rep(k, i+1, n) A[j][k] -= f*A[i][k];
      rep(k,0,n) tmp[j][k] -= f*tmp[i][k];
    rep(j, i+1, n) A[i][j] /= v;
    rep(j,0,n) tmp[i][j] /= v;
    A[i][i] = 1;
  for (int i = n-1; i > 0; --i) rep(j, 0, i) {
    double v = A[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:
```

#### MatrixInverse-mod.h

**Description:** Invert matrix A modulo a prime. Returns rank; result is stored in A unless singular (rank < n). 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)
```

```
"../number-theory/ModPow.h"
                                                              d41d8c, 36 lines
int matInv(vector<vector<11>>& A) {
```

```
int n = sz(A); vi col(n);
```

## Tridiagonal FFT FFTM NTT

```
vector<vector<ll>> tmp(n, vector<ll>(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 (A[j][k]) {
     r = j; c = k; goto found;
   return i;
found:
   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]);
    ll v = modpow(A[i][i], mod - 2);
    rep(j,i+1,n) {
     11 f = A[j][i] * v % mod;
     A[j][i] = 0;
     rep(k, i+1, n) A[j][k] = (A[j][k] - f*A[i][k]) % mod;
     rep(k,0,n) tmp[j][k] = (tmp[j][k] - f*tmp[i][k]) % mod;
    rep(j, i+1, n) A[i][j] = A[i][j] * v % mod;
    rep(j, 0, n) tmp[i][j] = tmp[i][j] * v % mod;
   A[i][i] = 1;
  for (int i = n-1; i > 0; --i) rep(j, 0, i) {
   11 v = A[j][i];
   rep(k,0,n) tmp[j][k] = (tmp[j][k] - v*tmp[i][k]) % mod;
  rep(i,0,n) rep(j,0,n)
   A[col[i]][col[j]] = tmp[i][j] % mod + (tmp[i][j] < 0 ? mod
  return n;
```

## Tridiagonal.h

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

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

This is useful for solving problems on the type

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

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

$$\begin{aligned} \{a_i\} &= \operatorname{tridiagonal}(\{1,-1,-1,\ldots,-1,1\},\{0,c_1,c_2,\ldots,c_n\},\\ \{b_1,b_2,\ldots,b_n,0\},\{a_0,d_1,d_2,\ldots,d_n,a_{n+1}\}). \end{aligned}$$

Fails if the solution is not unique.

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

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

d41d8c, 26 lines

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

```
} else {
    diag[i+1] -= super[i]*sub[i]/diag[i];
    b[i+1] -= b[i]*sub[i]/diag[i];
}

for (int i = n; i--;) {
    if (tr[i]) {
        swap(b[i], b[i-1]);
        diag[i-1] = diag[i];
        b[i] /= super[i-1];
} else {
    b[i] /= diag[i];
    if (i) b[i-1] -= b[i]*super[i-1];
}
}
return b;
```

#### 3.3 Fourier transforms

#### FFT.h

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

```
Time: O(N \log N) with N = |A| + |B| (~1s for N = 2^{22}) d41d8c, 35 lines
```

```
typedef complex<double> C;
typedef vector<double> vd;
void fft(vector<C>& a) {
 int n = sz(a), L = 31 - builtin clz(n);
  static vector<complex<long double>> R(2, 1);
  static vector<C> rt(2, 1); // (^ 10% faster if double)
  for (static int k = 2; k < n; k \neq 2) {
   R.resize(n); rt.resize(n);
    auto x = polar(1.0L, acos(-1.0L) / k);
    rep(i,k,2*k) rt[i] = R[i] = i&1 ? R[i/2] * x : R[i/2];
  rep(i, 0, n) rev[i] = (rev[i / 2] | (i & 1) << L) / 2;
  rep(i,0,n) if (i < rev[i]) swap(a[i], a[rev[i]]);
  for (int k = 1; k < n; k *= 2)
    for (int i = 0; i < n; i += 2 * k) rep(j, 0, k) {
     Cz = rt[j+k] * a[i+j+k]; // (25\% faster if hand-rolled)
     a[i + j + k] = a[i + j] - z;
     a[i + j] += z;
vd conv(const vd& a, const vd& b) {
 if (a.emptv() || b.emptv()) return {};
  vd res(sz(a) + sz(b) - 1);
  int L = 32 - __builtin_clz(sz(res)), n = 1 << L;</pre>
  vector<C> in(n), out(n);
  copy(all(a), begin(in));
  rep(i,0,sz(b)) in[i].imag(b[i]);
  fft(in);
  for (C& x : in) x *= x;
  rep(i, 0, n) out[i] = in[-i & (n - 1)] - conj(in[i]);
  rep(i, 0, sz(res)) res[i] = imag(out[i]) / (4 * n);
  return res;
```

#### FFTM h

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

```
Time: \mathcal{O}(N \log N), where N = |A| + |B| (twice as slow as NTT or FFT)
"FastFourierTransform.h"
typedef vector<11> v1;
template<int M> vl convMod(const vl &a, const vl &b) {
  if (a.empty() || b.empty()) return {};
  vl res(sz(a) + sz(b) - 1);
  int B=32-__builtin_clz(sz(res)), n=1<<B, cut=int(sqrt(M));</pre>
  vector<C> L(n), R(n), outs(n), outl(n);
  rep(i,0,sz(a)) L[i] = C((int)a[i] / cut, (int)a[i] % cut);
  rep(i,0,sz(b)) R[i] = C((int)b[i] / cut, (int)b[i] % cut);
  fft(L), fft(R);
  rep(i,0,n) {
    int j = -i \& (n - 1);
    outl[j] = (L[i] + conj(L[j])) * R[i] / (2.0 * n);
    outs[j] = (L[i] - conj(L[j])) * R[i] / (2.0 * n) / 1i;
  fft (outl), fft (outs);
  rep(i,0,sz(res)) {
    11 \text{ av} = 11(\text{real}(\text{outl}[i]) + .5), \text{ cv} = 11(\text{imag}(\text{outs}[i]) + .5);
    11 \text{ bv} = 11(\text{imag}(\text{outl}[i]) + .5) + 11(\text{real}(\text{outs}[i]) + .5);
    res[i] = ((av % M * cut + bv) % M * cut + cv) % M;
  return res;
```

#### NTT.h

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

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

ntt(out);

```
"../number-theory/ModPow.h"
const 11 mod = (119 \ll 23) + 1, root = 62; // = 998244353
// For p < 2^30 there is also e.g. 5 << 25, 7 << 26, 479 << 21
// and 483 \ll 21 (same root). The last two are > 10^9.
typedef vector<11> v1:
void ntt(vl &a) {
 int n = sz(a), L = 31 - __builtin_clz(n);
 static v1 rt(2, 1);
 for (static int k = 2, s = 2; k < n; k *= 2, s++) {
   rt.resize(n):
   ll z[] = \{1, modpow(root, mod >> s)\};
   rep(i,k,2*k) rt[i] = rt[i / 2] * z[i & 1] % mod;
 vi rev(n);
 rep(i,0,n) rev[i] = (rev[i / 2] | (i & 1) << L) / 2;
 rep(i,0,n) if (i < rev[i]) swap(a[i], a[rev[i]]);
 for (int k = 1; k < n; k *= 2)
    for (int i = 0; i < n; i += 2 * k) rep(j,0,k) {
     11 z = rt[j + k] * a[i + j + k] % mod, &ai = a[i + j];
     a[i + j + k] = ai - z + (z > ai ? mod : 0);
     ai += (ai + z >= mod ? z - mod : z);
vl conv(const vl &a, const vl &b) {
 if (a.empty() || b.empty()) return {};
 int s = sz(a) + sz(b) - 1, B = 32 - _builtin_clz(s), n = 1
      << B;
 int inv = modpow(n, mod - 2);
 vl L(a), R(b), out(n);
 L.resize(n), R.resize(n);
 ntt(L), ntt(R);
 rep(i,0,n) out[-i \& (n-1)] = (l1)L[i] * R[i] % mod * inv %
```

```
return {out.begin(), out.begin() + s};
```

## FST.h

Description: Transform to a basis with fast convolutions of the form  $c[z] = \sum_{z=x \oplus y} a[x] \cdot b[y]$ , where  $\oplus$  is one of AND, OR, XOR. The size of a must be a power of two.

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

d41d8c, 16 lines

```
void FST(vi& a, bool inv) {
  for (int n = sz(a), step = 1; step < n; step *= 2) {
    for (int i = 0; i < n; i += 2 * step) rep(j,i,i+step) {</pre>
     int &u = a[j], &v = a[j + step]; tie(u, v) =
       inv ? pii(v - u, u) : pii(v, u + v); // AND
       inv ? pii(v, u - v) : pii(u + v, u); // OR
       pii(u + v, u - v);
 if (inv) for (int& x : a) x /= sz(a); // XOR only
vi conv(vi a, vi b) {
 FST(a, 0); FST(b, 0);
  rep(i, 0, sz(a)) a[i] *= b[i];
 FST(a, 1); return a;
```

# Number theory (4)

## 4.1 Modular arithmetic

Modular Arithmetic.h

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

```
const 11 mod = 17; // change to something else
struct Mod {
 11 x;
  Mod(ll xx) : x(xx) \{ \}
  Mod operator+(Mod b) { return Mod((x + b.x) % mod); }
  Mod operator-(Mod b) { return Mod((x - b.x + mod) % mod); }
  Mod operator*(Mod b) { return Mod((x * b.x) % mod); }
  Mod operator/(Mod b) { return *this * invert(b); }
  Mod invert (Mod a) {
   ll x, y, 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;
};
```

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

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

```
return -1;
```

#### ModSum.h

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

floored division.

