iTMO

ITMO

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

overpressured but trained

The 2024 ICPC Northern Eurasia Finals
December 15, 2024

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 θ , area A and magic flux $F = b^2 + d^2 - a^2 - c^2$:

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

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

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 σ is the standard deviation. If X is instead continuous it will have a probability density function $f_X(x)$ and the sums above will instead be integrals with $p_X(x)$ replaced by $f_X(x)$.

Expectation is linear:

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

For independent X and Y,

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

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.

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

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

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

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

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 $\sim 3x$ faster. Uses 1.5x memory. Initial capacity must be a power of 2 (if provided).

```
#include <bits/extc++.h>
struct custom_int_hash {
```

```
static uint64 t splitmix64(uint64 t x) {
   x += 0x9e3779b97f4a7c15; x = (x ^ (x >> 30)) * 0
        xbf58476d1ce4e5b9; x = (x ^ (x >> 27)) * 0
        x94d049bb133111eb;
   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);
};
struct chash { // large odd number for C
 const uint 64_t C = 11(4e18 * acos(0)) | 71;
 11 operator()(11 x) const { return __builtin_bswap64(x*C); }
__gnu_pbds::gp_hash_table<11,int,chash /* custom_int_hash */> h
    ({},{},{},{},{1<<16});
LazySegmentTree.h
Description: Segment tree boilerplate
Time: \mathcal{O}(\log N).
                                                     88a0ab, 50 lines
class segtree {
 public:
 struct node
    // don't forget to set default value (used for leaves) not
         necessarily neutral element!
    void apply(int 1, int r, ... v) { }
 };
 node unite (const node &a, const node &b) const { node res; /*
       res = combine(a, b) */ return res; }
 inline void push (int x, int 1, int r) { int y = (1 + r) >> 1;
       int z = x + ((y - 1 + 1) << 1); }
 inline void pull(int x, int z) { tree[x] = unite(tree[x + 1],
       tree[z]); }
 int n; vector<node> tree;
 template <typename M> void build(int x, int 1, int r, const
      vector<M> &v) {
    if (1 == r) { tree[x].apply(1, r, v[1]); return; }
   int y = (1 + r) >> 1; int z = x + ((y - 1 + 1) << 1);
   build(x + 1, 1, y, v); build(z, y + 1, r, v); pull(x, z);
 node get(int x, int 1, int r, int 11, int rr) {
   if (ll <= 1 && r <= rr) { return tree[x]; }</pre>
   int y = (1 + r) >> 1; int z = x + ((y - 1 + 1) << 1);
   push(x, l, r);
   node res{};
   if (rr \le y) res = get (x + 1, 1, y, 11, rr);
    else if (11 > y) res = get(z, y + 1, r, 11, rr);
    else res = unite(get(x + 1, 1, y, 11, rr), get(z, y + 1, r,
         11, rr));
    pull(x, z);
    return res:
  template <typename... M>
 void modify(int x, int 1, int r, int 11, int rr, const M&...
    if (11 <= 1 && r <= rr) { tree[x].apply(1, r, v...); return</pre>
    int y = (1 + r) >> 1; int z = x + ((y - 1 + 1) << 1);
   push(x, l, r);
    if (l1 <= y) modify(x + 1, 1, y, 11, rr, v...);</pre>
```

if (rr > y) modify(z, y + 1, r, ll, rr, v...);

pull(x, z);

UnionFindRollback.h

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

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

(V)) 4aff3d, 21 lines

```
struct RollbackUF
  vi e; vector<pii> st;
 RollbackUF(int n) : e(n, -1) {}
  int size(int x) { return -e[find(x)]; }
  int find(int x) { return e[x] < 0 ? x : find(e[x]); }</pre>
  int time() { return sz(st); }
  void rollback(int t) {
    for (int i = time(); i --> t;)
      e[st[i].first] = st[i].second;
    st.resize(t);
 bool join(int a, int b) {
    a = find(a), b = find(b);
    if (a == b) return false;
    if (e[a] > e[b]) swap(a, b);
    st.push_back({a, e[a]});
    st.push back({b, e[b]});
    e[a] += e[b]; e[b] = a;
    return true;
};
```

SubMatrix.h

Description: Calculate submatrix sums quickly, given upper-left and lower-right corners (half-open).

Usage: SubMatrix<int> m(matrix); m.sum(0, 0, 2, 2); // top left 4 elements Time: $\mathcal{O}\left(N^2+Q\right)$

 $: \mathcal{O}\left(N^2 + Q\right)$ afd259, 11 lines

Matrix.h

Description: Basic operations on matrices. NEVER FORGET TRANS-POSING BEFORE MULTIPLYING.

LineContainer RationalLineContainer RMQ

```
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;
                                                        338ea3, 33 lines
template < class T, int N, int M> struct Matrix {
  using M1 = Matrix<T, M, N>; using M2 = Matrix<T, N, M>;
  M1 d{};
  M2 transpose() const {
    rep(i, 0, N) rep(j, 0, M) a.d[j][i] = d[i][j];
    return a;
  template<int K>
  Matrix<T, N, K> operator*(const Matrix<T, M, K>& b) const {
    Matrix<T, N, K> a;
    Matrix<T, K, M> b t = b.transpose();
    rep(i, 0, N) rep(j, 0, K) rep(k, 0, M) a.d[i][j] += d[i][k]*b_t.d
         [j][k];
    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;
  Matrix<T, N, N> operator^(ll p) const {
    if constexpr (N != M) throw std::invalid argument("Matrix
         is not square");
    assert (p >= 0);
    Matrix<T, N, N> 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 ("con-
vex hull trick").
Time: \mathcal{O}(\log N)
```

```
ad1f46, 29 lines
struct Line {
 mutable ll 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;
  11 div(11 a, 11 b) { return a / b - ((a ^ b) < 0 && a % b); }</pre>
  bool isect(iterator x, iterator y) {
   if (y == end()) return x \rightarrow p = inf, 0;
   if (x->k == y->k) x->p = x->m > y->m ? inf : -inf;
   else x->p = div(y->m - x->m, x->k - y->k);
   return x->p >= y->p;
  void add(ll k, ll m) {
    auto z = insert(\{k, m, 0\}), y = z++, x = y;
   while (isect(y, z)) z = erase(z);
   if (x != begin() \&\& isect(--x, y)) isect(x, y = erase(y));
    while ((y = x) != begin() && (--x)->p >= y->p)
      isect(x, erase(y));
```

```
ll query(ll x) {
    assert(!empty());
    auto 1 = *lower bound(x);
    return 1.k * x + 1.m;
};
RationalLineContainer.h
Description: linear CHT and persistent linear CHT for monotonic slopes
Time: \mathcal{O}(\log N)
"../../stress-tests/utilities/template.h"
constexpr int64 t CHT LINE INF = std::numeric limits<int64 t>::
    max();
struct line {
 int a{0};
 int64_t b{CHT_LINE_INF};
  line() = default;
  line(int a, int64_t b) noexcept : a(a), b(b) {}
  friend __int128 cross(const line &left, const line &right) {
    return ( int128)left.a * right.b - ( int128)left.b *
  friend line operator-(const line &left, const line &right) {
    return {left.a - right.a, left.b - right.b};
  [[nodiscard]] int64_t evaluate(int x) const { return (int64_t
      )a * x + b; }
};
// query for min. inserting for increasing/decreasing slope
template <bool increasing_slope> struct incremental_CHT {
 vector<line> lines:
 void insert(const line &new line) {
    while (lines.size() >= 2) {
      const line &line1 = lines[lines.size() - 2], line2 =
          lines.back(), line3 = new line;
      if ((cross(line3 - line2, line2 - line1) < 0) ^</pre>
           increasing_slope) { // > 0 for max
        lines.pop_back();
      } else break;
    if (!lines.empty()) {
      const line &line1 = lines.back(), line2 = new_line;
      if (line1.a == line2.a) {
        if (line1.b > line2.b) lines.pop_back(); // < for max</pre>
        else return;
    lines.push back(new line);
  // have to perform queries with decreasing x
  int64 t query(int x) {
    assert(!lines.empty());
    while (lines.size() >= 2) {
      const line &line1 = lines[lines.size() - 2], line2 =
           lines.back();
      if (line1.evaluate(x) < line2.evaluate(x)) lines.pop_back</pre>
           (); // > for max
      else break:
    return lines.back().evaluate(x);
template <bool increasing_slope> struct
    persistent incremental CHT {
```

```
vector<line> lines;
  vector<int> parents;
  int max_query_alive_vertex{};
  persistent_incremental_CHT() = default;
  int insert(const line &new line) {
    parents.push_back(-1); lines.push_back(new_line);
         max_query_alive_vertex = lines.size() - 1;
    if (lines.size() == 1) return 0;
    int k = lines.size() - 1, i = lines.size() - 2, i = parents
    parents[k] = j;
    while (i !=-1) {
      const line &line1 = lines[i], line2 = lines[j], line3 =
           new line;
      if ((cross(line3 - line2, line2 - line1) > 0) ^ !
           increasing_slope) { // < 0 \text{ for max}
        parents[k] = i; j = i; i = parents[i];
      } else break;
    while (j != -1) {
      const line &line1 = lines[j], line2 = new_line;
      if (line1.a == line2.a) {
        if (line1.b > line2.b) parents[k] = j = parents[j]; //
             < for max
        else return k;
      } else break;
    return k;
  int get(int root, int x) {
    if (root == -1 || parents[root] == -1) return root;
    const line &line1 = lines[parents[root]], line2 = lines[
    if (line1.evaluate(x) < line2.evaluate(x)) return parents[</pre>
         root] = get(parents[root], x);
    else return root;
  // have to perform queries with increasing x if slope
       increases
  int64 t query(int root, int x) {
    assert(!lines.empty());
    int64_t answer = lines[root].evaluate(x);
    return min(answer, lines[get(root, x)].evaluate(x));
};
RMQ.h
Description: Sparse tables.
Time: \mathcal{O}(1) for query, \mathcal{O}(n \log n) for build
                                                       27865b, 65 lines
    auto fun = [\mathcal{E}](int \ i, \ int \ j) \ \{ \ return \ min(i, \ j); \ \};
    SparseTable < int, decltype(fun) > st(a, fun);
// Sparse Table < int> st(a, [\mathcal{C}] (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) {
```

