

KTH Royal Institute of Technology

# Omogen Heap

Youngmin Park, Vedant Kumud, Ernest Lu

1	Contest
T	Contest
2	Mathematics
3	Data structures
4	Numerical
5	Number theory
6	Combinatorial
7	Graph
8	Geometry
9	Strings
10	Various
C	ontost (1)
	$\underline{\text{ontest}} \ (1)$
.ba	ashrc 4 li
ali xmo	-fsanitize=undefined,address'  las gc='g++ -Wall -Wconversion -Wfatal-errors -g -std=c++17  bdmap -e 'clear lock' -e 'keycode 66=less greater' #caps = coubleshoot.txt
_	52 li:
Wri Are Is Cou	e-submit:  te a few simple test cases if sample is not enough.  the limits close? If so, generate max cases.  the memory usage fine?  ald anything overflow?  te sure to submit the right file.
Pridate Are Car Rea Do Hav Any Cor Are Add Cre Go Go Go Exg As}	ang answer:  Int your solution! Print debug output, as well.  In your clearing all data structures between test cases?  In your algorithm handle the whole range of input?  In you handle all corner cases correctly?  If you understood the problem correctly?  If uninitialized variables?  If overflows?  If using N and M, i and j, etc.?  If you sure your algorithm works?  If special cases have you not thought of?  If you sure the STL functions you use work as you think?  If some assertions, maybe resubmit.  If some assertions, maybe resubmit.  If you sure the stcases to run your algorithm on.  If through the algorithm for a simple case.  If through this list again.  If your algorithm to a teammate.  If the teammate to look at your code.  If your output format correct? (including whitespace)  If your solution from the start or let a teammate do it.

Runtime error:

Have you tested all corner cases locally?

Any uninitialized variables?

Are you reading or writing outside the range of any vector? Any assertions that might fail?

Any possible division by 0? (mod 0 for example)

Any possible infinite recursion? Invalidated pointers or iterators?

Are you using too much memory?

Debug with resubmits (e.g. remapped signals, see Various).

Time limit exceeded:

Do you have any possible infinite loops?

What is the complexity of your algorithm?

Are you copying a lot of unnecessary data? (References)

How big is the input and output? (consider scanf) Avoid vector, map. (use arrays/unordered\_map)

What do your teammates think about your algorithm?

Memory limit exceeded:

What is the max amount of memory your algorithm should need? Are you clearing all data structures between test cases?

## Mathematics (2)

## 2.1 Geometry

## 2.1.1 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:  $r = \frac{A}{r}$ 

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

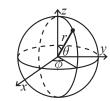
## 2.1.2 Quadrilaterals

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

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

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

#### 2.1.3 Spherical coordinates



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

#### 2.2Sums

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

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

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

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

#### 2.3

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

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

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

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

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

## Data structures (3)

OrderStatisticTree.h

**Description:** A set (not multiset!) with support for finding the n'th element, and finding the index of an element. To get a map, change null\_type. Time:  $\mathcal{O}(\log N)$ d41d8c, 16 lines

#include <bits/extc++.h> using namespace \_\_gnu\_pbds; template < class T> using Tree = tree<T, null\_type, less<T>, rb\_tree\_tag, tree\_order\_statistics\_node\_update>; void example() { Tree<int> t, t2; t.insert(8); auto it = t.insert(10).first;

```
assert(it == t.lower bound(9));
assert(t.order_of_key(10) == 1);
assert(t.order_of_key(11) == 2);
assert(*t.find_by_order(0) == 8);
t.join(t2); // assuming T < T2 or T > T2, merge t2 into t
```

### HashMap.h

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

#include <bits/extc++.h> // To use most bits rather than just the lowest ones: struct chash { // large odd number for C const uint64\_t C = 11(4e18 \* acos(0)) | 71; 11 operator()(11 x) const { return \_\_builtin\_bswap64(x \* C);

\_\_gnu\_pbds::gp\_hash\_table<11, int, chash> h({}, {}, {}, {}, {1

#### SegmentTree.h

<< 16});

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

d41d8c, 19 lines struct Tree { typedef int T; static constexpr T unit = INT\_MIN; T f(T a, T b) { return max(a, b); } //  $(any \ associative \ fn)$ vector<T> s; int n; Tree(int n = 0, T def = unit) : s(2\*n, def), n(n) {} void update(int pos, T val) { for (s[pos += n] = val; pos /= 2;) s[pos] = f(s[pos \* 2], s[pos \* 2 + 1]);T query (int b, int e) { // query [b, e)T ra = unit, rb = unit; for  $(b += n, e += n; b < e; b /= 2, e /= 2) {$ **if** (b % 2) ra = f(ra, s[b++]);**if** (e % 2) rb = f(s[--e], rb);return f(ra, rb); };

#### LazySeg.h

**Description:** Lazy segtree with templated info

d41d8c, 115 lines

```
template <class Info, class Tag> struct LazySegmentTree {
  std::vector<Info> info;
  std::vector<Tag> tag;
  LazySegmentTree(): n(0) {}
  LazySegmentTree(int n_, Info v_ = Info()) { init(n_, v_); }
  template <class T> LazySegmentTree(std::vector<T> init_) {
      init(init_); }
  void init(int n_, Info v_ = Info()) { init(std::vector(n_, v_
  template <class T> void init(std::vector<T> init_) {
   n = init_.size();
    info.assign(4 << std::__lg(n), Info());
    tag.assign(4 << std::__lg(n), Tag());
    std::function<void(int, int, int)> build = [&](int p, int 1
        , int r) {
     if (r - 1 == 1) {
       info[p] = init_[1];
        return;
```

```
int m = (1 + r) / 2;
    build(2 * p, 1, m);
    build(2 * p + 1, m, r);
    pull(p);
  };
  build(1, 0, n);
void pull(int p) { info[p] = info[2 * p] + info[2 * p + 1]; }
void apply(int p, const Tag &v) {
  info[p].apply(v);
  tag[p].apply(v);
void push(int p) {
  apply(2 * p, tag[p]);
  apply (2 * p + 1, tag[p]);
  tag[p] = Tag();
void modify(int p, int 1, int r, int x, const Info &v) {
  if (r - 1 == 1) {
    info[p] = v;
    return;
  int m = (1 + r) / 2;
  push(p);
  if (x < m)  {
    modify(2 * p, 1, m, x, v);
    modify(2 * p + 1, m, r, x, v);
  pull(p);
void modify(int p, const Info &v) { modify(1, 0, n, p, v); }
Info rangeQuery(int p, int l, int r, int x, int y) {
  if (1 >= y || r <= x) {
    return Info();
  if (1 >= x && r <= y) {
    return info[p];
  int m = (1 + r) / 2;
  push (p);
  return rangeQuery(2 * p, 1, m, x, y) + rangeQuery(2 * p +
      1, m, r, x, v);
Info rangeQuery(int 1, int r) { return rangeQuery(1, 0, n, 1,
void rangeApply(int p, int 1, int r, int x, int y, const Tag
  if (1 >= y || r <= x) {
    return;
  if (1 >= x && r <= y) {</pre>
    apply(p, v);
    return;
  int m = (1 + r) / 2;
  push (p);
  rangeApply(2 * p, 1, m, x, y, v);
  rangeApply(2 * p + 1, m, r, x, y, v);
  pull(p);
void rangeApply(int 1, int r, const Tag &v) {
  return rangeApply(1, 0, n, 1, r, v);
template <class F> int findFirst(int p, int 1, int r, int x,
     int y, F pred) {
  if (1 >= y || r <= x || !pred(info[p])) {</pre>
    return -1;
```

```
if (r - 1 == 1) {
      return 1;
    int m = (1 + r) / 2;
   push(p);
    int res = findFirst(2 * p, 1, m, x, y, pred);
    if (res == -1) {
      res = findFirst(2 * p + 1, m, r, x, y, pred);
    return res;
 template <class F> int findFirst(int 1, int r, F pred) {
    return findFirst(1, 0, n, 1, r, pred);
  template <class F> int findLast(int p, int 1, int r, int x,
      int y, F pred) {
    if (1 >= y || r <= x || !pred(info[p])) {</pre>
      return -1;
    if (r - 1 == 1) {
      return 1;
    int m = (1 + r) / 2;
    push(p);
    int res = findLast(2 * p + 1, m, r, x, y, pred);
    if (res == -1) {
      res = findLast(2 * p, 1, m, x, y, pred);
    return res:
  template <class F> int findLast(int 1, int r, F pred) {
    return findLast(1, 0, n, 1, r, pred);
};
```

```
PersistentSegmentTree.h
Description: Persistent segtree with allocation arena
                                                      d41d8c, 65 lines
template <typename T, int SZ> struct PST {
 struct Node {
   T val;
   int c[2];
   Node() {
     val = INF;
      c[0] = c[1] = 0;
 static const int LIM = 1.5e7;
 Node d[LTMl:
 int nxt = 0;
 int copy(int t) {
   d[nxt] = d[t];
   return nxt++;
 T comb (const T &a, const T &b) { return min(a, b); }
 void pull(int c) { d[c].val = comb(d[d[c].c[0]].val, d[d[c].c
       [1]].val); }
 T query(int lo, int hi, int t, int l, int r) {
   if (lo >= r || hi <= l)
      return INF;
   if (lo <= l && r <= hi)
      return d[t].val;
    int m = (1 + r) / 2;
   T lef = query(lo, hi, d[t].c[0], l, m);
   T \text{ rig} = \text{query(lo, hi, d[t].c[1], m, r)};
    return comb(lef, rig);
 int upd(int i, const T &v, int t, int 1, int r) {
```

return true;

};