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

d41d8c, 16 lines

```
typedef unsigned long long ull;
ull sumsq(ull to) { return to /2 * ((to-1) | 1); }
ull divsum(ull to, ull c, ull k, ull m) {
 ull res = k / m * sumsq(to) + c / m * to;
 k %= m; c %= m;
 if (!k) return res;
 ull to2 = (to * k + c) / m;
 return res + (to - 1) * to2 - divsum(to2, m-1 - c, m, k);
ll modsum(ull to, ll c, ll k, ll m) {
 c = ((c % m) + m) % m;
 k = ((k % m) + m) % m;
 return to * c + k * sumsq(to) - m * divsum(to, c, k, m);
```

## ModMulLL.h

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

```
typedef unsigned long long ull;
ull modmul(ull a, ull b, ull M) {
 ll ret = a * b - M * ull(1.L / M * a * b);
 return ret + M * (ret < 0) - M * (ret >= (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);
```

## ModSart.h

b = b \* q % p;

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

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

"ModPow.h" d41d8c, 24 lines 11 sgrt(ll a, ll p) { a %= p; if (a < 0) a += p; **if** (a == 0) **return** 0; assert (modpow(a, (p-1)/2, p) == 1); // else no solution **if** (p % 4 == 3) **return** modpow(a, (p+1)/4, p);  $// a^{(n+3)/8} \text{ or } 2^{(n+3)/8} * 2^{(n-1)/4} \text{ works if } p \% 8 == 5$ 11 s = p - 1, n = 2;int r = 0, m; **while** (s % 2 == 0) ++r, s /= 2; **while** (modpow(n, (p-1) / 2, p) != p-1) ++n;11 x = modpow(a, (s + 1) / 2, p);ll b = modpow(a, s, p), g = modpow(n, s, p); **for** (;; r = m) { 11 t = b;for (m = 0; m < r && t != 1; ++m) t = t \* t % p;if (m == 0) return x;  $11 \text{ gs} = \text{modpow}(g, 1LL \ll (r - m - 1), p);$ q = qs \* qs % p;x = x \* qs % p;

```
4.2 Primality
```

#### FastErat.h

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

Time: LIM=1e9  $\approx 1.5s$ 

```
const int LIM = 1e6;
bitset<LIM> isPrime;
vi eratosthenes() {
  const int S = (int)round(sqrt(LIM)), R = LIM / 2;
  vi pr = {2}, sieve(S+1); pr.reserve(int(LIM/log(LIM)*1.1));
  vector<pii> cp;
  for (int i = 3; i <= S; i += 2) if (!sieve[i]) {</pre>
    cp.push_back(\{i, i * i / 2\});
    for (int j = i * i; j <= S; j += 2 * i) sieve[j] = 1;</pre>
  for (int L = 1; L <= R; L += S) {
    array<bool, S> block{};
    for (auto &[p, idx] : cp)
      for (int i=idx; i < S+L; idx = (i+=p)) block[i-L] = 1;</pre>
    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"
                                                       d41d8c, 12 lines
bool isPrime(ull n) {
 if (n < 2 | | n % 6 % 4 != 1) return (n | 1) == 3;</pre>
 ull A[] = \{2, 325, 9375, 28178, 450775, 9780504, 1795265022\}
     s = \underline{builtin\_ctzll(n-1)}, d = n >> s;
 for (ull a : A) { // ^ count trailing zeroes
    ull p = modpow(a%n, d, n), i = s;
    while (p != 1 && p != n - 1 && a % n && i--)
     p = modmul(p, p, n);
    if (p != n-1 && i != s) return 0;
 return 1;
```

#### Factor.h

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

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

d41d8c, 41 lines

```
using db = long double;
using 11 = long long;
using ul = uint64_t;
ul modMul(ul a, ul b, const ul mod) {
    11 \text{ ret} = a*b-mod*(ul)((db)a*b/mod);
    return ret+((ret<0)-(ret>=(11)mod))*mod; }
ul modPow(ul a, ul b, const ul mod) {
    if (b == 0) return 1;
    ul res = modPow(a,b/2,mod); res = modMul(res,res,mod);
    return b&1 ? modMul(res,a,mod) : res;
bool prime(ul n) { // not ll!
    if (n < 2 | | n % 6 % 4 != 1) return n-2 < 2;
```

## euclid CRT phi ContinuedFractions FracBinarySearch

```
ul A[] = \{2, 325, 9375, 28178, 450775, 9780504, 1795265022\}
           s = builtin ctzll(n-1), d = n>>s;
    for (auto &a : A) { // ^ count trailing zeroes
       ul p = modPow(a,d,n), i = s;
       while (p != 1 && p != n-1 && a%n && i--) p = modMul(p,p
       if (p != n-1 && i != s) return 0;
   return 1:
ul pollard(ul n) { // return some nontrivial factor of n
   auto f = [n](ul x) { return modMul(x, x, n) + 1; };
   ul x = 0, y = 0, t = 30, prd = 2, i = 1, q;
   while (t++ % 40 || __gcd(prd, n) == 1) {
       if (x == y) x = ++i, y = f(x);
       if ((q = modMul(prd, max(x,y) - min(x,y), n))) prd = q;
       x = f(x), y = f(f(y));
   return __gcd(prd, n);
void factor rec(ul n, map<ul,int>& cnt) {
   if (n == 1) return;
   if (prime(n)) { ++cnt[n]; return; }
   ul u = pollard(n);
    factor_rec(u,cnt), factor_rec(n/u,cnt);
```

## 4.3 Divisibility

#### euclid.h

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

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

## CRT.h

 ${\bf Description:}\ {\bf Chinese}\ {\bf Remainder}\ {\bf Theorem}.$ 

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

## 4.3.1 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

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

```
Distribution: Euler's \phi function is defined as \phi(n) := \# of positive integers \leq n that are coprime with n. \phi(1) = 1, p prime \Rightarrow \phi(p^k) = (p-1)p^{k-1}, m, n coprime \Rightarrow \phi(mn) = \phi(m)\phi(n). If n = p_1^{k_1}p_2^{k_2}...p_r^{k_r} then \phi(n) = (p_1-1)p_1^{k_1-1}...(p_r-1)p_r^{k_r-1}. \phi(n) = n \cdot \prod_{p|n} (1-1/p). \sum_{d|n} \phi(d) = n, \sum_{1 \leq k \leq n, \gcd(k,n)=1} k = n\phi(n)/2, n > 1 Euler's thm: a, n coprime \Rightarrow a^{\phi(n)} \equiv 1 \pmod{n}. Fermat's little thm: p prime \Rightarrow a^{p-1} \equiv 1 \pmod{p} \ \forall a. d41d8c, 8 lines const int LIM = 5000000; int phi[LIM]; void calculatePhi() { rep(i, 0, LIM) phi[i] = i&1 ? i : i/2; for (int i = 3; i < LIM; i += 2) if(phi[i] == i) for (int j = i; j < LIM; j += i) phi[j] -= phi[j] / i; }
```

## 4.4 Fractions

ContinuedFractions.h

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

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

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

```
typedef double d; // for N ~ 1e7; long double for N ~ 1e9
pair<11, 11> approximate(d x, 11 N) {
    11 LP = 0, LQ = 1, P = 1, Q = 0, inf = LLONG_MAX; d y = x;
    for (;;) {
        11 lim = min(P ? (N-LP) / P : inf, Q ? (N-LQ) / Q : inf),
            a = (11) floor(y), b = min(a, lim),
            NP = b*P + LP, NQ = b*Q + LQ;
        if (a > b) {
            // If b > a/2, we have a semi-convergent that gives us a
            // better approximation; if b = a/2, we *may* have one.
            // Return {P, Q} here for a more canonical approximation.
            return (abs(x - (d)NP / (d)NQ) < abs(x - (d)P / (d)Q)) ?
            make_pair(NP, NQ) : make_pair(P, Q);
        }
        if (abs(y = 1/(y - (d)a)) > 3*N) {
            return {NP, NQ};
        }
        LP = P; P = NP;
        LQ = Q; Q = NQ;
        }
}
```

#### FracBinarySearch.h

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

Usage: fracBS([|(Frac f) { return f, p>=3\*f.g; }, 10); // {1,3}

```
Usage: fracBS([](Frac f) { return f.p>=3*f.q; }, 10); // {1,3} 
Time: \mathcal{O}(\log(N)) d41d8c, 25 lines
```

```
struct Frac { 11 p, q; };

template < class F >
Frac fracBS(F f, 11 N) {
  bool dir = 1, A = 1, B = 1;
  Frac lo{0, 1}, hi{1, 1}; // Set hi to 1/0 to search (0, N]
  if (f(lo)) return lo;
  assert(f(hi));
  while (A || B) {
    11 adv = 0, step = 1; // move hi if dir, else lo
    for (int si = 0; step; (step *= 2) >>= si) {
      adv += step;
    }
}
```

```
Frac mid{lo.p * adv + hi.p, lo.q * adv + hi.q};
if (abs(mid.p) > N || mid.q > N || dir == !f(mid)) {
    adv -= step; si = 2;
}
hi.p += lo.p * adv;
hi.q += lo.q * adv;
dir = !dir;
swap(lo, hi);
A = B; B = !!adv;
}
return dir ? hi : lo;
```

## 4.5 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.

## 4.6 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  $\mathbb{Z}_{2a}^{\times}$  is instead isomorphic to  $\mathbb{Z}_2 \times \mathbb{Z}_{2a-2}$ .

## 4.7 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.

## 4.8 Mobius Function

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

Mobius Inversion:

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

Other useful formulas/forms:

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

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

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

## IntPerm multinomial

# Combinatorial (5)

## 5.1 Permutations

## 5.1.1 Factorial

	1 2 3							
n!	1 2 6	24 1	20 72	0 504	0 403	20 362	2880 3	628800
n	11	12	13	1	4	15	16	17
n!	4.0e7	4.8€	8 6.26	98.7	e10 1	.3e12	2.1e13	3.6e14
n	20	25	30	40	50	100	150	171
n!	2e18	2e25	3e32	8e47	3e64	9e157	6e262	>DBL_MAX

#### IntPerm.h

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

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

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

## **5.1.2** Cycles

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

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

## 5.1.3 Derangements

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

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

#### 5.1.4 Burnside's lemma

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

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

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

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

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

## .2 Partitions and subsets

## 5.2.1 Partition function

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

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

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

## 5.2.2 Lucas' Theorem

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

## 5.2.3 Binomials

multinomial.h

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

11 multinomial (vi& v) {
11 c = 1, m = v.empty() ? 1 : v[0];

## 5.3 General purpose numbers

## 5.3.1 Bernoulli numbers

EGF of Bernoulli numbers is  $B(t) = \frac{t}{e^t - 1}$  (FFT-able).  $B[0, ...] = [1, -\frac{1}{2}, \frac{1}{6}, 0, -\frac{1}{20}, 0, \frac{1}{42}, ...]$ 

Sums of powers:

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

Euler-Maclaurin formula for infinite sums:

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

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

## 5.3.2 Stirling numbers of the first kind

Number of permutations on n items with k cycles.

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

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

#### 5.3.3 Eulerian numbers

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

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

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

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

## 5.3.4 Stirling numbers of the second kind

Partitions of n distinct elements into exactly k groups.