MoQueries Polynomial

```
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))]);
 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 +
        _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
  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 approximate 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:** $\mathcal{O}\left(N\sqrt{Q}\right)$

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

2.0.1 Segment Tree Beats

Ji Driver Segment Tree Beats. min = max = sum?, min?max? – store max and second max

```
min = + = , qcd.
```

store differences on segment, a[i] - a[j], but only BST of differences.

mod = set, sum?

break condition max; mod tag condition max == min.

```
sqrt = + = sum?max?min?
```

store sum, max, min, break condition just standard segtree qr <= l||r <= qr tag condition max - min <= 1

```
div = + sum?max?min?
```

```
store sum, max, min, break condition just standard segtree qr <= l||r <= qr \text{ tag condition } max - min <= 1
```

```
\& = , | =, max?
```

 $C=2^k$ is important the upper bound for numbers. store $max, pushand, pushor, and_on_seg, or_on_seg$ break condition: standard segtree $qr <= l||r <= ql \text{ tag condition for and } = ql <= l\&\&r <= qr\&\&((and_on_seg[v]^or_on_seg[v])\&x) == 0$ tag condition for or $= ql <= l\&\&r <= qr\&\&((and_on_seg[v]^or_on_seg[v])\&y) == 0$

Numerical (3)

3.1 Polynomials and recurrences

Polynomial.h

```
Description: Polynomial operations
```

d18448, 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) {</pre>
            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) {
            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++)
            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) {</pre>
               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++)
                a[i] = a[i] * k;
    }
    template<typename base>
   vector<base> mul(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++)
                R[i] = (R[i >> 1] >> 1) | ((i & 1) << (L -
                    1));
       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<base> mul(vector<base> a, vector<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<typename 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<base> inverse(const vector<base>& a, int size) {
    assert(size > 0 && a[0] != 0);
   vector<base> 0{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)  {
          return {0};
      if (size == 1)  {
          return \{1/a/0\};
      auto \ op = plug\_minus\_x(a);
      auto T = mul(a, op);
      T = onlu_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> 0{1};
    for (int sz = 2;; sz *= 2) {
        Q = mul(Q, sub(add(a, {1}), ln(Q, sz)), sz);
        if (sz >= size)
            break:
   Q.resize(size);
    return O;
```

PolyInterpolate BerlekampMassey

```
// O(n*log(n))
    template<typename base>
    vector<base> pow(vector<base> a, ll p, int size) {
        int i = 0;
        while (i < a.size()) {</pre>
            if (a[i] != 0)
                break:
            ++i;
        if (i == a.size()) {
            auto res = vector<base>(size, 0);
            if (p == 0)
                res[0] = 1;
            return res;
        a.erase(a.begin(), a.begin() + i);
        auto f = a[0];
        for (auto& x : a) x /= f;
        a = \exp(mul(ln(a, size), (base)p), size);
        for (int j = size - 1; j >= 0; --j) {
            if ((i > 0 \&\& p >= size) || j - p * i < 0)
                a[j] = 0;
            el se
                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>
```

PolyInterpolate.h

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

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

BerlekampMassey.h

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

```
Usage: berlekampMassey({0, 1, 1, 3, 5, 11}) // {1, 2}
Time: \mathcal{O}(N^2) and \mathcal{O}(nlog^2(k))
                                                     f796a5, 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 \star = 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++) {
        mint d = S[i];
        for (int j = 1; j \le L; j++) d += C[j] * S[i - j];
        if (d == 0) continue;
        mint coef = d * b.inv();
        for (int j = m; j < n; j++) C[j] -= coef * B[j - m];
        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 \neq -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++) {
        for (int j = 0; j < n + 1; j++) res[i + j] += a[i] * b[</pre>
            j];
    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
    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 *= -1;
    res *= a[i][i];
   if (res == 0) return 0;
    rep(j,i+1,n) {
     double v = a[j][i] / a[i][i];
     if (v != 0) rep(k, i+1, n) a[j][k] -= v * a[i][k];
  return res;
```

IntDeterminant.h

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

```
Time: \mathcal{O}\left(N^3\right)
const 11 mod = 12345;
```

3a8e36, 18 lines

```
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)
                                                        fbf9d8, 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"
                                                       71598f, 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:; }
```

SolveLinearBinary.h

Description: Solves Ax = b over \mathbb{F}_2 . If there are multiple solutions, one is returned arbitrarily. Returns rank, or -1 if no solutions. Destroys A and b. Time: $\mathcal{O}\left(n^2m\right)$ 9c2cbe, 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: $\mathcal{O}\left(n^3\right)$ 616852, 35 lines

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

01e75e, 16 lines

MatrixInverse-mod Tridiagonal FFT FFTM FST

Fails if the solution is not unique.

```
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"

63b429, 36 lines

```
int matInv(vector<vector<ll>>& A) {
 int n = sz(A); vi col(n);
  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]);
    11 v = modpow(A[i][i], mod - 2);
    rep(j,i+1,n) {
     ll 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 $\{a_i\} = \text{tridiagonal}(\{1, -1, -1, ..., -1, 1\}, \{0, c_1, c_2, ..., c_n\},\$ $\{b_1, b_2, \ldots, b_n, 0\}, \{a_0, d_1, d_2, \ldots, d_n, a_{n+1}\}\}$

```
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)
```

```
typedef double T;
vector<T> tridiagonal(vector<T> diag, const vector<T>& super,
   const vector<T>& sub, vector<T> b) {
 int n = sz(b); vi tr(n);
 rep(i, 0, n-1) {
    if (abs(diag[i]) < 1e-9 * abs(super[i])) { // diag[i] == 0
     b[i+1] = b[i] * diag[i+1] / super[i];
     if (i+2 < n) b[i+2] -= b[i] * sub[i+1] / super[i];</pre>
     diag[i+1] = sub[i]; tr[++i] = 1;
     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: conv(a, b) = c, where $c[x] = \sum a[i]b[x-i]$. For convolution of complex numbers or more than two vectors: FFT, multiply pointwise, divide by n, reverse(start+1, end), FFT back. Rounding is safe if $(\sum a_i^2 + \sum b_i^2) \log_2 N < 9 \cdot 10^{14}$ (in practice 10^{16}); higher for random inputs). Otherwise, use NTT/FFTMod. **Time:** $O(N \log N)$ with N = |A| + |B| (~1s for $N = 2^{22}$)

