

Massachusetts Institute of Technology

Benjamin Qi, Spencer Compton, Zhezheng Luo

adapted from KACTL and MIT NULL 2021-10-17

1 Contest

2 Mathematics

3 Data Structures

## TemplateShort .bashrc hash stress troubleshoot

| 4   | Number Theory   | 3                |
|---|---|------------------|
| 5   | Combinatorial   | 5                |
| 6   | Numerical   | 7                |
| 7   | Graphs  | 10               |
| 8   | Geometry  | 18               |
| 9   | Strings   | 22               |
| 10  | Various   | 24               |
| Contest (1)   |   |                  |
| Te  | mplateShort.cpp   | b3b6dc, 53 lines |
|   | clude <bits stdc++.h=""> ng namespace std;</bits>   |                  |
| <pre>using 11 = long long;<br/>using db = long double; // or double if tight TL<br/>using str = string;</pre> |   |                  |
| <pre>using pi = pair<int, int="">; #define mp make_pair #define f first #define s second</int,></pre>         |   |                  |
| tcT<br>tcT<br>usi<br>usi  | <pre>fine tcT template<class t=""> using V = vector<t>; , size_t SZ&gt; using AR = array<t,sz>; ng vi = V<int>; ng vb = V<bool>; ng vpi = V<pi>;</pi></bool></int></t,sz></t></class></pre>           |                  |
| #de<br>#de<br>#de<br>#de  | <pre>fine sz(x) int((x).size()) fine all(x) begin(x), end(x) fine sor(x) sort(all(x)) fine rsz resize fine pb push_back fine ft front() fine bk back()</pre>  |                  |
| #de<br>#de<br>#de<br>#de  | fine FOR(i,a,b) for (int i = (a); i < (b); ++i) fine FOR(i,a) FOR(i,0,a) fine ROF(i,a,b) for (int i = (b)-1; i >= (a);: fine ROF(i,a) ROF(i,0,a) fine rep(a) FOR(_,a) fine each(a,x) for (auto& a: x) | i)               |
| con   | st int MOD = 1e9+7;<br>st db PI = acos((db)-1);<br>9937 rng(0); // or mt19937_64  |                  |

```
1
   tcT> bool ckmin(T& a, const T& b) {
     return b < a ? a = b, 1 : 0; } // set a = min(a,b)
1
   tcT> bool ckmax(T& a, const T& b) {
     return a < b ? a = b, 1 : 0; } // set a = max(a,b)
\mathbf{2}
    tcT, class U> T fstTrue(T lo, T hi, U f) {
     ++hi; assert(lo <= hi); // assuming f is increasing
     while (lo < hi) { // find first index such that f is true
       T \text{ mid} = lo+(hi-lo)/2;
       f(mid) ? hi = mid : lo = mid+1;
     return lo;
   int main() { cin.tie(0)->sync_with_stdio(0); }
    .bashrc
   alias clr="printf '\33c'"
   co() { g++ -std=c++17 -02 -Wall -Wextra -Wshadow -Wconversion -
     \hookrightarrowo $1 $1.cpp; }
   run() { co $1 && ./$1; }
   hash.sh
    # Hash file ignoring whitespace and comments. Verifies that
    # code was correctly typed. Usage: 'sh hash.sh < A.cpp'
    cpp -dD -P -fpreprocessed|tr -d '[:space:]'|md5sum|cut -c-6
   stress.sh
    # A and B are executables you want to compare, gen takes int
    # as command line arg. Usage: 'sh stress.sh'
   for((i = 1; ; ++i)); do
       echo $i
        ./gen $i > int
       ./A < int > out1
        ./B < int > out2
       diff -w out1 out2 || break
       \# diff -w < (./A < int) < (./B < int) || break
    done
   troubleshoot.txt
   Write down most of your thoughts, even if you're not sure
   whether they're useful.
   Give your variables (and files) meaningful names.
   Stay organized and don't leave papers all over the place!
   You should know what your code is doing ...
   Pre-submit:
   Write a few simple test cases if sample is not enough.
   Are time limits close? If so, generate max cases.
   Is the memory usage fine?
   Could anything overflow?
   Remove debug output.
   Make sure to submit the right file.
   Wrong answer:
   Print your solution! Print debug output as well.
   Read the full problem statement again.
   Have you understood the problem correctly?
   Are you sure your algorithm works?
   Try writing a slow (but correct) solution.
```

Can your algorithm handle the whole range of input?

Did you consider corner cases (ex. n=1)?

```
Is your output format correct? (including whitespace)
Are you clearing all data structures between test cases?
Any uninitialized variables?
Any undefined behavior (array out of bounds)?
Any overflows or NaNs (or shifting 11 by >=64 bits)?
Confusing N and M, i and j, etc.?
Confusing ++i and i++?
Return vs continue vs break?
Are you sure the STL functions you use work as you think?
Add some assertions, maybe resubmit.
Create some test cases to run your algorithm on.
Go through the algorithm for a simple case.
Go through this list again.
Explain your algorithm to a teammate.
Ask the teammate to look at your code.
Go for a small walk, e.g. to the toilet.
Rewrite your solution from the start or let a teammate do it.
Geometry:
Work with ints if possible.
Correctly account for numbers close to (but not) zero. Related:
for functions like acos make sure absolute val of input is not
(slightly) greater than one.
Correctly deal with vertices that are collinear, concyclic,
```

#### Runtime error:

coplanar (in 3D), etc.

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?

Subtracting a point from every other (but not itself)?

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's your complexity? Large TL does not mean that something simple (like NlogN) isn't intended. Are you copying a lot of unnecessary data? (References) Avoid vector, map. (use arrays/unordered\_map) How big is the input and output? (consider FastIO) What do your teammates think about your algorithm?

#### Memory limit exceeded:

Calling count() on multiset?

What is the max amount of memory your algorithm should need? Are you clearing all data structures between test cases? If using pointers try BumpAllocator.

fe2351, 12 lines

## Mathematics (2)

## Trigonometry

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

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

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

$$\sin v + \sin w = 2\sin \frac{v+w}{2}\cos \frac{v-w}{2}$$

$$\cos v + \cos w = 2\cos \frac{v+w}{2}\cos \frac{v-w}{2}$$

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

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

## Geometry

## 2.2.1 Triangles

Side lengths: a, b, c

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

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

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

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

Length of median (divides triangle into two equal-area

triangles):  $m_a = \frac{1}{2}\sqrt{2b^2 + 2c^2 - a^2}$ 

Length of bisector (divides angles in two):

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

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

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

## 2.3 Derivatives/Integrals

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

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

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

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

Integration by parts:

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

## 2.4 Sums/Series

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

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

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

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

## Data Structures (3)

#### STL3.1

MapComparator.h

Description: example of function object (functor) for map or set

Usage: set<int, cmp> s; map<int, int, cmp> m; struct cmp{bool operator()(int l,int r)const{return l>r;}};

HashMap.h

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

Usage: ht<int,int> h({},{},{},{},{1<<16});

```
05a86f, 11 lines
<ext/pb_ds/assoc_container.hpp>
using namespace __gnu_pbds;
struct chash {
 const uint64_t C = 11(2e18*PI)+71; // large odd number
  const int RANDOM = rng();
  11 operator()(11 x) const {
    return __builtin_bswap64((x^RANDOM)*C); }
template<class K,class V> using um = unordered_map<K,V,chash>;
template<class K, class V> using ht = gp_hash_table<K, V, chash>;
template < class K, class V > V get (ht < K, V > & u, K x) {
 auto it = u.find(x); return it == end(u) ? 0 : it->s; }
```

#### IndexedSet.h

<ext/pb\_ds/assoc\_container.hpp>

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

```
using namespace __gnu_pbds;
tcT> using Tree = tree<T, null_type, less<T>,
 rb_tree_tag, tree_order_statistics_node_update>;
#define ook order_of_key
#define fbo find_by_order
void treeExample() {
 Tree<int> t, t2; t.insert(8);
 auto it = t.insert(10).f; assert(it == t.lb(9));
 assert(t.ook(10) == 1 && t.ook(11) == 2 && *t.fbo(0) == 8);
 t.join(t2); // assuming T < T2 or T > T2, merge t2 into t
```

## LCold.h

**Description:** Add lines of the form ax + b, query maximum y-coordinate for any x.

Time:  $\mathcal{O}(\log N)$ using T = 11; const T INF = LLONG\_MAX; // a/b rounded down

```
// 11 fdiv(11 a, 11 b) { return a/b-((a^b)<0&&a%b); }
bool _Q = 0;
struct Line {
T a, b; mutable T lst;
 T eval(T x) const { return a*x+b; }
 bool operator<(const Line&o)const{return _Q?lst<o.lst:a<o.a;}</pre>
 T last_gre(const Line& o) const { assert(a <= o.a);</pre>
    // greatest x s.t. a*x+b >= o.a*x+o.b
    return lst=(a==o.a?(b>=o.b?INF:-INF):fdiv(b-o.b,o.a-a));}
struct LineContainer: multiset<Line> {
 bool isect(iterator it) { auto n_it = next(it);
    if (n_it == end()) return it->lst = INF, 0;
    return it->last gre(*n it) >= n it->lst; }
  void add(T a, T b) { // remove lines after
    auto it = ins(\{a,b,0\}); while (isect(it)) erase(next(it));
    if (it == begin()) return;
    if (isect(--it)) erase(next(it)), isect(it);
    while (it != begin()) { // remove lines before
      --it; if (it->lst < next(it)->lst) break;
      erase(next(it)); isect(it); }
 T qmax(T x) { assert(!empty());
    _Q = 1; T res = lb(\{0,0,x\}) \rightarrow eval(x); _Q = 0;
    return res; }
};
```

## 1D Range Queries

## RMQ.h

Description: 1D range minimum query. Can also do queries for any associative operation in O(1) with D&C

Memory:  $\mathcal{O}(N \log N)$ Time:  $\mathcal{O}(1)$ 

```
tcT> struct RMQ {
  int level(int x) { return 31-__builtin_clz(x); }
 V<T> v; V<vi> jmp;
 int cmb(int a, int b) {
    return v[a] == v[b]?min(a,b):(v[a] < v[b]?a:b); }
  void init(const V<T>& _v) {
    v = v; jmp = \{vi(sz(v))\};
```

## SegTree LazySeg PSeg Treap ModIntShort

```
iota(all(jmp[0]),0);
  for (int j = 1; 1 << j <= sz(v); ++j) {
    jmp.pb(vi(sz(v) - (1 << j) +1));
   FOR(i,sz(jmp[j])) jmp[j][i] = cmb(jmp[j-1][i],
      jmp[j-1][i+(1<<(j-1))]);
int index(int 1, int r) {
  assert (1 \le r); int d = level(r-l+1);
  return cmb(jmp[d][1],jmp[d][r-(1<<d)+1]); }
T query(int 1, int r) { return v[index(1,r)]; }
```

## SegTree.h

Description: 1D point update, range query where cmb is any associative operation. If  $N=2^p$  then seg[1]==query(0,N-1).

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

tcT> struct SegTree { // cmb(ID,b) = b const T ID{}; T cmb(T a, T b) { return a+b; }

```
int n; V<T> seq;
  void init(int n) { // upd, query also work if n = n
   for (n = 1; n < _n; ) n *= 2;
   seg.assign(2*n,ID);}
  void pull(int p) { seq[p] = cmb(seq[2*p], seq[2*p+1]); }
  void upd(int p, T val) { // set val at position p
    seg[p += n] = val; for (p /= 2; p; p /= 2) pull(p); }
  T query(int 1, int r) { // associative op on [1, r]
   T ra = ID, rb = ID;
    for (1 += n, r += n+1; 1 < r; 1 /= 2, r /= 2) {
     if (1&1) ra = cmb(ra, seq[1++]);
     if (r&1) rb = cmb(seq[--r],rb);
    return cmb(ra,rb);
};
```

## LazySeg.h

**Description:** 1D range increment and sum query.

1630f3, 18 lines

```
8946fc, 26 lines
tcT, int SZ> struct LazvSeq {
  static assert(pct(SZ) == 1); // SZ must be power of 2
  const T ID = 0; T cmb(T a, T b) { return a+b; }
  T seg[2*SZ], lazy[2*SZ];
  LazySeg() { FOR(i,2*SZ) seg[i] = lazy[i] = ID; }
  void push(int ind, int L, int R) {
    seg[ind] += (R-L+1)*lazy[ind]; // dependent on operation
   if (L != R) F0R(i,2) lazy[2*ind+i] += lazy[ind];
   lazy[ind] = 0;
  } // recalc values for current node
  void pull(int ind) {seg[ind]=cmb(seg[2*ind], seg[2*ind+1]);}
  void build() { ROF(i,1,SZ) pull(i); }
  void upd(int lo,int hi,T inc,int ind=1,int L=0, int R=SZ-1) {
   push(ind,L,R); if (hi < L || R < lo) return;</pre>
    if (lo <= L && R <= hi) {
     lazy[ind] = inc; push(ind, L, R); return; }
    int M = (L+R)/2; upd(lo,hi,inc,2*ind,L,M);
    upd(lo,hi,inc,2*ind+1,M+1,R); pull(ind);
  T query(int lo, int hi, int ind=1, int L=0, int R=SZ-1) {
    push(ind,L,R); if (lo > R || L > hi) return ID;
    if (lo <= L && R <= hi) return seg[ind];
    int M = (L+R)/2; return cmb(query(lo,hi,2*ind,L,M),
      query(lo,hi,2*ind+1,M+1,R));
};
```

**Description:** Persistent min segtree with lazy updates, no propagation. If making d a vector then save the results of upd and build in local variables first to avoid issues when vector resizes in C++14 or lower. **Memory:**  $\mathcal{O}(N + Q \log N)$ 

```
8f37fa, 46 lines
tcT, int SZ> struct pseq {
 static const int LIM = 2e7:
 struct node {
   int 1, r; T val = 0, lazy = 0;
   void inc(T x) { lazy += x; }
   T get() { return val+lazy; }
 node d[LIM]; int nex = 0;
 int copy(int c) { d[nex] = d[c]; return nex++; }
 T cmb(T a, T b) { return min(a,b); }
 void pull(int c) { d[c].val =
    cmb(d[d[c].1].get(), d[d[c].r].get()); }
  //// MAIN FUNCTIONS
 T query(int c, int lo, int hi, int L, int R) {
   if (lo <= L && R <= hi) return d[c].get();</pre>
   if (R < lo || hi < L) return MOD;
   int M = (L+R)/2;
    return d[c].lazy+cmb(query(d[c].l,lo,hi,L,M),
              query(d[c].r,lo,hi,M+1,R));
 int upd(int c, int lo, int hi, T v, int L, int R) {
   if (R < lo || hi < L) return c;
   int x = copv(c);
   if (lo <= L && R <= hi) { d[x].inc(v); return x; }
   int M = (L+R)/2;
   d[x].l = upd(d[x].l, lo, hi, v, L, M);
   d[x].r = upd(d[x].r, lo, hi, v, M+1, R);
   pull(x); return x;
 int build(const V<T>& arr, int L, int R) {
   int c = nex++;
   if (L == R) {
     if (L < sz(arr)) d[c].val = arr[L];</pre>
    int M = (L+R)/2;
   d[c].l = build(arr, L, M), d[c].r = build(arr, M+1, R);
   pull(c); return c;
 vi loc; //// PUBLIC
 void upd(int lo, int hi, T v) {
   loc.pb(upd(loc.bk,lo,hi,v,0,SZ-1)); }
 T query(int ti, int lo, int hi) {
    return query(loc[ti],lo,hi,0,SZ-1); }
 void build(const V<T>&arr) {loc.pb(build(arr, 0, SZ-1));}
```