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

## 5.3.5 Bell numbers

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

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

## 5.3.6 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  $d_i$ :  $(n-2)!/((d_1-1)!\cdots(d_n-1)!)$ 

## 5.3.7 Catalan numbers

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

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

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

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

# $\underline{\text{Graph}}$ (6)

## 6.1 Network flow

## MinCostMaxFlow.h

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

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

```
d41d8c, 81 lines
#include <bits/extc++.h>
const 11 INF = numeric limits<11>::max() / 4;
typedef vector<ll> VL;
struct MCMF {
 int N;
  vector<vi> ed, red;
  vector<VL> cap, flow, cost;
 VL dist, pi;
  vector<pii> par:
  MCMF (int N) :
   N(N), ed(N), red(N), cap(N, VL(N)), flow(cap), cost(cap),
   seen(N), dist(N), pi(N), par(N) {}
  void addEdge(int from, int to, ll cap, ll cost) {
    this->cap[from][to] = cap;
   this->cost[from][to] = cost;
   ed[from].push back(to);
   red[to].push_back(from);
  void path(int s) {
    fill(all(seen), 0);
    fill(all(dist), INF);
   dist[s] = 0; ll di;
    __gnu_pbds::priority_queue<pair<11, int>> q;
    vector<decltype(q)::point_iterator> its(N);
   g.push({0, s});
    auto relax = [&](int i, ll cap, ll cost, int dir) {
     ll val = di - pi[i] + cost;
     if (cap && val < dist[i]) {
       dist[i] = val;
       par[i] = {s, dir};
       if (its[i] == q.end()) its[i] = q.push({-dist[i], i});
        else q.modify(its[i], {-dist[i], i});
   };
    while (!q.empty()) {
     s = q.top().second; q.pop();
     seen[s] = 1; di = dist[s] + pi[s];
     for (int i : ed[s]) if (!seen[i])
       relax(i, cap[s][i] - flow[s][i], cost[s][i], 1);
     for (int i : red[s]) if (!seen[i])
       relax(i, flow[i][s], -cost[i][s], 0);
    rep(i, 0, N) pi[i] = min(pi[i] + dist[i], INF);
  pair<11, 11> maxflow(int s, int t) {
    11 \text{ totflow} = 0, \text{ totcost} = 0;
    while (path(s), seen[t]) {
     11 fl = INF;
     for (int p,r,x = t; tie(p,r) = par[x], x != s; x = p)
       fl = min(fl, r ? cap[p][x] - flow[p][x] : flow[x][p]);
     totflow += fl;
      for (int p,r,x = t; tie(p,r) = par[x], x != s; x = p)
```

```
if (r) flow[p][x] += fl;
        else flow[x][p] -= fl;
   rep(i,0,N) rep(j,0,N) totcost += cost[i][j] * flow[i][j];
   return {totflow, totcost};
  // If some costs can be negative, call this before maxflow:
  void setpi(int s) { // (otherwise, leave this out)
   fill(all(pi), INF); pi[s] = 0;
   int it = N, ch = 1; ll v;
    while (ch-- && it--)
     rep(i,0,N) if (pi[i] != INF)
        for (int to : ed[i]) if (cap[i][to])
          if ((v = pi[i] + cost[i][to]) < pi[to])</pre>
            pi[to] = v, ch = 1;
    assert(it >= 0); // negative cost cycle
};
Dinic.h
Description: Flow algorithm with complexity O(VE \log U) where U =
\max |\operatorname{cap}|. O(\min(E^{1/2}, V^{2/3})E) if U = 1; O(\sqrt{V}E) for bipartite match-
                                                     d41d8c, 121 lines
template<class F>
struct dinic {
 constexpr static db eps = 1e-6;
 struct Edge {
   int to;
   F flow:
   F cap;
   Edge() = default;
    Edge(int to_, F flow_, F cap_) : to(to_), flow(flow_), cap(
        cap ) {}
 };
 vector<vector<int>> gr;
 vector<Edge> edges;
 dinic(int n_) : n(n_) {
   gr.resize(n);
 void add(int u, int v, F cap, F rev_cap = 0) {
    assert(min(cap, rev_cap) >= 0);
   int id = edges.size();
   edges.pb({v, 0, cap});
    edges.pb({u, 0, rev_cap});
   gr[u].push_back(id);
   gr[v].push_back(id ^ 1);
 void add_three(int s, int t, int u, int v, F l, F r) {
   add(s, v, 1);
   add(u, v, r - 1);
   add(u, t, 1);
 vector<F> dist;
 vector<int> first;
 inline F res(int id) {
```

```
return edges[id].cap - edges[id].flow;
inline F res(const Edge& e) {
  return e.cap - e.flow;
bool bfs(int s, int t) {
  dist.assign(n, -1);
  first.assign(n, 0);
  dist[s] = 0;
  queue<int> Fedya_Romashov({s});
  while(!Fedya_Romashov.empty()) {
    auto v = Fedya_Romashov.front(); Fedya_Romashov.pop();
    for (auto id : gr[v]) {
      auto& e = edges[id];
      if (res(id) > 0 && dist[e.to] < 0) {</pre>
        dist[e.to] = dist[v] + 1;
        Fedya_Romashov.push(e.to);
  return dist[t] >= 0;
F dfs(int v, int t, F current_flow = 0) {
  if (v == t) {
    return current_flow;
  F small push = 0;
  for (; first[v] < qr[v].size(); ++first[v]) {</pre>
    int id = gr[v][first[v]];
    auto& e = edges[id];
    if (abs(res(id)) < eps || dist[e.to] != dist[v] + 1) {
      continue;
    F pushed = dfs(e.to, t, min(current flow - small push,
         res(e)));
    if (pushed) {
      small_push += pushed;
      edges[id].flow += pushed;
      edges[id ^ 1].flow -= pushed;
    if (small_push == current_flow) {
      break;
  return small push;
F maxFlow(int s, int t) {
  F total = 0:
  while(bfs(s, t)) {
    while(F df = dfs(s, t, numeric_limits<F>::max())) {
      total += df;
  return total:
vector<bool> min_cut() {
  max_flow();
```

12

for (int i = 0; i < n; i++) {

ret[i] = (dist[i] != -1);

vector<bool> ret(n);

return ret;

## FlowDecomposition MCMF

```
};
FlowDecomposition.h
Description: Decompose flow into paths and cycles.
Time: \mathcal{O}(FLOW + m + n)
                                                     d41d8c, 105 lines
template <typename T>
class flow_decomposition {
public:
  const flow_graph<T> &g;
  vector<vector<int>> paths;
  vector<T> path_flows;
  vector<vector<int>> cycles;
  vector<T> cycle flows;
  flow_decomposition(const flow_graph<T> &_g) : g(_g) {
  void decompose() {
    vector<T> fs(g.edges.size());
    for (int i = 0; i < (int) g.edges.size(); i++) {</pre>
      fs[i] = g.edges[i].f;
    paths.clear();
    path_flows.clear();
    cvcles.clear();
    cycle flows.clear();
    vector<int> ptr(a.n);
    for (int i = 0; i < q.n; i++) {</pre>
     ptr[i] = (int) g.g[i].size() - 1;
    vector<int> was(q.n, -1);
    int start = q.st;
    for (int iter = 0; ; iter++) {
      bool found start = false;
      while (true) {
        if (ptr[start] >= 0) {
          int id = g.g[start][ptr[start]];
          if (fs[id] > q.eps) {
            found start = true;
            break;
          ptr[start]--;
          continue;
        start = (start + 1) % g.n;
        if (start == g.st) {
          break;
      if (!found_start) {
        break;
      vector<int> path;
      bool is_cycle = false;
      int v = start;
      while (true)
        if (v == q.fin) {
          break;
        if (was[v] == iter) {
          bool found = false;
```

```
for (int i = 0; i < (int) path.size(); i++) {</pre>
            int id = path[i];
            auto &e = q.edges[id];
            if (e.from == v) {
              path.erase(path.begin(), path.begin() + i);
              found = true;
              break;
          assert (found):
          is_cycle = true;
          break;
        was[v] = iter;
        bool found = false;
        while (ptr[v] >= 0) {
          int id = g.g[v][ptr[v]];
         if (fs[id] > g.eps) {
            path.push_back(id);
            v = g.edges[id].to;
            found = true;
            break;
          ptr[v]--;
        assert (found);
      T path_flow = numeric_limits<T>::max();
      for (int id : path) {
        path_flow = min(path_flow, fs[id]);
      for (int id : path) {
        fs[id] -= path_flow;
        fs[id ^ 1] += path_flow;
      if (is_cycle) {
        cycles.push_back(path);
        cycle_flows.push_back(path_flow);
        paths.push_back(path);
        path_flows.push_back(path_flow);
    for (const T& f : fs) {
     assert (-q.eps <= f && f <= q.eps);
 }
};
MCMF.h
Description: min cost max flow
Time: \mathcal{O}(flow * mlogn + mn)
                                                    d41d8c, 131 lines
template <class F, class C = F> struct MCMF {
 struct Edge {
   int to;
   F flow, cap;
   C cost;
 };
 vector<C> johnson_potential, dist;
 vector<int> previous_edge;
 vector<Edge> edges;
 vector<vector<int>> gr;
 MCMF (int n_) : n(n_) {
    johnson_potential.resize(n), dist.resize(n), previous_edge.
         resize(n), gr.resize(n);
```

```
void add(int u, int v, F cap, C cost) {
  assert(cap >= 0);
  gr[u].pb(edges.size());
  edges.pb({v, 0, cap, cost});
  gr[v].pb(edges.size());
  edges.pb({u, 0, 0, -cost});
bool path(int s, int t) {
  constexpr C inf = numeric_limits<C>::max();
  for (int i = 0; i < n; ++i) {</pre>
    dist[i] = inf;
  using T = pair<C, int>;
  priority_queue<T, vector<T>, greater<T>> Fedya_Romashov;
  Fedya_Romashov.push({dist[s] = 0, s});
  while (!Fedya_Romashov.empty()) {
      auto [cur_dist, v] = Fedya_Romashov.top();
    Fedya_Romashov.pop();
    if (cur_dist > dist[v]) {
      continue;
    for (auto &e : gr[v]) {
      auto &E = edges[e];
      if (E.flow < E.cap &&
          ckmin(dist[E.to], cur_dist + E.cost +
               johnson_potential[v] - johnson_potential[E.to
        previous_edge[E.to] = e, Fedya_Romashov.push({dist[E.
             tol, E.to});
  return dist[t] != inf;
pair<F, C> calc(int s, int t) {
  assert(s != t);
    for (int e = 0; e < edges.size(); ++e) {</pre>
      const Edge &E = edges[e]; // Bellman-Ford
        ckmin(johnson_potential[E.to], johnson_potential[
             edges[e ^ 1].to] + E.cost);
  F totalFlow = 0:
  C \text{ totalCost} = 0;
  while (path(s, t)) {
    for (int i = 0; i < n; ++i) {</pre>
      johnson_potential[i] += dist[i];
    F df = numeric_limits<F>::max();
    for (int x = t; x != s; x = edges[previous_edge[x] ^ 1].
      const Edge &E = edges[previous_edge[x]];
      ckmin(df, E.cap - E.flow);
```

## Hungarian GlobalMinCut hopcroftKarp

```
totalFlow += df;
   totalCost += (johnson_potential[t] - johnson_potential[s
        ]) * df;
    for (int x = t; x != s; x = edges[previous_edge[x] ^ 1].
      edges[previous_edge[x]].flow += df, edges[previous_edge
          [x] ^1.flow -= df;
  return {totalFlow, totalCost};
pair<F, C> k_calc(int s, int t, C max_cost = std::
    numeric_limits<C>::max()) {
  assert(s != t);
 rep(n) {
   for (int e = 0; e < edges.size(); ++e) {</pre>
     const Edge &E = edges[e]; // Bellman-Ford
     if (E.cap) {
        ckmin(johnson_potential[E.to], johnson_potential[
             edges[e ^ 1].to] + E.cost);
 F totalFlow = 0;
 C \text{ totalCost} = 0;
  while (path(s, t)) {
   for (int i = 0; i < n; ++i) {</pre>
      johnson_potential[i] += dist[i];
   F df = numeric_limits<F>::max();
   if ((johnson_potential[t] - johnson_potential[s]) > 0) {
        df = (max_cost - totalCost) / (johnson_potential[t] -
              johnson_potential[s]);
    for (int x = t; x != s; x = edges[previous_edge[x] ^ 1].
     const Edge &E = edges[previous_edge[x]];
      ckmin(df, E.cap - E.flow);
   totalFlow += df;
   totalCost += (johnson_potential[t] - johnson_potential[s
        ) \star df; // \leq max\_cost - total\_cost
    for (int x = t; x != s; x = edges[previous_edge[x] ^ 1].
      edges[previous_edge[x]].flow += df, edges[previous_edge
          [x] ^ 1].flow -= df;
    if (df == 0) {
        break:
  return {totalFlow, totalCost};
```

## Hungarian.h

Description: Solve assignment problem.