#### UnionFindRollback SubMatrix Matrix LineContainer Treap

```
int x = copy(t);
    if (r - 1 == 1) {
      d[x].val = min(d[x].val, v);
      return x;
    int m = (1 + r) / 2;
    if (i < m) {
     d[x].c[0] = upd(i, v, d[x].c[0], l, m);
     d[x].c[1] = upd(i, v, d[x].c[1], m, r);
   pull(x);
    return x;
  int build(const vector<T> &a, int 1, int r) {
    int c = nxt++;
    if (r - 1 == 1) {
     if (1 < (int)a.size())</pre>
       d[c].val = a[1];
      return c;
    int m = (1 + r) / 2;
    d[c].c[0] = build(a, 1, m);
    d[c].c[1] = build(a, m, r);
   pull(c);
    return c;
  vector<int> rts;
  void update_time(int i, const T &v) {
    debug(rts);
    rts.pb(upd(i, v, rts.back(), 0, SZ));
  void build(const vector<T> &a) { rts.pb(build(a, 0, SZ)); }
  T query_time(int ti, int lo, int hi) { return query(lo, hi,
       rts[ti], 0, SZ); }
PST<int, 1 << 19> pst;
UnionFindRollback.h
Description: Disjoint-set data structure with undo. If undo is not needed,
skip st, time() and rollback().
Usage: int t = uf.time(); ...; uf.rollback(t);
Time: \mathcal{O}(\log(N))
                                                      d41d8c, 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]); }
  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;
```

```
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)
                                                         d41d8c, 13 lines
template<class T>
struct SubMatrix {
  vector<vector<T>> p;
  SubMatrix(vector<vector<T>>& v) {
    int R = sz(v), C = sz(v[0]);
    p.assign(R+1, vector<T>(C+1));
    rep(r, 0, R) rep(c, 0, C)
      p[r+1][c+1] = v[r][c] + p[r][c+1] + p[r+1][c] - p[r][c];
 T sum(int u, int 1, int d, int r) {
    return p[d][r] - p[d][l] - p[u][r] + p[u][l];
};
Matrix.h
Description: Basic operations on square matrices.
Usage: Matrix<int, 3> A;
A.d = \{\{\{1,2,3\}\}, \{\{4,5,6\}\}, \{\{7,8,9\}\}\}\};
vector < int > vec = \{1, 2, 3\};
vec = (A^N) * vec;
                                                         d41d8c, 26 lines
template < class T, int N> struct Matrix {
  typedef Matrix M;
  arrav<arrav<T, N>, N> d{};
  M operator*(const M& m) const {
    rep(i,0,N) rep(j,0,N)
      rep(k, 0, N) \ a.d[i][j] += d[i][k]*m.d[k][j];
    return a;
  vector<T> operator*(const vector<T>& vec) const {
    vector<T> ret(N);
    rep(i, 0, N) rep(j, 0, N) ret[i] += d[i][j] * vec[j];
    return ret;
 M operator^(ll p) const {
    assert (p >= 0);
    M a, b(*this);
    rep(i, 0, N) \ a.d[i][i] = 1;
    while (p) {
      if (p&1) a = a*b;
      b = b*b;
      p >>= 1;
    return a;
};
LineContainer.h
Description: Container where you can add lines of the form kx+m, and
query maximum values at points x. Useful for dynamic programming ("con-
vex hull trick").
Time: \mathcal{O}(\log N)
                                                         d41d8c, 30 lines
struct Line {
 mutable 11 k, m, p;
  bool operator<(const Line& o) const { return k < o.k; }</pre>
  bool operator<(ll x) const { return p < x; }</pre>
struct LineContainer : multiset<Line, less<>>> {
  // (for doubles, use inf = 1/.0, div(a,b) = a/b)
  static const 11 inf = LLONG MAX;
```

```
ll div(ll a, ll b) { // floored division
    return a / b - ((a ^ b) < 0 && a % b); }
  bool isect(iterator x, iterator y) {
    if (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(v));
 11 query(11 x) {
    assert(!empty());
    auto 1 = *lower_bound(x);
    return 1.k * x + 1.m;
};
Treap.h
Description: A short self-balancing tree. It acts as a sequential container
with log-time splits/joins, and is easy to augment with additional data.
Time: \mathcal{O}(\log N)
                                                      d41d8c, 55 lines
struct Node {
 Node *1 = 0, *r = 0;
 int val, v, c = 1;
 Node(int val) : val(val), y(rand()) {}
 void recalc();
int cnt(Node* n) { return n ? n->c : 0; }
void Node::recalc() { c = cnt(1) + cnt(r) + 1; }
template < class F > void each (Node * n, F f) {
 if (n) { each(n->1, f); f(n->val); each(n->r, f); }
pair<Node*, Node*> split(Node* n, int k) {
 if (!n) return {};
 if (cnt(n->1) >= k) { // "n=>val>= k" for lower_bound(k)
    auto pa = split(n->1, k);
    n->1 = pa.second;
    n->recalc();
    return {pa.first, n};
    auto pa = split(n->r, k - cnt(n->1) - 1); // and just "k"
    n->r = pa.first;
    n->recalc();
    return {n, pa.second};
Node* merge(Node* 1, Node* r) {
 if (!1) return r;
 if (!r) return 1;
 if (1->y > r->y) {
   1->r = merge(1->r, r);
   l->recalc();
    return 1;
  } else {
    r->1 = merge(1, r->1);
    r->recalc();
    return r;
```

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

#### FenwickTree.h

**Description:** Computes partial sums a[0] + a[1] + ... + a[pos - 1], and updates single elements a[i], taking the difference between the old and new

**Time:** Both operations are  $\mathcal{O}(\log N)$ .

```
d41d8c, 22 lines
struct FT {
 vector<ll> s;
  FT(int n) : s(n) {}
  void update(int pos, 11 dif) { // a[pos] \leftarrow dif
    for (; pos < sz(s); pos |= pos + 1) s[pos] += dif;</pre>
  11 query (int pos) { // sum of values in [0, pos)
    11 \text{ res} = 0;
    for (; pos > 0; pos &= pos - 1) res += s[pos-1];
    return res;
  int lower_bound(11 sum) \{// min \ pos \ st \ sum \ of \ [0, \ pos] >= sum
    // Returns n if no sum is >= sum, or -1 if empty sum is.
    if (sum \le 0) return -1;
   int pos = 0;
    for (int pw = 1 << 25; pw; pw >>= 1) {
      if (pos + pw <= sz(s) && s[pos + pw-1] < sum)</pre>
        pos += pw, sum -= s[pos-1];
    return pos;
```

#### FenwickTree2d.h

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

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

```
"FenwickTree.h"
                                                     d41d8c, 22 lines
struct FT2 {
  vector<vi> ys; vector<FT> ft;
 FT2(int limx) : ys(limx) {}
  void fakeUpdate(int x, int y) {
   for (; x < sz(ys); x |= x + 1) ys[x].push_back(y);
  void init() {
   for (vi& v : ys) sort(all(v)), ft.emplace_back(sz(v));
  int ind(int x, int y) {
   return (int) (lower_bound(all(ys[x]), y) - ys[x].begin()); }
  void update(int x, int y, ll dif) {
   for (; x < sz(ys); x | = x + 1)
      ft[x].update(ind(x, y), dif);
  11 query(int x, int y) {
   11 sum = 0;
    for (; x; x &= x - 1)
     sum += ft[x-1].query(ind(x-1, y));
    return sum;
```

```
};
RMQ.h
Description: Range Minimum Queries on an array. Returns min(V[a], V[a
+1], ... V[b - 1]) in constant time.
Usage: RMO rmg(values);
rmq.query(inclusive, exclusive);
Time: \mathcal{O}(|V|\log|V|+Q)
                                                         d41d8c, 16 lines
template<class T>
struct RMQ {
 vector<vector<T>> jmp;
 RMQ(const vector<T>& V) : jmp(1, V) {
    for (int pw = 1, k = 1; pw * 2 <= sz(V); pw *= 2, ++k) {
      jmp.emplace_back(sz(V) - pw * 2 + 1);
      rep(j, 0, sz(jmp[k]))
        jmp[k][j] = min(jmp[k - 1][j], jmp[k - 1][j + pw]);
 T query(int a, int b) {
    assert (a < b); // or return inf if a == b
    int dep = 31 - __builtin_clz(b - a);
    return min(jmp[dep][a], jmp[dep][b - (1 << dep)]);</pre>
};
MoQueries.h
Description: Answer interval or tree path queries by finding an approxi-
mate TSP through the queries, and moving from one query to the next by
adding/removing points at the ends. If values are on tree edges, change step
to add/remove the edge (a, c) and remove the initial add call (but keep in).
Time: \mathcal{O}\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> 0) {
```

```
int L = 0, R = 0, blk = 350; // \sim N/sqrt(Q)
 vi s(sz(Q)), res = s;
#define K(x) pii(x.first/blk, x.second ^ -(x.first/blk & 1))
 iota(all(s), 0);
 sort(all(s), [\&](int s, int t){ return K(Q[s]) < K(Q[t]); });
 for (int qi : s) {
   pii q = Q[qi];
    while (L > q.first) add(--L, 0);
    while (R < q.second) add(R++, 1);</pre>
    while (L < q.first) del(L++, 0);
    while (R > q.second) del(--R, 1);
    res[qi] = calc();
 return res;
vi moTree(vector<array<int, 2>> Q, vector<vi>& 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;
```

## Numerical (4)

## 4.1 Polynomials and recurrences

#### Polynomial.h

d41d8c, 17 lines

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

#### PolyRoots.h

```
Description: Finds the real roots to a polynomial.
Usage: polyRoots(\{\{2,-3,1\}\},-1e9,1e9) // solve x^2-3x+2=0
Time: \mathcal{O}\left(n^2\log(1/\epsilon)\right)
"Polynomial.h"
                                                                     d41d8c, 23 lines
```

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

#### PolyInterpolate.h

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

d41d8c, 13 lines

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

#### BerlekampMassev.h

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

```
Usage: berlekampMassey({0, 1, 1, 3, 5, 11}) // {1, 2}
Time: \mathcal{O}\left(N^2\right)
```

```
"../number-theory/ModPow.h"
                                                     d41d8c, 20 lines
vector<ll> berlekampMassey(vector<ll> s) {
 int n = sz(s), L = 0, m = 0;
 vector<ll> C(n), B(n), T;
 C[0] = B[0] = 1;
 11 b = 1;
  rep(i, 0, n) \{ ++m;
   11 d = s[i] % mod;
   rep(j, 1, L+1) d = (d + C[j] * s[i - j]) % mod;
   if (!d) continue;
   T = C; 11 coef = d * modpow(b, mod-2) % mod;
   rep(j,m,n) C[j] = (C[j] - coef * B[j - m]) % mod;
   if (2 * L > i) continue;
   L = i + 1 - L; B = T; b = d; m = 0;
 C.resize(L + 1); C.erase(C.begin());
  for (11& x : C) x = (mod - x) % mod;
 return C;
```

#### LinearRecurrence.h

**Description:** Generates the k'th term of an n-order linear recurrence  $S[i] = \sum_{j} S[i-j-1]tr[j]$ , given  $S[0... \ge n-1]$  and tr[0...n-1]. Faster than matrix multiplication. Useful together with Berlekamp-Massey. Usage: linearRec( $\{0, 1\}, \{1, 1\}, k$ ) // k'th Fibonacci number Time:  $\mathcal{O}\left(n^2 \log k\right)$ d41d8c, 26 lines

```
typedef vector<ll> Poly;
11 linearRec(Poly S, Poly tr, 11 k) {
  int n = sz(tr);
  auto combine = [&](Poly a, Poly b) {
   Poly res(n \star 2 + 1);
   rep(i, 0, n+1) rep(j, 0, n+1)
     res[i + j] = (res[i + j] + a[i] * b[j]) % mod;
   for (int i = 2 * n; i > n; --i) rep(j,0,n)
     res[i - 1 - j] = (res[i - 1 - j] + res[i] * tr[j]) % mod;
    res.resize(n + 1);
   return res;
```

```
Poly pol(n + 1), e(pol);
pol[0] = e[1] = 1;
for (++k; k; k /= 2) {
  if (k % 2) pol = combine(pol, e);
  e = combine(e, e);
11 \text{ res} = 0:
rep(i, 0, n) res = (res + pol[i + 1] * S[i]) % mod;
return res;
```

## 4.2 Optimization

#### GoldenSectionSearch.h

**Description:** Finds the argument minimizing the function f in the interval [a,b] assuming f is unimodal on the interval, i.e. has only one local minimum and no local maximum. The maximum error in the result is eps. Works equally well for maximization with a small change in the code. See Ternary-Search.h in the Various chapter for a discrete version.

```
Usage: double func(double x) { return 4+x+.3*x*x; }
double xmin = qss(-1000, 1000, func);
Time: \mathcal{O}\left(\log((b-a)/\epsilon)\right)
```

d41d8c, 14 lines

```
double qss(double a, double b, double (*f)(double)) {
 double r = (sqrt(5)-1)/2, eps = 1e-7;
 double x1 = b - r*(b-a), x2 = a + r*(b-a);
 double f1 = f(x1), f2 = f(x2);
 while (b-a > eps)
   if (f1 < f2) { //change to > to find maximum
     b = x2; x2 = x1; f2 = f1;
     x1 = b - r*(b-a); f1 = f(x1);
     a = x1; x1 = x2; f1 = f2;
     x2 = a + r*(b-a); f2 = f(x2);
 return a:
```