```
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];
 vi rev(n);
  rep(i,0,n) \ rev[i] = (rev[i / 2] | (i & 1) << L) / 2;
  rep(i,0,n) if (i < rev[i]) swap(a[i], a[rev[i]]);
 for (int k = 1; k < n; k *= 2)
    for (int i = 0; i < n; i += 2 * k) rep(j,0,k) {
     Cz = rt[j+k] * a[i+j+k]; // (25\% faster if hand-rolled)
     a[i + j + k] = a[i + j] - z;
     a[i + j] += z;
vd conv(const vd& a, const vd& b) {
 if (a.empty() || b.empty()) return {};
  vd res(sz(a) + sz(b) - 1);
  int L = 32 - \underline{\quad} builtin_clz(sz(res)), n = 1 << L;
 vector<C> in(n), out(n);
```

```
copy(all(a), begin(in));
rep(i,0,sz(b)) in[i].imag(b[i]);
fft(in);
for (C& x : in) x *= x;
rep(i, 0, n) out[i] = in[-i & (n - 1)] - conj(in[i]);
fft(out);
rep(i, 0, sz(res)) res[i] = imag(out[i]) / (4 * n);
return res;
```

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;
```

FST.h

5609f3, 35 lines

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

```
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);
                                             // XOR
 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(g == 1); return Mod((x + mod) % mod);
  Mod operator^(11 e) {
   if (!e) return Mod(1);
   Mod r = *this ^ (e / 2); r = r * r;
    return e&1 ? *this * r : r;
};
```

ModLog.h

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

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

Time: O(\sqrt{m})

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

ModSum.h

Description: Sums of mod'ed arithmetic progressions.

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

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

4c49b1, 16 lines

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

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

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

ModMulLL.h

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

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

ModSqrt.h

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

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

```
"ModPow.h"
                                                        b548fc, 24 lines
ll sgrt(ll a, ll p) {
 a %= p; if (a < 0) a += p;
 if (a == 0) return 0;
 assert (modpow(a, (p-1)/2, p) == 1); // else no solution
 if (p % 4 == 3) return modpow(a, (p+1)/4, p);
 // a^{(n+3)/8} \text{ or } 2^{(n+3)/8} * 2^{(n-1)/4} \text{ works if } p \% 8 == 5
 11 s = p - 1, n = 2;
 int r = 0, m;
 while (s % 2 == 0)
    ++r, s /= 2;
  while (modpow(n, (p-1) / 2, p) != p-1) ++n;
 11 x = modpow(a, (s + 1) / 2, p);
 11 b = modpow(a, s, p), g = modpow(n, s, p);
 for (;; r = m) {
    11 t = b;
    for (m = 0; m < r && t != 1; ++m)
     t = t * t % p;
    if (m == 0) return x;
    11 \text{ qs} = \text{modpow}(q, 1LL \ll (r - m - 1), p);
   g = gs * gs % p;
   x = x * qs % p;
   b = b * q % p;
```

4.2 Primality

FastErat.h

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

```
ee2482, 20 lines
const int LIM = 1e6;
bitset<LIM> isPrime;
vi eratosthenes() {
  const int S = (int)round(sqrt(LIM)), R = LIM / 2;
 vi pr = \{2\}, sieve(S+1); pr.reserve(int(LIM/log(LIM) \star1.1));
  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 $a^b \mod c$.

${ m Factor.h}$

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.
                                                        6f4f4f, 34 lines
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;</pre>
    ul A[] = \{2, 325, 9375, 28178, 450775, 9780504, 1795265022\}
         , s = \underline{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
             ,n);
        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);
```

euclid CRT phi ContinuedFractions FracBinarySearch IntPerm

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 a-gcd instead. If a and b are coprime, then a is the inverse of a (mod 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

Description: Chinese Remainder Theorem.

crt (a, m, b, n) computes x such that $x\equiv a\pmod m$, $x\equiv b\pmod n$. If |a|< m and |b|< n, x will obey $0\le x< \mathrm{lcm}(m,n)$. 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

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

phi.h

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

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

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

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

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

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

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

d34567, 21 lines

```
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.q; }, 10); // {1,3} Time: $\mathcal{O}\left(\log(N)\right)$ 94107a, 25 lines

```
struct Frac { ll p, q; };
template<class F>
Frac fracBS(F f, ll N) {
 bool dir = 1, A = 1, B = 1;
 Frac lo{0, 1}, hi{1, 1}; // Set hi to 1/0 to search (0, N]
 if (f(lo)) return lo;
 assert(f(hi));
 while (A | | B) {
    11 adv = 0, step = 1; // move hi if dir, else lo
    for (int si = 0; step; (step *= 2) >>= si) {
     Frac mid{lo.p * adv + hi.p, lo.q * adv + hi.q};
      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 $\mathbb{Z}_{2^a}^{\times}$ is instead isomorphic to $\mathbb{Z}_2 \times \mathbb{Z}_{2^{a-2}}$.

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:

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

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

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

Combinatorial (5)

5.1 Permutations

5.1.1 Factorial

	n	1 2 3	4	5 6	7	8	9	10	
-	n!	1 2 6	24 1	20 720	5040	40320	362880	3628800	
	n						16		
	n!	4.0e7	′ 4.8e	8 6.2e9	9 8.7el	10 1.3e	12 2.1e1	3 3.6e14	
	n							171	
	n!	2e18	2e25	3e32 8	$8e47 \ 3$	e64 9e	157 6e26	$32 > DBL_M$	AX

IntPerm.h

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

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

72368d, 6 lines

```
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>
```

ITMO(Ivchenko, Konovalov, Matveev)

multinomial Dinic

11

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

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

5.2 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

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

General purpose numbers

5.3.1 Bernoulli numbers

EGF of Bernoulli numbers is $B(t)=\frac{t}{e^t-1}$ (FFT-able). $B[0,\ldots]=[1,-\frac12,\frac16,0,-\frac1{30},0,\frac1{42},\ldots]$

Sums of powers:

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

Euler-Maclaurin formula for infinite sums:

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

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

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 <math>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

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

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

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

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

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

6.1 Network flow

Dinic.h

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

```
"../../stress-tests/utilities/template.h"
template<class F>
struct dinic {
 static constexpr db eps = 1e-6;
  struct Edge {
    int to; F flow, cap;
    Edge() = default;
    Edge(int to_, F flow_, F cap_) : to(to_), flow(flow_), cap(
         cap_) {}
 int n; vvi gr; vector<Edge> edges;
 vector<F> dist; vector<int> first;
 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 1, F r) { add(s,
       v, 1); add(u, v, r - 1); add(u, t, 1); }
 F res(int id) { return edges[id].cap - edges[id].flow; }
 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});
```

auto v = Fedya_Romashov.front(); Fedya_Romashov.pop();