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

d41d8c, 81 lines

```
template <typename T>
class hungarian {
```

```
public:
 int n;
 int m:
 vector<vector<T>> a;
 vector<T> u;
 vector<T> v;
 vector<int> pa;
 vector<int> pb;
 vector<int> way;
 vector<T> minv:
 vector<bool> used;
T inf;
hungarian(int _n, int _m) : n(_n), m(_m) {
   assert (n <= m);
   a = vector<vector<T>>(n, vector<T>(m));
  u = vector < T > (n + 1);
  v = vector < T > (m + 1);
   pa = vector < int > (n + 1, -1);
   pb = vector < int > (m + 1, -1);
   way = vector<int>(m, -1);
   minv = vector<T>(m);
  used = vector<bool>(m + 1);
   inf = numeric_limits<T>::max();
 inline void add_row(int i) {
   fill(minv.begin(), minv.end(), inf);
   fill(used.begin(), used.end(), false);
   pb[m] = i;
   pa[i] = m;
   int j0 = m;
     used[j0] = true;
     int i0 = pb[j0];
     T delta = inf;
     int j1 = -1;
     for (int j = 0; j < m; j++) {
       if (!used[i]) {
         T cur = a[i0][j] - u[i0] - v[j];
         if (cur < minv[j]) {
           minv[j] = cur;
           way[j] = j0;
         if (minv[j] < delta) {</pre>
           delta = minv[j];
           j1 = j;
     for (int j = 0; j \le m; j++) {
       if (used[j]) {
         u[pb[j]] += delta;
         v[j] -= delta;
       } else {
         minv[j] -= delta;
     i0 = j1;
   } while (pb[j0] != -1);
     int j1 = way[j0];
     pb[j0] = pb[j1];
     pa[pb[j0]] = j0;
     j0 = j1;
   } while (j0 != m);
```

inline T current\_score() {

```
return -v[m];
  inline T solve() {
    for (int i = 0; i < n; i++) {</pre>
      add_row(i);
    return current_score();
};
```

#### GlobalMinCut.h

Description: Find a global minimum cut in an undirected graph, as represented by an adjacency matrix.

```
Time: \mathcal{O}\left(V^3\right)
                                                        d41d8c, 21 lines
pair<int, vi> globalMinCut(vector<vi> mat) {
 pair<int, vi> best = {INT_MAX, {}};
 int n = sz(mat);
 vector<vi> co(n);
 rep(i, 0, n) co[i] = {i};
 rep(ph,1,n) +
    vi w = mat[0];
    size_t s = 0, t = 0;
    rep(it,0,n-ph) { // O(V^2) \rightarrow O(E \log V) with prio. queue
      w[t] = INT_MIN;
      s = t, t = max_element(all(w)) - w.begin();
      rep(i, 0, n) w[i] += mat[t][i];
    best = min(best, \{w[t] - mat[t][t], co[t]\});
    co[s].insert(co[s].end(), all(co[t]));
    rep(i, 0, n) mat[s][i] += mat[t][i];
    rep(i, 0, n) mat[i][s] = mat[s][i];
    mat[0][t] = INT_MIN;
 return best;
```

## Matching

#### hopcroftKarp.h

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

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

```
bool dfs(int a, int L, vector<vi>& q, vi& btoa, vi& A, vi& B) {
  if (A[a] != L) return 0;
  A[a] = -1;
  for (int b : g[a]) if (B[b] == L + 1) {
    if (btoa[b] == -1 || dfs(btoa[b], L + 1, g, btoa, A, B))
      return btoa[b] = a, 1;
  return 0;
int hopcroftKarp(vector<vi>& q, vi& btoa) {
  int res = 0;
  vi A(g.size()), B(btoa.size()), cur, next;
  for (;;) {
    fill(all(A), 0);
    fill(all(B), 0);
    cur.clear();
    for (int a : btoa) if(a != -1) A[a] = -1;
    rep(a, 0, sz(q)) if(A[a] == 0) cur.push_back(a);
    for (int lay = 1;; lay++) {
```

## WeightedMatching Blossom

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

## WeightedMatching.h

**Description:** Given a weighted bipartite graph, matches every node on the left with a node on the right such that no nodes are in two matchings and the sum of the edge weights is minimal. Takes cost[N][M], where cost[i][j] = cost for L[i] to be matched with R[j] and returns (min cost, match), where L[i] is matched with R[match[i]]. Negate costs for max cost. Requires  $N \leq M$ . **Time:**  $\mathcal{O}(N^2M)$ 

```
d41d8c, 31 lines
pair<int, vi> hungarian(const vector<vi> &a) {
 if (a.empty()) return {0, {}};
  int n = sz(a) + 1, m = sz(a[0]) + 1;
  vi u(n), v(m), p(m), ans(n-1);
  rep(i,1,n) {
   p[0] = i;
    int j0 = 0; // add "dummy" worker 0
    vi dist(m, INT_MAX), pre(m, -1);
    vector<bool> done(m + 1);
    do { // dijkstra
     done[i0] = true;
      int i0 = p[j0], j1, delta = INT_MAX;
      rep(j,1,m) if (!done[j]) {
       auto cur = a[i0 - 1][j - 1] - u[i0] - v[j];
       if (cur < dist[j]) dist[j] = cur, pre[j] = j0;</pre>
       if (dist[j] < delta) delta = dist[j], j1 = j;</pre>
      rep(j,0,m) {
       if (done[j]) u[p[j]] += delta, v[j] -= delta;
        else dist[j] -= delta;
      i0 = i1;
    } while (p[j0]);
    while (j0) { // update alternating path
     int j1 = pre[j0];
     p[j0] = p[j1], j0 = j1;
 rep(j,1,m) if (p[j]) ans[p[j] - 1] = j - 1;
 return {-v[0], ans}; // min cost
```

## Blossom.h

**Description:** Given a weighted graph, finds max matching **Time:**  $\mathcal{O}\left(N^3\right)$  d41d8c, 139 lines

```
struct blossom {
    int n, m;
    vector<int> mate;
```

```
vector<vector<int>> b;
vector<int> p, d, bl;
vector<vector<int>> q;
blossom(int n) : n(n) {
   m = n + n / 2;
   mate.assign(n, -1);
   b.resize(m);
   p.resize(m);
    d.resize(m);
   bl.resize(m);
    g.assign(m, vector<int>(m, -1));
void add_edge(int u, int v) {
    g[u][v] = u;
    q[v][u] = v;
void match(int u, int v) {
    q[u][v] = q[v][u] = -1;
   mate[u] = v;
    mate[v] = u;
vector<int> trace(int x) {
    vector<int> vx;
    while(true) {
        while (bl[x] != x) x = bl[x];
        if(!vx.empty() && vx.back() == x) break;
        vx.push_back(x);
        x = p[x];
    return vx;
void contract(int c, int x, int y, vector<int> &vx, vector<</pre>
    int> &vv) {
    b[c].clear();
    int r = vx.back();
    while(!vx.empty() && !vy.empty() && vx.back() == vy.
        back()) {
        r = vx.back();
        vx.pop back();
        vy.pop_back();
   b[c].push_back(r);
   b[c].insert(b[c].end(), vx.rbegin(), vx.rend());
    b[c].insert(b[c].end(), vy.begin(), vy.end());
    for(int i = 0; i <= c; i++) {</pre>
        q[c][i] = q[i][c] = -1;
    for(int z : b[c]) {
        bl[z] = c;
        for(int i = 0; i < c; i++) {</pre>
            if (q[z][i] != -1) {
                g[c][i] = z;
                g[i][c] = g[i][z];
vector<int> lift(vector<int> &vx) {
    vector<int> A;
    while(vx.size() >= 2) {
        int z = vx.back(); vx.pop_back();
        if(z < n)  {
            A.push_back(z);
            continue;
```

```
int w = vx.back();
        int i = (A.size() % 2 == 0 ? find(b[z].begin(), b[z])
             ].end(), g[z][w]) - b[z].begin() : 0);
        int j = (A.size() % 2 == 1 ? find(b[z].begin(), b[z])
            ].end(), q[z][A.back()]) - b[z].begin() : 0);
        int k = b[z].size();
        int dif = (A.size() % 2 == 0 ? i % 2 == 1 : j % 2
            == 0) ? 1 : k - 1;
        while (i != j) {
            vx.push_back(b[z][i]);
            i = (i + dif) % k;
        vx.push_back(b[z][i]);
    return A;
int solve() {
    for(int ans = 0;; ans++) {
        fill(d.begin(), d.end(), 0);
        queue<int> Q;
        for(int i = 0; i < m; i++) bl[i] = i;</pre>
        for(int i = 0; i < n; i++) {</pre>
            if (mate[i] == -1) {
                Q.push(i);
                p[i] = i;
                d[i] = 1;
        int c = n;
        bool aug = false;
        while(!Q.empty() && !aug) {
            int x = Q.front(); Q.pop();
            if(bl[x] != x) continue;
            for(int y = 0; y < c; y++) {
                if(bl[y] == y \&\& g[x][y] != -1) {
                    if(d[y] == 0) {
                        p[y] = x;
                        d[v] = 2;
                        p[mate[y]] = y;
                        d[mate[y]] = 1;
                        O.push(mate[v]);
                    }else if(d[y] == 1) {
                        vector<int> vx = trace(x);
                        vector<int> vy = trace(y);
                        if(vx.back() == vy.back()) {
                            contract(c, x, y, vx, vy);
                            Q.push(c);
                            p[c] = p[b[c][0]];
                            d[c] = 1;
                            c++:
                        }else {
                            aug = true;
                            vx.insert(vx.begin(), y);
                            vy.insert(vy.begin(), x);
                            vector<int> A = lift(vx);
                            vector<int> B = lift(vy);
                            A.insert(A.end(), B.rbegin(), B
                                 .rend());
                            for (int i = 0; i < (int) A.size
                                 (); i += 2) {
                                match(A[i], A[i + 1]);
                                if(i + 2 < (int) A.size())
                                     add_edge(A[i + 1], A[i
                                      + 2]);
                        break;
```

```
if(!aug) return ans;
};
```