#### HillClimbing.h

Description: Poor man's optimization for unimodal functions<sub>d41d8c, 14 lines</sub>

```
typedef array<double, 2> P;
template < class F > pair < double, P > hillClimb (P start, F f) {
 pair<double, P> cur(f(start), start);
 for (double jmp = 1e9; jmp > 1e-20; jmp /= 2) {
    rep(j, 0, 100) rep(dx, -1, 2) rep(dy, -1, 2) {
      P p = cur.second;
      p[0] += dx * jmp;
      p[1] += dy * jmp;
      cur = min(cur, make_pair(f(p), p));
 return cur;
```

#### Integrate.h

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

```
template<class F>
double quad(double a, double b, F f, const int n = 1000) {
  double h = (b - a) / 2 / n, v = f(a) + f(b);
  rep(i,1,n*2)
   v += f(a + i*h) * (i&1 ? 4 : 2);
```

```
return v * h / 3;
IntegrateAdaptive.h
Description: Fast integration using an adaptive Simpson's rule.
Usage: double sphereVolume = quad(-1, 1, [](double x) {
return quad(-1, 1, [&] (double y)
return quad(-1, 1, [\&] (double z) {
return \bar{x} \times x + y \times y + z \times z < 1; \}); \}); \}); \}
                                                           d41d8c, 15 lines
typedef double d;
#define S(a,b) (f(a) + 4*f((a+b) / 2) + f(b)) * (b-a) / 6
template <class F>
d rec(F& f, da, db, deps, dS) {
  dc = (a + b) / 2;
  d S1 = S(a, c), S2 = S(c, b), T = S1 + S2;
  if (abs(T - S) <= 15 * eps || b - a < 1e-10)</pre>
    return T + (T - S) / 15;
  return rec(f, a, c, eps / 2, S1) + rec(f, c, b, eps / 2, S2);
template < class F>
d \text{ quad}(d \text{ a, } d \text{ b, } F \text{ f, } d \text{ eps} = 1e-8)  {
  return rec(f, a, b, eps, S(a, b));
```

#### Simplex.h

**Description:** Solves a general linear maximization problem: maximize  $c^T x$ subject to  $Ax \leq b$ ,  $x \geq 0$ . Returns -inf if there is no solution, inf if there are arbitrarily good solutions, or the maximum value of  $c^T x$  otherwise. The input vector is set to an optimal x (or in the unbounded case, an arbitrary solution fulfilling the constraints). Numerical stability is not guaranteed. For better performance, define variables such that x = 0 is viable.

```
Usage: vvd A = \{\{1, -1\}, \{-1, 1\}, \{-1, -2\}\};
vd b = \{1, 1, -4\}, c = \{-1, -1\}, x;
T val = LPSolver(A, b, c).solve(x);
```

b[s] = a[s] \* inv2;

rep(j,0,n+2) **if** (j != s) D[r][j] \*= inv;rep(i,0,m+2) if (i != r)  $D[i][s] \star = -inv;$ 

**Time:**  $\mathcal{O}(NM * \#pivots)$ , where a pivot may be e.g. an edge relaxation.

```
\mathcal{O}(2^n) in the general case.
                                                      d41d8c, 68 lines
typedef double T; // long double, Rational, double + mod<P>...
typedef vector<T> vd;
typedef vector<vd> vvd;
const T eps = 1e-8, inf = 1/.0;
#define MP make pair
#define ltj(X) if(s == -1 || MP(X[j], N[j]) < MP(X[s], N[s])) s=j
struct LPSolver {
  int m, n;
  vi N, B;
  vvd D:
  LPSolver (const vvd& A, const vd& b, const vd& c) :
    m(sz(b)), n(sz(c)), N(n+1), B(m), D(m+2), vd(n+2)) {
      rep(i, 0, m) rep(j, 0, n) D[i][j] = A[i][j];
      rep(i,0,m) { B[i] = n+i; D[i][n] = -1; D[i][n+1] = b[i]; }
      rep(j,0,n) \{ N[j] = j; D[m][j] = -c[j]; \}
      N[n] = -1; D[m+1][n] = 1;
  void pivot(int r, int s) {
    T *a = D[r].data(), inv = 1 / a[s];
    rep(i, 0, m+2) if (i != r && abs(D[i][s]) > eps) {
      T *b = D[i].data(), inv2 = b[s] * inv;
      rep(j,0,n+2) b[j] -= a[j] * inv2;
```

```
D[r][s] = inv;
    swap(B[r], N[s]);
  bool simplex(int phase) {
    int x = m + phase - 1;
    for (;;) {
     int s = -1;
      rep(j,0,n+1) if (N[j] != -phase) ltj(D[x]);
      if (D[x][s] >= -eps) return true;
      int r = -1:
      rep(i,0,m) {
       if (D[i][s] <= eps) continue;</pre>
       if (r == -1 || MP(D[i][n+1] / D[i][s], B[i])
                     < MP(D[r][n+1] / D[r][s], B[r])) r = i;
      if (r == -1) return false;
     pivot(r, s);
 T solve(vd &x) {
    int r = 0;
    rep(i,1,m) if (D[i][n+1] < D[r][n+1]) r = i;
    if (D[r][n+1] < -eps) {
     pivot(r, n);
      if (!simplex(2) || D[m+1][n+1] < -eps) return -inf;</pre>
      rep(i, 0, m) if (B[i] == -1) {
       int s = 0;
        rep(j,1,n+1) ltj(D[i]);
        pivot(i, s);
   bool ok = simplex(1); x = vd(n);
    rep(i,0,m) if (B[i] < n) x[B[i]] = D[i][n+1];
    return ok ? D[m][n+1] : inf;
};
```

#### 4.3 Matrices

#### Determinant.h

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

```
duble 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)
```

d41d8c, 18 lines

```
const 11 mod = 12345;
11 det(vector<vector<11>>& a) {
  int n = sz(a); 11 ans = 1;
  rep(i,0,n) {
```

```
rep(j,i+1,n) {
    while (a[j][i] != 0) { // gcd step
        ll t = a[i][i] / a[j][i];
        if (t) rep(k,i,n)
            a[i][k] = (a[i][k] - a[j][k] * t) % mod;
        swap(a[i], a[j]);
        ans *= -1;
        }
        ans = ans * a[i][i] % mod;
        if (!ans) return 0;
    }
    return (ans + mod) % mod;
}
```

#### SolveLinear.h

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

```
typedef vector<double> vd;
const double eps = 1e-12;
int solveLinear(vector<vd>& A, vd& b, vd& x) {
 int n = sz(A), m = sz(x), rank = 0, br, bc;
 if (n) assert(sz(A[0]) == m);
 vi col(m); iota(all(col), 0);
 rep(i,0,n) {
   double v, bv = 0;
   rep(r,i,n) rep(c,i,m)
     if ((v = fabs(A[r][c])) > bv)
       br = r, bc = c, bv = v;
   if (bv <= eps) {
     rep(j,i,n) if (fabs(b[j]) > eps) return -1;
    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[i][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:

```
fail:; }
```

#### SolveLinearBinary.h

typedef bitset<1000> bs;

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

```
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:  $\mathcal{O}\left(n^3\right)$ 

```
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];
```

d41d8c, 16 lines

```
rep(k,0,n) tmp[j][k] \rightarrow f*tmp[i][k];
  rep(j,i+1,n) A[i][j] /= v;
 rep(j,0,n) tmp[i][j] /= v;
 A[i][i] = 1;
for (int i = n-1; i > 0; --i) rep(j, 0, i) {
 double v = A[j][i];
 rep(k,0,n) tmp[j][k] -= v*tmp[i][k];
rep(i,0,n) \ rep(j,0,n) \ A[col[i]][col[j]] = tmp[i][j];
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

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

Fails if the solution is not unique.

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

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

d41d8c, 26 lines

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

#### 4.4 Fourier transforms

FastFourierTransform.h

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

```
d41d8c, 35 lines
typedef complex<double> C:
typedef vector<double> vd;
void fft(vector<C>& a) {
 int n = sz(a), L = 31 - builtin clz(n);
 static vector<complex<long double>> R(2, 1);
 static vector<C> rt(2, 1); // (^ 10% faster if double)
 for (static int k = 2; k < n; k *= 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 - __builtin_clz(sz(res)), n = 1 << L;</pre>
 vector<C> in(n), out(n);
 copy(all(a), begin(in));
 rep(i,0,sz(b)) in[i].imag(b[i]);
 fft(in);
 for (C& x : in) x *= x;
 rep(i, 0, n) out[i] = in[-i & (n - 1)] - conj(in[i]);
 rep(i, 0, sz(res)) res[i] = imag(out[i]) / (4 * n);
 return res;
```

#### FastFourierTransformMod.h

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

**Time:**  $\mathcal{O}(N \log N)$ , where N = |A| + |B| (twice as slow as NTT or FFT) "FastFourierTransform.h" d41d8c, 22 lines

```
typedef vector<ll> v1;
template<int M> v1 convMod(const v1 &a, const v1 &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)) {
```

```
ll av = ll(real(outl[i])+.5), cv = ll(imag(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;
```

NumberTheoreticTransform.h

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

```
"../number-theory/ModPow.h"
const 11 mod = (119 << 23) + 1, root = 62; // = 998244353
```

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

#### FastSubsetTransform.h

Description: Transform to a basis with fast convolutions of the form  $c[z] \, = \, \sum_{z=x \oplus y} a[x] \cdot b[y], \text{ where } \oplus \text{ is one of AND, OR, XOR.}$  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) {
     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
                                             // XOR
       pii(u + v, u - v);
```

```
if (inv) for (int& x : a) \times /= sz(a); // XOR only
vi conv(vi a, vi b) {
 FST(a, 0); FST(b, 0);
 rep(i, 0, sz(a)) a[i] *= b[i];
 FST(a, 1); return a;
```

## Number theory (5)

## 5.1 Modular arithmetic

#### Modular Arithmetic.h

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

```
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^(ll e) {
    if (!e) return Mod(1);
   Mod r = *this ^ (e / 2); r = r * r;
   return e&1 ? *this * r : r;
};
```

#### ModInverse.h

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

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

#### ModPow.h

d41d8c, 8 lines

```
const 11 mod = 1000000007; // faster if const
ll modpow(ll b, ll e) {
 11 \text{ ans} = 1;
  for (; e; b = b * b % mod, e /= 2)
   if (e & 1) ans = ans * b % mod;
  return ans;
```

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

```
d41d8c, 11 lines
11 modLog(ll a, ll b, ll m) {
 unordered_map<11, 11> A;
  while (j <= n && (e = f = e * a % m) != b % m)
  A[e * b % m] = j++;
  if (e == b % m) return j;
 if (__gcd(m, e) == __gcd(m, b))
   rep(i,2,n+2) if (A.count(e = e * f % m))
     return n * i - A[e];
```

```
return -1;
```

#### ModSum.h

**Description:** Sums of mod'ed arithmetic progressions.

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

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

```
d41d8c, 16 lines
typedef unsigned long long ull;
ull sumsq(ull to) { return to /2 * ((to-1) | 1); }
ull divsum(ull to, ull c, ull k, ull m) {
 ull res = k / m * sumsq(to) + c / m * to;
 k %= m; c %= m;
 if (!k) return res;
 ull to2 = (to * k + c) / m;
 return res + (to - 1) * to2 - divsum(to2, m-1 - c, m, k);
11 modsum(ull to, 11 c, 11 k, 11 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) {
 11 ret = a * b - M * ull(1.L / M * a * b);
 return ret + M * (ret < 0) - M * (ret >= (11)M);
ull modpow(ull b, ull e, ull mod) {
 ull ans = 1;
 for (; e; b = modmul(b, b, mod), e /= 2)
   if (e & 1) ans = modmul(ans, b, mod);
```