Description: Easy BBST. Use split and merge to implement insert and

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

bdb758, 65 lines

```
using pt = struct tnode*;
struct tnode {
 int pri, val; pt c[2]; // essential
 int sz; 11 sum; // for range queries
 bool flip = 0; // lazy update
 tnode (int val) {
   pri = rnq(); sum = val = _val;
   sz = 1; c[0] = c[1] = nullptr;
 ~tnode() { FOR(i,2) delete c[i]; }
```

```
int getsz(pt x) { return x?x->sz:0; }
11 getsum(pt x) { return x?x->sum:0; }
pt prop(pt x) { // lazy propagation
 if (!x || !x->flip) return x;
 swap (x->c[0], x->c[1]);
  x->flip = 0; F0R(i,2) if (x->c[i]) x->c[i]->flip ^= 1;
pt calc(pt x) {
 pt a = x->c[0], b = x->c[1];
  assert(!x->flip); prop(a), prop(b);
 x->sz = 1+getsz(a)+getsz(b);
 x->sum = x->val+getsum(a)+getsum(b);
 return x:
void tour(pt x, vi& v) { // print values of nodes,
 if (!x) return; // inorder traversal
 prop(x); tour(x->c[0],v); v.pb(x->val); tour(x->c[1],v);
pair<pt, pt> split(pt t, int v) { // >= v goes to the right
 if (!t) return {t,t};
 prop(t);
 if (t->val >= v) {
    auto p = split(t->c[0], v); t->c[0] = p.s;
    return {p.f,calc(t)};
 } else {
    auto p = split(t->c[1], v); t->c[1] = p.f;
    return {calc(t),p.s};
pair<pt, pt> splitsz(pt t, int sz) { // sz nodes go to left
 if (!t) return {t,t};
 prop(t);
 if (qetsz(t->c[0]) >= sz) {
    auto p = splitsz(t->c[0],sz); t->c[0] = p.s;
    return {p.f,calc(t)};
    auto p=splitsz(t->c[1],sz-qetsz(t->c[0])-1); t->c[1]=p.f;
    return {calc(t),p.s};
pt merge(pt 1, pt r) { // keys in 1 < keys in r
 if (!1 || !r) return 1?:r;
 prop(l), prop(r); pt t;
 if (1->pri > r->pri) 1->c[1] = merge(1->c[1],r), t = 1;
 else r - c[0] = merge(1, r - c[0]), t = r;
 return calc(t);
pt ins(pt x, int v) { // insert v
 auto a = split(x,v), b = split(a.s,v+1);
 return merge(a.f, merge(new tnode(v),b.s)); }
pt del(pt x, int v) { // delete v
 auto a = split(x,v), b = split(a.s,v+1);
 return merge(a.f,b.s); }
```

## Number Theory (4)

## 4.1 Modular Arithmetic

## ModIntShort.h

```
Description: Modular arithmetic.
```

```
Usage: mi a = MOD+5; cout << (int)inv(a); // 400000003 63d209, 20 lines
struct mi { // WARNING: needs some adjustment to work with FFT
  int v; explicit operator int() const { return v; }
  mi():v(0) {}
  mi(11 _v):v(int(_v%MOD)) { v += (v<0)*MOD; }
```

```
mi& operator+= (mi& a, mi b) {
  if ((a.v += b.v) >= MOD) a.v -= MOD;
  return a; }
mi& operator = (mi& a, mi b) {
  if ((a.v -= b.v) < 0) a.v += MOD;
  return a; }
mi operator+(mi a, mi b) { return a += b; }
mi operator-(mi a, mi b) { return a -= b; }
mi operator*(mi a, mi b) { return mi((ll)a.v*b.v); }
mi& operator *= (mi& a, mi b) { return a = a*b; }
mi pow(mi a, ll p) { assert(p >= 0); // won't work for negative
  return p==0?1:pow(a*a,p/2)*(p&1?a:1); }
mi inv(mi a) { assert(a.v != 0); return pow(a, MOD-2); }
mi operator/(mi a, mi b) { return a*inv(b); }
bool operator == (mi a, mi b) { return a.v == b.v; }
```

## ModFact.h

**Description:** pre-compute factorial mod inverses, assumes MOD is prime and SZ < MOD.

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

```
vmi invs, fac, ifac;
void genFac(int SZ) {
  invs.rsz(SZ), fac.rsz(SZ), ifac.rsz(SZ);
  invs[1] = fac[0] = ifac[0] = 1;
  FOR (i, 2, SZ) invs [i] = mi(-(11) MOD/i*(int) invs [MOD%i]);
  FOR(i,1,SZ) fac[i] = fac[i-1]*i, ifac[i] = ifac[i-1]*invs[i];
mi comb(int a, int b) {
  if (a < b || b < 0) return 0;
  return fac[a] * ifac[b] * ifac[a-b]; }
```

## ModMulLL.h

Description: Multiply two 64-bit integers mod another if 128-bit is not available. modMul is equivalent to (ul) (\_int128(a) \*b%mod). Works for  $0 \le a, b < mod < 2^{63}$ 530181, 9 lines

```
using ul = uint64 t;
ul modMul(ul a, ul b, const ul mod) {
 11 \text{ ret} = a*b-mod*(ul)((db)a*b/mod);
 return ret+((ret<0)-(ret>=(11)mod))*mod; }
ul modPow(ul a, ul b, const ul mod) {
 if (b == 0) return 1;
 ul res = modPow(a,b/2,mod); res = modMul(res,res,mod);
 return b&1 ? modMul(res,a,mod) : res;
```

#### FastMod.h

**Description:** Barrett reduction computes a%b about 4 times faster than usual where b > 1 is constant but not known at compile time. Division by bis replaced by multiplication by m and shifting right 64 bits.

```
using ul = uint64_t; using L = __uint128_t;
struct FastMod {
  ul b, m; FastMod(ul b) : b(b), m(-1ULL / b) {}
  ul reduce(ul a) {
    ul q = (ul) ((\underline{uint128_t (m)} * a) >> 64), r = a - q * b;
    return r - (r >= b) * b; }
};
```

### ModSart.h

**Description:** square root of integer mod a prime, -1 if doesn't exist.

```
"ModInt.h"
                                                              4ece48, 14 lines
using T = int;
T sgrt(mi a) {
```

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

```
mi p = pow(a, (MOD-1)/2);
if (p.v != 1) return p.v == 0 ? 0 : -1;
T s = MOD-1; int r = 0; while (s%2 == 0) s /= 2, ++r;
mi n = 1; while (pow(n, (MOD-1)/2).v != 1) n = T(n)+1;
// n non-square, ord(g)=2^r, ord(b)=2^m, ord(g)=2^r, m<r
for (mi \ x = pow(a, (s+1)/2), b = pow(a, s), q = pow(n, s);;) {
  if (b.v == 1) return min(x.v, MOD-x.v); // x^2=ab
  int m = 0; for (mi t = b; t.v != 1; t *= t) ++m;
  rep(r-m-1) q *= q; // ord(q) = 2^{m+1}
  x \neq q, q \neq q, b \neq q, r = m; // ord(q) = 2^m, ord(b) < 2^m
```

## ModSum.h

**Description:** Counts # of lattice points (x, y) s.t.  $1 \le x, 1 \le y, ax + by \le s$ (mod 2)<sup>64</sup> and related quantities.

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

23cbf6, 20 lines

```
using ul = uint64 t;
ul sum2(ul n) { return n/2*((n-1)|1); } // sum(0..n-1)
// \text{ return } | \{(x,y) \mid 1 \le x, 1 \le y, a*x+b*y \le S\} |
        = sum_{i=1}^{qs} (S-a*i)/b
ul triSum(ul a, ul b, ul s) { assert(a > 0 && b > 0);
 ul gs = s/a, rs = s%a; //ans = sum \{i=0\}^{gs-1}(i*a+rs)/b
 ul ad = a/b*sum2(qs)+rs/b*qs; a %= b, rs %= b;
 return ad+(a?triSum(b,a,a*gs+rs):0); // reduce if a \ge b
} // then swap x and y axes and recurse
// \text{ return sum}_{x=0}^{n-1} (a*x+b)/m
       = | f(x,y) | 0 < m*y <= a*x+b < a*n+b} |
// assuming a*n+b does not overflow
ul divSum(ul n, ul a, ul b, ul m) { assert(m > 0);
 ul extra = b/m*n; b %= m;
 return extra+(a?triSum(m,a,a*n+b):0); }
// \return sum {x=0}^{n-1} (a*x+b)%m
ul modSum(ul n, ll a, ll b, ul m) { assert(m > 0);
 a = (a\%m+m)\%m, b = (b\%m+m)\%m;
 return a*sum2(n)+b*n-m*divSum(n,a,b,m); }
```

## Primality

## **4.2.1** 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 1000000.

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

## 4.2.2 Divisors

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

The number of divisors of n is at most around 100 for n < 5e4, 500 for n < 1e7, 2000 for n < 1e10, 200000 for n < 1e19.

**Dirichlet Convolution:** Given a function f(x), let

```
(f * g)(x) = \sum_{d \mid x} g(d) f(x/d).
```

If the partial sums  $s_{f*q}(n)$ ,  $s_q(n)$  can be computed in O(1)and  $s_f(1...n^{2/3})$  can be computed in  $O(n^{2/3})$  then all  $s_f\left(\frac{n}{d}\right)$  can as well. Use

$$s_{f*g}(n) = \sum_{d=1}^{n} g(d)s_f(n/d).$$

If  $f(x) = \mu(x)$  then g(x) = 1, (f \* g)(x) = (x == 1), and  $s_f(n) = 1 - \sum_{i=2}^{n} s_f(n/i).$ 

If 
$$f(x) = \phi(x)$$
 then  $g(x) = 1$ ,  $(f * g)(x) = x$ , and  $s_f(n) = \frac{n(n+1)}{2} - \sum_{i=2}^n s_f(n/i)$ .

**Description:** Tests primality up to SZ. Runs faster if only odd indices are stored.

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

41c6ed, 20 lines

```
template<int SZ> struct Sieve {
 bitset<SZ> is prime; vi primes;
 Sieve() {
   is_prime.set(); is_prime[0] = is_prime[1] = 0;
   for (int i = 4; i < SZ; i += 2) is prime[i] = 0;
    for (int i = 3; i*i < SZ; i += 2) if (is_prime[i])
     for (int j = i*i; j < SZ; j += i*2) is_prime[j] = 0;
   FOR(i,SZ) if (is prime[i]) primes.pb(i);
 // int sp[SZ]{}; // smallest prime that divides
 // Sieve() { // above is faster
 // FOR(i,2,SZ) f
      if (sp[i] == 0) sp[i] = i, primes.pb(i);
       for (int p: primes) {
         if (p > sp[i] \mid | i * p >= SZ) break;
         sp[i*p] = p;
 // }
 // }
```

## MultiplicativePrefix.h

**Description:** Prefix sums of some multiplicative functions. Solve for all square-free numbers in range, then fix.

**Time:** faster than  $\mathcal{O}\left(N^{2/3}\right)$ ?

```
ll pre0(ll n, int i) { // prod(primes) in factorization of x
 ll res = n*(n+1)/2; // assume all square-free
 for (;;++i) {
   ll p = S.pr[i], nn = n/p/p; if (!nn) break;
    for(ll coef=p*(p-1);nn;nn/=p)res-=coef*pre0(nn,i+1);
 return res;
ll prel(ll n, int i) { // gcd of x and arithmetic derivative
 // p^e contributes p^e if e^p == 0 and p^e = 1 otherwise
 11 res = n; // assume all square-free
 for (;;++i) {
   ll p = S.pr[i], nn = n/p/p; if (!nn) break;
   11 lst = 1, mul = p*p;
    for (int e = 2; nn; mul *= p, nn /= p, ++e) {
```

574c0e, 10 lines

```
ll nex = mul; if (e%p) nex /= p;
   if (lst != nex) res += (nex-lst) *pre1(nn, i+1);
return res;
```

## PrimeCnt.h

**Description:** Counts number of primes up to N. Can also count sum of

```
Time: \mathcal{O}\left(N^{3/4}/\log N\right), 60ms for N=10^{11}, 2.5s for N=10^{13} co4e96, 20 lines
11 count primes(11 N) { // count primes(1e13) == 346065536839
 if (N <= 1) return 0;
  int sq = (int)sqrt(N);
  vl big_ans((sq+1)/2), small_ans(sq+1);
  FOR(i, 1, sq+1) small_ans[i] = (i-1)/2;
  FOR(i, sz(big\_ans)) big_ans[i] = (N/(2*i+1)-1)/2;
  vb skip(sq+1); int prime_cnt = 0;
  for (int p = 3; p \le sq; p += 2) if (!skip[p]) { // primes
    for (int j = p; j \le sq; j += 2*p) skip[j] = 1;
    FOR(j,min((ll)sz(big_ans),(N/p/p+1)/2)) {
      11 \text{ prod} = (11)(2*j+1)*p;
      big_ans[j] -= (prod > sq ? small_ans[(double)N/prod]
             : big_ans[prod/2])-prime_cnt;
    for (int j = sq, q = sq/p; q >= p; --q) for (; j >= q*p; --j)
      small_ans[j] -= small_ans[q]-prime_cnt;
    ++prime_cnt;
  return big_ans[0]+1;
```

#### MillerRabin.h

**Description:** Deterministic primality test, works up to 2<sup>64</sup>. For larger numbers, extend A randomly.

```
"ModMulLL.h"
                                                      89df33, 11 lines
bool prime (ul n) { // not 11!
 if (n < 2 | | n % 6 % 4 != 1) return n-2 < 2;
  ul A[] = \{2, 325, 9375, 28178, 450775, 9780504, 1795265022\},\
     s = builtin ctzll(n-1), d = n>>s;
  each(a,A) { // ^ count trailing zeroes
   ul p = modPow(a,d,n), i = s;
   while (p != 1 \&\& p != n-1 \&\& a%n \&\& i--) p = modMul(p,p,n);
   if (p != n-1 && i != s) return 0;
 return 1:
```

#### FactorFast.h

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

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

```
"MillerRabin.h", "ModMulLL.h"
ul pollard(ul n) { // return some nontrivial factor of n
  auto f = [n](ul x) \{ return modMul(x, x, n) + 1; \};
  ul x = 0, y = 0, t = 30, prd = 2, i = 1, q;
  while (t++ % 40 || gcd(prd, n) == 1) {
   if (x == y) x = ++i, y = f(x);
   if ((q = modMul(prd, max(x,y)-min(x,y), n))) prd = q;
   x = f(x), y = f(f(y));
  return gcd(prd, n);
void factor_rec(ul n, map<ul,int>& cnt) {
```

```
if (n == 1) return;
if (prime(n)) { ++cnt[n]; return; }
ul u = pollard(n);
factor_rec(u,cnt), factor_rec(n/u,cnt);
```

## Euclidean Algorithm

## FracInterval.h

**Description:** Given fractions a < b with non-negative numerators and denominators, finds fraction f with lowest denominator such that a < f < b. Should work with all numbers less than 2<sup>62</sup>.

```
pl bet(pl a, pl b) {
 11 num = a.f/a.s; a.f = num*a.s, b.f = num*b.s;
 if (b.f > b.s) return {1+num,1};
 auto x = bet(\{b.s, b.f\}, \{a.s, a.f\});
 return {x.s+num*x.f,x.f};
```

#### Euclid.h

Description: Generalized Euclidean algorithm. euclid and invGeneral work for  $A, B < 2^{62}$ . minBetween assumes that  $0 \le L \le R < B$ , works for  $AB < 2^{62}$  (same for min\_rem)

Time:  $\mathcal{O}(\log AB)$ 5ddb26, 49 lines // ceil(a/b)

```
// 11 cdiv(11 a, 11 b) { return a/b+((a^b)>0&a^b); }
pl euclid(11 A, 11 B) { // For A, B>=0, finds (x, y) s.t.
  // Ax+By=gcd(A,B), |Ax|, |By| \le AB/gcd(A,B)
 if (!B) return {1,0};
  pl p = euclid(B,A%B); return {p.s,p.f-A/B*p.s}; }
ll invGeneral(11 A, ll B) { // find x in [0,B) such that Ax=1
  pl p = euclid(A,B); assert(p.f*A+p.s*B == 1);
  return p.f+(p.f<0)*B; \} // must have gcd(A,B)=1
11 minBetween(11 A, 11 B, 11 L, 11 R) {
  // min x s.t. exists v s.t. L \le A*x-B*y \le R
  A %= B:
 if (L == 0) return 0;
  if (A == 0) return -1;
  ll k = cdiv(L,A); if (A*k \le R) return k;
  ll x = minBetween (B, A, A-R%A, A-L%A); // min x s.t. exists y
  // s.t. -R <= Bx-Av <= -L
  return x == -1 ? x : cdiv(B*x+L,A); // solve for y
// find min((Ax+C)%B) for 0 <= x <= M
// aka find minimum non-negative value of A*x-B*y+C
// where 0 <= x <= M, 0 <= y
11 minRemainder(11 A, 11 B, 11 C, 11 M) {
  assert (A >= 0 && B > 0 && C >= 0 && M >= 0);
 A %= B, C %= B; ckmin(M,B-1);
  if (A == 0) return C;
  if (C >= A) { // make sure C<A
    ll ad = cdiv(B-C,A);
    M \rightarrow ad; if (M < 0) return C;
    C += ad*A-B;
  11 q = B/A, new_B = B%A; // new_B < A
  if (new_B == 0) return C; // B-q*A
  // now minimize A*x-new B*y+C
  // where \theta \le x, y and x+q*y \le M, \theta \le C \le new_B \le A
  // q*y \rightarrow C-new_B*y
  if (C/new_B > M/q) return C-M/q*new_B;
  M -= C/new_B*q; C %= new_B; // now C < new_B
```

```
// given y, we can compute x = ceil[((B-q*A)*y-C)/A]
// so x+q*y = ceil((B*y-C)/A) <= M
11 \text{ max}_Y = (M*A+C)/B; // \text{ must have } y \leq max_Y
11 max_X = cdiv(new_B*max_Y-C,A); // must have x <= max_X</pre>
if (\max_X \times A - \text{new}_B \times \max_Y + C >= \text{new}_B) - -\max_X;
// now we can remove upper bound on y
return minRemainder(A, new_B, C, max_X);
```