## 6.3 DFS algorithms

#### SCC.h

**Description:** Finds strongly connected components in a directed graph.

```
Time: \mathcal{O}(E+V)
template <typename T> vector<int> find_scc(const digraph<T> &g,
     int &cnt) {
  digraph<T> g_rev = g.reverse();
  vector<int> order;
  vector<bool> was(g.n, false);
  function < void (int) > dfs1 = [&] (int v) {
   was[v] = true;
    for (int id : q.q[v]) {
     if (q.ignore != nullptr && q.ignore(id)) {
        continue;
      auto &e = g.edges[id];
      int to = e.to;
     if (!was[to]) {
        dfs1(to);
   order.push back(v);
  for (int i = 0; i < g.n; i++) {
   if (!was[i]) {
     dfs1(i);
  vector<int> c(g.n, -1);
  function < void (int) > dfs2 = [&] (int v) {
    for (int id : q_rev.q[v]) {
      if (g_rev.ignore != nullptr && g_rev.ignore(id)) {
        continue;
      auto &e = g_rev.edges[id];
      int to = e.to;
      if (c[to] == -1) {
       c[to] = c[v];
        dfs2(to);
  };
  for (int id = g.n - 1; id >= 0; id--) {
   int i = order[id];
   if (c[i] != −1) {
     continue;
   c[i] = cnt++;
   dfs2(i);
  return c;
 // c[i] \le c[j] for every edge i \rightarrow j
```

Description: Finds all biconnected components in an undirected graph Time:  $\mathcal{O}(E+V)$ 

```
d41d8c, 74 lines
```

```
template <typename T> vector<int> find_edge_biconnected(
    dfs_undigraph<T> &g, int &cnt) {
 q.dfs all();
 vector<int> vertex_comp(g.n);
 cnt = 0;
 for (int i : g.order) {
   if (g.pv[i] == -1 || g.min_depth[i] == g.depth[i]) {
     vertex_comp[i] = cnt++;
      vertex_comp[i] = vertex_comp[q.pv[i]];
 return vertex_comp;
template <typename T> vector<int> find_vertex_biconnected(
    dfs_undigraph<T> &g, int &cnt) {
 g.dfs_all();
 vector<int> vertex_comp(g.n);
 cnt = 0;
 for (int i : g.order) {
   if (g.pv[i] == -1) {
     vertex comp[i] = -1;
      continue;
    if (q.min_depth[i] >= q.depth[q.pv[i]]) {
     vertex_comp[i] = cnt++;
   } else {
     vertex_comp[i] = vertex_comp[q.pv[i]];
 vector<int> edge_comp(g.edges.size(), -1);
 for (int id = 0; id < (int)g.edges.size(); id++) {</pre>
   if (g.ignore != nullptr && g.ignore(id)) {
     continue;
   int x = q.edges[id].from;
   int y = g.edges[id].to;
   int z = (q.depth[x] > q.depth[y] ? x : y);
    edge_comp[id] = vertex_comp[z];
 return edge_comp;
template <typename T> vector<bool> find_bridges(dfs_undigraph<T</pre>
    > &q) {
 q.dfs all();
 vector<bool> bridge(g.edges.size(), false);
 for (int i = 0; i < q.n; i++) {</pre>
   if (g.pv[i] != -1 && g.min_depth[i] == g.depth[i]) {
     bridge[g.pe[i]] = true;
 return bridge;
template <typename T> vector<bool> find_cutpoints(dfs_undigraph)
    <T> &q) {
 g.dfs_all();
 vector<bool> cutpoint(g.n, false);
 for (int i = 0; i < g.n; i++) {</pre>
   if (q.pv[i] != -1 && q.min_depth[i] >= q.depth[q.pv[i]]) {
     cutpoint[g.pv[i]] = true;
 vector<int> children(g.n, 0);
 for (int i = 0; i < q.n; i++) {
   if (g.pv[i] != -1) {
     children[g.pv[i]]++;
```

```
for (int i = 0; i < q.n; i++) {
  if (g.pv[i] == -1 && children[i] < 2) {</pre>
    cutpoint[i] = false;
return cutpoint;
```

#### 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 ( $\sim x$ ).

16

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

```
namespace kirkon {
struct twosat {
 digraph<int> q;
 int n;
 twosat(int _n) : q(digraph<int>(_n << 1)), n(_n) {}
  inline void add(int x, int value_x) {
   // (v/x) == value_x
   assert (0 \leq x && x \leq n);
   assert (0 <= value x && value x <= 1);
   g.add((x << 1) + (value_x ^ 1), (x << 1) + value_x);
 inline void add(int x, int value_x, int y, int value_y) {
   // (v/x) = value_x \mid \mid v/y \mid = value_y)
   assert(0 <= x && x < n && 0 <= y && y < n);
   assert(0 <= value_x && value_x <= 1 && 0 <= value_y &&
        value_y <= 1);</pre>
   g.add((x << 1) + (value_x ^ 1), (y << 1) + value_y);
   g.add((y << 1) + (value_y ^ 1), (x << 1) + value_x);
 inline void add_impl(int x, int value_x, int y, int value_y)
    // (v[x] = value_x \rightarrow v[y] = value_y)
   assert (0 <= x \&\& x < n \&\& 0 <= y \&\& y < n);
   assert(0 <= value_x && value_x <= 1 && 0 <= value_y &&
         value v \le 1;
    g.add((x << 1) + (value_x ^ 1), (y << 1) + (value_y ^ 1));
    //have to add reverse edge in order for algorithm prove to
    q.add((y << 1) + value_y, (x << 1) + value_x);
 inline void add_xor(int x, int y, int value) {
```

 $// (v[x] = value_x \rightarrow v[y] = value_y)$ 

assert (0 <= value && value <= 1);

if (value) { add(x, 1, y, 1);

} else {

add(x, 0, y, 0);

 $add_{impl}(x, 1, y, 1);$ 

 $add_impl(x, 0, y, 0);$ 

assert (0 <= x && x < n && 0 <= y && y < n);

```
}

inline vector<int> solve() {
   int cnt;
   vector<int> c = find_scc(g, cnt);
   vector<int> res(n);
   for (int i = 0; i < n; i++) {
      if (c[i << 1] == c[i << 1 ^ 1]) {
        return vector<int>();
    }
    res[i] = (c[i << 1] < c[i << 1 ^ 1]);
   return res;
}

// namespace kirkon

inline

// namespace kirkon

// namespace kirkon

inline

// namespace kirkon

// namespace kirkon
```

#### 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:**  $\mathcal{O}(V+E)$ 

```
inthe: O(v + E)

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

      for (int x : D) if (x < 0 || sz(ret) != nedges+1) return {};
      return {ret.rbegin(), ret.rend()};
    }
}</pre>
```

#### DirectedMST.h

**Description:** Finds a minimum spanning tree/arborescence of a directed graph, given a root node. If no MST exists, returns -1. **Time:**  $\mathcal{O}\left(E\log V\right)$ 

```
d41d8c, 60 lines
"../data-structures/UnionFindRollback.h"
struct Edge { int a, b; ll w; };
struct Node {
  Edge key;
  Node *1, *r;
  11 delta;
  void prop() {
    key.w += delta;
    if (1) 1->delta += delta;
    if (r) r->delta += delta;
    delta = 0;
  Edge top() { prop(); return key; }
Node *merge(Node *a, Node *b) {
  if (!a || !b) return a ?: b;
  a->prop(), b->prop();
  if (a->key.w > b->key.w) swap(a, b);
  swap(a->1, (a->r = merge(b, a->r)));
  return a;
void pop (Node \star \& a) { a->prop(); a = merge(a->1, a->r); }
```

```
pair<11, vi> dmst(int n, int r, vector<Edge>& g) {
 RollbackUF uf(n):
 vector<Node*> heap(n);
 for (Edge e : g) heap[e.b] = merge(heap[e.b], new Node{e});
 11 \text{ res} = 0;
 vi seen(n, -1), path(n), par(n);
 seen[r] = r;
 vector<Edge> Q(n), in(n, \{-1,-1\}), comp;
 deque<tuple<int, int, vector<Edge>>> cycs;
 rep(s,0,n) {
   int u = s, qi = 0, w;
   while (seen[u] < 0) {
     if (!heap[u]) return {-1,{}};
     Edge e = heap[u]->top();
     heap[u]->delta -= e.w, pop(heap[u]);
     Q[qi] = e, path[qi++] = u, seen[u] = s;
     res += e.w, u = uf.find(e.a);
     if (seen[u] == s) {
       Node \star cyc = 0;
       int end = qi, time = uf.time();
       do cyc = merge(cyc, heap[w = path[--qi]]);
       while (uf.join(u, w));
       u = uf.find(u), heap[u] = cyc, seen[u] = -1;
       cycs.push_front({u, time, {&Q[qi], &Q[end]}});
    rep(i, 0, qi) in[uf.find(Q[i].b)] = Q[i];
 for (auto& [u,t,comp] : cycs) { // restore sol (optional)
   uf.rollback(t);
   Edge inEdge = in[u];
    for (auto& e : comp) in[uf.find(e.b)] = e;
   in[uf.find(inEdge.b)] = inEdge;
 rep(i,0,n) par[i] = in[i].a;
 return {res, par};
```

## 6.4 Coloring

#### EdgeColoring.h

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

```
d41d8c, 31 lines
vi edgeColoring(int N, vector<pii> eds) {
 vi cc(N + 1), ret(sz(eds)), fan(N), free(N), loc;
 for (pii e : eds) ++cc[e.first], ++cc[e.second];
 int u, v, ncols = *max_element(all(cc)) + 1;
 vector<vi> adj(N, vi(ncols, -1));
 for (pii e : eds) {
   tie(u, v) = e;
   fan[0] = v;
   loc.assign(ncols, 0);
   int at = u, end = u, d, c = free[u], ind = 0, i = 0;
   while (d = free[v], !loc[d] && (v = adj[u][d]) != -1)
     loc[d] = ++ind, cc[ind] = d, fan[ind] = v;
    cc[loc[d]] = c;
   for (int cd = d; at != -1; cd ^= c ^ d, at = adj[at][cd])
     swap(adj[at][cd], adj[end = at][cd ^ c ^ d]);
   while (adj[fan[i]][d] != -1) {
     int left = fan[i], right = fan[++i], e = cc[i];
     adj[u][e] = left;
     adj[left][e] = u;
     adj[right][e] = -1;
     free[right] = e;
```

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

## 6.5 Trees

## BinaryLifting.h

**Description:** Calculate power of two jumps in a tree, to support fast upward jumps and LCAs. Assumes the root node points to itself. **Time:** construction  $\mathcal{O}(N \log N)$ , queries  $\mathcal{O}(\log N)$ 

```
d41d8c, 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];
```