while(!Fedya_Romashov.empty()) {

FlowDecomposition MCMF

```
for (auto id : gr[v]) {
        auto& e = edges[id];
        if (res(id) > 0 && dist[e.to] < 0) { dist[e.to] = dist[</pre>
             v] + 1; Fedya_Romashov.push(e.to); }
    return dist[t] >= 0;
  F dfs(int v, int t, F current flow = numeric limits<F>::max()
    if (v == t) return current flow;
    F small push = 0:
    for (; first[v] < gr[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 max flow(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(int s, int t) {
    max flow(s, t);
    vector<bool> ret(n);
    for (int i = 0; i < n; i++) { ret[i] = (dist[i] != -1); }</pre>
};
FlowDecomposition.h
Description: Decompose flow into paths and cycles.
Time: \mathcal{O}(FLOW + m + n)
                                                     72e3cb, 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(q.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(q.n);
for (int i = 0; i < g.n; i++) {
 ptr[i] = (int) q.q[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) % q.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 = g.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 = q.q[v][ptr[v]];
     if (fs[id] > q.eps) {
       path.push back(id);
       v = q.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)
                                                     7afab8, 47 lines
template <class F, class C = F> struct MCMF {
  struct Edge { int to; F flow, cap; C cost; };
 int n; vector<C> pot, dist;
 vector<int> previous_edge; vector<Edge> edges; vector<vector<</pre>
      int>> gr;
 MCMF (int n ) : n(n ) {
    pot.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();
    dist.assign(n, inf);
    using T = pair<C, int>; priority_queue<T, vector<T>,
        greater<T>> bfs:
    bfs.push({dist[s] = 0, s});
    while (!bfs.emptv()) {
      auto [cur_dist, v] = bfs.top(); bfs.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 + pot[v] - pot[E.to]))
          previous_edge[E.to] = e, bfs.push({dist[E.to], E.to})
    return dist[t] != inf;
 pair<F, C> calc(int s, int t) {
    assert(s != t);
    rep(n) for (int e = 0; e < edges.size(); ++e) {
      const Edge &E = edges[e]; if (E.cap) ckmin(pot[E.to], pot
          [edges[e ^ 1].to] + E.cost); // Bellman-Ford
    F totalFlow = 0; C totalCost = 0;
    while (path(s, t)) {
      for (int i = 0; i < n; ++i) pot[i] += (dist[i] ==</pre>
          numeric_limits<C>::max() ? 0 : dist[i]);
      F df = numeric limits<F>::max();
```

```
for (int x = t; x != s; x = edges[previous_edge[x] ^ 1].
          to) { const Edge &E = edges[previous_edge[x]]; ckmin
          (df, E.cap - E.flow); }
     totalFlow += df; totalCost += (pot[t] - pot[s]) * df;
     for (int x = t; x != s; x = edges[previous_edge[x] ^ 1].
          to) edges[previous_edge[x]].flow += df, edges[
          previous_edge[x] ^ 1].flow -= df;
   return {totalFlow, totalCost};
};
```

MCMFPushRelabel.h

Description: Min-cost max-flow. Supports lower bounds and negative costs (and even cycles!)

```
Time: \mathcal{O}\left(\tilde{V}^3 \log(VC)\right)
"../../stress-tests/utilities/template.h"
                                                      4ef0ee, 275 lines
template <class F> struct HLPP {
  struct Edge {
   int to, inv;
   F rem, cap;
  vector<vector<Edge>> G;
  vector<F> excess:
  vector<int> hei, arc, prv, nxt, act, bot;
  queue<int> 0;
  int n, high, cut, work;
  HLPP(int k) : G(k) {}
  int addEdge(int u, int v, F cap, F rcap = 0) {
    assert(u != v);
    G[u].push_back({v, sz(G[v]), cap, cap});
    G[v].push_back({u, sz(G[u]) - 1, rcap, rcap});
    return sz(G[u]) - 1;
  void raise(int v, int h) {
    prv[nxt[prv[v]] = nxt[v]] = prv[v];
   hei[v] = h;
    if (excess[v] > 0) {
     bot[v] = act[h];
     act[h] = v;
     high = max(high, h);
    if (h < n)
     cut = max(cut, h + 1);
    nxt[v] = nxt[prv[v] = h += n];
   prv[nxt[nxt[h] = v]] = v;
  void global(int s, int t) {
   hei.assign(n, n \star 2);
    act.assign(n \star 2, -1);
   iota(all(prv), 0);
   iota(all(nxt), 0);
   hei[t] = high = cut = work = 0;
   hei[s] = n;
    for (int x : \{t, s\})
      for (Q.push(x); !Q.empty(); Q.pop()) {
        int v = Q.front();
        for (auto &e : G[v])
          if (hei[e.to] == n * 2 \&\& G[e.to][e.inv].rem)
            Q.push(e.to), raise(e.to, hei[v] + 1);
  void push (int v, Edge &e, bool z) {
    auto f = min(excess[v], e.rem);
   if (f > 0) {
     if (z && !excess[e.to]) {
       bot[e.to] = act[hei[e.to]];
        act[hei[e.to]] = e.to;
```

```
e.rem -= f;
     G[e.to][e.inv].rem += f;
     excess[v] -= f;
     excess[e.to] += f;
 void discharge(int v) {
   int h = n * 2, k = hei[v];
   for (int j = 0; j < sz(G[v]); j++) {</pre>
     auto &e = G[v][arc[v]];
     if (e.rem) {
       if (k == hei[e.to] + 1) {
         push(v, e, 1);
         if (excess[v] <= 0)
           return:
       } else
         h = min(h, hei[e.to] + 1);
     if (++arc[v] >= sz(G[v]))
       arc[v] = 0;
   if (k < n \&\& nxt[k + n] == prv[k + n]) {
     for (int j = k; j < cut; j++)</pre>
       while (nxt[j + n] < n)
         raise(nxt[j + n], n);
     cut = k;
   } else
     raise(v, h), work++;
 // Compute maximum flow from src to dst
 F flow(int src, int dst) {
   excess.assign(n = sz(G), 0);
   arc.assign(n, 0);
   prv.assign(n * 3, 0);
   nxt.assign(n * 3, 0);
   bot.assign(n, 0);
   for (auto &e : G[src])
     excess[src] = e.rem, push(src, e, 0);
   global(src, dst);
   for (; high; high--)
     while (act[high] !=-1) {
       int v = act[high];
       act[high] = bot[v];
       if (v != src && hei[v] == high) {
         discharge(v);
         if (work > 4 * n)
           global(src, dst);
   global(src, dst);
   return excess[dst];
 // Get flow through e-th edge of vertex v
 F getFlow(int v, int e) { return G[v][e].cap - G[v][e].rem; }
 // Get if v belongs to cut component with src
 bool cutSide(int v) { return hei[v] >= n; }
template <class T> struct Circulation {
 const T INF = numeric_limits<T>::max() / 2;
 T lowerBoundSum = 0;
 HLPP<T> mf;
 // Initialize for n vertices
 Circulation(int k): mf(k + 2) {}
 void addEdge(int s, int e, T l, T r) {
   mf.addEdge(s + 2, e + 2, r - 1);
   if (1 > 0) {
     mf.addEdge(0, e + 2, 1);
```

```
mf.addEdge(s + 2, 1, 1);
      lowerBoundSum += 1;
    } else {
      mf.addEdge(0, s + 2, -1);
      mf.addEdge(e + 2, 1, -1);
      lowerBoundSum += -1:
 bool solve(int s, int e) {
    // \ mf.addEdge(e+2, s+2, INF); // to reduce as maxflow with
         lower bounds, in circulation problem skip this line
    return lowerBoundSum == mf.flow(0, 1);
    // to get maximum LR flow, run maxflow from s+2 to e+2
};
template <class T> struct MinCostCirculation {
  const int SCALE = 3; // scale by 1/(1 << SCALE)</pre>
  const T INF = numeric_limits<T>::max() / 2;
 struct EdgeStack {
   int s, e;
   T l, r, cost;
 };
  struct Edge {
   int pos, rev;
   T rem, cap, cost;
 };
 int n:
 vector<EdgeStack> estk;
 Circulation<T> circ;
 vector<vector<Edge>> gph;
 vector<T> p;
 MinCostCirculation(int k) : circ(k), gph(k), p(k) { n = k; }
  void addEdge(int s, int e, T l, T r, T cost){
    estk.push_back({s, e, l, r, cost});
 pair<bool, T> solve() {
    for(auto &i : estk) {
      if(i.s != i.e) circ.addEdge(i.s, i.e, i.l, i.r);
    if(!circ.solve(-1, -1)){
      return make_pair(false, T(0));
    vector<int> ptr(n);
    T eps = 0;
    for(auto &i : estk){
     T curFlow;
      if(i.s != i.e) curFlow = i.r - circ.mf.G[i.s + 2][ptr[i.s
          11.rem;
      else curFlow = i.r;
      int srev = sz(gph[i.e]);
      int erev = sz(gph[i.s]);
      if(i.s == i.e) srev++;
      gph[i.s].push_back({i.e, srev, i.r - curFlow, i.r, i.cost
            * (n + 1) \});
      gph[i.e].push_back({i.s, erev, -i.l + curFlow, -i.l, -i.
          cost * (n + 1) });
      eps = max(eps, abs(i.cost) * (n + 1));
      if(i.s != i.e){
       ptr[i.s] += 2;
        ptr[i.e] += 2;
    while(true){
      auto cost = [&](Edge &e, int s, int t){
       return e.cost + p[s] - p[t];
      eps = 0;
```