## CRT.h

**Description:** Chinese Remainder Theorem.  $a.f \pmod{a.s}, b.f \pmod{b.s}$  $\implies$ ? (mod lcm(a.s, b.s)). Should work for  $ab < 2^6$ 

```
"Euclid.h"
pl CRT(pl a, pl b) {
 if (a.s < b.s) swap(a,b);
  11 x, y; tie(x, y) = euclid(a.s, b.s);
  11 q = a.s*x+b.s*y, 1 = a.s/q*b.s;
  if ((b.f-a.f)%g) return {-1,-1}; // no solution
  // ?*a.s+a.f \equiv b.f \pmod{b.s}
  // ?= (b.f-a.f)/g*(a.s/g)^{-1} \pmod{b.s/g}
  x = (b.f-a.f) %b.s \times x %b.s/q \times a.s + a.f;
  return \{x+(x<0)*1,1\};
```

## 4.4 Pythagorean Triples

The Pythagorean triples are uniquely generated by

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

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

## 4.5 Lifting the Exponent

For n > 0, p prime, and ints x, y s.t.  $p \nmid x, y$  and p|x-y:

```
• p \neq 2 or p = 2, 4|x - y|
  \implies v_n(x^n - y^n) = v_n(x - y) + v_n(n).
```

• 
$$p = 2, 2|n \implies v_2(x^n - y^n) = v_2((x^2)^{n/2} - (y^2)^{n/2}).$$

## Combinatorial (5)

## Permutations

## **5.1.1** Cycles

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

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

### 5.1.2 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 g (g.x = x).

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

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

## 5.2 Partitions and subsets

## 5.2.1 Partition function

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

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

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

## 5.2.2 Lucas' Theorem

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

## 5.3 General purpose numbers

## 5.3.1 Bernoulli numbers

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

Sums of powers:

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

Euler-Maclaurin formula for infinite sums:

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

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

## 5.3.2 Stirling numbers of the first kind

Number of permutations on n items with k cycles.

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

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

## 5.3.3 Eulerian numbers

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

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

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

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

## 5.3.4 Stirling numbers of the second kind

Partitions of n distinct elements into exactly k groups.

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

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

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

## 5.3.5 Bell numbers

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

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

## 5.3.6 Labeled unrooted trees

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

## 5.3.7 Catalan numbers

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

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

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

- 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.

## 5.4 Young Tableaux

Let a **Young diagram** have shape  $\lambda = (\lambda_1 \ge \cdots \ge \lambda_k)$ , where  $\lambda_i$  equals the number of cells in the *i*-th (left-justified) row from the top. A **Young tableau** of shape  $\lambda$  is a filling of the  $n = \sum \lambda_i$  cells with a permutation of  $1 \dots n$  such that each row and column is increasing.

**Hook-Length Formula**: For the cell in position (i, j), let  $h_{\lambda}(i, j) = |\{(I, J)|i \leq I, j \leq J, (I = i \text{ or } J = j)\}|$ . The number of Young tableaux of shape  $\lambda$  is equal to  $f^{\lambda} = \frac{n!}{\prod h_{\lambda}(i, j)}$ .

Schensted's Algorithm: converts a permutation  $\sigma$  of length n into a pair of Young Tableaux  $(S(\sigma), T(\sigma))$  of the same shape. When inserting  $x = \sigma_i$ ,

- 1. Add x to the first row of S by inserting x in place of the largest y with x < y. If y doesn't exist, push x to the end of the row, set the value of T at that position to be i, and stop.
- 2. Add y to the second row using the same rule, keep repeating as necessary.

All pairs  $(S(\sigma), T(\sigma))$  of the same shape correspond to a unique  $\sigma$ , so  $n! = \sum (f^{\lambda})^2$ . Also,  $S(\sigma^R) = S(\sigma)^T$ .

Let  $d_k(\sigma)$ ,  $a_k(\sigma)$  be the lengths of the longest subseqs which are a union of k decreasing/ascending subseqs, respectively. Then  $a_k(\sigma) = \sum_{i=1}^k \lambda_i, d_k(\sigma) = \sum_{i=1}^k \lambda_i^*$ , where  $\lambda_i^*$  is size of the i-th column.

## 5.5 Other

### DeBruijnSeg.h

**Description:** Given alphabet [0, k) constructs a cyclic string of length  $k^n$  that contains every length n string as substr.

a6961b. 13 lines

```
vi deBruijnSeq(int k, int n) {
   if (k == 1) return {0};
   vi seq, aux(n+1);
   function<void(int,int)> gen = [&](int t, int p) {
      if (t > n) { // +lyndon word of len p
         if (n*p == 0) FOR(i,1,p+1) seq.pb(aux[i]);
    } else {
      aux[t] = aux[t-p]; gen(t+1,p);
      while (++aux[t] < k) gen(t+1,t);
   }
};
gen(1,1); return seq;</pre>
```

## NimProduct.h

**Description:** Product of nimbers is associative, commutative, and distributive over addition (xor). Forms finite field of size  $2^{2^k}$ . Defined by  $ab = \max(\{a'b + ab' + a'b' : a' < a, b' < b\})$ . Application: Given 1D coin turning games  $G_1, G_2, G_1 \times G_2$  is the 2D coin turning game defined as follows. If turning coins at  $x_1, x_2, \ldots, x_m$  is legal in  $G_1$  and  $y_1, y_2, \ldots, y_n$  is legal in  $G_2$ , then turning coins at all positions  $(x_i, y_j)$  is legal assuming that the coin at  $(x_m, y_n)$  goes from heads to tails. Then the grundy function g(x, y) of  $G_1 \times G_2$  is  $g_1(x) \times g_2(y)$ .

Time: 64<sup>2</sup> xors per multiplication, memorize to speed up. 5afe17, 46 lines

```
using ul = uint64_t;
struct Precalc {
  ul tmp[64][64], y[8][8][256];
  unsigned char x[256][256];
  Precalc() { // small nim products, all < 256
    FOR(i, 256) FOR(j, 256) x[i][j] = mult < 8 > (i, j);
    FOR(i,8) FOR(j,i+1) FOR(k,256)
     y[i][j][k] = mult<64>(prod2(8*i,8*j),k);
  ul prod2(int i, int j) { // nim prod of 2^i, 2^j
   ul& u = tmp[i][j]; if (u) return u;
   if (!(i&j)) return u = 1ULL<<(i|j);</pre>
    int a = (i\&j)\&-(i\&j); // a=2^k, consider 2^{2^k}
   return u=prod2(i^a,j)^prod2((i^a)|(a-1),(j^a)|(i&(a-1)));
    // 2^{2^k} *2^{2^k} = 2^{2^k} +2^{2^k-1}
   }  // 2^{2^i}*2^{2^j} = 2^{2^i+2^j}   if i < j 
  template<int L> ul mult(ul a, ul b) {
    ul c = 0; F0R(i,L) if (a>>i&1)
     FOR(j,L) if (b>>j&1) c ^= prod2(i,j);
    return c;
  // 2^{8*i}*(a>>(8*i)&255) * 2^{8*j}*(b>>(8*j)&255)
  // \rightarrow (2^{8*i}*2^{8*i})*((a>>(8*i)&255)*(b>>(8*i)&255))
  ul multFast(ul a, ul b) const { // faster nim product
   ul res = 0; auto f=[](ul c,int d) {return c>>(8*d)&255;};
   F0R(i,8) {
     FOR(j,i) res ^= y[i][j][x[f(a,i)][f(b,j)]
              x[f(a,j)][f(b,i)];
      res ^= y[i][i][x[f(a,i)][f(b,i)]];
```

```
return res;
}
};
const Precalc P;

struct nb { // nimber
  ul x; nb() { x = 0; }
  nb(ul _x): x(_x) {}
  explicit operator ul() { return x; }
  nb operator+(nb y) { return nb(x^y.x); }
  nb operator+(nb y) { return nb(P.multFast(x,y.x)); }
  friend nb pow(nb b, ul p) {
    nb res = 1; for (;p;p/=2,b=b*b) if (p&1) res = res*b; return res; } // b^{2^2-2^2-1}=1 where 2^{2^2-2} > b
  friend nb inv(nb b) { return pow(b,-2); }
};
```

## MatroidIsect.h

return 0;

**Description:** Computes a set of maximum size which is independent in both graphic and colorful matroids, aka a spanning forest where no two edges are of the same color. In general, construct the exchange graph and find a shortest path. Can apply similar concept to partition matroid.

**Usage:** MatroidIsect<Gmat, Cmat> M(sz (ed), Gmat (ed), Cmat (col)) **Time:**  $\mathcal{O}\left(GI^{1.5}\right)$  calls to oracles, where G is size of ground set and I is size of independent set.

```
"DSU.h"
                                                     d0051c, 51 lines
struct Gmat { // graphic matroid
 int V = 0; vpi ed; DSU D;
 Gmat(vpi ed):ed(ed) {
   map < int, int > m; each(t,ed) m[t.f] = m[t.s] = 0;
    each(t,m) t.s = V++;
    each (t,ed) t.f = m[t.f], t.s = m[t.s];
 void clear() { D.init(V); }
 void ins(int i) { assert(D.unite(ed[i].f,ed[i].s)); }
 bool indep(int i) { return !D.sameSet(ed[i].f,ed[i].s); }
struct Cmat { // colorful matroid
 int C = 0; vi col; V<bool> used;
 Cmat(vi col):col(col) {each(t,col) ckmax(C,t+1); }
 void clear() { used.assign(C,0); }
 void ins(int i) { used[col[i]] = 1; }
 bool indep(int i) { return !used[col[i]]; }
template<class M1, class M2> struct MatroidIsect {
 int n; V<bool> iset; M1 m1; M2 m2;
 bool augment() {
   vi pre(n+1,-1); queue<int> q({n});
   while (sz(q)) {
     int x = q.ft; q.pop();
     if (iset[x]) {
       ml.clear(); F0R(i,n) if (iset[i] && i != x) ml.ins(i);
       FOR(i,n) if (!iset[i] && pre[i] == -1 && ml.indep(i))
          pre[i] = x, q.push(i);
     } else {
       auto backE = [&]() { // back edge
         FOR(c, 2) FOR(i, n) if((x==i | | iset[i]) &&(pre[i]==-1) ==c) {
           if (!m2.indep(i))return c?pre[i]=x,q.push(i),i:-1;
           m2.ins(i); }
         return n:
        for (int y; (y = backE()) != -1;) if (y == n) {
         for(; x != n; x = pre[x]) iset[x] = !iset[x];
         return 1; }
```

```
}
MatroidIsect(int n, M1 m1, M2 m2):n(n), m1(m1), m2(m2) {
    iset.assign(n+1,0); iset[n] = 1;
    m1.clear(); m2.clear(); // greedily add to basis
    R0F(i,n) if (m1.indep(i) && m2.indep(i))
    iset[i] = 1, m1.ins(i), m2.ins(i);
    while (augment());
}
```

## Numerical (6)

## 6.1 Matrix

### Matrix.h

**Description:** 2D matrix operations.

```
"ModInt.h"
                                                      c2e27f, 34 lines
using T = mi;
using Mat = V<V<T>>; // use array instead if tight TL
Mat makeMat(int r, int c) { return Mat(r, V<T>(c)); }
Mat makeId(int n) {
  Mat m = makeMat(n,n);
  FOR(i,n) m[i][i] = 1;
  return m;
Mat& operator+=(Mat& a, const Mat& b) {
  assert (sz(a) == sz(b) && sz(a[0]) == sz(b[0]));
  FOR(i, sz(a)) FOR(j, sz(a[0])) a[i][j] += b[i][j];
  return a;
Mat& operator = (Mat& a, const Mat& b) {
  assert (sz(a) == sz(b) && sz(a[0]) == sz(b[0]));
  FOR(i, sz(a)) FOR(j, sz(a[0])) a[i][j] -= b[i][j];
Mat operator*(const Mat& a, const Mat& b) {
  int x = sz(a), y = sz(a[0]), z = sz(b[0]);
  assert (y == sz(b)); Mat c = makeMat(x,z);
  FOR(i,x) FOR(j,y) FOR(k,z) c[i][k] += a[i][j]*b[j][k];
Mat operator+(Mat a, const Mat& b) { return a += b; }
Mat operator-(Mat a, const Mat& b) { return a -= b; }
Mat& operator*=(Mat& a, const Mat& b) { return a = a*b; }
Mat pow(Mat m, 11 p) {
  int n = sz(m); assert (n == sz(m[0]) \&\& p >= 0);
  Mat res = makeId(n);
  for (; p; p /= 2, m \star= m) if (p&1) res \star= m;
  return res;
```

#### | MatrixInv.h

**Description:** Uses gaussian elimination to convert into reduced row echelon form and calculates determinant. For determinant via arbitrary modulos, use a modified form of the Euclidean algorithm because modular inverse may not exist. If you have computed  $A^{-1} \pmod{p^k}$ , then the inverse  $\pmod{p^{2k}}$  is  $A^{-1}(2I-AA^{-1})$ .