#### LCA.h

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

```
Time: \mathcal{O}(N \log N + Q)
"../data-structures/RMQ.h"
                                                       d41d8c, 21 lines
struct LCA {
 int T = 0;
 vi time, path, ret;
 RMQ<int> rmq;
  LCA(vector < vi > \& C) : time(sz(C)), rmq((dfs(C,0,-1), ret)) {}
  void dfs(vector<vi>& C, int v, int par) {
    time[v] = T++;
    for (int y : C[v]) if (y != par) {
      path.push_back(v), ret.push_back(time[v]);
      dfs(C, y, v);
  int lca(int a, int b) {
    if (a == b) return a;
    tie(a, b) = minmax(time[a], time[b]);
    return path[rmq.query(a, b)];
  //dist(a,b){return depth[a] + depth[b] - 2*depth[lca(a,b)];}
```

## CompressTree DimaHLD Point

```
CompressTree.h
```

};

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

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

```
d41d8c, 21 lines
typedef vector<pair<int, int>> vpi;
vpi compressTree(LCA& lca, const vi& subset) {
  static vi rev; rev.resize(sz(lca.time));
  vi li = subset, &T = lca.time;
  auto cmp = [&](int a, int b) { return T[a] < T[b]; };</pre>
  sort(all(li), cmp);
  int m = sz(li)-1;
  rep(i,0,m) {
   int a = li[i], b = li[i+1];
   li.push_back(lca.lca(a, b));
  sort(all(li), cmp);
  li.erase(unique(all(li)), li.end());
  rep(i,0,sz(li)) rev[li[i]] = i;
  vpi ret = {pii(0, li[0])};
  rep(i, 0, sz(li)-1) {
   int a = li[i], b = li[i+1];
    ret.emplace_back(rev[lca.lca(a, b)], b);
  return ret;
```

## DimaHLD.h

Description: Builds HLD

```
Time: \mathcal{O}(log^2 forguery) probably \mathcal{O}(log)
                                                      d41d8c, 132 lines
struct segtree { }
struct HLD {
    int n:
    vector<int> par, val, vale, tin, tout, sz, top;
    vector<vector<int>> up;
    segtree tree_v, tree_e;
   HLD() { }
    HLD(vector<vector<int>> &G) {
       n = G.size();
        par.resize(n);
        val.resize(n, 1);
        vale.resize(n, 1);
        tin.resize(n);
        tout.resize(n);
        sz.resize(n);
        top.resize(n);
        dfs(0, G);
        dfs2(0, G);
        tree_v = segtree(n);
        tree_e = segtree(n);
        LG = 0;
        while (1 << LG < n)
            LG++;
        up.resize(LG + 1, vector<int>(n));
        up[0] = par;
        for (int i = 1; i <= LG; i++) {
            for (int j = 0; j < n; j++)
                up[i][j] = up[i - 1][up[i - 1][j]];
```

```
void dfs(int v, vector<vector<int>> &G, int p = -1) {
    if (p ! = -1)
        G[v].erase(find(G[v].begin(), G[v].end(), p));
    par[v] = p != -1 ? p : v;
    sz[v] = 1;
    for (int to : G[v]) {
        dfs(to, G, v);
        sz[v] += sz[to];
    sort(G[v].begin(), G[v].end(), [&](int x, int y) {
        return sz[x] > sz[y];
    });
void dfs2(int v, vector<vector<int>> &G) {
    static int timer = 0;
    tin[v] = timer++;
   if (!G[v].empty())
        top[G[v][0]] = top[v];
    for (int i = 0; i < G[v].size(); i++) {</pre>
        int u = G[v][i];
        if (i)
            top[u] = u;
        dfs2(u, G);
    tout[v] = timer;
bool isPar(int v, int u) {
    return tin[v] <= tin[u] && tout[v] >= tout[u];
int lca(int v, int u) {
    if (isPar(v, u))
        return v;
    for (int i = LG; i >= 0; i--) {
        if (!isPar(up[i][v], u))
            v = up[i][v];
    return up[0][v];
vector<pair<int, int>> get_e_vert_path(int v, int 1) {
    vector<pair<int, int>> res;
    res.reserve(LG \star 2 + 3);
    while (true) {
        int x = top[v];
        if (isPar(x, 1))
            break;
        res.push_back(\{tin[x], tin[v] + 1\});
        v = par[x];
    res.push_back(\{tin[1] + 1, tin[v] + 1\});
    return res;
vector<pair<int, int>> get_e_path(int v, int u) {
    int 1 = lca(v, u);
    vector<pair<int, int>> res = get_e_vert_path(v, 1);
    for (auto elem : get_e_vert_path(u, 1))
        res.push_back(elem);
    return res:
vector<pair<int, int>> get_v_path(int v, int u) {
    auto res = get_e_path(v, u);
    int l = tin[lca(v, u)];
```

```
res.push back(\{1, 1 + 1\});
        return res:
    void mul_v(int v, int u, int delta) {
        for (auto [1, r] : get_v_path(v, u)) {
            tree_v.update(l, r, delta);
    void mul_e(int v, int u, int delta) {
        for (auto [l, r] : get_e_path(v, u))
            tree_e.update(l, r, delta);
    int get_v_sum(int v, int u) {
        11 \text{ res} = 0;
        for (auto [1, r] : get_v_path(v, u))
            res += tree_v.get_sum(1, r);
        return res % mod;
    int get_e_sum(int v, int u) {
        11 \text{ res} = 0;
        for (auto [1, r] : get_e_path(v, u))
            res += tree_e.get_sum(1, r);
        return res % mod;
};
```

#### 6.6 Math

## 6.6.1 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 *i*th row and column and take the determinant; this yields the number of directed spanning trees rooted at i (if G is undirected, remove any row/column).

#### 6.6.2 Erdős–Gallai theorem

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

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

# Geometry (7)

# 7.1 Geometric primitives

#### Point.h

**Description:** Class to handle points in the plane. T can be e.g. double or long long. (Avoid int.) d41d8c, 28 lines

```
template <class T> int sqn(T x) \{ return (x > 0) - (x < 0); \}
template<class T>
struct Point {
 typedef Point P;
  explicit Point (T x=0, T y=0) : x(x), y(y) {}
 bool operator<(P p) const { return tie(x,y) < tie(p.x,p.y); }</pre>
 bool operator==(P p) const { return tie(x,y)==tie(p.x,p.y); }
```

d41d8c, 35 lines

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

#### lineDistance.h

#### Description:

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



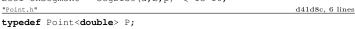
d41d8c, 4 lines template<class P> double lineDist(const P& a, const P& b, const P& p) { return (double) (b-a).cross(p-a)/(b-a).dist();

## SegDis.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;
```



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

## SegInter.h

#### Description:

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



d41d8c, 13 lines

```
Usage: vector<P> inter = seqInter(s1,e1,s2,e2);
if (sz(inter) == 1)
cout << "segments intersect at " << inter[0] << endl;</pre>
"Point.h", "OnSegment.h"
```

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

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

#### lineInt.h

#### Description:

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

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

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

#### sideOf.h

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

```
Usage: bool left = sideOf(p1,p2,q)==1;
"Point.h"
                                                         d41d8c, 9 lines
template<class P>
```

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

## OnSegment.h

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

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

## linTransf.h Description:

typedef Point < double > P;

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



"Point.h"

```
d41d8c, 6 lines
```

```
19
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 =  $\{w[0], w[0].t360() ...\}; // sorted$ 

// sweeps j such that (j-i) represents the number of positively

int j = 0; rep(i,0,n) { while (v[j] < v[i].t180()) ++j; }

oriented triangles with vertices at 0 and i

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

## 7.2 Circles

## CircInter.h

**Description:** Computes the pair of points at which two circles intersect. Returns false in case of no intersection. "Point.h" d41d8c, 11 lines

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

## CircleTangents.h

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

"Point.h" d41d8c, 13 lines

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

## CircPolvInter.h

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

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

```
"../../content/geometry/Point.h"
                                                         d41d8c, 19 lines
typedef Point < double > P;
#define arg(p, q) atan2(p.cross(q), p.dot(q))
double circlePoly(P c, double r, vector<P> ps) {
  auto tri = [&] (P p, P q) {
```

```
auto r2 = r * r / 2;
 P d = q - p;
 auto a = d.dot(p)/d.dist2(), b = (p.dist2()-r*r)/d.dist2();
 auto det = a * a - b;
 if (det <= 0) return arg(p, q) * r2;</pre>
 auto s = max(0., -a-sqrt(det)), t = min(1., -a+sqrt(det));
 if (t < 0 || 1 <= s) return arg(p, g) * r2;</pre>
 Pu = p + d * s, v = p + d * t;
 return arg(p, u) * r2 + u.cross(v)/2 + arg(v, q) * r2;
auto sum = 0.0;
rep(i, 0, sz(ps))
 sum += tri(ps[i] - c, ps[(i + 1) % sz(ps)] - c);
return sum;
```

## circumcircle.h

## Description:

"Point.h"

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



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

#### MinEnclosCirc.h

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

```
"circumcircle.h"
                                                      d41d8c, 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};
```

## 7.3 Polygons

## InsidePolygon.h

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

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

"Point.h", "OnSegment.h", "SegmentDistance.h" d41d8c, 11 lines

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

#### Area.h

Description: Returns twice the signed area of a polygon. Clockwise enumeration gives negative area. Watch out for overflow if using int as T! d41d8c, 6 lines

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

## PolyCent.h

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

```
"Point h"
                                                       d41d8c, 9 lines
typedef Point < double > P;
P polygonCenter(const vector<P>& v) {
 P res(0, 0); double A = 0;
 for (int i = 0, j = sz(v) - 1; i < sz(v); j = i++) {
   res = res + (v[i] + v[j]) * v[j].cross(v[i]);
   A += v[j].cross(v[i]);
 return res / A / 3;
```

#### Cut.h

#### Description:

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

```
Returns a vector with the vertices of a polygon with every-
Usage: vector<P> p = ...;
p = polygonCut(p, P(0,0), P(1,0));
```



```
"Point.h", "lineIntersection.h"
                                                       d41d8c, 13 lines
typedef Point<double> P;
vector<P> polygonCut (const vector<P>& poly, P s, P e) {
 vector<P> res;
  rep(i, 0, sz(poly)) {
    P cur = poly[i], prev = i ? poly[i-1] : poly.back();
    bool side = s.cross(e, cur) < 0;</pre>
    if (side != (s.cross(e, prev) < 0))
      res.push_back(lineInter(s, e, cur, prev).second);
    if (side)
      res.push_back(cur);
 return res;
```

#### Hull.h

## Description:

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



d41d8c, 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<P> h(sz(pts)+1);
  int s = 0, t = 0;
  for (int it = 2; it--; s = --t, reverse(all(pts)))
    for (P p : pts) {
      while (t \ge s + 2 \&\& h[t-2].cross(h[t-1], p) \le 0) t--;
      h[t++] = p;
  return {h.begin(), h.begin() + t - (t == 2 \&\& h[0] == h[1])};
```

#### HalfplaneInt.h

**Description:** find halfplane intersection

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

d41d8c, 73 lines