#### ModSart.h

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

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

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

```
5.2 Primality
```

#### FastEratosthenes.h

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

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

#### MillerRabin.h

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

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

```
"ModMulLL.h"
                                                       d41d8c, 12 lines
bool isPrime(ull n) {
 if (n < 2 || n % 6 % 4 != 1) return (n | 1) == 3;</pre>
 ull A[] = \{2, 325, 9375, 28178, 450775, 9780504, 1795265022\},
      s = \underline{builtin\_ctzll(n-1)}, d = n >> s;
  for (ull a : A) { // ^ count trailing zeroes
    ull p = modpow(a%n, d, n), i = s;
    while (p != 1 && p != n - 1 && a % n && i--)
      p = modmul(p, p, n);
    if (p != n-1 && i != s) return 0;
 return 1;
```

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

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

```
"ModMulLL.h", "MillerRabin.h"
                                                     d41d8c, 18 lines
ull pollard(ull n) {
 auto f = [n] (ull x) { return modmul(x, x, n) + 1; };
 ull x = 0, y = 0, t = 30, prd = 2, i = 1, q;
  while (t++ % 40 || __gcd(prd, n) == 1) {
   if (x == y) x = ++i, y = f(x);
    if ((q = modmul(prd, max(x,y) - min(x,y), n))) prd = q;
    x = f(x), y = f(f(y));
 return __gcd(prd, n);
vector<ull> factor(ull n) {
 if (n == 1) return {};
 if (isPrime(n)) return {n};
 ull x = pollard(n);
```

```
auto 1 = factor(x), r = factor(n / x);
1.insert(1.end(), all(r));
return 1;
```

## 5.3 Divisibility

euclid.h

KTH

Description: Extended GCD

d41d8c, 7 lines

d41d8c, 8 lines

```
array<int, 3> exgcd(int a, int b) {
  if (b == 0) {
    return {a, 1, 0};
  }
  auto [g, x, y] = exgcd(b, a % b);
  return {g, y, x - a / b * y};
}
```

#### CRT.h

Description: Chinese Remainder Theorem.

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

#### 5.3.1 Bézout's identity

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

$$ax + by = d$$

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

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

#### phiFunction.h

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

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

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

void calculatePhi() {
 row(i, 0, LIM) phi(i) = icl 2 i + i/2;
}

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

## 5.4 Fractions

ContinuedFractions.h

**Description:** Given N and a real number  $x \ge 0$ , 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  $\infty$ ; if x is the root of a degree 2 polynomial the a's eventually become cyclic. Time:  $\mathcal{O}(\log N)$ 

typedef double d; // for  $N \sim 1e7$ ; long double for  $N \sim 1e9$ pair<11, 11> approximate(d x, 11 N) { 11 LP = 0, LQ = 1, P = 1, Q = 0, inf =  $LLONG_MAX$ ; d y = x; 11  $\lim = \min(P ? (N-LP) / P : \inf, Q ? (N-LQ) / Q : \inf),$ a = (ll) floor(y), b = min(a, lim),NP = b\*P + LP, NO = b\*O + LO;**if** (a > b) { // If b > a/2, we have a semi-convergent that gives us a // better approximation; if b = a/2, we \*may\* have one. // Return {P, Q} here for a more canonical approximation. return (abs(x - (d)NP / (d)NO) < abs(x - (d)P / (d)O)) ? make\_pair(NP, NQ) : make\_pair(P, Q); **if**  $(abs(y = 1/(y - (d)a)) > 3*N) {$ return {NP, NQ}; LP = P; P = NP;LQ = Q; Q = NQ;

#### FracBinarySearch.h

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

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

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

## 5.5 Pythagorean Triples

The Pythagorean triples are uniquely generated by

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

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

## 5.6 Primes

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

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

### 5.7 Estimates

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

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

#### 5.8 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:

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

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

$$g(n) = \sum_{1 \le m \le n} f(\left|\frac{n}{m}\right|) \Leftrightarrow f(n) = \sum_{1 \le m \le n} \mu(m)g(\left|\frac{n}{m}\right|)$$

## Combinatorial (6)

### 6.1 Permutations

#### 6.1.1 Factorial

#### IntPerm.h

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

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

d41d8c, 6 lines

#### multinomial PushRelabel MinCostMaxFlow

#### 6.1.2 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 q (q.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

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

#### Partitions and subsets

#### 6.2.1 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  $\binom{n}{m} \equiv \prod_{i=0}^{k} \binom{n_i}{m_i} \pmod{p}$ .

#### 6.2.2 Binomials

multinomial.h

```
Description: Computes \binom{k_1 + \dots + k_n}{k_1, k_2, \dots, k_n} = \frac{(\sum k_i)!}{k_1!k_2!\dots k_n!}
11 multinomial(vi& v) {
  11 c = 1, m = v.empty() ? 1 : v[0];
  rep(i,1,sz(v)) rep(j,0,v[i])
    c = c * ++m / (j+1);
  return c;
```

## General purpose numbers

#### 6.3.1 Bell numbers

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

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

#### 6.3.2 Labeled unrooted trees

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

#### 6.3.3 Catalan numbers

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

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

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

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

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

# Graph (7)

#### 7.1 Network flow

PushRelabel.h

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

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

```
struct PushRelabel {
 struct Edge {
   int dest, back;
   11 f, c;
 };
 vector<vector<Edge>> g;
 vector<11> ec:
 vector<Edge*> cur;
 vector<vi> hs; vi H;
 PushRelabel(int n): q(n), ec(n), cur(n), hs(2*n), H(n) {}
 void addEdge(int s, int t, ll cap, ll rcap=0) {
    if (s == t) return;
   g[s].push_back({t, sz(g[t]), 0, cap});
   g[t].push_back({s, sz(g[s])-1, 0, rcap});
 void addFlow(Edge& e, ll f) {
   Edge &back = g[e.dest][e.back];
    if (!ec[e.dest] && f) hs[H[e.dest]].push_back(e.dest);
   e.f += f; e.c -= f; ec[e.dest] += f;
   back.f -= f; back.c += f; ec[back.dest] -= f;
 ll calc(int s, int t) {
    int v = sz(q); H[s] = v; ec[t] = 1;
   vi co(2*v); co[0] = v-1;
    rep(i,0,v) cur[i] = q[i].data();
    for (Edge& e : q[s]) addFlow(e, e.c);
    for (int hi = 0;;) {
     while (hs[hi].empty()) if (!hi--) return -ec[s];
     int u = hs[hi].back(); hs[hi].pop_back();
      while (ec[u] > 0) // discharge u
       if (cur[u] == g[u].data() + sz(g[u])) {
          H[u] = 1e9;
          for (Edge& e : g[u]) if (e.c && H[u] > H[e.dest]+1)
            H[u] = H[e.dest]+1, cur[u] = &e;
          if (++co[H[u]], !--co[hi] && hi < v)</pre>
            rep(i, 0, v) if (hi < H[i] && H[i] < v)
              --co[H[i]], H[i] = v + 1;
          hi = H[u];
        } else if (cur[u] \rightarrow c \&\& H[u] == H[cur[u] \rightarrow dest]+1)
          addFlow(*cur[u], min(ec[u], cur[u]->c));
        else ++cur[u];
 bool leftOfMinCut(int a) { return H[a] >= sz(q); }
```

#### MinCostMaxFlow.h

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

10

```
Time: \mathcal{O}\left(FE\log(V)\right) where F is max flow. \mathcal{O}\left(VE\right) for setpi. d41d8c, 79 lines
#include <bits/extc++.h>
const 11 INF = numeric_limits<11>::max() / 4;
struct MCMF {
  struct edge {
    int from, to, rev;
    11 cap, cost, flow;
  int N;
  vector<vector<edge>> ed;
  vi seen;
  vector<ll> dist, pi;
  vector<edge*> par;
  MCMF(int N) : N(N), ed(N), seen(N), dist(N), pi(N), par(N) {}
  void addEdge(int from, int to, ll cap, ll cost) {
    if (from == to) return;
    ed[from].push_back(edge{ from, to, sz(ed[to]), cap, cost, 0 });
    ed[to].push_back(edge{ to,from,sz(ed[from])-1,0,-cost,0 });
  void path(int s) {
    fill(all(seen), 0);
    fill(all(dist), INF);
    dist[s] = 0; ll di;
    __gnu_pbds::priority_queue<pair<ll, int>> q;
    vector<decltype(q)::point_iterator> its(N);
    q.push({ 0, s });
    while (!q.empty()) {
      s = q.top().second; q.pop();
      seen[s] = 1; di = dist[s] + pi[s];
      for (edge& e : ed[s]) if (!seen[e.to]) {
        11 val = di - pi[e.to] + e.cost;
        if (e.cap - e.flow > 0 && val < dist[e.to]) {</pre>
          dist[e.to] = val;
          par[e.to] = &e;
           if (its[e.to] == q.end())
             its[e.to] = q.push({ -dist[e.to], e.to });
             q.modify(its[e.to], { -dist[e.to], e.to });
    rep(i,0,N) pi[i] = min(pi[i] + dist[i], INF);
  pair<11, 11> maxflow(int s, int t) {
    11 \text{ totflow} = 0, \text{ totcost} = 0;
    while (path(s), seen[t]) {
      11 fl = INF;
      for (edge* x = par[t]; x; x = par[x->from])
        fl = min(fl, x->cap - x->flow);
      totflow += fl;
      for (edge* x = par[t]; x; x = par[x->from]) {
        x \rightarrow flow += fl;
        ed[x->to][x->rev].flow -= fl;
```

```
rep(i,0,N) for(edge& e : ed[i]) totcost += e.cost * e.flow;
    return {totflow, totcost/2};
  // If some costs can be negative, call this before maxflow:
  void setpi(int s) { // (otherwise, leave this out)
    fill(all(pi), INF); pi[s] = 0;
    int it = N, ch = 1; l1 v;
    while (ch-- && it--)
     rep(i,0,N) if (pi[i] != INF)
        for (edge& e : ed[i]) if (e.cap)
          if ((v = pi[i] + e.cost) < pi[e.to])</pre>
            pi[e.to] = v, ch = 1;
    assert(it >= 0); // negative cost cycle
};
```

#### EdmondsKarp.h

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

template<class T> T edmondsKarp(vector<unordered\_map<int, T>>& graph, int source, int sink) { assert (source != sink); T flow = 0: vi par(sz(graph)), q = par; for (;;) {