Hungarian GlobalMinCut

```
for(int i = 0; i < n; i++) {</pre>
    for(auto &e : gph[i]){
      if(e.rem > 0) eps = max(eps, -cost(e, i, e.pos));
  if(eps <= T(1)) break;</pre>
  eps = max(T(1), eps >> SCALE);
  bool upd = 1;
  for(int it = 0; it < 5 && upd; it++){</pre>
    upd = false:
    for(int i = 0; i < n; i++) {</pre>
      for(auto &e : gph[i]){
        if(e.rem > 0 && p[e.pos] > p[i] + e.cost + eps){
          p[e.pos] = p[i] + e.cost + eps;
          upd = true;
      }
    if(!upd) break;
  if(!upd) continue;
  vector<T> excess(n);
  queue<int> que;
  auto push = [&](Edge &e, int src, T flow) {
   e.rem -= flow;
    qph[e.pos][e.rev].rem += flow;
    excess[src] -= flow;
    excess[e.pos] += flow;
   if(excess[e.pos] <= flow && excess[e.pos] > 0){
      que.push (e.pos);
  };
  vector<int> ptr(n);
  auto relabel = [&](int v){
   ptr[v] = 0;
   p[v] = -INF;
    for(auto &e : gph[v]){
      if(e.rem > 0){
        p[v] = max(p[v], p[e.pos] - e.cost - eps);
   }
  };
  for(int i = 0; i < n; i++) {</pre>
    for(auto &j : gph[i]) {
      if(j.rem > 0 && cost(j, i, j.pos) < 0){</pre>
        push(j, i, j.rem);
  while (sz (que)) {
    int x = que.front();
    que.pop();
    while(excess[x] > 0){
      for(; ptr[x] < sz(qph[x]); ptr[x]++){</pre>
        Edge &e = gph[x][ptr[x]];
        if(e.rem > 0 && cost(e, x, e.pos) < 0){</pre>
          push(e, x, min(e.rem, excess[x]));
          if(excess[x] == 0) break;
      if(excess[x] == 0) break;
      relabel(x);
T ans = 0;
for(int i = 0; i < n; i++) {</pre>
  for(auto &j : gph[i]){
    j.cost /= (n + 1);
```

```
ans += j.cost * (j.cap - j.rem);
   return make_pair(true, ans / 2);
 void bellmanFord() {
   fill(all(p), T(0));
   bool upd = 1;
    while (upd) {
      upd = 0:
      for(int i = 0; i < n; i++) {</pre>
        for(auto &j : gph[i]){
          if(j.rem > 0 && p[j.pos] > p[i] + j.cost){
            p[j.pos] = p[i] + j.cost;
            upd = 1;
};
Hungarian.h
Description: Solve assignment problem.
Time: \mathcal{O}\left(n^2 * m\right)
                                                      a470a9, 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> wav;
 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;
    do {
     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]) {</pre>

```
minv[j] = cur;
            way[j] = j0;
          if (minv[j] < delta) {</pre>
            delta = minv[j];
            j1 = j;
      for (int j = 0; j <= m; j++) {
        if (used[j]) {
          u[pb[j]] += delta;
          v[j] -= delta;
        } else {
          minv[j] -= delta;
      j0 = 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: $\mathcal{O}(V^3)$

```
a53e91, 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:
```

hopcroftKarp WeightedMatching Blossom

6.2 Matching

hopcroftKarp.h

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

```
Time: \mathcal{O}\left(\sqrt{V}E\right)
bool dfs(int a, int L, vector<vi>& g, vi& btoa, vi& A, vi& B) {
  if (A[a] != L) return 0;
  A[a] = -1;
  for (int b : g[a]) if (B[b] == L + 1) {
   B[b] = 0:
    if (btoa[b] == -1 || dfs(btoa[b], L + 1, q, 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(g)) if(A[a] == 0) cur.push_back(a);
    for (int lay = 1;; lay++) {
     bool islast = 0;
      next.clear();
      for (int a : cur) for (int b : g[a]) {
        if (btoa[b] == -1) {
          B[b] = lay;
          islast = 1;
        else if (btoa[b] != a && !B[b]) {
          B[b] = lav;
          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:** $\mathcal{O}(N^2M)$

```
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 "dumny" worker 0
   vi dist(m, INT_MAX), pre(m, -1);
   vector\bool> done(m + 1);
```

```
do { // dijkstra
      done[j0] = true;
      int i0 = p[j0], j1, delta = INT_MAX;
      rep(j,1,m) if (!done[j]) {
        auto cur = a[i0 - 1][j - 1] - u[i0] - v[j];
        if (cur < dist[j]) dist[j] = cur, pre[j] = j0;
        if (dist[j] < delta) delta = dist[j], j1 = j;</pre>
      rep(j,0,m) {
        if (done[j]) u[p[j]] += delta, v[j] -= delta;
        else dist[j] -= delta;
      j0 = j1;
    } while (p[j0]);
    while (j0) { // update alternating path
      int j1 = pre[j0];
      p[j0] = p[j1], j0 = j1;
 rep(j,1,m) if (p[j]) ans[p[j] - 1] = j - 1;
 return {-v[0], ans}; // min cost
Blossom.h
Description: Given a weighted graph, finds max matching
Time: \mathcal{O}(N^3)
"../../stress-tests/utilities/template.h"
                                                     0015e1, 129 lines
struct blossom {
    int n, m;
    vi mate: vvi b:
    vi p, d, bl; vvi 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); q
         .assign(m, vi(m, -1)); }
    void add_edge(int u, int v) {
        q[u][v] = u;
        g[v][u] = v;
    void match(int u, int v) {
        q[u][v] = q[v][u] = -1;
        mate[u] = v;
        mate[v] = u;
    vi trace(int x) {
        vi 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, vi &vx, vi &vy) {
        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++) {

```
g[c][i] = g[i][c] = -1;
    for(int z : b[c]) {
        bl[z] = c;
        for(int i = 0; i < c; i++) {</pre>
            if(g[z][i] != -1) {
                g[c][i] = z;
                g[i][c] = g[i][z];
vi lift(vi &vx) {
    vi 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;
}
    for(int ans = 0; ans++) {
        fill(d.begin(), d.end(), 0);
        queue<int> 0;
        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 && q[x][y] != -1) {
                    if(d[v] == 0) {
                        p[y] = x;
                        d[v] = 2;
                        p[mate[y]] = y;
                        d[mate[y]] = 1;
                        Q.push(mate[y]);
                    }else if(d[y] == 1) {
                        vi vx = trace(x);
                        vi vy = trace(y);
                        if(vx.back() == vy.back()) {
                             contract(c, x, y, vx, vy);
                             Q.push(c);
```

return bridge;

```
16
```

```
p[c] = p[b[c][0]];
                                d[c] = 1;
                                C++;
                             }else {
                                 aug = true;
                                vx.insert(vx.begin(), y);
                                vy.insert(vy.begin(), x);
                                vi A = lift(vx);
                                vi B = lift(vy);
                                A.insert(A.end(), B.rbegin(), B
                                      .rend());
                                for(int i = 0; i < (int) A.size</pre>
                                     (); 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: $\mathcal{O}(E+V)$