**Time:**  $\mathcal{O}(N^3)$ , determinant of  $1000 \times 1000$  matrix of modints in 1 second if you reduce # of operations by half

```
const db EPS = 1e-9; // adjust?
int getRow(V<V<db>>& m, int R, int i, int nex) {
   pair<db,int> bes{0,-1}; // find row with max abs value
   FOR(j,nex,R) ckmax(bes,{abs(m[j][i]),j});
   return bes.f < EPS ? -1 : bes.s; }</pre>
```

```
int getRow(V<vmi>& m, int R, int i, int nex) {
   FOR(j, nex, R) if (m[j][i] != 0) return j;
    return -1; }
pair<T,int> gauss(Mat& m) { // convert to reduced row echelon
    if (!sz(m)) return {1,0};
   int R = sz(m), C = sz(m[0]), rank = 0, nex = 0;
   T prod = 1; // determinant
   F0R(i,C) {
        int row = getRow(m,R,i,nex);
        if (row == -1) { prod = 0; continue; }
       if (row != nex) prod *= -1, swap(m[row], m[nex]);
       prod *= m[nex][i]; rank++;
       T x = 1/m[nex][i]; FOR(k,i,C) m[nex][k] *= x;
       FOR(j,R) if (j != nex) {
           T v = m[j][i]; if (v == 0) continue;
            FOR(k,i,C) m[j][k] -= v*m[nex][k];
       }
       nex++;
    return {prod, rank};
Mat inv(Mat m) {
    int R = sz(m); assert(R == sz(m[0]));
   Mat x = makeMat(R, 2*R);
   F0R(i,R) {
       x[i][i+R] = 1;
       FOR(j,R) x[i][j] = m[i][j];
    if (gauss(x).s != R) return Mat();
   Mat res = makeMat(R,R);
   FOR(i,R) FOR(j,R) res[i][j] = x[i][j+R];
    return res;
```

## MatrixTree.h

Description: Kirchhoff's Matrix Tree Theorem. Given adjacency matrix, calculates # of spanning trees.

```
066e59, 11 lines
"MatrixInv.h"
T numSpan(Mat m) {
  int n = sz(m); Mat res(n-1, n-1);
  FOR(i,n) FOR(j,i+1,n) {
   mi ed = m[i][j]; res[i][i] += ed;
   if (j != n-1) {
     res[j][j] += ed;
      res[i][j] -= ed, res[j][i] -= ed;
  return gauss (res).f;
```

## ShermanMorrison.h

**Description:** Calculates  $(A + uv^T)^{-1}$  given  $B = A^{-1}$ . Not invertible if sum=0.

```
"MatrixInv.h"
                                                       3a3f34, 7 lines
void ad (Mat& B, const V<T>& u, const V<T>& v) {
 int n = sz(A); V < T > x(n), y(n);
 FOR(i,n) FOR(j,n)
   x[i] += B[i][j]*u[j], y[j] += v[i]*B[i][j];
  T sum = 1; FOR(i,n) FOR(j,n) sum += v[i]*B[i][j]*u[j];
 FOR(i,n) FOR(j,n) B[i][j] -= x[i]*y[j]/sum;
```

## 6.2 Polynomials

using T = mi; using poly = V<T>;

```
Polv.h
```

Description: Basic poly ops including division. Can replace T with double, complex.

```
void remz(poly& p) { while (sz(p)\&\&p.bk==T(0)) p.pop back(); }
poly REMZ(poly p) { remz(p); return p; }
poly rev(poly p) { reverse(all(p)); return p; }
poly shift (poly p, int x) {
 if (x \ge 0) p.insert(begin(p), x, 0);
 else assert(sz(p)+x \ge 0), p.erase(begin(p),begin(p)-x);
poly RSZ(const poly& p, int x) {
 if (x <= sz(p)) return poly(begin(p), begin(p)+x);
 poly q = p; q.rsz(x); return q; }
T eval(const poly& p, T x) { // evaluate at point x
 T res = 0; R0F(i,sz(p)) res = x*res+p[i];
poly dif(const poly& p) { // differentiate
 poly res; FOR(i,1,sz(p)) res.pb(T(i)*p[i]);
 return res; }
poly integ(const poly& p) { // integrate
 static poly invs{0,1};
 for (int i = sz(invs); i \le sz(p); ++i)
   invs.pb(-MOD/i*invs[MOD%i]);
 poly res(sz(p)+1); F0R(i,sz(p)) res[i+1] = p[i]*invs[i+1];
 return res;
poly& operator+=(poly& 1, const poly& r) {
 1.rsz(max(sz(1),sz(r))); FOR(i,sz(r)) 1[i] += r[i];
poly& operator -= (poly& 1, const poly& r) {
 1.rsz(max(sz(1),sz(r))); FOR(i,sz(r)) 1[i] -= r[i];
poly& operator *= (poly& 1, const T& r) { each(t,1) t *= r;
 return 1; }
poly& operator/=(poly& 1, const T& r) { each(t,1) t /= r;
 return 1; }
poly operator+(poly 1, const poly& r) { return 1 += r; }
poly operator-(poly 1, const poly& r) { return 1 -= r; }
poly operator-(poly 1) { each(t,1) t *= -1; return 1; }
poly operator*(poly 1, const T& r) { return 1 *= r; }
poly operator*(const T& r, const poly& 1) { return 1*r; }
poly operator/(poly 1, const T& r) { return 1 /= r; }
poly operator*(const poly& 1, const poly& r) {
 if (!min(sz(l),sz(r))) return {};
 poly x(sz(1)+sz(r)-1);
 FOR(i, sz(1)) FOR(j, sz(r)) x[i+j] += l[i]*r[j];
 return x:
poly& operator*=(poly& 1, const poly& r) { return 1 = 1*r; }
pair<poly, poly> quoRem(poly a, poly b) {
 remz(a); remz(b); assert(sz(b));
 T lst = b.bk, B = T(1)/lst; each(t,a) t *= B;
  each(t,b) t \star= B;
  poly q(max(sz(a)-sz(b)+1,0));
 for (int dif; (dif=sz(a)-sz(b)) >= 0; remz(a)) {
   q[dif] = a.bk; F0R(i,sz(b)) a[i+dif] -= q[dif]*b[i]; }
  each(t,a) t *= lst;
 return {q,a}; // quotient, remainder
poly operator% (const poly& a, const poly& b) {
 return quoRem(a,b).s; }
```

```
T resultant (poly a, poly b) { // R(A,B)
 // =b_m^n*prod_{j=1}^mA(mu_j)
 // =b_m^na_n^m*prod_{i=1}^nprod_{j=1}^m (mu_j-lambda_i)
 // = (-1)^{mn}a_n^m*prod_{i=1}^nB(lambda_i)
 // = (-1) ^{nm}R(B, A)
 // Also, R(A,B)=b_m^{deg(A)-deg(A-CB)}R(A-CB,B)
 int ad = sz(a)-1, bd = sz(b)-1;
 if (bd <= 0) return bd < 0 ? 0 : pow(b.bk,ad);
 int pw = ad; a = a\%b; pw -= (ad = sz(a) -1);
 return resultant(b,a)*pow(b.bk,pw)*T((bd&ad&1)?-1:1);
```

### PolvInterpolate.h

44d0dd, 73 lines

**Description:** n points determine unique polynomial of degree < n-1. For numerical precision pick v[k]  $f = c * \cos(k/(n-1) * \pi), k = 0 \dots n-1$ . Time:  $\mathcal{O}\left(n^2\right)$ 

```
"Poly.h"
                                                      aada3a, 8 lines
polv interpolate(V<pair<T,T>> v) {
 poly res, tmp{1};
 FOR(i,sz(v)) { T prod = 1; // add one point at a time
    FOR(j,i) v[i].s = prod*v[j].s, prod *= v[i].f-v[j].f;
   v[i].s \neq prod; res += v[i].s*tmp; tmp *= poly{-v[i].f,1};
  \} // add multiple of (x-v[0].f)*(x-v[1].f)*...*(x-v[i-1].f)
 return res;
```

## FFT.h

**Description:** Multiply polynomials of ints for any modulus  $< 2^{31}$ . For XOR convolution ignore m within fft.

```
Time: \mathcal{O}(N \log N)
"ModInt.h"
                                                      9d4a1a, 41 lines
// const int MOD = 998244353;
tcT> void fft(V<T>& A, bool inv = 0) { // NTT
  int n = sz(A); assert((T::mod-1)%n == 0); V<T> B(n);
  for (int b = n/2; b; b /= 2, swap (A, B)) { // w = n/b'th root
    T w = pow(T::rt(), (T::mod-1)/n*b), m = 1;
    for (int i = 0; i < n; i += b*2, m*= w) FOR (j, b) {
      T u = A[i+j], v = A[i+j+b] *m;
      B[i/2+j] = u+v; B[i/2+j+n/2] = u-v;
  if (inv) { reverse(1+all(A));
    Tz = T(1)/T(n); each(t,A) t *= z; }
} // for NTT-able moduli
tcT> V<T> mul(V<T> A, V<T> B) {
  if (!min(sz(A),sz(B))) return {};
  int s = sz(A) + sz(B) - 1, n = 1; for (; n < s; n *= 2);
  bool eq = A == B; A.rsz(n), fft(A);
  if (eq) B = A; // squaring A, reuse result
  else B.rsz(n), fft(B);
  FOR(i,n) A[i] \star= B[i];
  fft(A,1); A.rsz(s); return A;
template<class M, class T> V<M> mulMod(V<T> x, V<T> y) {
  auto con = [](const V<T>& v) {
    V < M > w(sz(v)); FOR(i,sz(v)) w[i] = (int)v[i];
    return w; };
  return mul(con(x),con(y));
} // arbitrary moduli
tcT> V<T> MUL(const V<T>& A, const V<T>& B) {
  using m0 = mint < (119 << 23) + 1,62 >; auto c0 = mulMod < m0 > (A,B);
  using m1 = mint < (5 << 25) + 1, 62>; auto c1 = mulMod < m1 > (A, B);
  using m2 = mint < (7 << 26) +1, 62>; auto c2 = mulMod < m2> (A,B);
  int n = sz(c0); V<T> res(n); m1 r01 = 1/m1(m0::mod);
  m2 r02 = 1/m2 (m0::mod), r12 = 1/m2 (m1::mod);
  F0R(i,n) { // a=remainder mod m0::mod, b fixes it mod m1::mod
    int a = c0[i].v, b = ((c1[i]-a)*r01).v,
```

```
c = (((c2[i]-a)*r02-b)*r12).v;
 res[i] = (T(c)*m1::mod+b)*m0::mod+a; // c fixes m2::mod
return res;
```

## PolyInvSimpler.h

**Description:** computes  $A^{-1}$  such that  $AA^{-1} \equiv 1 \pmod{x^n}$ . Newton's method: If you want F(x) = 0 and  $F(Q_k) \equiv 0 \pmod{x^a}$  then  $Q_{k+1} = Q_k - \frac{F(Q_k)}{F'(Q_k)} \pmod{x^{2a}}$  satisfies  $F(Q_{k+1}) \equiv 0 \pmod{x^{2a}}$ . Application: if f(n), g(n) are the #s of forests and trees on n nodes then  $\sum_{n=0}^{\infty} f(n)x^n = \exp\left(\sum_{n=1}^{\infty} \frac{g(n)}{n!}\right)$ 

**Usage:** vmi v= $\{1,5,2,3,4\}$ ; ps $(\exp(2*\log(v,9),9))$ ; // squares v Time:  $\mathcal{O}(N \log N)$ 

```
"PolyConv.h"
poly inv(poly A, int n) { // Q-(1/Q-A)/(-Q^{-2})}
  poly B{1/A[0]};
  for (int x = 2; x/2 < n; x *= 2)
   B = 2*B-RSZ (conv(RSZ(A,x),conv(B,B)),x);
  return RSZ(B,n);
poly sqrt(const poly& A, int n) { //Q-(Q^2-A)/(2Q)
  assert(A[0] == 1); poly B\{1\};
  for (int x = 2; x/2 < n; x *= 2)
   B = T(1)/T(2) *RSZ(B+conv(RSZ(A,x),inv(B,x)),x);
  return RSZ(B,n);
// return {quotient, remainder}
pair<poly, poly> divi(const poly& f, const poly& g) {
  if (sz(f) < sz(g)) return {{},f};</pre>
  poly q = conv(inv(rev(q), sz(f) - sz(q) + 1), rev(f));
  q = rev(RSZ(q, sz(f) - sz(g) + 1));
  poly r = RSZ(f-conv(q,q),sz(q)-1); return \{q,r\};
poly log(poly A, int n) { assert(A[0] == 1); // (In A)' = A'/A
  A.rsz(n); return integ(RSZ(conv(dif(A),inv(A,n-1)),n-1)); }
poly exp(poly A, int n) { assert(A[0] == 0);
  polv B{1}, IB{1};
  for (int x = 1; x < n; x *= 2) {
    IB = 2*IB-RSZ(conv(B,conv(IB,IB)),x); // inverse of B to x
    poly Q = dif(RSZ(A,x)); Q += RSZ(conv(IB, dif(B) - conv(B,Q))
       \hookrightarrow, 2 * x - 1);
    // first x-1 terms of dif(B)-conv(B,O) are zero
    B = B+RSZ(conv(B,RSZ(A,2*x)-integ(0)),2*x);
  } // We know that Q=A' is B'/B to x-1 places, we want to find
     \hookrightarrow B'/B to 2x-1 places
  return RSZ(B,n);
// poly expOld(poly A, int n) \{ // Q-(lnQ-A)/(1/Q) \}
// assert(A[0] == 0); poly B = {1};
// while (sz(B) < n) f int x = 2*sz(B);
// B = RSZ(B+conv(B,RSZ(A,x)-log(B,x)),x); }
// return RSZ(B,n);
```

## $\operatorname{Misc}$

#### LinRec.h

**Description:** Berlekamp-Massey, computes linear recurrence C of order Nfor sequence s of 2N terms.

```
Usage: LinRec L; L.init({0,1,1,2,3,5,8}); // Fibonacci
Time: init \Rightarrow \mathcal{O}(N|C|), eval \Rightarrow \mathcal{O}(|C|^2 \log p).
```

"Poly.h" e63980, 32 lines struct LinRec {

```
void BM() { // find smallest C such that C[0]=1 and
  // for all i \ge sz(C)-1, sum_{j=0}^{sz(C)-1}C[j]*s[i-j]=0
  // If we treat C and s as polynomials in D, then
  // for all i \ge sz(C)-1, [D^i]C*s=0
  int x = 0; T b = 1;
  poly B; B = C = \{1\}; // B is fail vector
  FOR(i,sz(s)) { // update C after adding a term of s
    ++x; int L = sz(C), M = i+3-L;
    T d = 0; FOR(j,L) d += C[j]*s[i-j]; // [D^i]C*s
    if (d == 0) continue; // [D^i]C \times s = 0
    poly _C = C; T coef = d/b;
    C.rsz(max(L,M)); FOR(j,sz(B)) C[j+x] -= coef*B[j];
    if (L < M) B = _C, b = d, x = 0;
void init(const poly& _s) {
  s = s; BM();
  rC = C; reverse(all(rC)); // poly for getPow
  C.erase(begin(C)); each(t,C) t \star = -1;
poly getPow(ll p) { // get x^p mod rC
  if (p == 0) return {1};
  poly r = getPow(p/2); r = (r*r) %rC;
  return p&1?(r*poly{0,1})%rC:r;
T dot(poly v) { // dot product with seq
  T ans = 0; FOR(i,sz(v)) ans += v[i]*s[i];
  return ans; } // get p-th term of rec
T eval(ll p) { assert(p >= 0); return dot(getPow(p)); }
```

## Integrate.h

**Description:** Integration of a function over an interval using Simpson's rule, exact for polynomials of degree up to 3. The error should be proportional to  $dif^4$ , although in practice you will want to verify that the result is stable to desired precision when epsilon changes. 3ebaab, 7 lines

```
// db f(db x) { return x*x+3*x+1; }
template < class F > db quad (F f, db a, db b) {
 const int n = 1000;
 db dif = (b-a)/2/n, tot = f(a)+f(b);
 FOR(i,1,2*n) tot += f(a+i*dif)*(i&1?4:2);
 return tot*dif/3;
```

### IntegrateAdaptive.h

Description: Unused. Fast integration using adaptive Simpson's rule, exact for polynomials of degree up to 5.

```
Usage: db z, y;
db h(db x) { return x*x + y*y + z*z \le 1; }
db g(db y) \{ :: y = y; \text{ return quad(h, } -1, 1); \}
db f(db z) \{ :: z = z; \text{ return quad}(q, -1, 1); \}
db sphereVol = quad(f,-1,1), pi = sphereVol*3/4;
                                                       3b316e, 10 lines
template < class F > db simpson(F f, db a, db b) {
 db c = (a+b)/2; return (f(a)+4*f(c)+f(b))*(b-a)/6; }
template < class F > db rec (F& f, db a, db b, db eps, db S) {
  db c = (a+b)/2;
  db S1 = simpson(f,a,c), S2 = simpson(f,c,b), T = S1+S2;
  if (abs(T-S)<=15*eps || b-a<1e-10) return T+(T-S)/15;
  return rec(f,a,c,eps/2,S1)+rec(f,c,b,eps/2,S2);
template < class F > db quad (F f, db a, db b, db eps = 1e-8) {
  return rec(f,a,b,eps,simpson(f,a,b)); }
```

## Simplex.h

**Description:** Solves a general linear maximization problem: maximize  $c^T x$ subject to Ax < b, x > 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 \text{ val} = LPSolver(A, b, c).solve(x);
```

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