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

## LowerBoundFlow.h

**Description:** Dinic to run flows, has a lower bound option too<sub>41d8c, 93 lines</sub>

```
template <typename C, typename R = C, bool lb = false> struct
    dinic {
  typedef C flow_type;
  typedef R result_type;
  static const flow_type oo = std::numeric_limits<flow_type>::
      max();
  struct edge {
```

```
int src, dst, rev;
  flow_type cap, flow;
  edge(int src, int dst, int rev, flow_type cap, flow_type
      : src(src), dst(dst), rev(rev), cap(cap), flow(flow) {}
};
dinic(int n)
    : adj(n + 2 * 1b), que(n + 2 * 1b), level(n + 2 * 1b),
      edge_pos(n + 2 * lb) {}
int add_edge(int src, int dst, flow_type cap,
             flow_type rcap = 0) { // if lb rcap is low
  adj[src].emplace_back(src, dst, (int)adj[dst].size(), cap,
      1b ? rcap : 0);
  if (src == dst)
    adj[src].back().rev++;
  adj[dst].emplace_back(dst, src, (int)adj[src].size() - 1,
      lb ? 0 : rcap, 0);
  return (int)adj[src].size() - 1 - (src == dst);
inline bool side_of_S(int u) { return level[u] == -1; }
result_type max_flow(int source, int sink) {
  result_type flow = 0;
  while (true) {
    int front = 0, back = 0;
    std::fill(level.begin(), level.end(), -1);
    for (level[que[back++] = sink] = 0; front < back && level</pre>
         [source] == -1;
         ++front) {
      int u = que[front];
      for (const edge &e : adj[u])
        if (level[e.dst] == -1 && rev(e).flow < rev(e).cap)</pre>
          level[que[back++] = e.dst] = 1 + level[u];
    if (level[source] == -1)
     break;
    std::fill(edge_pos.begin(), edge_pos.end(), 0);
    auto find_path = [&] (int from, flow_type res) {
     if (from == sink)
      for (int &ept = edge_pos[from]; ept < (int)adj[from].</pre>
          size(); ++ept) {
        edge &e = adj[from][ept];
        if (e.flow == e.cap || level[e.dst] + 1 != level[from
            1)
        flow_type push = find_path(e.dst, std::min(res, e.cap
              - e.flow));
        if (push > 0) {
          e.flow += push;
          rev(e).flow -= push;
          if (e.flow == e.cap)
            ++ept;
          return push;
     return static_cast<flow_type>(0);
    for (flow_type f; (f = find_path(source, oo)) > 0;)
      flow += f:
  return flow;
result_type max_flow_lb(int source, int sink) {
  int n = adj.size() - 2;
  vector<flow_type> delta(n + 2);
  for (int u = 0; u < n; ++u)
    for (auto &e : adj[u]) {
      delta[u] -= e.flow;
      delta[e.dst] += e.flow;
```

```
result_type sum = 0;
    int s = n, t = n + 1;
    for (int u = 0; u < n; ++u) {</pre>
      if (delta[u] > 0) {
        add_edge(s, u, delta[u], 0);
        sum += delta[u];
      } else if (delta[u] < 0)</pre>
        add_edge(u, t, -delta[u], 0);
    add_edge(sink, source, oo, 0);
    if (max_flow(s, t) != sum)
      return -1; // no solution
    result_type flow = adj[sink].back().flow;
    adj[sink].pop_back();
    adj[source].pop_back();
    return flow + max_flow(source, sink);
  std::vector<std::vector<edge>> adj;
private:
 std::vector<int> que;
 std::vector<int> level;
 std::vector<int> edge_pos;
 inline edge &rev(const edge &e) { return adj[e.dst][e.rev]; }
```

#### MinCut.h

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

#### GlobalMinCut.h

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

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

11

```
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;
```

#### GomoryHu.h

**Description:** Given a list of edges representing an undirected flow graph, returns edges of the Gomory-Hu tree. The max flow between any pair of vertices is given by minimum edge weight along the Gomory-Hu tree path. **Time:**  $\mathcal{O}(V)$  Flow Computations

```
"PushRelabel.h"
```

```
d41d8c, 13 lines
typedef array<11, 3> Edge;
vector<Edge> gomoryHu(int N, vector<Edge> ed) {
```

**Description:** Simple bipartite matching algorithm. Graph q should be a list

of neighbors of the left partition, and btoa should be a vector full of -1's of

the same size as the right partition. Returns the size of the matching. btoa[i]

will be the match for vertex i on the right side, or -1 if it's not matched.

```
vector<Edge> tree;
vi par(N);
rep(i,1,N) {
 PushRelabel D(N); // Dinic also works
 for (Edge t : ed) D.addEdge(t[0], t[1], t[2], t[2]);
 tree.push_back({i, par[i], D.calc(i, par[i])});
 rep(j,i+1,N)
   if (par[j] == par[i] && D.leftOfMinCut(j)) par[j] = i;
return tree:
```

## Matching

#### hopcroftKarp.h

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

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

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

DFSMatching.h

cur.swap(next);

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

rep(a, 0, sz(g))

```
Usage: vi btoa(m, -1); dfsMatching(q, btoa);
Time: \mathcal{O}(VE)
bool find(int j, vector<vi>& g, vi& btoa, vi& vis) {
 if (btoa[j] == -1) return 1;
 vis[j] = 1; int di = btoa[j];
  for (int e : q[di])
    if (!vis[e] && find(e, g, btoa, vis)) {
      btoa[e] = di;
      return 1;
  return 0;
int dfsMatching(vector<vi>& g, vi& btoa) {
  rep(i, 0, sz(q)) {
    vis.assign(sz(btoa), 0);
    for (int j : q[i])
      if (find(j, g, btoa, vis)) {
        btoa[j] = i;
        break;
  return sz(btoa) - (int)count(all(btoa), -1);
MinimumVertexCover.h
Description: Finds a minimum vertex cover in a bipartite graph. The size
is the same as the size of a maximum matching, and the complement is a
maximum independent set.
"DFSMatching.h"
                                                      d41d8c, 20 lines
vi cover(vector<vi>& g, int n, int m) {
```

```
vi match(m, -1);
int res = dfsMatching(g, match);
vector<bool> lfound(n, true), seen(m);
for (int it : match) if (it != -1) lfound[it] = false;
vi q, cover;
rep(i,0,n) if (lfound[i]) q.push_back(i);
while (!q.emptv()) {
  int i = q.back(); q.pop_back();
  lfound[i] = 1;
  for (int e : g[i]) if (!seen[e] && match[e] != -1) {
    seen[e] = true;
    q.push_back(match[e]);
rep(i,0,n) if (!lfound[i]) cover.push_back(i);
rep(i,0,m) if (seen[i]) cover.push_back(n+i);
assert(sz(cover) == res);
return cover;
```

#### WeightedMatching.h

**Description:** Given a weighted bipartite graph, matches every node on the left with a node on the right such that no nodes are in two matchings and the sum of the edge weights is minimal. Takes 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 "dummy" worker 0
  vi dist(m, INT_MAX), pre(m, -1);
  vector<bool> done(m + 1);
  do { // dijkstra
    done[j0] = true;
    int i0 = p[j0], j1, delta = INT_MAX;
    rep(j,1,m) if (!done[j]) {
      auto cur = a[i0 - 1][j - 1] - u[i0] - v[j];
      if (cur < dist[j]) dist[j] = cur, pre[j] = j0;
      if (dist[j] < delta) delta = dist[j], j1 = j;</pre>
    rep(j,0,m) {
      if (done[j]) u[p[j]] += delta, v[j] -= delta;
      else dist[j] -= delta;
    j0 = j1;
  } while (p[j0]);
  while (j0) { // update alternating path
    int j1 = pre[j0];
    p[j0] = p[j1], j0 = j1;
rep(j,1,m) if (p[j]) ans[p[j] - 1] = j - 1;
return {-v[0], ans}; // min cost
```

#### GeneralMatching.h

**Description:** Matching for general graphs. Fails with probability N/mod. Time:  $\mathcal{O}(N^3)$ 

```
"../numerical/MatrixInverse-mod.h"
                                                     d41d8c, 40 lines
vector<pii> generalMatching(int N, vector<pii>& ed) {
 vector<vector<ll>> mat(N, vector<ll>(N)), A;
 for (pii pa : ed) {
   int a = pa.first, b = pa.second, r = rand() % mod;
   mat[a][b] = r, mat[b][a] = (mod - r) % mod;
 int r = matInv(A = mat), M = 2*N - r, fi, f;
 assert (r % 2 == 0);
 if (M != N) do {
   mat.resize(M, vector<11>(M));
   rep(i,0,N) {
     mat[i].resize(M);
     rep(j,N,M) {
       int r = rand() % mod;
       mat[i][j] = r, mat[j][i] = (mod - r) % mod;
 } while (matInv(A = mat) != M);
 vi has (M, 1); vector<pii> ret;
 rep(it,0,M/2) {
   rep(i,0,M) if (has[i])
      rep(j,i+1,M) if (A[i][j] && mat[i][j]) {
       fi = i; fj = j; goto done;
    } assert(0); done:
    if (fj < N) ret.emplace back(fi, fj);</pre>
   has[fi] = has[fj] = 0;
    rep(sw,0,2) {
     11 a = modpow(A[fi][fi], mod-2);
      rep(i,0,M) if (has[i] && A[i][fj]) {
       ll b = A[i][fi] * a % mod;
        rep(j, 0, M) A[i][j] = (A[i][j] - A[fi][j] * b) % mod;
      swap(fi,fj);
```

```
return ret;
```

## DFS algorithms

#### SCC.h

Description: Finds strongly connected components in a directed graph. If vertices u, v belong to the same component, we can reach u from v and vice

Usage:  $scc(graph, [\&](vi\& v) \{ ... \})$  visits all components in reverse topological order. comp[i] holds the component index of a node (a component only has edges to components with lower index). ncomps will contain the number of components. Time:  $\mathcal{O}\left(E+V\right)$ 

vi val, comp, z, cont; int Time, ncomps; template < class G, class F> int dfs (int j, G& g, F& f) { int low = val[j] = ++Time, x; z.push back(j); for (auto e : q[j]) if (comp[e] < 0) low = min(low, val[e] ?: dfs(e,g,f));**if** (low == val[j]) {  $x = z.back(); z.pop_back();$ comp[x] = ncomps;cont.push back(x); } while (x != j);f(cont); cont.clear(); ncomps++; return val[j] = low; template < class G, class F> void scc(G& q, F f) { int n = sz(q);val.assign(n, 0); comp.assign(n, -1); Time = ncomps = 0;rep(i,0,n) if (comp[i] < 0) dfs(i, g, f);

#### BiconnectedComponents.h

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

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

d41d8c, 33 lines

```
vi num, st;
vector<vector<pii>> ed;
int Time;
template<class F>
int dfs(int at, int par, F& f) {
  int me = num[at] = ++Time, e, y, top = me;
  for (auto pa : ed[at]) if (pa.second != par) {
    tie(y, e) = pa;
    if (num[y]) {
     top = min(top, num[y]);
     if (num[y] < me)
        st.push back(e);
    } else {
      int si = sz(st);
      int up = dfs(y, e, f);
     top = min(top, up);
     if (up == me) {
```

```
st.push_back(e);
       f(vi(st.begin() + si, st.end()));
       st.resize(si);
      else if (up < me) st.push_back(e);</pre>
      else { /* e is a bridge */ }
 return top;
template<class F>
void bicomps(F f) {
 num.assign(sz(ed), 0);
 rep(i,0,sz(ed)) if (!num[i]) dfs(i, -1, f);
```

#### 2sat.h

struct TwoSat {

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

```
Usage: TwoSat ts(number of boolean variables);
ts.either(0, \sim3); // Var 0 is true or var 3 is false
ts.setValue(2); // Var 2 is true
ts.atMostOne(\{0, \sim 1, 2\}); // <= 1 of vars 0, \sim 1 and 2 are true
ts.solve(); // Returns true iff it is solvable
ts.values[0..N-1] holds the assigned values to the vars
```

**Time:**  $\mathcal{O}(N+E)$ , where N is the number of boolean variables, and E is the number of clauses. d41d8c, 56 lines

```
int N;
vector<vi> gr;
vi values; // 0 = false, 1 = true
TwoSat(int n = 0) : N(n), gr(2*n) {}
int addVar() { // (optional)
  gr.emplace_back();
 gr.emplace_back();
```