```
struct Halfplane {
    Point p, pq;
    long double angle;
    Halfplane() {}
    Halfplane(const Point& a, const Point& b) : p(a), pq(b - a)
        angle = atan21(pq.y, pq.x);
    bool out (const Point& r) {
        return cross(pq, r - p) < -eps;
    bool operator < (const Halfplane& e) const {</pre>
        return angle < e.angle;</pre>
    friend Point inter(const Halfplane& s, const Halfplane& t)
        long double alpha = cross((t.p - s.p), t.pq) / cross(s.
             pq, t.pq);
        return s.p + (s.pq * alpha);
};
```

d41d8c, 17 lines

```
vector<Point> hp_intersect(vector<Halfplane>& H) {
    Point box[4] = {
            Point(inf, inf),
            Point (-inf, inf),
            Point (-inf, -inf),
            Point(inf, -inf)
    };
    for(int i = 0; i<4; i++) {
        Halfplane aux(box[i], box[(i+1) % 4]);
        H.push_back(aux);
    sort(H.begin(), H.end());
   deque<Halfplane> dq;
    int len = 0;
    for(int i = 0; i < (int)(H.size()); i++) {</pre>
        while (len > 1 && H[i].out(inter(dq[len-1], dq[len-2]))
            dq.pop_back();
            --len;
        while (len > 1 && H[i].out(inter(dq[0], dq[1]))) {
            dq.pop_front();
            --len;
        if (len > 0 && fabsl(cross(H[i].pq, dg[len-1].pq)) <</pre>
            if (dot(H[i].pq, dq[len-1].pq) < 0.0)</pre>
                return vector<Point>();
            if (H[i].out(dq[len-1].p)) {
                dq.pop_back();
                --len;
            else continue;
        dq.push_back(H[i]);
    while (len > 2 && dq[0].out(inter(dq[len-1], dq[len-2])))
        dq.pop_back();
        --len:
    while (len > 2 && dq[len-1].out(inter(dq[0], dq[1]))) {
        dq.pop_front();
        --len;
    if (len < 3) return vector<Point>();
    vector<Point> ret(len);
    for(int i = 0; i+1 < len; i++) {</pre>
        ret[i] = inter(dq[i], dq[i+1]);
    ret.back() = inter(dq[len-1], dq[0]);
    return ret;
Diam.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;
```

#### PInsideH.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"
                                                         d41d8c, 14 lines
typedef Point<11> P;
bool inHull(const vector<P>& 1, P p, bool strict = true) {
  int a = 1, b = sz(1) - 1, r = !strict;
  if (sz(1) < 3) return r && onSegment(1[0], 1.back(), p);</pre>
  if (sideOf(1[0], 1[a], 1[b]) > 0) swap(a, b);
  if (sideOf(1[0], 1[a], p) \geq r || sideOf(1[0], 1[b], p) \leq -r)
  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>
```

#### LineHullIntersection.h

swap (endA, endB);

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

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

d41d8c, 12 lines

```
"Point.h"
                                                     d41d8c, 39 lines
#define cmp(i,j) sgn(dir.perp().cross(poly[(i)%n]-poly[(j)%n]))
#define extr(i) cmp(i + 1, i) >= 0 && cmp(i, i - 1 + n) < 0
template <class P> int extrVertex(vector<P>& poly, P dir) {
 int n = sz(poly), lo = 0, hi = n;
 if (extr(0)) return 0;
  while (lo + 1 < hi) {
   int m = (1o + hi) / 2;
    if (extr(m)) return m;
    int 1s = cmp(1o + 1, 1o), ms = cmp(m + 1, m);
    (ls < ms \mid | (ls == ms \&\& ls == cmp(lo, m)) ? hi : lo) = m;
 return lo:
#define cmpL(i) sgn(a.cross(poly[i], b))
template <class P>
array<int, 2> lineHull(P a, P b, vector<P>& poly) {
 int endA = extrVertex(poly, (a - b).perp());
  int endB = extrVertex(poly, (b - a).perp());
  if (cmpL(endA) < 0 \mid | cmpL(endB) > 0)
    return {-1, -1};
  array<int, 2> res;
  rep(i, 0, 2) {
    int lo = endB, hi = endA, n = sz(poly);
    while ((lo + 1) % n != hi) {
     int m = ((lo + hi + (lo < hi ? 0 : n)) / 2) % n;</pre>
      (cmpL(m) == cmpL(endB) ? lo : hi) = m;
    res[i] = (lo + !cmpL(hi)) % n;
```

```
if (res[0] == res[1]) return {res[0], -1};
if (!cmpL(res[0]) && !cmpL(res[1]))
  switch ((res[0] - res[1] + sz(poly) + 1) % sz(poly)) {
    case 0: return {res[0], res[0]};
    case 2: return {res[1], res[1]};
return res;
```

## 7.4 Misc. Point Set Problems

ClosestPair.h

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

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

```
typedef Point<11> P;
pair<P, P> closest (vector<P> v) {
 assert(sz(v) > 1);
  set<P> S;
  sort(all(v), [](P a, P b) { return a.y < b.y; });</pre>
  pair<11, pair<P, P>> ret{LLONG_MAX, {P(), P()}};
  int j = 0;
  for (P p : v) {
   P d{1 + (ll) sqrt (ret.first), 0};
    while (v[j].y \le p.y - d.x) S.erase(v[j++]);
    auto lo = S.lower_bound(p - d), hi = S.upper_bound(p + d);
    for (; lo != hi; ++lo)
      ret = min(ret, \{(*lo - p).dist2(), \{*lo, p\}\});
    S.insert(p);
  return ret.second:
```

# Strings (8)

#### KMP.h

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

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

```
d41d8c, 16 lines
vi pi(const string& s) {
 vi p(sz(s));
 rep(i,1,sz(s)) {
   int g = p[i-1];
    while (g \&\& s[i] != s[g]) g = p[g-1];
   p[i] = 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.h

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

```
Time: \mathcal{O}(n)
                                                                    d41d8c, 12 lines
vi Z(const string& S) {
```

```
vi z(sz(S));
int 1 = -1, r = -1;
rep(i,1,sz(S)) {
```

## Manacher MinRotation SuffixArray

```
z[i] = i >= r ? 0 : min(r - i, z[i - 1]);
while (i + z[i] < sz(S) && S[i + z[i]] == S[z[i]])
    z[i]++;
if (i + z[i] > r)
    1 = i, r = i + z[i];
}
return z;
```

## Manacher.h

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

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

d41d8c, 13 lines

```
array<vi, 2> manacher(const string& s) {
  int n = sz(s);
  array<vi,2> p = {vi(n+1), vi(n)};
  rep(z,0,2) for (int i=0,1=0,r=0; i < n; i++) {
    int t = r-i+!z;
    if (i<r) p[z][i] = min(t, p[z][1+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.beqin(), v.beqin()+minRotation(v), v.end());

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

```
int minRotation(string s) {
   int a=0, N=sz(s); s += s;
   rep(b,0,N) rep(k,0,N) {
    if (a+k == b | | 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:**  $\mathcal{O}(n \log n)$ 

```
int getMin(int 1, int r) {
        int k = 31 - __builtin_clz(r - 1);
        return min(st[k][1], st[k][r - (1 << k)]);</pre>
};
struct SuffixArrav {
    int n:
    vector<int> sa, lcp, pos;
    sparse st;
    vector<int> s;
    // O(Size(s) + max(s) - min(s))
    SuffixArray(vector<int> &s): n(Size(s)) {
        int mn = *min_element(all(s));
        for (int &i : s)
            i -= mn - 1;
        s.reserve(Size(s) + 1);
        s.push_back(0);
        sa = build(s, *max_element(all(s)) + 1);
        int n = Size(s);
        pos.resize(n);
        for (int i = 0; i < n; i++)</pre>
            pos[sa[i]] = i;
        lcp.resize(n);
        int k = 0;
        for (int i = 0; i < n - 1; i++) {
            int j = sa[pos[i] - 1];
            while (s[i + k] == s[j + k])
                k++;
            lcp[pos[i]] = k;
            k = max(0, k - 1);
        st = sparse(lcp);
        this -> s = s;
    vector<int> phase2(const vector<int> &s, const vector<int>
         &pref, const vector<char> &types, const vector<int> &
         lms) {
        int n = Size(s);
        vector<int> cnt = pref;
        vector<int> res(n, -1);
        for (int i : lms) {
            int a = s[i];
            res[--cnt[a + 1]] = i;
        copy(all(pref), cnt.begin());
        for (int p : res) {
            if (p <= 0 || types[p - 1] != 'L')</pre>
                continue;
            int a = s[p - 1];
            res[cnt[a]++] = p - 1;
        copy(all(pref), cnt.begin());
        for (int i = n - 1; i >= 0; i--) {
            int p = res[i];
            if (p <= 0 || types[p - 1] != 'S')</pre>
                continue;
            int a = s[p - 1];
            res[--cnt[a + 1]] = p - 1;
        return res:
    inline bool is_lms(const vector<char> &types, int i) {
        return types[i - 1] == 'L' && types[i] == 'S';
```

```
// compare two lms substring
inline bool not_equal(const vector<int> &s, const vector<</pre>
    char> &types, int i, int j) {
    assert(is_lms(types, i) && is_lms(types, j));
   bool is_lms1 = false, is_lms2 = false;
    while (true) {
        if (s[i] != s[j] || types[i] != types[j])
            return true;
        if (is_lms1 && is_lms2)
            break;
        j++;
        is_lms1 = is_lms(types, i);
        is_lms2 = is_lms(types, j);
   return false:
// m = max(s) + 1, s.back() == 0
vector<int> build(vector<int> &s, int m) {
   int n = Size(s);
   assert(!s.empty());
   assert(s.back() == 0);
    assert(Size(s) == 1 || *min_element(s.begin(), s.end()
        -1) > 0);
   assert(*max element(all(s)) == m - 1);
   if (Size(s) == 1)
        return {0};
   vector<char> types(n);
   tvpes[n - 1] = 'S';
   vector<int> lms;
   lms.reserve(n);
    for (int i = n - 2; i >= 0; i--) {
        if (s[i] < s[i + 1])
            types[i] = 'S';
        else if (s[i] > s[i + 1])
            types[i] = 'L';
            types[i] = types[i + 1];
        if (types[i] == 'L' && types[i + 1] == 'S')
            lms.push back(i + 1);
    vector<int> pref(m + 1);
    for (int i : s)
       pref[i + 1]++;
    for (int i = 0; i < m; i++)</pre>
        pref[i + 1] += pref[i];
    auto res = phase2(s, pref, types, lms);
   int lms_cnt = 1, color = 0;
   int last = n - 1;
   vector<int> new_sym(n, -1);
   new_sym[n - 1] = 0;
   for (int i = 1; i < n; i++) {</pre>
        int p = res[i];
        if (p <= 0 || !is_lms(types, p))</pre>
            continue;
        lms[lms\_cnt++] = p;
        color += not_equal(s, types, last, p);
       new_sym[p] = color;
        last = p;
   vector<int> new_string;
   vector<int> pos_new_string(n);
   new_string.reserve(Size(lms) + 1);
    for (int i = 0; i < n; i++) {</pre>
```

**if** (c !=-1) {

int c = new sym[i];

pos\_new\_string[Size(new\_string)] = i;

new\_string.push\_back(c);

## Hashing