```
Time: \mathcal{O}(E+V)
"../../stress-tests/utilities/template.h", "graphs_structures.h"
                                                      3103a2, 43 lines
template <typename T> vi find_scc(const digraph<T> &q, int &cnt
  digraph<T> g_rev = g.reverse();
  vi order;
  vector<bool> was(g.n, false);
  function < void (int) > dfs1 = [&] (int v) {
    was[v] = true;
    for (int id : g.g[v]) {
     auto &e = g.edges[id];
     int to = e.to;
     if (!was[to]) {
        dfs1(to);
   order.push_back(v);
  for (int i = 0; i < q.n; i++) {
   if (!was[i]) {
     dfs1(i);
  vector<int> c(g.n, -1);
  function<void(int) > dfs2 = [&](int v) {
    for (int id : g rev.g[v]) {
      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
BiComp.h
Description: Finds all biconnected components in an undirected graph
Time: \mathcal{O}(E+V)
template <typename T> vector<int> find_edge_biconnected(
    dfs_undigraph<T> &g, int &cnt) {
 vector<int> vertex_comp(q.n);
 cnt = 0;
 for (int i : q.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(q.n);
 cnt = 0;
 for (int i : g.order) {
    if (q.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[g.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 = (g.depth[x] > g.depth[y] ? x : y);
    edge_comp[id] = vertex_comp[z];
 return edge_comp;
template <typename T> vector<bool> find_bridges(dfs_undigraph<T</pre>
    > &a) {
 g.dfs_all();
 vector<bool> bridge(g.edges.size(), false);
 for (int i = 0; i < g.n; i++) {</pre>
   if (g.pv[i] != -1 && g.min_depth[i] == g.depth[i]) {
     bridge[g.pe[i]] = true;
```

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.

"../../stress-tests/utilities/template.h", "graphs.structures.h", "SCC.h" b9f520, 23 lines

```
struct twosat {
 digraph<int> g; int n;
  twosat(int _n) : g(digraph<int>(_n << 1)), n(_n) {}
  inline void add(int x, int value_x) { // (v[x] = value_x)
    assert(0 <= x && x < n); assert(0 <= value_x && value_x <=
    q.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 = value_y)
    assert(0 <= x && x < n && 0 <= y && y < n); assert(0 <=
         value_x && value_x <= 1 && 0 <= value_y && value_y <=</pre>
    q.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 \le x \&\& x \le n \&\& 0 \le y \&\& y \le n); assert(0 \le
         value x && value x <= 1 && 0 <= value y && value y <=
    g.add((x << 1) + (value_x ^ 1), (y << 1) + (value_y ^ 1));
         g.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 <= x && x < n && 0 <= y && y < n); assert(0 <=
         value && value <= 1);</pre>
    if (value) { add(x, 1, y, 1); add(x, 0, y, 0); }
    else { add_impl(x, 1, y, 1); add_impl(x, 0, y, 0); } }
  vi solve() { int cnt; vi c = find_scc(g, cnt); vi res(n);
    for (int i = 0; i < n; i++) {</pre>
      if (c[i << 1] == c[i << 1 ^ 1]) return vi();</pre>
      res[i] = (c[i << 1] < c[i << 1 ^ 1]);
    return res: }
};
```

EulerWalk.h

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

```
vi eulerWalk(vector<vector<pii>>>& gr, int nedges, int src=0) {
 int n = sz(qr);
  vi D(n), its(n), eu(nedges), ret, s = \{src\};
  D[src]++; // to allow Euler paths, not just cycles
  while (!s.empty()) {
   int x = s.back(), y, e, &it = its[x], end = sz(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 \mid | sz(ret) != nedges+1) return {};
  return {ret.rbegin(), ret.rend()};
```

DirectedMST.h

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

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

8a344f, 60 lines

```
struct Edge { int a, b; ll w; };
struct Node {
  Edge key;
 Node *1, *r;
  ll delta;
  void prop() {
   key.w += delta;
   if (1) 1->delta += delta;
   if (r) r->delta += delta;
   delta = 0;
  Edge top() { prop(); return key; }
Node *merge(Node *a, Node *b) {
 if (!a || !b) return a ?: b;
  a->prop(), b->prop();
  if (a->key.w > b->key.w) swap(a, b);
  swap(a->1, (a->r = merge(b, a->r)));
  return a:
void pop(Node*\& 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 * 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};
```

Diikstra.h

Description: Calculates shortest paths from s in a graph that might NOT have negative edge weights.

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

```
"../../stress-tests/utilities/template.h", "graphs_structures.h"
template <typename T> vector<T> dijkstra(const graph<T> &g, int
  assert(0 <= start && start < g.n);
 constexpr T MAXVALUE = numeric_limits<T>::max(); vector<T>
       dist(g.n, MAXVALUE);
 priority queue<pair<T, int>, vector<pair<T, int>>, greater
      pair<T, int>>> s; dist[start] = 0; s.emplace(dist[start
       1, start);
 while (!s.emptv()) {
   T expected = s.top().first; int i = s.top().second; s.pop()
    if (dist[i] != expected) continue;
    for (int id : q.q[i]) {
     auto &e = g.edges[id]; int to = e.from ^ e.to ^ i;
     if (dist[i] + e.cost < dist[to]) {</pre>
        dist[to] = dist[i] + e.cost; s.emplace(dist[to], to);
 return dist;
```

BellmanFord.h

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

```
"../../stress-tests/utilities/template.h", "graphs_structures.h"
template <typename T> vector<T> fordbellman(const graph<T> &g,
    int start) {
 assert(0 <= start && start < g.n);
 constexpr T MAXVALUE = numeric_limits<T>::max(); vector<T>
       dist(g.n, MAXVALUE);
 dist[start] = 0; vector<int> pv(g.n, -1);
 auto relax = [&]() {
    for (int id = 0; id < g.edges.size(); ++id) {</pre>
      auto& e = g.edges[id];
     if (dist[e.from] < MAXVALUE) {</pre>
        if (dist[e.to] > dist[e.from] + e.cost) {
          dist[e.to] = dist[e.from] + e.cost;
          pv[e.to] = e.from;
```

```
};
for (int i = 0; i + 1 < g.n; ++i) relax();</pre>
return dist:
/* find cycle with negative value
auto cur = dist; relax();
int\ broken = -1;
for (int \ i = 0; \ i < g.n; ++i) {
  if (cur[i] != dist[i]) \{ broken = i; break; \}
if (broken = -1) \{ return \{ \}; \}
for (int i = 0; i < g.n; ++i) broken = pv[broken];
vector<int> path{broken}; broken = pv[broken];
while (broken != path [0]) {
  path.pb(broken);
  broken = pv/broken;
reverse(all(path));
return path;
*/
```

Johnson.h

Description: calculates all-pairs shortest paths in a graph that might have negative edge weights.

```
Time: \mathcal{O}(N^3)
"../../stress-tests/utilities/template.h", "../../content/graph/BellmanFord.h",
"../../content/graph/dijkstra.h"
vector<vector<ll>> johnson(graph<ll> &g) {
 int n = q.n;
  ++q.n; q.q.resize(q.n);
  for (int i = 0; i < n; ++i) g.add(n, i, 0);
  vector<ll> d = fordbellman(g, n);
  g.edges.erase(g.edges.end() - n, g.edges.end());
  for (auto& adj : q.q) while (!adj.empty() && adj.back() >= q.
       edges.size()) { adj.pop_back(); }
  --q.n; q.q.resize(q.n);
  for(auto& e : q.edges) e.cost += d[e.from] - d[e.to];
  vector<vector<ll>> ans(n, vector<ll>(n, numeric_limits<ll>::
  for (int v = 0; v < n; ++v) {
    ans[v] = dijkstra(q, v);
    for (int u = 0; u < n; ++u) if (ans[v][u] != numeric_limits</pre>
         <11>::max()) ans[v][u] -= d[v] - d[u];
  return ans;
```

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: $\mathcal{O}(NM)$