```
\mathcal{O}\left(2^{N}\right) in the general case.
```

c99f9c, 67 lines

```
using T = db; // double probably suffices
using vd = V<T>; using vvd = V<vd>;
const T eps = 1e-8, inf = 1/.0;
#define ltj(X) if (s=-1 \mid | mp(X[j],N[j]) < mp(X[s],N[s])) s=j
struct LPSolver {
 int m, n; // # m = contraints, # n = variables
 vi N, B; // N[i] = non-basic variable (i-th column), = 0
  vvd D; // B[j] = basic variable (j-th row)
  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)) {
    FOR(i,m) FOR(j,n) D[i][j] = A[i][j];
    FOR(i,m) B[i] = n+i, D[i][n] = -1, D[i][n+1] = b[i];
    // B[i]: basic variable for each constraint
    // D[i][n]: artificial variable for testing feasibility
    FOR(j,n) N[j] = j, D[m][j] = -c[j];
    // D[m] stores negation of objective,
    // which we want to minimize
    N[n] = -1; D[m+1][n] = 1; // to find initial feasible
  } // solution, minimize artificial variable
  void pivot(int r, int s) { // swap B[r] (row)
    T inv = 1/D[r][s]; // with N[r] (column)
    FOR(i, m+2) if (i != r && abs(D[i][s]) > eps) {
     T binv = D[i][s]*inv;
      FOR(j, n+2) if (j != s) D[i][j] -= D[r][j]*binv;
      D[i][s] = -binv;
    D[r][s] = 1; FOR(j, n+2) D[r][j] *= inv; // scale r-th row
    swap(B[r],N[s]);
 bool simplex(int phase) {
   int x = m+phase-1;
    while (1) { // if phase=1, ignore artificial variable
     int s = -1; FOR(j, n+1) if (N[j] != -phase) ltj(D[x]);
      // find most negative col for nonbasic (NB) variable
      if (D[x][s] >= -eps) return 1;
      // can't get better sol by increasing NB variable
      int r = -1;
      F0R(i,m) {
       if (D[i][s] <= eps) continue;</pre>
        if (r == -1 \mid | mp(D[i][n+1] / D[i][s], B[i])
              < mp(D[r][n+1] / D[r][s], B[r])) r = i;
       // find smallest positive ratio
      } // -> max increase in NB variable
      if (r == -1) return 0; // objective is unbounded
      pivot(r,s);
 T solve(vd& x) { // 1. check if x=0 feasible
    int r = 0; FOR(i,1,m) if (D[i][n+1] < D[r][n+1]) r = i;
    if (D[r][n+1] < -eps) { // if not, find feasible start}
      pivot(r,n); // make artificial variable basic
      assert(simplex(2)); // I think this will always be true??
      if (D[m+1][n+1] < -eps) return -inf;</pre>
      // D[m+1][n+1] is max possible value of the negation of
      // artificial variable, optimal value should be zero
```

```
// if exists feasible solution
FOR(i,m) if (B[i] == -1) { // artificial var basic
    int s = 0; FOR(j,1,n+1) ltj(D[i]); // -> nonbasic
    pivot(i,s);
}
bool ok = simplex(1); x = vd(n);
FOR(i,m) if (B[i] < n) x[B[i]] = D[i][n+1];
return ok ? D[m][n+1] : inf;
};</pre>
```

# Graphs (7)

**Erdos-Gallai:**  $d_1 \ge \cdots \ge d_n$  can be degree sequence of simple graph on n vertices iff their sum is even and  $\sum_{i=1}^k d_i \le k(k-1) + \sum_{i=k+1}^n \min(d_i, k), \forall 1 \le k \le n$ .

## 7.1 Cycles

# DirectedCycle.h Description: skack

3504a5, 17 lines

```
template<int SZ> struct DirCyc {
    vi adj[SZ], stk, cyc; vb inStk, vis;
    void dfs(int x) {
        stk.pb(x); inStk[x] = vis[x] = 1;
        each(i,adj[x]) {
            if (inStk[i]) cyc = {find(all(stk),i),end(stk)};
            else if (!vis[i]) dfs(i);
            if (sz(cyc)) return;
        }
        stk.pop_back(); inStk[x] = 0;
    }
    vi init(int N) {
        inStk.rsz(N), vis.rsz(N);
        FOR(i,N) if (!vis[i] && !sz(cyc)) dfs(i);
        return cyc;
    }
};
```

#### NegativeCvcle.h

**Description:** use Bellman-Ford (make sure no underflow)

688ec8, 11 lines

```
vi negCyc(int N, V<pair<pi,int>> ed) {
  vl d(N); vi p(N); int x = -1;
  rep(N) {
    x = -1; each(t,ed) if (ckmin(d[t.f.s],d[t.f.f]+t.s))
    p[t.f.s] = t.f.f, x = t.f.s;
    if (x == -1) return {};
}
  rep(N) x = p[x]; // enter cycle
  vi cyc{x}; while (p[cyc.bk] != x) cyc.pb(p[cyc.bk]);
  reverse(all(cyc)); return cyc;
}
```

## 7.2 DSU

#### DSII b

**Description:** Disjoint Set Union with path compression and union by size. Add edges and test connectivity. Use for Kruskal's or Boruvka's minimum spanning tree.

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

e42a83, 11 lines

```
struct DSU {
    vi e; void init(int N) { e = vi(N,-1); }
    int get(int x) { return e[x] < 0 ? x : e[x] = get(e[x]); }
    bool sameSet(int a, int b) { return get(a) == get(b); }
    int size(int x) { return -e[get(x)]; }
    bool unite(int x, int y) { // union by size
        x = get(x), y = get(y); if (x == y) return 0;
        if (e[x] > e[y]) swap(x,y);
        e[x] += e[y]; e[y] = x; return 1;
    }
};
```

## 7.3 Trees

## LCAjump.h

**Description:** Calculates least common ancestor in tree with verts  $0 \dots N-1$  and root R using binary jumping.

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

**Time:**  $\mathcal{O}(N \log N)$  build,  $\mathcal{O}(\log N)$  query

6b0ee9, 28 lines

```
struct LCA {
 int N; V<vi> par, adj; vi depth;
 void init(int _N) { N = _N;
   int d = 1; while ((1 << d) < N) ++d;
   par.assign(d, vi(N)); adj.rsz(N); depth.rsz(N);
 void ae(int x, int y) { adj[x].pb(y), adj[y].pb(x); }
 void gen(int R = 0) { par[0][R] = R; dfs(R); }
 void dfs(int x = 0) {
   FOR(i, 1, sz(par)) par[i][x] = par[i-1][par[i-1][x]];
   each(y,adj[x]) if (y != par[0][x])
      depth[y] = depth[par[0][y]=x]+1, dfs(y);
 int jmp(int x, int d) {
   FOR(i, sz(par)) if ((d>>i)&1) x = par[i][x];
    return x; }
 int lca(int x, int y) {
    if (depth[x] < depth[y]) swap(x,y);</pre>
    x = jmp(x, depth[x] - depth[y]); if (x == y) return x;
   R0F(i,sz(par)) {
     int X = par[i][x], Y = par[i][y];
     if (X != Y) x = X, y = Y;
    return par[0][x];
 int dist(int x, int y) { // # edges on path
    return depth[x]+depth[y]-2*depth[lca(x,y)]; }
};
```

## LCArmq.h

**Description:** Euler Tour LCA. Compress takes a subset S of nodes and computes the minimal subtree that contains all the nodes pairwise LCAs and compressing edges. Returns a list of (par, orig\_index) representing a tree rooted at 0. The root points to itself.

```
struct LCA {
  int N; V<vi> adj;
  vi depth, pos, par, rev; // rev is for compress
  vpi tmp; RMQ<pi> r;
  void init(int _N) { N = _N; adj.rsz(N);
    depth = pos = par = rev = vi(N); }
  void ae(int x, int y) { adj[x].pb(y), adj[y].pb(x); }
  void dfs(int x) {
    pos[x] = sz(tmp); tmp.eb(depth[x],x);
    each(y,adj[x]) if (y != par[x]) {
        depth[y] = depth[par[y]=x]+1, dfs(y);
    }
}
```

```
tmp.eb(depth[x],x); }

void gen(int R = 0) { par[R] = R; dfs(R); r.init(tmp); }
int lca(int u, int v){
    u = pos[u], v = pos[v]; if (u > v) swap(u,v);
    return r.query(u,v).s; }
int dist(int u, int v) {
    return depth[u]+depth[v]-2*depth[lca(u,v)]; }

vpi compress(vi S) {
    auto cmp = [&](int a, int b) { return pos[a] < pos[b]; };
    sort(all(S),cmp); ROF(i,sz(S)-1) S.pb(lca(S[i],S[i+1]));
    sort(all(S),cmp); S.erase(unique(all(S)),end(S));
    vpi ret{{0,S[0]}}; FOR(i,sz(S)) rev[S[i]] = i;
    FOR(i,1,sz(S)) ret.eb(rev[lca(S[i-1],S[i])],S[i]);
    return ret;
}
</pre>
```

#### HLD.h

**Description:** Heavy-Light Decomposition, add val to verts and query sum in path/subtree.

**Time:** any tree path is split into  $\mathcal{O}(\log N)$  parts

```
"LazySeg.h"
                                                    1802e2, 48 lines
template<int SZ, bool VALS_IN_EDGES> struct HLD {
 int N; vi adi[SZ];
 int par[SZ], root[SZ], depth[SZ], sz[SZ], ti;
 int pos[SZ]; vi rpos; // rpos not used but could be useful
 void ae(int x, int y) { adj[x].pb(y), adj[y].pb(x); }
 void dfsSz(int x) {
   sz[x] = 1;
    each(y,adj[x]) {
     par[y] = x; depth[y] = depth[x]+1;
     adj[v].erase(find(all(adj[v]),x));
     dfsSz(y); sz[x] += sz[y];
     if (sz[y] > sz[adj[x][0]]) swap(y,adj[x][0]);
 void dfsHld(int x) {
   pos[x] = ti++; rpos.pb(x);
    each(y,adj[x]) {
     root[y] = (y == adj[x][0] ? root[x] : y);
     dfsHld(y); }
 void init(int N, int R = 0) { N = N;
   par[R] = depth[R] = ti = 0; dfsSz(R);
   root[R] = R; dfsHld(R);
 int lca(int x, int y) {
    for (; root[x] != root[y]; y = par[root[y]])
     if (depth[root[x]] > depth[root[y]]) swap(x,y);
    return depth[x] < depth[y] ? x : y;</pre>
 LazySeg<11,SZ> tree; // segtree for sum
 template <class BinaryOp>
 void processPath(int x, int y, BinaryOp op) {
    for (; root[x] != root[y]; y = par[root[y]]) {
     if (depth[root[x]] > depth[root[y]]) swap(x,y);
     op(pos[root[y]],pos[y]); }
    if (depth[x] > depth[y]) swap(x,y);
    op (pos[x]+VALS_IN_EDGES, pos[y]);
 void modifyPath(int x, int y, int v) {
   processPath(x,y,[this,&v](int 1, int r) {
     tree.upd(l,r,v); }); }
 11 queryPath(int x, int y) {
   11 res = 0; processPath(x,y,[this,&res](int 1, int r) {
      res += tree.query(1,r); });
    return res; }
```

```
void modifySubtree(int x, int v) {
 tree.upd(pos[x]+VALS_IN_EDGES, pos[x]+sz[x]-1,v); }
```

## Centroid.h

**Description:** The centroid of a tree of size N is a vertex such that after removing it, all resulting subtrees have size at most  $\frac{N}{2}$ . Supports updates in the form "add 1 to all verts v such that dist(x, v) < y."

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

```
Time: \mathcal{O}(N \log N) build, \mathcal{O}(\log N) update and query
                                                     907e21, 54 lines
void ad(vi& a, int b) { ckmin(b, sz(a)-1); if (b>=0) a[b]++; }
void prop(vi& a) { R0F(i,sz(a)-1) a[i] += a[i+1]; }
template<int SZ> struct Centroid {
  vi adj[SZ]; void ae(int a,int b){adj[a].pb(b),adj[b].pb(a);}
 bool done[SZ]; // processed as centroid yet
  int N, sub[SZ], cen[SZ], lev[SZ]; // subtree size, centroid anc
  int dist[32-__builtin_clz(SZ)][SZ]; // dists to all ancs
  vi stor[SZ], STOR[SZ];
  void dfs(int x, int p) { sub[x] = 1;
   each(y,adj[x]) if (!done[y] && y != p)
     dfs(y,x), sub[x] += sub[y];
  int centroid(int x) {
    dfs(x,-1);
    for (int sz = sub[x];;) {
     pi mx = \{0, 0\};
     each(y,adj[x]) if (!done[y] && sub[y] < sub[x])
       ckmax(mx, {sub[y], y});
     if (mx.f*2 \le sz) return x;
     x = mx.s;
  void genDist(int x, int p, int lev) {
   dist[lev][x] = dist[lev][p]+1;
   each(y,adj[x]) if (!done[y] && y != p) genDist(y,x,lev);
  void gen(int CEN, int _x) { // CEN = centroid above x
   int x = centroid(_x); done[x] = 1; cen[x] = CEN;
    sub[x] = sub[x]; lev[x] = (CEN == -1 ? 0 : lev[CEN]+1);
   dist[lev[x]][x] = 0;
    stor[x].rsz(sub[x]),STOR[x].rsz(sub[x]+1);
   each(y,adj[x]) if (!done[y]) genDist(y,x,lev[x]);
   each(y,adj[x]) if (!done[y]) gen(x,y);
  void init(int N) { N = N; FOR(i, 1, N+1) done[i] = 0;
   gen(-1,1); } // start at vert 1
  void upd(int x, int y) {
   int cur = x, pre = -1;
   R0F(i, lev[x]+1) {
     ad(stor[cur],y-dist[i][x]);
     if (pre != -1) ad(STOR[pre], y-dist[i][x]);
     if (i > 0) pre = cur, cur = cen[cur];
  } // call propAll() after all updates
  void propAll() { FOR(i,1,N+1) prop(stor[i]), prop(STOR[i]); }
  int query(int x) { // get value at vertex x
   int cur = x, pre = -1, ans = 0;
   ROF(i, lev[x]+1) { // if pre != -1, subtract those from
     ans += stor[cur][dist[i][x]]; // same subtree
     if (pre != -1) ans -= STOR[pre][dist[i][x]];
     if (i > 0) pre = cur, cur = cen[cur];
    return ans;
```

## 7.3.1 SqrtDecompton

HLD generally suffices. If not, here are some common strategies:

- Rebuild the tree after every  $\sqrt{N}$  queries.
- Consider vertices with > or  $<\sqrt{N}$  degree separately.
- For subtree updates, note that there are  $O(\sqrt{N})$ distinct sizes among child subtrees of any node.

Block Tree: Use a DFS to split edges into contiguous groups of size  $\sqrt{N}$  to  $2\sqrt{N}$ .

Mo's Algorithm for Tree Paths: Maintain an array of vertices where each one appears twice, once when a DFS enters the vertex (st) and one when the DFS exists (en). For a tree path  $u \leftrightarrow v$  such that st[u] < st[v],

- If u is an ancestor of v, query [st[u], st[v]].
- Otherwise, query [en[u], st[v]] and consider LCA(u, v) separately.

Solutions with worse complexities can be faster if you optimize the operations that are performed most frequently. Use arrays instead of vectors whenever possible. Iterating over an array in order is faster than iterating through the same array in some other order (ex. one given by a random permutation) or DFSing on a tree of the same size. Also, the difference between  $\sqrt{N}$  and the optimal block (or buffer) size can be quite large. Try up to 5x smaller or larger (at least).

## 7.4 DFS Algorithms

## EulerPath.h

Description: Eulerian path starting at src if it exists, visits all edges exactly once. Works for both directed and undirected. Returns vector of  $\{\text{vertex}, \text{label of edge to vertex}\}$ . Second element of first pair is always -1. Time:  $\mathcal{O}(N+M)$ 

template<bool directed> struct Euler { int N; V<vpi> adj; V<vpi::iterator> its; vb used; void init(int \_N) { N = \_N; adj.rsz(N); } void ae(int a, int b) { int M = sz(used); used.pb(0); adj[a].eb(b,M); if (!directed) adj[b].eb(a,M); } vpi solve(int src = 0) { its.rsz(N); F0R(i,N) its[i] = begin(adj[i]); vpi ans, s{{src,-1}}; // {{vert,prev vert},edge label} int lst = -1; // ans generated in reverse order while (sz(s)) { int x = s.bk.f; auto& it=its[x], en=end(adj[x]); while (it != en && used[it->s]) ++it; if (it == en) { // no more edges out of vertex if (lst != -1 && lst != x) return {};

```
// not a path, no tour exists
                ans.pb(s.bk); s.pop_back(); if (sz(s)) lst=s.bk
            } else s.pb(*it), used[it->s] = 1;
       } // must use all edges
        if (sz(ans) != sz(used)+1) return {};
        reverse(all(ans)); return ans;
};
```

11

#### SCCT.h

**Description:** Tarjan's, DFS once to generate strongly connected components in topological order. a, b in same component if both  $a \to b$  and  $b \to a$ exist. Uses less memory than Kosaraju b/c doesn't store reverse edges.

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