```
return N++;
void either(int f, int j) {
  f = \max(2*f, -1-2*f);
  j = \max(2 * j, -1 - 2 * j);
  gr[f].push_back(j^1);
  gr[j].push_back(f^1);
void setValue(int x) { either(x, x); }
void atMostOne(const vi& li) { // (optional)
  if (sz(li) <= 1) return;</pre>
  int cur = \simli[0];
  rep(i,2,sz(li)) {
    int next = addVar();
    either(cur, ~li[i]);
    either(cur, next);
    either(~li[i], next);
    cur = ~next;
  either(cur, ~li[1]);
vi val, comp, z; int time = 0;
int dfs(int i) {
```

int low = val[i] = ++time, x; z.push\_back(i);

for(int e : gr[i]) if (!comp[e])

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

13

d41d8c, 31 lines

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

## 7.4 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:  $\mathcal{O}(NM)$ 

adj[u][e] = left;

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];
```

d41d8c, 21 lines

```
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;
```

### 7.5 Heuristics

#### MaximalCliques.h

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

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

d41d8c, 12 lines

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

#### MaximumClique.h

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

Time: Runs in about 1s for n=155 and worst case random graphs (p=.90). Runs faster for sparse graphs. d41d8c, 49 lines

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

```
C[1].clear(), C[2].clear();
       for (auto v : T) {
         int k = 1;
         auto f = [&](int i) { return e[v.i][i]; };
         while (any_of(all(C[k]), f)) k++;
         if (k > mxk) mxk = k, C[mxk + 1].clear();
         if (k < mnk) T[j++].i = v.i;
         C[k].push_back(v.i);
       if (j > 0) T[j - 1].d = 0;
       rep(k, mnk, mxk + 1) for (int i : C[k])
         T[j].i = i, T[j++].d = k;
       expand(T, lev + 1);
     } else if (sz(q) > sz(qmax)) qmax = q;
     q.pop_back(), R.pop_back();
 vi maxClique() { init(V), expand(V); return qmax; }
 Maxclique(vb conn) : e(conn), C(sz(e)+1), S(sz(C)), old(S) {
    rep(i,0,sz(e)) V.push_back({i});
};
```

#### MaximumIndependentSet.h

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

## 7.6 Trees

#### BinaryLifting.h

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

**Time:** construction  $\mathcal{O}(N \log N)$ , queries  $\mathcal{O}(\log N)$ 

d41d8c, 25 lines

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

#### LCA.h

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

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

```
"../data-structures/RMQ.h"
                                                                  d41d8c, 21 lines
struct LCA {
```

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

#### CompressTree.h

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

Time:  $\mathcal{O}(|S| \log |S|)$ "LCA.h"

typedef vector<pair<int, int>> vpi; vpi compressTree(LCA& lca, const vi& subset) { static vi rev; rev.resize(sz(lca.time)); vi li = subset, &T = lca.time;

```
auto cmp = [&](int a, int b) { return T[a] < T[b]; };</pre>
sort(all(li), cmp);
int m = sz(1i)-1;
rep(i,0,m) {
  int a = li[i], b = li[i+1];
  li.push_back(lca.lca(a, b));
sort(all(li), cmp);
li.erase(unique(all(li)), li.end());
rep(i, 0, sz(li)) rev[li[i]] = i;
vpi ret = {pii(0, li[0])};
rep(i, 0, sz(li)-1) {
  int a = li[i], b = li[i+1];
  ret.emplace_back(rev[lca.lca(a, b)], b);
return ret;
```

#### HLD.h

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

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

```
"../data-structures/LazySegmentTree.h"
                                                            d41d8c, 46 lines
template <bool VALS_EDGES> struct HLD {
  int N, tim = 0;
```

```
vector<vi> adj;
vi par, siz, depth, rt, pos;
Node *tree;
HLD (vector<vi> adj )
  : N(sz(adj_)), adj(adj_), par(N, -1), siz(N, 1), depth(N),
    rt(N),pos(N),tree(new Node(0, N)) { dfsSz(0); dfsHld(0); }
```

```
void dfsSz(int v) {
   if (par[v] != -1) adj[v].erase(find(all(adj[v]), par[v]));
   for (int& u : adj[v]) {
     par[u] = v, depth[u] = depth[v] + 1;
     dfsSz(u);
     siz[v] += siz[u];
     if (siz[u] > siz[adj[v][0]]) swap(u, adj[v][0]);
 void dfsHld(int v) {
   pos[v] = tim++;
   for (int u : adj[v]) {
     rt[u] = (u == adj[v][0] ? rt[v] : u);
     dfsHld(u);
  template <class B> void process(int u, int v, B op) {
   for (; rt[u] != rt[v]; v = par[rt[v]]) {
     if (depth[rt[u]] > depth[rt[v]]) swap(u, v);
     op(pos[rt[v]], pos[v] + 1);
   if (depth[u] > depth[v]) swap(u, v);
   op(pos[u] + VALS_EDGES, pos[v] + 1);
  void modifyPath(int u, int v, int val) {
   process(u, v, [&](int 1, int r) { tree->add(1, r, val); });
  int queryPath(int u, int v) { // Modify depending on problem
   int res = -1e9;
   process(u, v, [&](int l, int r) {
       res = max(res, tree->query(1, r));
   return res;
 int querySubtree(int v) { // modifySubtree is similar
   return tree->query(pos[v] + VALS_EDGES, pos[v] + siz[v]);
};
```

#### 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"
struct Edge { int a, b; ll w; };
struct Node {
  Edge key;
  Node *1, *r;
  11 delta:
  void prop() {
   key.w += delta;
   if (1) 1->delta += delta;
   if (r) r->delta += delta;
   delta = 0:
  Edge top() { prop(); return key; }
Node *merge(Node *a, Node *b) {
  if (!a || !b) return a ?: b;
  a->prop(), b->prop();
  if (a->key.w > b->key.w) swap(a, b);
  swap(a->1, (a->r = merge(b, a->r)));
  return a;
void pop(Node*\& a) { a->prop(); a = merge(a->1, a->r); }
pair<ll, vi> dmst(int n, int r, vector<Edge>& g) {
  RollbackUF uf(n);
  vector<Node*> heap(n);
```

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

## 7.7 Math

## 7.7.1 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).

#### 7.7.2 Erdős–Gallai theorem

A simple graph with node degrees  $d_1 > \cdots > d_n$  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 (8)

## 8.1 Geometric primitives

#### Point.h

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

```
template <class T> int sgn(T x) \{ return (x > 0) - (x < 0); \}
```

```
template<class T>
struct Point {
 typedef Point P;
 Тх, у;
  // operators imply that point should be IN FRONT of the class
        it passes in
  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 std::ostream& operator<<(std::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.



d41d8c, 6 lines

d41d8c, 4 lines

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

#### SegmentDistance.h

#### Description:

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

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

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

SegmentIntersection.h

#### Description:

If a unique intersection point between the line segments going from s1 to e1 and from s2 to e2 exists then it is returned. If no intersection point exists an empty vector is returned. If infinitely many exist a vector with 2 elements is returned, containing the endpoints of the common line segment. The wrong position will be returned if P is Point<|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"
                                                         d41d8c, 13 lines
```

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

#### lineIntersection.h

#### Description:

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  $\{-1, e^2\}$ (0,0)} is returned. The wrong position will be returned if P is Point<|l> and the intersection point does not have integer coordinates. Products of three coordinates are used in inter- 1 mediate steps so watch out for overflow if using int or ll.



```
Usage: auto res = lineInter(s1,e1,s2,e2);
if (res.first == 1)
cout << "intersection point at " << res.second << endl;</pre>
"Point.h"
                                                        d41d8c, 8 lines
```

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

#### sideOf.h

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

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

```
d41d8c, 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 l = (e-s).dist()*eps;
  return (a > 1) - (a < -1);
```

#### linearTransformation.h Description:

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



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

#### Angle.h

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

```
Usage: vector<Angle> v = \{w[0], w[0].t360() ...\}; // sorted
int j = 0; rep(i,0,n) { while (v[j] < v[i].t180()) ++j; }
// sweeps j such that (j-i) represents the number of positively
oriented triangles with vertices at 0 and i
                                                     d41d8c, 35 lines
```

```
struct Angle {
 int x, y;
 Angle(int x, int y, int t=0) : x(x), y(y), t(t) {}
 Angle operator-(Angle b) const { return {x-b.x, y-b.y, t}; }
 int half() const {
   assert(x || y);
    return y < 0 || (y == 0 && x < 0);
 Angle t90() const { return \{-y, x, t + (half() \&\& x >= 0)\}; \}
 Angle t180() const { return {-x, -y, t + half()}; }
 Angle t360() const { return {x, y, t + 1}; }
bool operator<(Angle a, Angle b) {</pre>
 // add a.dist2() and b.dist2() to also compare distances
  return make_tuple(a.t, a.half(), a.y * (ll)b.x) <</pre>
         make_tuple(b.t, b.half(), a.x * (11)b.y);
// Given two points, this calculates the smallest angle between
// them, i.e., the angle that covers the defined line segment.
pair<Angle, Angle> segmentAngles(Angle a, Angle b) {
 if (b < a) swap(a, b);
  return (b < a.t180() ?
          make_pair(a, b) : make_pair(b, a.t360()));
Angle operator+(Angle a, Angle b) { // point \ a + vector \ b
 Angle r(a.x + b.x, a.y + b.y, a.t);
 if (a.t180() < r) r.t--;
 return r.t180() < a ? r.t360() : r;</pre>
Angle angleDiff(Angle a, Angle b) { // angle b- angle a
 int tu = b.t - a.t; a.t = b.t;
  return \{a.x*b.x + a.y*b.y, a.x*b.y - a.y*b.x, tu - (b < a)\};
```

## 8.2 Circles

#### CircleTangents.h

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

```
"Point.h"
                                                                         d41d8c, 13 lines
```

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

#### CircleLine.h

Description: Finds the intersection between a circle and a line. Returns a vector of either 0, 1, or 2 intersection points. P is intended to be Point < double >. <vector>, "Point.h" d41d8c, 9 lines

```
template<class P>
std::vector<P> circleLine(P c, double r, P a, P b) {
 P ab = b - a, p = a + ab * (c-a).dot(ab) / ab.dist2();
 double s = a.cross(b, c), h2 = r*r - s*s / ab.dist2();
 if (h2 < 0) return {};
 if (h2 == 0) return {p};
 P h = ab.unit() * sqrt(h2);
 return {p - h, p + h};
```

#### CirclePolygonIntersection.h

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

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

```
"../../content/geometry/Point.h"
// make sure to consider the value of infinity when doing this
     ... if it's too big it'll blow up! (from halfplane
     intersection and stuff)
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;
    Pd = q - p;
    auto a = d.dot(p)/d.dist2(), b = (p.dist2()-r*r)/d.dist2();
    auto det = a * a - b;
    if (det <= 0) return arg(p, q) * r2;</pre>
    auto s = max(0., -a-sgrt(det)), t = min(1., -a+sgrt(det));
    if (t < 0 || 1 <= s) return arg(p, q) * r2;</pre>
    P u = p + d * s, v = p + d * t;
    return arg(p,u) * r2 + u.cross(v)/2 + arg(v,g) * r2;
  auto sum = 0.0;
  rep(i, 0, sz(ps))
   sum += tri(ps[i] - c, ps[(i + 1) % sz(ps)] - c);
  return sum;
```

#### circumcircle.h

#### Description:

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



typedef Point < double > P; double ccRadius (const P& A, const P& B, const P& C) { **return** (B-A).dist() \* (C-B).dist() \* (A-C).dist() /

d41d8c, 12 lines