```
7d20ae, 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;
DCPOffline.h
Description: DCP offline algorithm. Actually could be generalized to any
offline queries.
Time: \mathcal{O}\left(V\log^2(V)\right)
"../../stress-tests/utilities/template.h"
                                                      93309a, 87 lines
struct dsu save {
    int v, rnkv, u, rnku;
    dsu_save() = default;
    dsu_save(int _v, int _rnkv, int _u, int _rnku) : v(_v),
         rnkv( rnkv), u( u), rnku( rnku) {}
};
struct dsu with rollbacks {
    vector<int> p, rnk; int comps;
    stack<dsu save> op;
    dsu with rollbacks() = default;
    dsu_with_rollbacks(int n) {
        p.resize(n); rnk.resize(n);
        iota(p.begin(), p.end(), 0); rnk.assign(n, 0); comps =
    int find set(int v) {
        return (v == p[v]) ? v : find_set(p[v]);
   bool unite(int v, int u) {
        v = find_set(v); u = find_set(u);
        if (v == u) return false;
        if (rnk[v] > rnk[u]) swap(v, u);
        op.push(dsu_save(v, rnk[v], u, rnk[u])); p[v] = u;
        if (rnk[u] == rnk[v]) rnk[u]++;
        return true;
    void rollback() {
        if (op.empty()) return;
        dsu_save x = op.top(); op.pop(); comps++;
        p[x.v] = x.v; rnk[x.v] = x.rnkv;
        p[x.u] = x.u; rnk[x.u] = x.rnku;
};
struct query {
```

```
int v, u;
    bool united = true;
    query(int _v, int _u) : v(_v), u(_u) {}
struct OuervTree {
    vector<vector<query>> t;
    dsu with rollbacks dsu;
    int T:
    QueryTree() = default;
    QueryTree(int _T, int n) : T(_T) {
        dsu = dsu with rollbacks(n);
        t.resize(4 * T + 4);
    void add_to_tree(int v, int l, int r, int ul, int ur, query
        if (ul > ur) return;
        if (1 == u1 && r == ur) {
            t[v].push_back(q);
            return;
        int mid = (1 + r) / 2;
        add_to_tree(2 * v, 1, mid, ul, min(ur, mid), q);
        add_to_tree(2 * v + 1, mid + 1, r, max(ul, mid + 1), ur
    void add_query(query q, int 1, int r) { // edge(q.u, q.v)
         lives on segment [l, r]
        add_to_tree(1, 0, T - 1, 1, r, q);
    void dfs(int v, int 1, int r, vector<int>& ans) {
        for (query& q : t[v]) {
            q.united = dsu.unite(q.v, q.u);
        if (1 == r) ans[1] = dsu.comps; // here you can
             customize answers on queries
        else { int mid = (1 + r) / 2; dfs(2 * v, 1, mid, ans);
             dfs(2 * v + 1, mid + 1, r, ans);
        for (auto q : t[v]) {
            if (q.united) dsu.rollback();
    vector<int> solve() {
        vector<int> ans(T);
        dfs(1, 0, T - 1, ans);
        return ans;
};
6.5 Trees
DimssHLD.h
Description: Builds HLD
Time: \mathcal{O}(\log^2) probably actually \mathcal{O}(\log)
"../../stress-tests/utilities/template.h"
                                                     b652bc, 97 lines
static constexpr int mod = 1e9 + 7;
struct HLD {
    int n;
    vector<int> par, val, vale, tin, tout, sz, top;
    vector<vector<int>> up;
    int LG;
    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 \le 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(),
        ; ((a
    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 1 = tin[lca(v, u)];
    res.push_back(\{1, 1 + 1\});
    return res;
void mul_v(int v, int u, int delta) {
    for (auto [l, r] : get_v_path(v, u)) tree_v.update(l, r
         , delta);
void mul_e(int v, int u, int delta) {
    for (auto [1, r] : get_e_path(v, u)) tree_e.update(l, r
         , delta);
int get_v_sum(int v, int u) {
   11 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 res = 0; for (auto [1, r] : get_e_path(v, u)) res +=
         tree_e.get_sum(1, r);
    return res % mod;
```

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 $d_1 \geq \cdots \geq d_n$ exists iff $d_1 + \cdots + d_n$ is even and for every $k = 1 \dots n$,

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

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.)

```
template \langle class T \rangle 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); }
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 /S on the result of the cross product.

```
5a9b8c, 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;

return ((p-s)*d-(e-s)*t).dist()/d;

f4553b, 6 lines typedef Point < double > P; double segDist(P& s, P& e, P& p) { if (s==e) return (p-s).dist(); auto d = (e-s).dist2(), t = min(d, max(.0, (p-s).dot(e-s)));

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.

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

template < class P > vector < P > seqInter(P a, P b, P c, P d) {

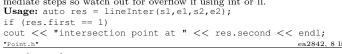


```
auto oa = c.cross(d, a), ob = c.cross(d, b),
     oc = a.cross(b, c), od = a.cross(b, d);
// Checks if intersection is single non-endpoint point.
if (sqn(oa) * sqn(ob) < 0 && sqn(oc) * sqn(od) < 0)
  return { (a * ob - b * oa) / (ob - oa) };
set<P> s:
if (onSegment(c, d, a)) s.insert(a);
if (onSegment(c, 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<ll> and the intersection point does not have integer coordinates. Products of three coordinates are used in inter- \sim \lambda mediate steps so watch out for overflow if using int or ll.



```
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"
                                                          856221, 9 lines
template<class P>
```

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

```
cc1cab, 3 lines
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:

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

2c1075, 6 lines

ccf6ab, 9 lines

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

Angle.h

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

 $\label{eq:Usage: vector Angle v = {w[0], w[0].t360() ...}; // sorted int j = 0; rep(i,0,n) { while (v[j] < v[i].t180()) ++j; } // sweeps j such that (j-i) represents the number of positively oriented triangles with vertices at 0 and i $6cf4ff, 35 lines$}$

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

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" f6bd8a, 11 lines typedef Point<double> P;

```
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 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.

CircPolvInter.h

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

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

"../../content/geometry/Point.h" 5661af, 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> P u = p + d * s, v = p + d * t;return arg(p,u) * r2 + u.cross(v)/2 + arg(v,q) * r2;**auto** sum = 0.0; rep(i, 0, sz(ps))sum += tri(ps[i] - c, ps[(i + 1) % sz(ps)] - c);return sum;

circumcircle.h

Description:

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



"Point.h" 6a6e48, 9 li

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

```
"circumcircle.h" bdf483, 17 lines
pair<P, double> mec(vector<P> ps) {
    shuffle(all(ps), mt19937(time(0)));
    P o = 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: vectorP v = {P{4,4}, P{1,2}, P{2,1}}; bool in = inPolygon(v, P{3, 3}, false); Time: O(n)
```

"Point.h", "OnSegment.h", "SegmentDistance.h" 52b242, 11 lines

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

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!

"Point.h"

7944c4, 6 lines

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

PolyCent.h

Description: Returns the center of mass for a polygon.

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

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;</pre>

Cut.h

Description:

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

Usage: vector<P> p = ...; p = polygonCut(p, P(0,0), P(1,0));"Point.h", "lineIntersection.h"



de6e85, 13 lines

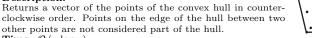
```
typedef Point<double> P;
vector<P> polygonCut(const vector<P>& poly, P s, P e) {
  vector<P> res;
  rep(i, 0, sz(poly)) {
   P cur = poly[i], prev = i ? poly[i-1] : poly.back();
   bool side = s.cross(e, cur) < 0;</pre>
   if (side != (s.cross(e, prev) < 0))
     res.push_back(lineInter(s, e, cur,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

Hull.h

Description:

clockwise order. Points on the edge of the hull between two other points are not considered part of the hull.

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





"Point.h"