```
struct SCC {
 int N, ti = 0; V<vi> adj;
 vi disc, comp, stk, comps;
 void init(int _N) { N = _N, adj.rsz(N);
    disc.rsz(N), comp.rsz(N,-1);}
  void ae(int x, int y) { adj[x].pb(y); }
  int dfs(int x) {
    int low = disc[x] = ++ti; stk.pb(x);
    each(y,adj[x]) if (comp[y] == -1) // comp[y] == -1,
      ckmin(low, disc[y]?:dfs(y)); // disc[y] != 0 -> in stack
    if (low == disc[x]) { // make new SCC
      // pop off stack until you find x
      comps.pb(x); for (int y = -1; y != x;)
        comp[y = stk.bk] = x, stk.pop_back();
    return low;
  void gen() {
    FOR(i, N) if (!disc[i]) dfs(i);
    reverse(all(comps));
};
```

## TwoSAT.h

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

Usage: TwoSat ts;

```
ts.either(0, \sim3); // Var 0 is true or var 3 is false
ts.setVal(2); // Var 2 is true
ts.atMostOne(\{0, \sim 1, 2\}); // <= 1 of vars 0, \sim 1 and 2 are true
ts.solve(N); // Returns true iff it is solvable
ts.ans[0..N-1] holds the assigned values to the vars
                                                       805e1c, 32 lines
struct TwoSAT {
  int N = 0; vpi edges;
  void init(int _N) { N = _N; }
  int addVar() { return N++; } // for atMostOne
  void either(int x, int y) {
    x = max(2*x, -1-2*x), y = max(2*y, -1-2*y);
    edges.eb(x,v); }
  void implies (int x, int y) { either (\sim x, y); }
  void must(int x) { either(x,x); }
  void atMostOne(const vi& li) {
    if (sz(li) <= 1) return;
    int cur = \simli[0];
    FOR(i, 2, sz(li)) {
      int next = addVar();
```

either(cur,~li[i]); either(cur,next);

either(~li[i],next); cur = ~next;

## BCC MaximalCliques Dinic GomoryHu MCMF

```
either(cur,~li[1]);
  vb solve(int N = -1) {
   if (N != -1) N = N;
   SCC S; S.init(2*N);
   each (e,edges) S.ae (e.f^1,e.s), S.ae (e.s^1,e.f);
   S.gen(); reverse(all(S.comps)); // reverse topo order
    for (int i = 0; i < 2*N; i += 2)
     if (S.comp[i] == S.comp[i^1]) return {};
   vi tmp(2*N); each(i,S.comps) if (!tmp[i])
     tmp[i] = 1, tmp[S.comp[i^1]] = -1;
    vb ans(N); FOR(i,N) ans[i] = tmp[S.comp[2*i]] == 1;
};
```

## BCC.h

Description: Biconnected components of edges. Removing any vertex in BCC doesn't disconnect it. To get block-cut tree, create a bipartite graph with the original vertices on the left and a vertex for each BCC on the right. Draw edge  $u \leftrightarrow v$  if u is contained within the BCC for v. Self-loops are not included in any BCC while BCCS of size 1 represent bridges.

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

0625a6, 35 lines

```
struct BCC {
  V<vpi> adj; vpi ed;
 V<vi> edgeSets, vertSets; // edges for each bcc
  int N, ti = 0; vi disc, stk;
  void init(int _N) { N = _N; disc.rsz(N), adj.rsz(N); }
  void ae(int x, int y) {
   adj[x].eb(y,sz(ed)), adj[y].eb(x,sz(ed)), ed.eb(x,y); }
  int dfs(int x, int p = -1) { // return lowest disc
   int low = disc[x] = ++ti;
   each(e,adj[x]) if (e.s != p) {
     if (!disc[e.f]) {
       stk.pb(e.s); // disc[x] < LOW -> bridge
       int LOW = dfs(e.f,e.s); ckmin(low,LOW);
       if (disc[x] <= LOW) { // get edges in bcc
          edgeSets.eb(); vi& tmp = edgeSets.bk; // new bcc
          for (int y = -1; y != e.s; )
            tmp.pb(y = stk.bk), stk.pop_back();
     } else if (disc[e.f] < disc[x]) // back-edge</pre>
       ckmin(low, disc[e.f]), stk.pb(e.s);
   return low;
  void gen() {
   F0R(i,N) if (!disc[i]) dfs(i);
   vb in(N);
    each(c,edgeSets) { // edges contained within each BCC
     vertSets.eb(); // so you can easily create block cut tree
     auto ad = [\&] (int x) {
       if (!in[x]) in[x] = 1, vertSets.bk.pb(x); };
      each(e,c) ad(ed[e].f), ad(ed[e].s);
      each(e,c) in[ed[e].f] = in[ed[e].s] = 0;
};
```

#### MaximalCliques.h

**Description:** Used only once. Finds all maximal cliques.

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

f5cd93, 16 lines

```
using B = bitset<128>; B adj[128];
// possibly in clique, not in clique, in clique
void cliques (B P = \simB(), B X={}, B R={}) {
```

```
if (!P.any()) {
  if (!X.any()) // do smth with R
  return;
int q = (P|X)._Find_first();
// clique must contain q or non-neighbor of q
B cands = P\&\sim adj[q];
FOR(i,N) if (cands[i]) {
  R[i] = 1; cliques(P&adj[i], X&adj[i], R);
  R[i] = P[i] = 0; X[i] = 1;
```

## 7.5 Flows

**Konig's Theorem:** In a bipartite graph, max matching = min vertex cover.

Dilworth's Theorem: For any partially ordered set, the sizes of the max antichain and of the min chain decomposition are equal. Equivalent to Konig's theorem on the bipartite graph (U, V, E) where U = V = S and (u, v) is an edge when u < v. Those vertices outside the min vertex cover in both U and V form a max antichain.

## Dinic.h

template<class F> struct Dinic {

**Description:** Fast flow. After computing flow, edges  $\{u, v\}$  such that  $lev[u] \neq 0$ , lev[v] = 0 are part of min cut. Time:  $\mathcal{O}(N^2M)$  flow

```
struct Edge { int to, rev; F cap; };
int N; V<V<Edge>> adj;
void init(int _N) { N = _N; adj.rsz(N); }
pi ae(int a, int b, F cap, F rcap = 0) {
  assert(min(cap,rcap) >= 0); // saved me > once
  adj[a].pb({b,sz(adj[b]),cap});
  adj[b].pb({a,sz(adj[a])-1,rcap});
  return {a,sz(adj[a])-1};
F edgeFlow(pi loc) { // get flow along original edge
  const Edge& e = adj.at(loc.f).at(loc.s);
  return adj.at(e.to).at(e.rev).cap;
vi lev, ptr;
bool bfs(int s, int t) { // level=shortest dist from source
  lev = ptr = vi(N);
  lev[s] = 1; queue < int > q({s});
  while (sz(q)) { int u = q.ft; q.pop();
    each(e,adj[u]) if (e.cap && !lev[e.to]) {
```

q.push(e.to), lev[e.to] = lev[u]+1;

for (int& i = ptr[v]; i < sz(adj[v]); i++) {</pre>

if (F df = dfs(e.to,t,min(flo,e.cap))) {

if (lev[e.to]!=lev[v]+1||!e.cap) continue;

return df; } // saturated >=1 one edge

e.cap -= df; adj[e.to][e.rev].cap += df;

if (e.to == t) return 1;

F dfs(int v, int t, F flo) {

if (v == t) return flo;

Edge& e = adj[v][i];

return 0;

```
return 0;
 F maxFlow(int s, int t) {
    F \text{ tot} = 0; while (bfs(s,t)) while (F df =
      dfs(s,t,numeric_limits<F>::max())) tot += df;
};
```

## GomorvHu.h

**Description:** Returns edges of Gomory-Hu tree (second element is weight). Max flow between pair of vertices of undirected graph is given by min edge weight along tree path. Uses the fact that for any  $i, j, k, \lambda_{ik} \geq \min(\lambda_{ij}, \lambda_{jk})$ , where  $\lambda_{ij}$  denotes the flow between i and j.

**Time:** N - 1 calls to Dinic

```
"Dinic.h"
                                                      0d712e, 16 lines
template < class F > V < pair < pi, F >> gomoryHu (int N,
    const V<pair<pi,F>>& ed) {
 vi par(N); Dinic<F> D; D.init(N);
  vpi ed_locs; each(t,ed)ed_locs.pb(D.ae(t.f.f,t.f.s,t.s,t.s));
  V<pair<pi,F>> ans;
  FOR(i,1,N) {
    each(p,ed_locs) { // reset capacities
      auto& e = D.adj.at(p.f).at(p.s);
      auto& e_rev = D.adj.at(e.to).at(e.rev);
      e.cap = e_rev.cap = (e.cap+e_rev.cap)/2;
    ans.pb({{i,par[i]},D.maxFlow(i,par[i])});
    FOR(j, i+1, N) if (par[j] == par[i] \&\& D.lev[j]) par[j] = i;
 return ans;
```

## MCMF.h

c76643, 43 lines

Description: Minimum-cost maximum flow, assumes no negative cycles. It is possible to choose negative edge costs such that the first run of Dijkstra is slow, but this hasn't been an issue in the past. Edge weights  $\geq 0$  for every subsequent run. To get flow through original edges, assign ID's during ae. **Time:** Ignoring first run of Dijkstra,  $\mathcal{O}(FM \log M)$  if caps are integers and F is max flow.

```
77bfb0, 46 lines
struct MCMF {
 using F = 11; using C = 11; // flow type, cost type
 struct Edge { int to, rev; F flo, cap; C cost; };
 int N; V<C> p, dist; vpi pre; V<V<Edge>> adj;
 void init(int _N) { N = _N;
    p.rsz(N), adj.rsz(N), dist.rsz(N), pre.rsz(N); }
  void ae(int u, int v, F cap, C cost) { assert(cap >= 0);
    adj[u].pb({v,sz(adj[v]),0,cap,cost});
    adj[v].pb({u,sz(adj[u])-1,0,0,-cost});
  } // use asserts, don't try smth dumb
 bool path(int s, int t) { // send flow through lowest cost
    const C inf = numeric_limits<C>::max(); dist.assign(N,inf);
    using T = pair<C, int>;
    priority_queue<T, V<T>, greater<T>> todo;
    todo.push(\{dist[s] = 0, s\});
    while (sz(todo)) { // Dijkstra
     T x = todo.top(); todo.pop();
     if (x.f > dist[x.s]) continue;
      each(e,adj[x.s]) { // all weights should be non-negative
       if (e.flo < e.cap && ckmin(dist[e.to],
            x.f+e.cost+p[x.s]-p[e.to]))
          pre[e.to]={x.s,e.rev}, todo.push({dist[e.to],e.to});
    } // if costs are doubles, add some EPS so you
    // don't traverse ~0-weight cycle repeatedly
```

## Hungarian UnweightedMatch WeightedMatch

```
return dist[t] != inf; // true if augmenting path
  pair<F,C> calc(int s, int t) { assert(s != t);
   FOR(_,N) FOR(i,N) each(e,adj[i]) // Bellman-Ford
     if (e.cap) ckmin(p[e.to],p[i]+e.cost);
    F \text{ totFlow} = 0; C \text{ totCost} = 0;
    while (path(s,t)) { // p -> potentials for Dijkstra
     FOR(i,N) p[i] += dist[i]; // don't matter for unreachable
     F df = numeric_limits<F>::max();
      for (int x = t; x != s; x = pre[x].f) {
        Edge& e = adj[pre[x].f][adj[x][pre[x].s].rev];
        ckmin(df,e.cap-e.flo); }
      totFlow += df; totCost += (p[t]-p[s]) *df;
      for (int x = t; x != s; x = pre[x].f) {
        Edge& e = adj[x][pre[x].s]; e.flo -= df;
        adj[pre[x].f][e.rev].flo += df;
    } // get max flow you can send along path
    return {totFlow,totCost};
};
```

## Matching

## Hungarian.h

**Description:** Given array of (possibly negative) costs to complete each of N(1-indexed) jobs w/ each of M workers  $(N \le M)$ , finds min cost to complete all jobs such that each worker is assigned to at most one job. Dijkstra with potentials works in almost the same way as MCMF.

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

09d0ec, 28 lines

```
using C = 11:
C hungarian(const V<V<C>>& a) {
  int N = sz(a)-1, M = sz(a[0])-1; assert (N \le M);
  V<C> u(N+1), v(M+1); // potentials to make edge weights >= \theta
  vi job (M+1);
  FOR(i,1,N+1) { // find alternating path with job i
    const C inf = numeric_limits<C>::max();
    int w = 0; job[w] = i; // add "dummy" worker 0
    V<C> dist(M+1, inf); vi pre(M+1, -1); vb done(M+1);
    while (job[w]) { // dijkstra
     done[w] = 1; int j = job[w], nexW; C delta = inf;
      // fix dist[j], update dists from j
     FOR(W,M+1) if (!done[W]) { // try all workers
       if (ckmin(dist[W],a[j][W]-u[j]-v[W])) pre[W] = w;
        if (ckmin(delta,dist[W])) nexW = W;
     FOR(W,M+1) { // subtract constant from all edges going
        // from done -> not done vertices, lowers all
        // remaining dists by constant
        if (done[W]) u[job[W]] += delta, v[W] -= delta;
       else dist[W] -= delta;
     w = nexW;
    } // potentials adjusted so all edge weights >= 0
    for (int W; w; w = W) job[w] = job[W = pre[w]];
  \} // job[w] = 0, found alternating path
  return -v[0]; // min cost
```

#### UnweightedMatch.h

Description: Edmond's Blossom Algorithm. General unweighted matching with 1-based indexing. If vis[v]=0 when bfs returns 0, v is not part of every max matching.

**Time:**  $\mathcal{O}(N^3)$ , faster in practice

513ca2, 54 lines

```
template<int SZ> struct UnweightedMatch {
 int match[SZ], N; vi adj[SZ];
```

```
void ae(int u, int v) { adj[u].pb(v), adj[v].pb(u); }
 queue<int> q;
 int par[SZ], vis[SZ], orig[SZ], aux[SZ];
 void augment(int u, int v) { // toggle edges on u-v path
   while (1) { // one more matched pair
      int pv = par[v], nv = match[pv];
      match[v] = pv; match[pv] = v;
     v = nv; if (u == pv) return;
 int lca(int u, int v) { // find LCA of supernodes in O(dist)
    static int t = 0;
    for (++t;;swap(u,v)) {
     if (!u) continue;
     if (aux[u] == t) return u; // found LCA
     aux[u] = t; u = orig[par[match[u]]];
 void blossom(int u, int v, int a) { // go other way
   for (; orig[u] != a; u = par[v]) { // around cycle
     par[u] = v; v = match[u]; // treat u as if vis[u] = 1
     if (vis[v] == 1) vis[v] = 0, q.push(v);
     orig[u] = orig[v] = a; // merge into supernode
 bool bfs(int u) { // u is initially unmatched
   FOR(i, N+1) par[i] = 0, vis[i] = -1, orig[i] = i;
   q = queue < int > (); vis[u] = 0, q.push(u);
    while (sz(q)) { // each node is pushed to q at most once
     int v = q.ft; q.pop(); // 0 -> unmatched vertex
      each(x,adj[v]) {
       if (vis[x] == -1) { // neither of x, match[x] visited
         vis[x] = 1; par[x] = v;
         if (!match[x]) return augment(u,x),1;
         vis[match[x]] = 0, q.push(match[x]);
        } else if (vis[x] == 0 && orig[v] != orig[x]) {
          int a = lca(orig[v],orig[x]); // odd cycle
         blossom(x,v,a), blossom(v,x,a);
        } // contract O(n) times
   return 0;
 int calc(int _N) { // rand matching -> constant improvement
   N = N; FOR(i, N+1) match[i] = aux[i] = 0;
    int ans = 0; vi V(N); iota(all(V),1); shuffle(all(V),rng);
      \hookrightarrow // find rand matching
    each(x,V) if (!match[x]) each(y,adj[x]) if (!match[y]) {
     match[x] = y, match[y] = x; ++ans; break; }
    FOR(i,1,N+1) if (!match[i] && bfs(i)) ++ans;
    return ans;
};
```

## WeightedMatch.h

Description: General max weight max matching with 1-based indexing. Edge weights must be positive, combo of UnweightedMatch and Hungarian. Time:  $\mathcal{O}(N^3)$ ? 120873, 145 lines