```
abs((B-A).cross(C-A))/2;
P ccCenter (const P& A, const P& B, const P& C) {
 P b = C-A, c = B-A;
 return A + (b*c.dist2()-c*b.dist2()).perp()/b.cross(c)/2;
```

#### MinimumEnclosingCircle.h

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

```
d41d8c, 17 lines
"circumcircle.h"
pair<P, double> mec(vector<P> ps) {
  shuffle(all(ps), mt19937(time(0)));
  P \circ = ps[0];
  double r = 0, EPS = 1 + 1e-8;
  rep(i, 0, sz(ps)) if ((o - ps[i]).dist() > r * EPS) {
   o = ps[i], r = 0;
    rep(j,0,i) if ((o - ps[j]).dist() > r * EPS) {
     o = (ps[i] + ps[j]) / 2;
     r = (o - ps[i]).dist();
      rep(k, 0, j) if ((o - ps[k]).dist() > r * EPS) {
       o = ccCenter(ps[i], ps[j], ps[k]);
        r = (o - ps[i]).dist();
  return {o, r};
```

## 8.3 Polygons

#### InsidePolygon.h

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

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

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

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

#### PolygonArea.h

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

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

#### PolygonCenter.h

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

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

d41d8c, 9 lines

```
typedef Point<double> P;
```

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

#### PolygonCut.h Description:

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

```
Usage: vector<P> p = ...;
p = polygonCut(p, P(0,0), P(1,0));
```

"Point.h", "lineIntersection.h" d41d8c, 13 lines

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

#### PolygonUnion.h

**Description:** Calculates the area of the union of n polygons (not necessarily convex). The points within each polygon must be given in CCW order. (Epsilon checks may optionally be added to sideOf/sgn, but shouldn't be

**Time:**  $\mathcal{O}(N^2)$ , where N is the total number of points

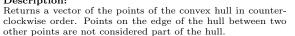
```
"Point.h", "sideOf.h"
                                                                         d41d8c, 33 lines
```

```
typedef Point<double> P;
double rat(P a, P b) { return sgn(b.x) ? a.x/b.x : a.y/b.y; }
double polyUnion(vector<vector<P>>& poly) {
 double ret = 0;
 rep(i, 0, sz(poly)) rep(v, 0, sz(poly[i])) {
   P A = poly[i][v], B = poly[i][(v + 1) % sz(poly[i])];
   vector<pair<double, int>> segs = {{0, 0}, {1, 0}};
    rep(j,0,sz(poly)) if (i != j) {
      rep(u, 0, sz(poly[j])) {
       P C = poly[j][u], D = poly[j][(u + 1) % sz(poly[j])];
       int sc = sideOf(A, B, C), sd = sideOf(A, B, D);
       if (sc != sd) {
         double sa = C.cross(D, A), sb = C.cross(D, B);
         if (min(sc, sd) < 0)
            segs.emplace_back(sa / (sa - sb), sgn(sc - sd));
        } else if (!sc && !sd && j<i && sgn((B-A).dot(D-C))>0){
          segs.emplace_back(rat(C - A, B - A), 1);
          segs.emplace_back(rat(D - A, B - A), -1);
    sort (all (segs));
    for (auto\& s : segs) s.first = min(max(s.first, 0.0), 1.0);
   double sum = 0;
   int cnt = segs[0].second;
    rep(j,1,sz(segs)) {
     if (!cnt) sum += seqs[j].first - seqs[j - 1].first;
     cnt += segs[j].second;
    ret += A.cross(B) * sum;
```

```
return ret / 2;
```

#### ConvexHull.h

#### Description: Returns a vector of the points of the convex hull in counterclockwise order. Points on the edge of the hull between two



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

```
<vector>, "Point.h"
template < class T>
std::vector<Point<T>> convexHull(std::vector<Point<T>> pts) {
 using P = Point<T>;
 if ((int) pts.size() <= 1) return pts;</pre>
  sort(pts.begin(), pts.end());
  std::vector<P> h((int) pts.size()+1);
  int s = 0, t = 0;
  for (int it = 2; it--; s = --t, reverse(pts.begin(), pts.end
    for (P p : pts) {
      while (t >= s + 2 \&\& h[t-2].cross(h[t-1], p) <= 0) t--;
 return {h.begin(), h.begin() + t - (t == 2 && h[0] == h[1])};
```

#### HullDiameter.h

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

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

"Point.h"

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

#### PointInsideHull.h

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

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

```
"Point.h", "sideOf.h", "OnSegment.h"
                                                       d41d8c, 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

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

#### HalfplaneIntersection.h

Description: NlogN Halfplane intersection Time:  $\mathcal{O}(n \log n)$ 

```
"Point.h", <vector>, <deque>
```

d41d8c, 94 lines

```
using namespace std;
const long double eps = 1e-9, inf = 1e9;
template <class T>
struct Halfplane {
    // the halfplane determined by p and pq takes all points
         that are _counterclockwise_
    // wrt p \Rightarrow (p + pq) vector
   Point<T> p, pq;
   long double angle;
   Halfplane() {}
```

```
Halfplane(const Point<T>& a, const Point<T>& b) : p(a), pq(
        b - a) {
        angle = atan21(pq.y, pq.x);
   bool out(const Point<T>& r) {
        return pq.cross(r - p) < -eps;</pre>
   bool operator < (const Halfplane& e) const {
        return angle < e.angle;</pre>
    friend Point<T> inter(const Halfplane& s, const Halfplane&
        long double alpha = (t.p - s.p).cross(t.pq) / s.pq.
             cross(t.pq);
        return s.p + (s.pq * alpha);
template <class T>
vector<Point<T>> hp_intersect(vector<Halfplane<T>> &H) {
    // important note: H is changed within this function so
         make a copy if you need it later!
    using P = Point<T>;
    using Halfplane = Halfplane<T>;
   P box[4] = \{
       P(inf, inf),
       P(-inf, inf),
       P(-inf, -inf),
       P(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(dg[len - 1], dg[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(H[i].pq.cross(dq[len - 1].pq)) <</pre>
            if (H[i].pq.dot(dq[len - 1].pq) < 0.0) {</pre>
                return vector<P>();
            if (H[i].out(dq[len - 1].p)) {
                dq.pop_back();
                --len;
            } else continue;
        dq.push_back(H[i]);
        ++len;
```

```
while (len > 2 && dq[0].out(inter(dq[len - 1], dq[len - 2])
    )) {
    dq.pop_back();
    --len;
while (len > 2 && dq[len - 1].out(inter(dq[0], dq[1]))) {
    dq.pop_front();
    --len;
if (len < 3) return vector<P>();
vector<P> 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:
```

#### Minkowski.h

Description: Minkowski sum of two convex polygons Assumes the polygons are sorted ccw Time Complexity: O(-P-+-Q-)

```
<vector>, "Point.h"
                                                      d41d8c, 27 lines
template <typename P>
void reorder_polygon(std::vector<P>& poly) {
    int pos = 0;
    for (int i = 1; i < poly.size(); i++) {</pre>
        if (poly[i].y < poly[pos].y || (poly[i].y == poly[pos].</pre>
            y && poly[i].x < poly[pos].x))
            pos = i;
    rotate(poly.begin(), poly.begin() + pos, poly.end());
template <typename P>
std::vector<P> minkowski(std::vector<P> poly1, std::vector<P>
    polv2) {
    reorder_polygon(poly1), reorder_polygon(poly2);
    poly1.push_back(poly1[0]);
    poly1.push_back(poly1[1]);
    poly2.push_back(poly2[0]);
    poly2.push back(poly2[1]);
    std::vector<P> result;
    int i = 0, j = 0;
    while (i < poly1.size() - 2 || j < poly2.size() - 2) {</pre>
        result.push_back(poly1[i] + poly2[j]);
        auto cr = (poly1[i + 1] - poly1[i]).cross(poly2[j + 1]
             - poly2[j]);
        if (cr >= 0 && i < poly1.size() - 2) i++;</pre>
        if (cr <= 0 && j < poly2.size() - 2) j++;</pre>
    return result;
```

## 8.4 Misc. Point Set Problems

ClosestPair.h

Description: Finds the closest pair of points.

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

```
"Point.h"
                                                        d41d8c, 17 lines
typedef Point<11> P;
pair<P, P> closest(vector<P> v) {
  assert (sz(v) > 1);
  sort(all(v), [](P a, P b) { return a.y < b.y; });
  pair<ll, pair<P, P>> ret{LLONG_MAX, {P(), P()}};
  int \dot{1} = 0;
```

```
for (P p : v) {
  P d{1 + (ll) sqrt(ret.first), 0};
  while (v[j].y <= 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;</pre>
```

#### 8.5 3D

#### PolyhedronVolume.h

**Description:** Magic formula for the volume of a polyhedron. Faces should point outwards.

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

#### Point3D.h

**Description:** Class to handle points in 3D space. T can be e.g. double or long long.  $\frac{\text{d41d8c, 32 lines}}{\text{d41d8c, 32 lines}}$ 

```
template<class T> struct Point3D {
  typedef Point3D P;
  typedef const P& R;
  T x, y, z;
  explicit Point3D(T x=0, T y=0, T z=0) : x(x), y(y), z(z) {}
  bool operator<(R p) const {</pre>
   return tie(x, y, z) < tie(p.x, p.y, p.z); }
  bool operator==(R p) const {
   return tie(x, y, z) == tie(p.x, p.y, p.z); }
  P operator+(R p) const { return P(x+p.x, y+p.y, z+p.z); }
  P operator-(R p) const { return P(x-p.x, y-p.y, z-p.z); }
  P operator*(T d) const { return P(x*d, y*d, z*d); }
  P operator/(T d) const { return P(x/d, y/d, z/d); }
  T dot(R p) const { return x*p.x + y*p.y + z*p.z; }
  P cross(R p) const {
    return P(y*p.z - z*p.y, z*p.x - x*p.z, x*p.y - y*p.x);
  T dist2() const { return x*x + y*y + z*z; }
  double dist() const { return sqrt((double)dist2()); }
  //Azimuthal angle (longitude) to x-axis in interval [-pi, pi]
  double phi() const { return atan2(y, x); }
  //Zenith angle (latitude) to the z-axis in interval [0, pi]
  double theta() const { return atan2(sqrt(x*x+y*y),z); }
  P unit() const { return *this/(T)dist(); } //makes dist()=1
  //returns unit vector normal to *this and p
  P normal(P p) const { return cross(p).unit(); }
  //returns point rotated 'angle' radians ccw around axis
  P rotate (double angle, P axis) const {
    double s = sin(angle), c = cos(angle); P u = axis.unit();
    return u*dot(u)*(1-c) + (*this)*c - cross(u)*s;
};
```

#### 3dHull.h

**Description:** Computes all faces of the 3-dimension hull of a point set. \*No four points must be coplanar\*, or else random results will be returned. All faces will point outwards.

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

```
"Point3D.h" d41d8c, 49 lines
```

typedef Point3D<double> P3;

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

#### sphericalDistance.h

**Description:** Returns the shortest distance on the sphere with radius radius between the points with azimuthal angles (longitude) f1  $(\phi_1)$  and f2  $(\phi_2)$  from x axis and zenith angles (latitude) t1  $(\theta_1)$  and t2  $(\theta_2)$  from z axis (0 = north pole). All angles measured in radians. The algorithm starts by converting the spherical coordinates to cartesian coordinates so if that is what you have you can use only the two last rows. dx\*radius is then the difference between the two points in the x direction and d\*radius is the total distance between the points.