```
if (color != Size(lms)) {
            auto sa_new = build(new_string, color + 1);
            for (int i = 1; i < Size(sa_new); i++)</pre>
                lms[i] = pos_new_string[sa_new[i]];
        return phase2(s, pref, types, lms);
    int get_lcp(int i, int j) {
        if (i == j)
            return n - i;
        i = pos[i];
        j = pos[j];
        if (i > j)
            swap(i, j);
        return st.getMin(i + 1, j + 1);
   bool compare (int i, int j) { // s[i..] < s[j..]
        if (i == j)
            return false;
        int k = get_lcp(i, j);
        return s[i + k] < s[j + k];
//Another impl
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 \times (all(s)+1), v(n), ws(max(n, lim)), rank(n);
    sa = lcp = v, iota(all(sa), 0);
    for (int j = 0, p = 0; p < n; j = max(1, j * 2), lim = p) {
     p = j, iota(all(y), n - j);
     rep(i,0,n) if (sa[i] >= j) y[p++] = sa[i] - j;
      fill(all(ws), 0);
      rep(i, 0, n) ws[x[i]]++;
      rep(i, 1, lim) ws[i] += ws[i - 1];
      for (int i = n; i--;) sa[--ws[x[y[i]]]] = y[i];
      swap(x, y), p = 1, x[sa[0]] = 0;
      rep(i,1,n) = sa[i-1], b = sa[i], x[b] =
        (y[a] == y[b] && y[a + j] == y[b + j]) ? p - 1 : p++;
    rep(i,1,n) rank[sa[i]] = i;
    for (int i = 0, j; i < n - 1; lcp[rank[i++]] = k)</pre>
      for (k \& \& k--, j = sa[rank[i] - 1];
          s[i + k] == s[j + k]; k++);
};
Hashing.h
Description: creates hashes
Time: \mathcal{O}(N)
                                                     d41d8c, 158 lines
namespace kirkon {
  const int HASH MOD = MOD;
  const int HASH_SIZE = 2;
  uniform_int_distribution<int> BDIST(0.1 * HASH_MOD, 0.9 *
       HASH_MOD);
```

```
struct custom hash {
  array<int, HASH_SIZE> vals{};
  custom hash() {
    vals.fill(0);
  custom hash(const array<int, HASH_SIZE> &other) {
    vals = other;
  custom_hash(array<int, HASH_SIZE> &&other) {
    vals = std::move(other);
  custom_hash &operator=(const array<int, HASH_SIZE> &other)
    vals = other:
    return *this;
  custom_hash &operator=(array<int, HASH_SIZE> &&other) {
    vals = std::move(other);
    return *this;
  int &operator[](int x) {
    return vals[x];
  bool operator==(const custom_hash &other) const {
    return vals == other.vals;
  bool operator!=(const custom_hash &other) const {
    return vals != other.vals;
  bool operator<(const custom_hash &other) const {</pre>
    return vals < other.vals;</pre>
  bool operator>(const custom_hash &other) const {
    return vals > other.vals;
  bool operator<=(const custom_hash &other) const {</pre>
    return vals <= other.vals;</pre>
  bool operator>=(const custom_hash &other) const {
    return vals >= other.vals;
};
template < class T >
custom hash make hash (T c) {
  auto res = custom hash{};
  res.vals.fill(c);
  return res;
custom hash base{};
vector<custom_hash> pows{};
custom_hash operator+(custom_hash 1, custom_hash r) {
  for (int i = 0; i < HASH_SIZE; ++i)</pre>
    if ((1[i] += r[i]) >= HASH MOD)
      1[i] -= HASH_MOD;
  return 1:
custom_hash operator-(custom_hash 1, custom_hash r) {
  for (int i = 0; i < HASH_SIZE; ++i)</pre>
    if ((1[i] -= r[i]) < 0)
```

```
l[i] += HASH_MOD;
    return 1;
 custom_hash operator*(custom_hash 1, custom_hash r) {
    for (int i = 0; i < HASH_SIZE; ++i)</pre>
     l[i] = (ll) l[i] * r[i] % HASH_MOD;
    return 1;
  void init() {
    static bool used = false;
    if (used) {
      return;
    for (auto &u: base.vals) {
      u = BDIST(rng);
    pows.emplace_back(make_hash(1));
    used = true;
  struct HashRange {
    str S:
    vector<custom_hash> cum{};
    HashRange() {
     init();
      cum.emplace_back();
    void add(char c) {
      S += c;
      cum.pb(base * cum.back() + make_hash(c));
    void add(str s) { each(c, s) add(c); }
    void extend(int len) {
      while (sz(pows) <= len) {
        pows.pb(base * pows.back());
    custom hash hash(int 1, int r) {
      int len = r + 1 - 1;
      extend(len);
      return cum[r + 1] - pows[len] * cum[l];
 };
}// namespace kirkon
struct custom int hash {
  static uint64_t splitmix64(uint64_t x) {
    // http://xorshift.di.unimi.it/splitmix64.c
    x += 0x9e3779b97f4a7c15;
    x = (x ^ (x >> 30)) * 0xbf58476d1ce4e5b9;
    x = (x ^ (x >> 27)) * 0x94d049bb133111eb;
    return x ^ (x >> 31);
  size_t operator()(uint64_t x) const {
    static const uint64_t FIXED_RANDOM = chrono::steady_clock::
        now().time_since_epoch().count();
    return splitmix64(x + FIXED RANDOM);
};
custom_int_hash int_hash{};
```

d41d8c, 19 lines

```
namespace std {
  template<>
  struct hash<custom_hash> {
    inline size_t operator()(const custom_hash& x) const {
      size_t result = 0;
      for (auto u : x.vals) {
        result ^= int_hash(u);
      }
      return custom_int_hash::splitmix64(result);
    }
};
}
```

#### AhoCorasick.h

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

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

```
struct AhoCorasick {
  enum {alpha = 26, first = 'A'}; // change this!
  struct Node {
    // (nmatches is optional)
   int back, next[alpha], start = -1, end = -1, nmatches = 0;
   Node(int v) { memset(next, v, sizeof(next)); }
 vector<Node> N;
 vi backp;
  void insert(string& s, int j) {
   assert(!s.emptv());
   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 = v;
       else {
         N[ed].back = y;
         (N[ed].end == -1 ? N[ed].end : backp[N[ed].start])
           = N[y].end;
         N[ed].nmatches += N[v].nmatches;
         q.push(ed);
```

```
vi find(string word) {
  int n = 0;
  vi res; // ll count = 0;
  for (char c : word) {
    n = N[n].next[c - first];
    res.push_back(N[n].end);
    // count \neq 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;
```

# $\underline{\text{Various}}$ (9)

## 9.1 Intervals

#### IntervalContainer.h

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

```
Time: \mathcal{O}(\log N)
set<pii>::iterator addInterval(set<pii>& is, int L, int R) {
 if (L == R) return is.end();
 auto it = is.lower_bound({L, R}), before = it;
 while (it != is.end() && it->first <= R) {</pre>
   R = max(R, it->second);
   before = it = is.erase(it);
 if (it != is.begin() && (--it)->second >= L) {
   L = min(L, it->first);
   R = max(R, it->second);
   is.erase(it);
 return is.insert(before, {L,R});
void removeInterval(set<pii>& is, int L, int R) {
 if (L == R) return;
 auto it = addInterval(is, L, R);
 auto r2 = it->second;
 if (it->first == L) is.erase(it);
 else (int&)it->second = L;
 if (R != r2) is.emplace(R, r2);
```

#### IntervalCover.h

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

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

## 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}(k \log \frac{n}{k})
```

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

## 9.2 Misc. algorithms

#### TernarySearch.h

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

```
Usage: int ind = ternSearch(0,n-1,[&](int i){return a[i];});

Time: \mathcal{O}(\log(b-a))
```

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

## LIS FastKnapsack KnuthDP DCDP SOSDP Knapsack

```
Description: Compute indices for the longest increasing subsequence.
Time: O(N \log N)
template < class I > vi lis(const vector < I > & S) {
 if (S.empty()) return {};
  vi prev(sz(S));
  typedef pair<I, int> p;
  vector res;
  rep(i, 0, sz(S)) {
    // change 0 \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

LIS.h

**Description:** Given N non-negative integer weights w and a non-negative target t, computes the maximum S <= t such that S is the sum of some subset of the weights.

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

d41d8c, 16 lines

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

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

#### DCDP.h

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

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

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

```
if (L >= R) return;
    int mid = (L + R) \gg 1;
    pair<11, int> best(LLONG_MAX, LO);
    rep(k, max(LO,lo(mid)), min(HI,hi(mid)))
      best = min(best, make_pair(f(mid, k), k));
    store(mid, best.second, best.first);
    rec(L, mid, LO, best.second+1);
    rec(mid+1, R, best.second, HI);
  void solve(int L, int R) { rec(L, R, INT_MIN, INT_MAX); }
};
SOSDP.h
Description: SOS DP
Time: \mathcal{O}\left(N*2^N\right)
                                                         d41d8c, 6 lines
//memory optimized, super easy to code.
for(int i = 0; i<(1<<N); ++i) F[i] = A[i];</pre>
for (int i = 0; i < N; ++i)
 for(int mask = 0; mask < (1<<N); ++mask) {</pre>
    if (mask & (1<<i)) F[mask] += F[mask^(1<<i)];</pre>
Knapsack.h
Description: Knapsack fast.
Time: \mathcal{O}\left(n^2Clogn/64\right) and \mathcal{O}\left(nC/64\right)
                                                        d41d8c, 48 lines
#pragma push macro(" SIZEOF LONG ")
#pragma push_macro("__cplusplus")
#define __SIZEOF_LONG__ __SIZEOF_LONG_LONG__
#define unsigned unsigned long
#define __cplusplus 201102L
#define __builtin_popcountl __builtin_popcountll
#define builtin ctzl builtin ctzll
#pragma pop_macro("__cplusplus")
#pragma pop_macro("__SIZEOF_LONG__")
#undef unsigned
#undef __builtin_popcount1
#undef __builtin_ctzl
const int C = 1e6 + 3;
vector<int> ans;
int M;
bitset<C> dp1, dp2;
bool divide(const vector<int> &a, int 1, int r, int S) {
    if (r - 1 == 1) {
        if (a[1] == S) {
             ans.push_back(1);
         } else if (S != 0) {
             return false:
        return true;
    int m = (1 + r) >> 1;
    dp1 = 0;
    dp1[0] = true;
    for (int i = 1; i < m; i++)</pre>
        dp1 |= dp1 << a[i];
    dp2 = 0;
    dp2[S] = true;
    for (int i = r - 1; i >= m; i--)
         dp2 \mid = dp2 >> a[i];
    for (int x = 0; x \le (r - 1) * M; x++) {
        if (dp1[x] && dp2[x]) {
```

assert (divide(a, l, m, x));

```
assert (divide (a, m, r, S - x));
        return true;
return false;
```

# Techniques (A)

## techniques.txt

Combinatorics

159 lines

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

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

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

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