```
template<int SZ> struct WeightedMatch {
 struct edge { int u,v,w; }; edge g[SZ*2][SZ*2];
 void ae(int u, int v, int w) { g[u][v].w = g[v][u].w = w; }
 int N,NX,lab[SZ*2],match[SZ*2],slack[SZ*2],st[SZ*2];
 int par[SZ*2],floFrom[SZ*2][SZ],S[SZ*2],aux[SZ*2];
 vi flo[SZ*2]; queue<int> q;
 void init(int _N) { N = _N; // init all edges
   FOR(u, 1, N+1) FOR(v, 1, N+1) g[u][v] = {u, v, 0};
 int eDelta(edge e) { // >= 0 at all times
    return lab[e.u]+lab[e.v]-g[e.u][e.v].w*2; }
```

```
void updSlack(int u, int x) { // smallest edge -> blossom x
  if (!slack[x] \mid | eDelta(g[u][x]) < eDelta(g[slack[x]][x]))
    slack[x] = u; }
void setSlack(int x) {
  slack[x] = 0; FOR(u, 1, N+1) if (g[u][x].w > 0
   && st[u] != x && S[st[u]] == 0) updSlack(u,x); }
void qPush(int x) {
  if (x \le N) q.push(x);
  else each(t,flo[x]) qPush(t); }
void setSt(int x, int b) {
  st[x] = b; if (x > N) each(t,flo[x]) setSt(t,b); }
int getPr(int b, int xr) { // get even position of xr
  int pr = find(all(flo[b]),xr)-begin(flo[b]);
  if (pr&1) { reverse(1+all(flo[b])); return sz(flo[b])-pr; }
  return pr; }
void setMatch(int u, int v) { // rearrange flo[u], matches
  edge e = g[u][v]; match[u] = e.v; if (u <= N) return;
  int xr = floFrom[u][e.u], pr = getPr(u,xr);
  FOR(i,pr) setMatch(flo[u][i],flo[u][i^1]);
  setMatch(xr,v); rotate(begin(flo[u]),pr+all(flo[u])); }
void augment (int u, int v) { // set matches including u->v
  while (1) { // and previous ones
    int xnv = st[match[u]]; setMatch(u,v);
    if (!xnv) return;
    setMatch(xnv,st[par[xnv]]);
    u = st[par[xnv]], v = xnv;
int lca(int u, int v) { // same as in unweighted
  static int t = 0; // except maybe return 0
  for (++t;u||v;swap(u,v)) {
   if (!u) continue;
    if (aux[u] == t) return u;
    aux[u] = t; u = st[match[u]];
    if (u) u = st[par[u]];
  return 0;
void addBlossom(int u, int anc, int v) {
  int b = N+1; while (b \le NX \&\& st[b]) ++b;
  if (b > NX) ++NX; // new blossom
  lab[b] = S[b] = 0; match[b] = match[anc]; flo[b] = {anc};
  auto blossom = [&](int x) {
    for (int y; x != anc; x = st[par[y]])
      flo[b].pb(x), flo[b].pb(y = st[match[x]]), qPush(y);
  blossom(u); reverse(1+all(flo[b])); blossom(v); setSt(b,b);
  // identify all nodes in current blossom
  FOR(x, 1, NX+1) q[b][x].w = q[x][b].w = 0;
  FOR(x, 1, N+1) floFrom[b][x] = 0;
  each(xs,flo[b]) { // find tightest constraints
    FOR(x,1,NX+1) if (g[b][x].w == 0 \mid \mid eDelta(g[xs][x]) <
      eDelta(g[b][x])) g[b][x]=g[xs][x], g[x][b]=g[x][xs];
    FOR(x, 1, N+1) if (floFrom[xs][x]) floFrom[b][x] = xs;
  } // floFrom to deconstruct blossom
  setSlack(b); // since didn't qPush everything
void expandBlossom(int b) {
  each(t,flo[b]) setSt(t,t); // undo setSt(b,b)
  int xr = floFrom[b][q[b][par[b]].u], pr = getPr(b,xr);
  for (int i = 0; i < pr; i += 2) {
    int xs = flo[b][i], xns = flo[b][i+1];
    par[xs] = q[xns][xs].u; S[xs] = 1; // no setSlack(xns)?
    S[xns] = slack[xs] = slack[xns] = 0; qPush(xns);
  S[xr] = 1, par[xr] = par[b];
  FOR(i,pr+1,sz(flo[b])) { // matches don't change
   int xs = flo[b][i]; S[xs] = -1, setSlack(xs); }
  st[b] = 0; // blossom killed
```

13

## MaxMatchLexMin MaxMatchFast ChordalGraphRecognition

```
bool onFoundEdge(edge e) {
    int u = st[e.u], v = st[e.v];
    if (S[v] == -1) { // v unvisited, matched with smth else
     par[v] = e.u, S[v] = 1; slack[v] = 0;
      int nu = st[match[v]]; S[nu] = slack[nu] = 0; qPush(nu);
    } else if (S[v] == 0) {
      int anc = lca(u,v); // if 0 then match found!
      if (!anc) return augment(u,v), augment(v,u),1;
     addBlossom(u,anc,v);
   return 0;
  bool matching() {
    q = queue<int>();
    FOR(x, 1, NX+1) {
     S[x] = -1, slack[x] = 0; // all initially unvisited
     if (st[x] == x \&\& !match[x]) par[x] = S[x] = 0, qPush(x);
    if (!sz(q)) return 0;
    while (1) {
      while (sz(q)) { // unweighted matching with tight edges
       int u = q.ft; q.pop(); if (S[st[u]] == 1) continue;
        FOR(v, 1, N+1) if (g[u][v].w > 0 && st[u] != st[v]) {
          if (eDelta(g[u][v]) == 0) { // condition is strict}
            if (onFoundEdge(g[u][v])) return 1;
          } else updSlack(u,st[v]);
      int d = INT MAX;
      FOR(b, N+1, NX+1) if (st[b] == b \&\& S[b] == 1)
       ckmin(d, lab[b]/2); // decrease lab[b]
      FOR(x,1,NX+1) if (st[x] == x \&\& slack[x]) {
       if (S[x] == -1) ckmin(d,eDelta(g[slack[x]][x]));
        else if (S[x] == 0) ckmin(d, eDelta(g[slack[x]][x])/2);
      } // edge weights shouldn't go below 0
      FOR(u,1,N+1) {
        if (S[st[u]] == 0) {
          if (lab[u] <= d) return 0; // why?</pre>
          lab[u] -= d;
       } else if (S[st[u]] == 1) lab[u] += d;
      } // lab has opposite meaning for verts and blossoms
     FOR(b, N+1, NX+1) if (st[b] == b \&\& S[b] != -1)
       lab[b] += (S[b] == 0 ? 1 : -1) *d*2;
      q = queue<int>();
      FOR(x,1,NX+1) if (st[x]==x \&\& slack[x] // new tight edge
        && st[slack[x]] != x && eDelta(g[slack[x]][x]) == 0)
          if (onFoundEdge(g[slack[x]][x])) return 1;
      FOR(b, N+1, NX+1) if (st[b] == b \&\& S[b] == 1 \&\& lab[b] == 0)
        expandBlossom(b); // odd dist blossom taken apart
    return 0;
  pair<ll, int> calc() {
    NX = N; st[0] = 0; FOR(i, 1, 2*N+1) aux[i] = 0;
    FOR(i,1,N+1) match[i] = 0, st[i] = i, flo[i].clear();
    int wMax = 0;
    FOR(u, 1, N+1) FOR(v, 1, N+1)
     floFrom[u][v] = (u == v ? u : 0), ckmax(wMax,g[u][v].w);
    FOR(u, 1, N+1) lab[u] = wMax; // start high and decrease
    int num = 0; 11 wei = 0; while (matching()) ++num;
    FOR(u, 1, N+1) if (match[u] \&\& match[u] < u)
     wei += g[u][match[u]].w; // edges in matching
    return {wei, num};
};
```

```
MaxMatchLexMin.h
```

Description: lexiographically least matching wrt left vertices Usage: solve (L, R, sz (L))

Time:  $\log |L|$  times sum of complexities of gen, maxMatch 44bb49, 26 lines

```
vpi maxMatch(vi L, vi R); // return pairs in max matching
pair<vi, vi> gen(vi L, vi R); // return {Lp, Rp}, vertices on
// left/right that can be reached by alternating path from
// unmatched node on left after finding max matching
vpi res: // stores answer
void solve(vi L, vi R, int x) { // first |L|-x elements of L
 if (x \le 1) { // are in matching, easy if x \le 1
    vpi v = maxMatch(L,R);
    if (sz(v) != sz(L)) L.pop_back(), v = maxMatch(L,R);
    assert(sz(v) == sz(L));
    res.insert(end(res),all(v)); return;
  vi Lp, Rp; tie(Lp, Rp) = gen(L, R); vi Lm = sub(L, Lp), Rm = sub(R, Rp);
  // Lp U Rm is max indep set, Lm U Rp is min vertex cover
  // Lp and Rm independent, edges from Lm to Rp can be ignored
  vpi v = maxMatch(Lm,Rm); assert(sz(v) == sz(Lm));
  res.insert(end(res),all(v));
  vi L2(all(L)-x/2); vi Lp2, Rp2; tie(Lp2, Rp2) = gen(L2, R);
  int cnt = 0; each(t,Lp2) cnt += t >= L[sz(L)-x];
  solve(Lp2, Rp2, cnt); // Rp2 covered by best matching
  vi LL = sub(Lp, Lp2), RR = sub(Rp, Rp2); // those in Lp but not
  // Lp2 that are < L[sz(L)-x/2] must be in answer, not cnt
  cnt = 0; each(t, LL) cnt += t >= L[sz(L)-x/2];
  solve(LL,RR,cnt); // do rest
\} // x reduced by factor of at least two
```

## MaxMatchFast.h

Description: Fast bipartite matching.

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

```
ec6c96, 31 lines
vpi maxMatch(int L, int R, const vpi& edges) {
 V < vi > adj = V < vi > (L);
 vi nxt(L,-1), prv(R,-1), lev, ptr;
 FOR(i,sz(edges)) adj.at(edges[i].f).pb(edges[i].s);
 while (true) {
   lev = ptr = vi(L); int max_lev = 0;
    queue<int> q; F0R(i,L) if (nxt[i]==-1) lev[i]=1, q.push(i);
   while (sz(q)) {
     int x = q.ft; q.pop();
     for (int v: adj[x]) {
       int z = prv[y];
       if (z == -1) \max_{e} [ev = lev[x];
       else if (!lev[z]) lev[z] = lev[x]+1, q.push(z);
     if (max_lev) break;
   if (!max_lev) break;
   FOR(i,L) if (lev[i] > max_lev) lev[i] = 0;
   auto dfs = [&](auto self, int x) -> bool {
     for (;ptr[x] < sz(adj[x]);++ptr[x]) {</pre>
        int y = adj[x][ptr[x]], z = prv[y];
        if (z == -1 \mid | (lev[z] == lev[x]+1 \&\& self(self,z)))
          return nxt[x]=y, prv[y]=x, ptr[x]=sz(adj[x]), 1;
     return 0;
   FOR(i,L) if (nxt[i] == -1) dfs(dfs,i);
 vpi ans; FOR(i,L) if (nxt[i] != -1) ans.pb(\{i,nxt[i]\});
 return ans:
```

## 7.7 Advanced

## ChordalGraphRecognition.h

**Description:** Recognizes graph where every induced cycle has length exactly 3 using maximum adjacency search.

```
int N, M;
set<int> adj[MX];
int cnt[MX];
vi ord, rord;
vi find_path(int x, int y, int z) {
 vi pre(N,-1);
  queue<int> q; q.push(x);
  while (sz(q)) {
    int t = q.ft; q.pop();
    if (adj[t].count(y)) {
      pre[y] = t; vi path = {y};
      while (path.bk != x) path.pb(pre[path.bk]);
      path.pb(z);
      return path;
    each(u,adj[t]) if (u != z \&\& !adj[u].count(z) \&\& pre[u] ==
       →-1) {
      pre[u] = t;
      q.push(u);
 assert (0);
int main() {
  setIO(); re(N,M);
  F0R(i,M) {
    int a,b; re(a,b);
    adj[a].insert(b), adj[b].insert(a);
  rord = vi(N, -1);
  priority_queue<pi> pq;
  FOR(i,N) pg.push(\{0,i\});
  while (sz(pq)) {
    pi p = pq.top(); pq.pop();
    if (rord[p.s] != -1) continue;
    rord[p.s] = sz(ord); ord.pb(p.s);
    each(t,adj[p.s]) pq.push({++cnt[t],t});
  assert(sz(ord) == N);
  each(z,ord) {
    pi biq = \{-1, -1\};
    each(y,adj[z]) if (rord[y] < rord[z])
      ckmax(big,mp(rord[y],y));
    if (big.f == -1) continue;
    int y = big.s;
    each(x,adj[z]) if (rord[x] < rord[y]) if (!adj[y].count(x))
      ps("NO");
      vi v = find_path(x, y, z);
      ps(sz(v));
      each(t,v) pr(t,'');
      exit(0);
  ps("YES");
 reverse(all(ord));
  each (z, ord) pr(z, '');
```

## DominatorTree EdgeColor DirectedMST LCT

### DominatorTree.h

**Description:** Used only a few times. Assuming that all nodes are reachable from root, a dominates b iff every path from root to b passes through a.

from root, a dominates b iff every path from root to b passes through Time:  $\mathcal{O}(M \log N)$ 

ime:  $O(M \log N)$  4b8836, 41 lines

```
template<int SZ> struct Dominator {
  vi adj[SZ], ans[SZ]; // input edges, edges of dominator tree
  vi radj[SZ], child[SZ], sdomChild[SZ];
  int label[SZ], rlabel[SZ], sdom[SZ], dom[SZ], co = 0;
  int par[SZ], bes[SZ];
  void ae(int a, int b) { adj[a].pb(b); }
  int get(int x) { // DSU with path compression
    // get vertex with smallest sdom on path to root
   if (par[x] != x) {
     int t = get(par[x]); par[x] = par[par[x]];
     if (sdom[t] < sdom[bes[x]]) bes[x] = t;</pre>
   return bes[x];
  void dfs(int x) { // create DFS tree
   label[x] = ++co; rlabel[co] = x;
    sdom[co] = par[co] = bes[co] = co;
   each(y,adj[x]) {
     if (!label[y]) {
       dfs(y); child[label[x]].pb(label[y]); }
     radj[label[y]].pb(label[x]);
  void init(int root) {
   dfs(root);
   ROF(i, 1, co+1) {
      each(j,radj[i]) ckmin(sdom[i],sdom[get(j)]);
     if (i > 1) sdomChild[sdom[i]].pb(i);
      each(j,sdomChild[i]) {
       int k = qet(i);
       if (sdom[j] == sdom[k]) dom[j] = sdom[j];
       else dom[j] = k;
     each(j,child[i]) par[j] = i;
    FOR(i, 2, co+1) {
     if (dom[i] != sdom[i]) dom[i] = dom[dom[i]];
     ans[rlabel[dom[i]]].pb(rlabel[i]);
};
```

## EdgeColor.h

**Description:** Used only once. Naive implementation of Misra & Gries edge coloring. By Vizing's Theorem, a simple graph with max degree d can be edge colored with at most d+1 colors

**Time:**  $\mathcal{O}(N^2M)$ , faster in practice

cc2b29, 40 lines

```
template<int SZ> struct EdgeColor {
    int N = 0, maxDeg = 0, adj[SZ][SZ], deg[SZ];
    void init(int _N) { N = _N;
        F0R(i,N) { deg[i] = 0; F0R(j,N) adj[i][j] = 0; } }

void ae(int a, int b, int c) {
    adj[a][b] = adj[b][a] = c; }

int delEdge(int a, int b) {
    int c = adj[a][b]; adj[a][b] = adj[b][a] = 0;
    return c; }