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

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

#### KMP

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

d41d8c 16 lin

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

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

#### Zfunc.h

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

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

d41d8c, 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, <math>p[1][i] = longest odd (half rounded down).

Time:  $\mathcal{O}\left(N\right)$  d41d8c, 13 lines

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

#### MinRotation.h

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

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

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

return a;

#### SuffixArray SuffixTree Hashing AhoCorasick

```
SuffixArray.h
Description: Suffix Array
Time: \mathcal{O}(n \log n)
                                                      d41d8c, 46 lines
vector<int> suffix array(string S) {
  int N = S.size();
  vector<int> sa(N), classes(N);
  for (int i = 0; i < N; i++) {</pre>
   sa[i] = N - 1 - i;
   classes[i] = S[i];
  stable_sort(sa.begin(), sa.end(), [&S](int i, int j) { return
       S[i] < S[i]; });
  for (int len = 1; len < N; len *= 2) {
   vector<int> c(classes);
    for (int i = 0; i < N; i++) {</pre>
     bool same = i && sa[i - 1] + len < N && c[sa[i]] == c[sa[
          i - 1]] &&
                  c[sa[i] + len / 2] == c[sa[i - 1] + len / 2];
     classes[sa[i]] = same ? classes[sa[i - 1]] : i;
    vector<int> cnt(N), s(sa);
    for (int i = 0; i < N; i++)
     cnt[i] = i;
    for (int i = 0; i < N; i++) {
     int s1 = s[i] - len;
     if (s1 >= 0)
        sa[cnt[classes[s1]]++] = s1;
  return sa;
vector<int> lcp array(const vector<int> &sa, string S) {
  int N = S.size();
  vector < int > rank(N), lcp(N - 1);
  for (int i = 0; i < N; i++)
   rank[sa[i]] = i;
  int pre = 0;
  for (int i = 0; i < N; i++) {
   if (rank[i] < N - 1) {
      int i = sa[rank[i] + 1];
      while (max(i, j) + pre < S.size() && S[i + pre] == S[j +
          prel)
        ++pre;
      lcp[rank[i]] = pre;
     if (pre > 0)
        --pre;
 return lcp;
SuffixTree.h
```

if (s[a+k] > s[b+k]) { a = b; break; }

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

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

d41d8c, 50 lines

struct SuffixTree {

```
enum { N = 200010, ALPHA = 26 }; // N \sim 2*maxlen+10
  int toi(char c) { return c - 'a'; }
  string a; //v = cur \ node, q = cur \ position
  int t[N][ALPHA], 1[N], r[N], p[N], s[N], v=0, q=0, m=2;
  void ukkadd(int i, int c) { suff:
    if (r[v]<=q) {
      if (t[v][c]==-1) { t[v][c]=m; l[m]=i;
        p[m++]=v; v=s[v]; q=r[v]; goto suff; }
      v=t[v][c]; q=l[v];
    if (q==-1 || c==toi(a[q])) q++; else {
      l[m+1]=i; p[m+1]=m; l[m]=l[v]; r[m]=q;
      p[m]=p[v]; t[m][c]=m+1; t[m][toi(a[q])]=v;
      l[v]=q; p[v]=m; t[p[m]][toi(a[l[m]])]=m;
      v=s[p[m]]; q=l[m];
      while (q<r[m]) { v=t[v][toi(a[q])]; q+=r[v]-l[v]; }</pre>
      if (q==r[m]) s[m]=v; else s[m]=m+2;
      q=r[v]-(q-r[m]); m+=2; qoto suff;
  SuffixTree(string a) : a(a) {
    fill(r,r+N,sz(a));
    memset(s, 0, sizeof s);
    memset(t, -1, sizeof t);
    fill(t[1],t[1]+ALPHA,0);
    s[0] = 1; 1[0] = 1[1] = -1; r[0] = r[1] = p[0] = p[1] = 0;
    rep(i,0,sz(a)) ukkadd(i, toi(a[i]));
  // example: find longest common substring (uses ALPHA = 28)
  pii best;
  int lcs(int node, int i1, int i2, int olen) {
    if (l[node] <= i1 && i1 < r[node]) return 1;</pre>
    if (1[node] <= i2 && i2 < r[node]) return 2;</pre>
    int mask = 0, len = node ? olen + (r[node] - 1[node]) : 0;
    rep(c, 0, ALPHA) if (t[node][c] != -1)
      mask |= lcs(t[node][c], i1, i2, len);
    if (mask == 3)
      best = max(best, {len, r[node] - len});
    return mask;
  static pii LCS(string s, string t) {
    SuffixTree st(s + (char) ('z' + 1) + t + (char) ('z' + 2));
    st.lcs(0, sz(s), sz(s) + 1 + sz(t), 0);
    return st.best;
Hashing.h
Description: Self-explanatory methods for string hashing.
// Arithmetic mod 2^64-1. 2x slower than mod 2^64 and more
// code, but works on evil test data (e.g. Thue-Morse, where
// ABBA... and BAAB... of length 2^10 hash the same mod 2^64).
// "typedef ull H;" instead if you think test data is random,
// or work mod 10^9+7 if the Birthday paradox is not a problem.
typedef uint64 t ull;
struct H {
  ull x; H(ull x=0) : x(x) {}
  H operator+(H \circ) { return x + \circ.x + (x + \circ.x < x); }
  H operator-(H o) { return *this + ~o.x; }
  H operator*(H o) { auto m = ( uint128 t)x * o.x;
    return H((ull)m) + (ull)(m >> 64); }
  ull get() const { return x + !~x; }
  bool operator==(H o) const { return get() == o.get(); }
  bool operator<(H o) const { return get() < o.get(); }</pre>
```

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

#### AhoCorasick.h

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

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

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

# Various (10)

## 10.1 Intervals

IntervalContainer.h

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

```
Time: \mathcal{O}(\log N)
                                                     d41d8c, 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) {
   R = max(R, it->second);
   before = it = is.erase(it);
  if (it != is.begin() && (--it)->second >= L) {
   L = min(L, it->first);
   R = max(R, it->second);
   is.erase(it);
  return is.insert(before, {L,R});
void removeInterval(set<pii>& is, int L, int R) {
  if (L == R) return;
 auto it = addInterval(is, L, R);
  auto r2 = it->second;
```

```
if (it->first == L) is.erase(it);
else (int&)it->second = L;
if (R != r2) is.emplace(R, r2);
```

#### IntervalCover.h

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

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

#### ConstantIntervals.h

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

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

```
d41d8c, 19 lines
template < class F, class G, class T>
void rec(int from, int to, F& f, G& g, int& i, T& p, T q) {
 if (p == q) return;
 if (from == to) {
   g(i, to, p);
    i = to; p = q;
    int mid = (from + to) >> 1;
    rec(from, mid, f, q, i, p, f(mid));
    rec(mid+1, to, f, g, i, p, q);
template < class F, class G>
void constantIntervals(int from, int to, F f, G g) {
 if (to <= from) return;</pre>
 int i = from; auto p = f(i), q = f(to-1);
 rec(from, to-1, f, q, i, p, q);
 g(i, to, q);
```

## 10.2 Misc. algorithms

#### 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))
                                                              d41d8c, 16 lines
int knapsack(vi w, int t) {
  int a = 0, b = 0, x;
```

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

## 10.3 Dynamic programming

#### KnuthDP.h

d41d8c, 19 lines

**Description:** When doing DP on intervals:  $a[i][j] = \min_{i < k < j} (a[i][k] + a[i][k])$ a[k][j] + f(i,j), where the (minimal) optimal k increases with both i and j, one can solve intervals in increasing order of length, and search k = p[i][j] for a[i][j] only between p[i][j-1] and p[i+1][j]. This is known as Knuth DP. Sufficient criteria for this are if  $f(b,c) \leq f(a,d)$  and  $f(a,c) + f(b,d) \le f(a,d) + f(b,c)$  for all  $a \le b \le c \le d$ . Consider also: LineContainer (ch. Data structures), monotone queues, ternary search. Time:  $\mathcal{O}(N^2)$ 

#### DivideAndConquerDP.h

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

```
Time: \mathcal{O}((N + (hi - lo)) \log N)
                                                                            d41d8c, 18 lines
```

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

## 10.4 Debugging tricks

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

## 10.5 Optimization tricks

\_builtin\_ia32\_ldmxcsr(40896); disables denormals (which make floats 20x slower near their minimum value).

#### 10.5.1 Bit hacks

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

#### 10.5.2 Pragmas

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

#### FastMod.h

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

```
typedef unsigned long long ull;
struct FastMod {
  ull b, m;
  FastMod(ull b) : b(b), m(-1ULL / b) {}
  ull reduce(ull a) { // a % b + (0 or b)
    return a - (ull)((__uint128_t(m) * a) >> 64) * b;
  }
};
```

#### FastInput.h

Description: debug

void dbg\_out() { cout << endl; }</pre>

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

Usage: ./a.out < input.txt

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

d41d8c, 17 lines

d41d8c, 37 lines

```
inline char gc() { // like getchar()
    static char buf[1 << 16];
    static size_t bc, be;
    if (bc >= be) {
        buf[0] = 0, bc = 0;
        be = fread(buf, 1, sizeof(buf), stdin);
    }
    return buf[bc++]; // returns 0 on EOF
}

int readInt() {
    int a, c;
    while ((a = gc()) < 40);
    if (a == '-') return -readInt();
    while ((c = gc()) >= 48) a = a * 10 + c - 480;
    return a - 48;
}
Debug.h
```

```
template <typename Head, typename... Tail> void dbg_out (Head H,
     Tail... T) {
  cout << ' ' << H;
 dbg_out(T...);
template <typename T, typename = void> struct is_iterable :
    false_type {};
template <typename T>
struct is iterable<
    T, void_t<decltype(begin(declval<T>())), decltype(end(
        declval<T>()))>>
    : true_type {};
template <typename T>
typename enable_if<is_iterable<T>::value && !is_same<T, string
    >::value,
                   ostream &>::type
operator << (ostream &cout, T const &v);
template <typename A, typename B>
ostream & operator << (ostream & cout, pair <A, B> const &p) {
 return cout << "(" << p.f << ", " << p.s << ")";
template <typename T>
typename enable_if<is_iterable<T>::value && !is_same<T, string
    >::value,
                   ostream &>::type
operator << (ostream &cout, T const &v) {
 cout << "[";
  for (auto it = v.begin(); it != v.end();) {
    cout << *it;
    if (++it != v.end())
      cout << ", ";
 return cout << "]";</pre>
#ifdef LOCAL
#define debug(...) cout << "(" << #__VA_ARGS__ << "):", dbg_out
     ( VA ARGS )
#else
#define debug(...) "zzz"
#endif
```

# 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 Ouadratic reciprocity Pollard-Rho Miller-Rabin Hensel lifting Vieta root jumping Game theory Combinatorial games Game trees Mini-max Nim Games on graphs Games on graphs with loops Grundy numbers Bipartite games without repetition General games without repetition Alpha-beta pruning Probability theory Optimization Binary search Ternary search Unimodality and convex functions Binary search on derivative Numerical methods Numeric integration Newton's method Root-finding with binary/ternary search Golden section search Matrices Gaussian elimination Exponentiation by squaring Sorting Radix sort Geometry Coordinates and vectors \* Cross product \* Scalar product Convex hull 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

23