```
typedef Point<11> 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.beqin(), h.beqin() + t - (t == 2 && h[0] == h[1])};
```

HalfplaneInt.h

Description: find halfplane intersection

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

c8d64d, 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;
    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);
};
```

```
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, dq[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;
        dg.push_back(H[i]);
    while (len > 2 && dq[0].out(inter(dq[len-1], dq[len-2]))) {
        dq.pop_back();
    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"
                                                       65e3f1, 12 lines
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)
```

res = $\max(\text{res}, \{(S[i] - S[j]).dist2(), \{S[i], S[j]\}\});$

for $(;; j = (j + 1) % n) {$

```
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"
                                                        fe002c, 14 lines
typedef Point<ll> P;
bool inHull(const vector<P>& 1, P p, bool strict = true) {
  int a = 1, b = sz(1) - 1, r = !strict;
 if (sz(1) < 3) return r && onSegment(1[0], 1.back(), p);</pre>
 if (sideOf(1[0], 1[a], 1[b]) > 0) swap(a, b);
 if (sideOf(1[0], 1[a], p) >= r || sideOf(1[0], 1[b], p) <= -r)
    return false;
  while (abs(a - b) > 1) {
    int c = (a + b) / 2;
    (sideOf(1[0], 1[c], p) > 0 ? b : a) = c;
  return 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)
```

```
"Point.h"
#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 (10 + 1 < hi) {
    int m = (lo + hi) / 2;
    if (extr(m)) return m;
    int 1s = cmp(1o + 1, 1o), ms = cmp(m + 1, m);
    (1s < ms \mid | (1s == ms \&\& 1s == cmp(1o, m)) ? hi : 1o) = 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;
      (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"
                                                      da1262, 17 lines
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; });
  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

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

```
Time: \mathcal{O}(n)
                                                                                   0c4338, 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: $\mathcal{O}(n)$ b21212, 12 lines

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

```
z[i] = i >= r ? 0 : min(r - i, z[i - 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, p[1][i] = longest odd (half rounded

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

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

MinRotation.h

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

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

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

```
"../../stress-tests/utilities/template.h"
                                                      870800, 211 lines
#define Size(x) (int)(x).size()
struct sparse {
    vector<vector<int>> st;
    sparse() { }
    sparse (const vector<int> &a) {
        int n = Size(a);
        int k = 0;
        while (1 << k < n)
            k++:
        st.resize(k + 1, vector<int>(n));
        copy(all(a), st[0].begin());
        for (int i = 1; i <= k; i++) {
            for (int j = 0; j + (1 << i) <= n; j++)</pre>
                st[i][j] = min(st[i-1][j], st[i-1][j+(1)]
                      << (i - 1)));
```

```
int getMin(int 1, int r) {
        int k = 31 - \underline{\text{builtin\_clz}(r - 1)};
        return min(st[k][l], st[k][r - (1 << k)]);</pre>
struct SuffixArray {
    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++) {</pre>
            int j = sa[pos[i] - 1];
            while (s[i + k] == s[j + 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';
```

23

Hashing

```
// compare two lms substring
inline bool not equal(const vector<int> &s, const vector<
    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:
        i++:
        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);
    types[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])
           tvpes[i] = 'L';
        el se
           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>
        int c = new_sym[i];
```

```
if (c !=-1) {
                pos_new_string[Size(new_string)] = i;
                new string.push back(c);
        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 \ x(all(s)+1), \ y(n), \ ws(max(n, lim)), \ rank(n);
    sa = lcp = y, iota(all(sa), 0);
    for (int j = 0, p = 0; p < n; j = max(1, j * 2), lim = p) {
      p = j, iota(all(y), n - j);
      rep(i,0,n) if (sa[i] >= j) y[p++] = sa[i] - j;
      fill(all(ws), 0);
      rep(i,0,n) ws[x[i]]++:
      rep(i,1,lim) ws[i] += ws[i-1];
      for(int \ i = n; \ i--;) \ sa[--ws[x[y[i]]]] = y[i];
      swap(x, y), p = 1, x[sa[0]] = 0;
      rep(i,1,n) \ a = sa[i-1], \ b = sa[i], \ x[b] =
        (y[a] = y[b] \otimes y[a+j] = y[b+j]) ? p-1 : p++;
    rep(i,1,n) \ rank[sa[i]] = i;
    for (int i = 0, j; i < n - 1; lcp[rank[i++]] = k)
      for (k \&\& k--, j = sa[rank[i]-1];
          s[i+k] = s[j+k]; k++);
};
Hashing.h
Description: creates hashes
Time: \mathcal{O}(N)
"../../stress-tests/utilities/template.h"
constexpr int HASH_MOD = MOD; constexpr 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]; }
  // if C++20 is available use auto operator\iff and bool
       operator = instead
  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 { return vals</pre>
       < other.vals; }
 bool operator>(const custom_hash &other) const { return vals
      > other.vals; }
 bool operator<=(const custom_hash &other) const { return vals</pre>
        <= other.vals; }
 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) 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) if ((1[i] -= r[i]) < 0) 1
       [i] += HASH_MOD; return 1;
custom hash operator*(custom hash 1, custom hash r) {
 for (int i = 0; i < HASH_SIZE; ++i) 1[i] = (11) 1[i] * r[i] %
       HASH MOD; return 1;
  static bool used = false; if (exchange(used, true)) { return;
  for (auto &u: base.vals) { u = BDIST(rng); }
 pows.emplace_back(make_hash(1));
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 *</pre>
        pows.back()); }
  custom_hash hash(int 1, int r) { int len = r + 1 - 1; extend(
      len); return cum[r + 1] - pows[len] * cum[1]; }
struct custom_int_hash {
 static uint64_t splitmix64(uint64_t x) {
```

```
x += 0x9e3779b97f4a7c15; x = (x ^ (x >> 30)) * 0
        xbf58476d1ce4e5b9; x = (x ^ (x >> 27)) * 0
        x94d049bb133111eb;
    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{};
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 Aho-Corasick 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 $\mathcal{O}(NM)$. _{13b7ef. 66 lines}

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

```
int &ed = N[n].next[i], y = N[prev].next[i];
      if (ed == -1) ed = y;
      else {
        N[ed].back = y;
        (N[ed].end == -1 ? N[ed].end : backp[N[ed].start])
          = N[y].end;
        N[ed].nmatches += N[y].nmatches;
        q.push(ed);
vi find(string word) {
  int n = 0;
  vi res; // ll count = 0;
  for (char c : word) {
   n = N[n].next[c - first];
    res.push_back(N[n].end);
    // count \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;
```

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: $\mathcal{O}(\log N)$ cf4189, 23 lines

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

IntervalCover.h

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

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

346<u>94c, 19 lines</u>

```
template < class T >
    vi cover(pair < T, T > G, vector < pair < T, T > I) {
        vi S(sz(I)), R;
        iota(all(S), 0);
        sort(all(S), [&](int a, int b) { return I[a] < I[b]; });
        T cur = G.first;
        int at = 0;
        while (cur < G.second) { // (A)
            pair < T, int > mx = make_pair(cur, -1);
        while (at < sz(I) && I[S[at]].first <= cur) {
                mx = max(mx, make_pair(I[S[at]].second, S[at]));
                at++;
        }
        if (mx.second == -1) return {};
        cur = mx.first;
        R.push_back(mx.second);
    }
    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) {
    q(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 q) {
 if (to <= from) return;</pre>
 int i = from; auto p = f(i), q = f(to-1);
 rec(from, to-1, f, g, i, p, q);
 g(i, to, q);
```

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);</pre>
```

```
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;
```

LIS.h

Description: Compute indices for the longest increasing subsequence.

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

c6efe4, 17 lines

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

FastKnapsack.h

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

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

c8bca4, 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++];</pre>
  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(i, 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;
```

9.3 Dynamic programming

KnuthDP.h

Description: When doing DP on intervals: $a[i][j] = \min_{i < k < j} (a[i][k] + a[i][j])$ a[k][i] + f(i,i), 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: $\mathcal{O}(N^2)$

LIS FastKnapsack KnuthDP DCDP SOSDP Knapsack

```
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 a[i] for i = L..R - 1.
Time: \mathcal{O}\left(\left(N + (hi - lo)\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) >> 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)
                                                            b219f4, 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)];
Knapsack.h
Description: Knapsack fast.
Time: \mathcal{O}\left(n^2Clogn/64\right) and \mathcal{O}\left(nC/64\right)
                                                           5e90f6, 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 \mid = 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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