V<bool> genCol(int x) {
    V<bool> col(N+1); F0R(i,N) col[adj[x][i]] = 1;
    return col; }

int freeCol(int u) {
    auto col = genCol(u); int x = 1;
    while (col[x]) ++x; return x; }

void invert(int x, int d, int c) {
```

```
FOR(i,N) if (adj[x][i] == d)
      delEdge(x,i), invert(i,c,d), ae(x,i,c); }
 void ae(int u, int v) {
   // check if you can add edge w/o doing any work
    assert(N); ckmax(maxDeg, max(++deg[u], ++deg[v]));
    auto a = genCol(u), b = genCol(v);
   FOR(i,1,maxDeg+2) if (!a[i] && !b[i])
      return ae(u, v, i);
   V < bool > use(N); vi fan = \{v\}; use[v] = 1;
    while (1) {
      auto col = genCol(fan.bk);
      if (sz(fan) > 1) col[adj[fan.bk][u]] = 0;
      int i=0; while (i \le N \&\& (use[i] \mid | col[adj[u][i]])) i++;
      if (i < N) fan.pb(i), use[i] = 1;</pre>
      else break;
    int c = freeCol(u), d = freeCol(fan.bk); invert(u,d,c);
    int i = 0; while (i < sz(fan) && genCol(fan[i])[d]</pre>
      && adj[u][fan[i]] != d) i ++;
    assert (i != sz(fan));
   FOR(j,i) ae(u,fan[j],delEdge(u,fan[j+1]));
    ae(u,fan[i],d);
};
```

## DirectedMST.h

**Description:** Chu-Liu-Edmonds algorithm. Computes minimum weight directed spanning tree rooted at r, edge from  $par[i] \to i$  for all  $i \neq r$ . Use DSU with rollback if need to return edges.

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

```
"DSUrb.h"
                                                     5d5c10, 61 lines
struct Edge { int a, b; ll w; };
struct Node { // lazy skew heap node
  Edge key; Node *1, *r; 11 delta;
  void prop() {
    kev.w += delta;
    if (1) 1->delta += delta;
    if (r) r->delta += delta;
    delta = 0;
 Edge top() { prop(); return key; }
Node *merge(Node *a, Node *b) {
 if (!a || !b) return a ?: b;
 a->prop(), b->prop();
  if (a->key.w > b->key.w) swap(a, b);
  swap(a->1, a->r = merge(b, a->r));
  return a;
void pop(Node*\& a) { a->prop(); a = merge(a->1, a->r); }
pair<ll, vi> dmst(int n, int r, const vector<Edge>& g) {
  DSUrb dsu; dsu.init(n);
  vector<Node*> heap(n); // store edges entering each vertex
  // in increasing order of weight
  each(e,q) heap[e.b] = merge(heap[e.b], new Node{e});
  ll res = 0; vi seen(n,-1); seen[r] = r;
  vpi in(n, {-1,-1}); // edge entering each vertex in MST
  vector<pair<int, vector<Edge>>> cycs;
  FOR(s,n) {
    int u = s, w;
    vector<pair<int, Edge>> path;
    while (seen[u] < 0) {</pre>
      if (!heap[u]) return {-1,{}};
      seen[u] = s;
      Edge e = heap[u] \rightarrow top(); path.pb({u,e});
      heap[u]->delta -= e.w, pop(heap[u]);
      res += e.w, u = dsu.get(e.a);
```

```
if (seen[u] == s) { // found cycle, contract
     Node * cvc = 0; cvcs.eb();
      do {
        cyc = merge(cyc, heap[w = path.bk.f]);
        cycs.bk.s.pb(path.bk.s);
        path.pop_back();
      } while (dsu.unite(u,w));
      u = dsu.qet(u); heap[u] = cyc, seen[u] = -1;
      cycs.bk.f = u;
  each(t,path) in[dsu.get(t.s.b)] = \{t.s.a,t.s.b\};
} // found path from root to s, done
while (sz(cycs)) { // expand cycs to restore sol
  auto c = cycs.bk; cycs.pop back();
  pi inEdge = in[c.f];
  each(t,c.s) dsu.rollback();
  each(t,c.s) in[dsu.get(t.b)] = \{t.a,t.b\};
  in[dsu.get(inEdge.s)] = inEdge;
vi par(n); FOR(i,n) par[i] = in[i].f;
// i == r ? in[i].s == -1 : in[i].s == i
return {res,par};
```

#### LCT.h

**Description:** Link-Cut Tree. Given a function  $f(1\dots N)\to 1\dots N$ , evaluates  $f^b(a)$  for any a,b. sz is for path queries; sub, vsub are for subtree queries. x->access() brings x to the top and propagates it; its left subtree will be the path from x to the root and its right subtree will be empty. Then sub will be the number of nodes in the connected component of x and vsub will be the number of nodes under x. Use makeRoot for arbitrary path queries.

Usage: FOR(i,1,N+1)LCT[i]=new snode(i); link(LCT[1],LCT[2],1); Time:  $\mathcal{O}(\log N)$  e24bf7.115 lines

```
typedef struct snode* sn;
struct snode { ////// VARIABLES
 sn p, c[2]; // parent, children
 sn extra; // extra cycle node for "The Applicant"
 bool flip = 0; // subtree flipped or not
 int val, sz; // value in node, # nodes in current splay tree
 int sub, vsub = 0; // vsub stores sum of virtual children
 snode(int _val) : val(_val) {
   p = c[0] = c[1] = extra = NULL; calc(); }
 friend int getSz(sn x) { return x?x->sz:0; }
 friend int getSub(sn x) { return x?x->sub:0; }
 void prop() { // lazy prop
   if (!flip) return;
   swap(c[0],c[1]); flip = 0;
   FOR(i,2) if (c[i]) c[i]->flip ^= 1;
 void calc() { // recalc vals
   F0R(i,2) if (c[i]) c[i]->prop();
   sz = 1+getSz(c[0])+getSz(c[1]);
   sub = 1+getSub(c[0])+getSub(c[1])+vsub;
 ////// SPLAY TREE OPERATIONS
 int dir() {
   if (!p) return -2;
   FOR(i,2) if (p->c[i] == this) return i;
   return -1; // p is path-parent pointer
 } // -> not in current splay tree
 // test if root of current splay tree
 bool isRoot() { return dir() < 0; }</pre>
 friend void setLink(sn x, sn y, int d) {
   if (y) y->p = x;
   if (d >= 0) x -> c[d] = y; }
 void rot() { // assume p and p->p propagated
```

```
assert(!isRoot()); int x = dir(); sn pa = p;
    setLink(pa->p, this, pa->dir());
    setLink(pa, c[x^1], x); setLink(this, pa, x^1);
   pa->calc();
  void splay() {
    while (!isRoot() && !p->isRoot()) {
     p->p->prop(), p->prop(), prop();
     dir() == p->dir() ? p->rot() : rot();
    if (!isRoot()) p->prop(), prop(), rot();
   prop(); calc();
  sn fbo(int b) { // find by order
   prop(); int z = getSz(c[0]); // of splay tree
    if (b == z) { splay(); return this; }
    return b < z ? c[0] -> fbo(b) : c[1] -> fbo(b-z-1);
  ////// BASE OPERATIONS
  void access() { // bring this to top of tree, propagate
    for (sn v = this, pre = NULL; v; v = v->p) {
     v->splay(); // now switch virtual children
     if (pre) v->vsub -= pre->sub;
     if (v->c[1]) v->vsub += v->c[1]->sub;
     v \rightarrow c[1] = pre; v \rightarrow calc(); pre = v;
   splay(); assert(!c[1]); // right subtree is empty
  void makeRoot() {
    access(); flip ^= 1; access(); assert(!c[0] && !c[1]); }
  ////// OUERIES
  friend sn lca(sn x, sn y) {
   if (x == y) return x;
   x->access(), y->access(); if (!x->p) return NULL;
   x->splay(); return x->p?:x; // y was below x in latter case
  } // access at y did not affect x -> not connected
  friend bool connected(sn x, sn y) { return lca(x,y); }
  // # nodes above
  int distRoot() { access(); return getSz(c[0]); }
  sn getRoot() { // get root of LCT component
   access(); sn a = this;
    while (a->c[0]) a = a->c[0], a->prop();
   a->access(); return a;
  sn getPar(int b) { // get b-th parent on path to root
   access(); b = getSz(c[0])-b; assert(b \ge 0);
   return fbo(b);
  } // can also get min, max on path to root, etc
  ////// MODIFICATIONS
  void set(int v) { access(); val = v; calc(); }
  friend void link(sn x, sn y, bool force = 0) {
    assert(!connected(x,y));
    if (force) y->makeRoot(); // make x par of y
   else { y->access(); assert(!y->c[0]); }
   x->access(); setLink(y,x,0); y->calc();
  friend void cut(sn y) { // cut y from its parent
   y->access(); assert(y->c[0]);
    y \rightarrow c[0] \rightarrow p = NULL; y \rightarrow c[0] = NULL; y \rightarrow calc(); }
  friend void cut(sn x, sn y) { // if x, y adj in tree
    x->makeRoot(); y->access();
    assert (y-c[0] == x && !x-c[0] && !x-c[1]); cut(y); }
sn LCT[MX];
////// THE APPLICANT SOLUTION
void setNex(sn a, sn b) { // set f[a] = b
 if (connected(a,b)) a->extra = b;
```

```
else link(b,a); }
void delNex(sn a) { // set f[a] = NULL
   auto t = a->getRoot();
   if (t == a) { t->extra = NULL; return; }
   cut(a); assert(t->extra);
   if (!connected(t,t->extra))
        link(t->extra,t), t->extra = NULL;
}
sn getPar(sn a, int b) { // get f^b[a]
   int d = a->distRoot(); if (b <= d) return a->getPar(b);
   b -= d+1; auto r = a->getRoot()->extra; assert(r);
   d = r->distRoot()+1; return r->getPar(b%d);
}
```

# TopTree.h Description: Top tree (generalization of LCT to support path & subtree

```
updates & queries)
* Usage:
   Implement
      void update()
       void downdate()
       void do_flip_path()
       void do_other_operation() ...
    When update() is called, you can assume downdate() has
   \hookrightarrowalready been called.
   In general, do_op() should eagerly apply the operation but
   \hookrightarrow not touch the
   children. In downdate(), you can push down to the children
   \hookrightarrow with ch->do_op().
   WARNING : if different operations do not trivially commute
   \hookrightarrow, you *must*
   implement a way to swap/alter them to compose in a
   \hookrightarrow consistent order, and you
   must use that order when implementing downdate(). This can

→ be nontrivial!

    Creating vertices:
       n->is_path = n->is_vert = true;
       n->update();
    Creating edges: no setup/update() needed, just call
       link(e, va, vb);
    Updates:
      auto cur = get_path(va, vb); // or get_subtree(va, vb)
      cur->do_stuff();
       cur->downdate();
       cur->update_all();
* Node types:
* path edges: compress(c[0], self, c[1])
      assert(is_path && !is_vert);
       assert(c[0] && c[1]);
      assert (c[0] \rightarrow is_path \&\& c[1] \rightarrow is_path);
       assert (!c[2]);
     (path) vertices: self + rake(c[0], c[1])
       assert (is_path && is_vert);
      assert(!c[2]);
       if (c[0]) assert(!c[0]->is_path);
       if (c[1]) assert(!c[1]->is_path);
     non-path edges: rake(c[0], self + c[2], c[1])
       assert (!is_path && !is_vert);
       assert(c[2]);
       assert(c[2]->is_path);
       if (c[0]) assert(!c[0]->is_path);
       if (c[1]) assert(!c[1]->is_path);
```

```
using pt = struct top_tree_node*;
struct top_tree_node {
private:
  mutable pt p = nullptr;
  AR<pt, 3> c{nullptr, nullptr, nullptr};
  // \return direction in which parent points to you
  int d() const {
    if (!p) return -1;
    FOR(i,3) if (this == p->c[i]) return i;
    assert (false);
  // \return true if this is root of rake or compress tree
  bool r() const { return !p || p->is path != is path; }
  // 3 types of verts: path edges, path verts, non-path edges
  bool is_path, is_vert;
  bool flip_path = false;
  // MODIFY STUFF BELOW
  int path len, best path;
  bool own parity, path parity;
  AR<int,2> best_down, best_up;
  void do flip path() {
    assert(is_path); flip_path ^= 1;
    swap(best_down, best_up);
  void downdate() {
    if (flip path) {
      assert(is path);
      if (!is_vert) F0R(i,2) if (c[i]) c[i]->do_flip_path(); //
         \hookrightarrow if vert, then you're at a leaf so don't propagate
      swap(c[0], c[1]);
      flip_path = false;
  void update() { // TODO: find longest path of each parity
    assert (!flip path);
    if (is_path && !is_vert) {
      assert(c[0] && c[1] && !c[2]);
      path_len = 1+c[0]->path_len+c[1]->path_len;
      path_parity = own_parity^c[0]->path_parity^c[1]->
         \hookrightarrowpath_parity;
      best_up = c[0] -> best_up;
      FOR(2.2)
        ckmax(best_up[c[0]->path_parity^own_parity^z],c[0]->
           \hookrightarrowpath_len+1+c[1]->best_up[z]);
      best_down = c[1]->best_down;
      F0R(z,2)
        ckmax(best_down[c[1]->path_parity^own_parity^z],c[1]->
           \hookrightarrowpath_len+1+c[0]->best_down[z]);
      best_path = max(c[0]->best_path,c[1]->best_path);
      FOR(z, 2) ckmax(best_path,c[0]->best_down[z]+1+c[1]->
         ⇔best_up[z^own_parity]);
    } else {
      path len = 0;
      path_parity = 0;
      best_up = \{0, -MOD\};
      best_path = 0;
      if (!is_vert) {
        best_path = c[2]->best_path;
        AR < int, 2 > tmp_up = c[2] - > best_up;
        F0R(i,2) ++tmp_up[i];
        if (own_parity) swap(tmp_up[0],tmp_up[1]);
```

16

```
F0R(i,2) {
          ckmax(best_up[i],tmp_up[i]);
          if (i == 0) ckmax(best_path,best_up[i]);
      F0R(d,2) {
        if (!c[d]) continue;
        ckmax(best_path,c[d]->best_path);
          ckmax(best path,c[d]->best up[z]+best up[z]);
          ckmax(best_up[z],c[d]->best_up[z]);
      // rake(c[0], c[1], self+c[2])
      best down = best up;
  void downdate all() {
    if (p) p->downdate_all();
    downdate();
  // Returns the root
  pt update all() {
   pt cur = this; cur->update();
    while (cur->p) { cur = cur->p; cur->update(); }
private:
  // sets v to be the d'th child of x
  friend void setLink(pt x, pt y, int d) {
   if (y) y \rightarrow p = x;
   if (d != -1) x -> c[d] = y;
  void rot() { // rotate this up, parent down
   assert(!is_vert && !r());
   pt pa = p;
    int x = d(); assert (x == 0 | | x == 1);
   pt ch = c[!x];
    setLink(pa->p,this,pa->d());
    setLink(pa,ch,x);
    setLink(this,pa,!x);
   pa->update();
  void rot_2(int c_d) { // rotate this up (along with c_d'th
    ⇔child), parent down
    assert(!is_vert && !r());
    assert(c[c_d] && !c[c_d]->is_vert);
    if (d() == c_d) { rot(); return; }
    pt pa = p;
    int x = d(); assert (x == 0 | | x == 1);
    assert(c d == !x);
    pt ch = c[c_d] - c[!x];
    setLink(pa->p,this,pa->d());
    setLink(pa,ch,x);
    setLink(this->c[c_d],pa,!x);
   pa->update();
  void splay_dir(int x) { // splay while direction is x
    while (!r() \&\& d() == x) {
     if (!p->r() && p->d() == x) p->rot();
      rot();
  // splay path edge along with child path edge
  void splay_2(int c_d) {
    assert (!is_vert && is_path);
    assert(c[c_d] && !c[c_d]->is_vert);
    while (!r()) {
     if (!p->r()) {
```

```
if (p->d() == d()) p->rot();
      else rot_2(c_d);
    rot_2(c_d);
// splay parent edge to top of tree, bringing this along for
void splay_2() {
  assert(!is vert && is path && !r());
  p->splay_2(d());
// splay vertex as close to top of tree as possible
void splay_vert() {
  assert (is vert && is path); if (r()) return; // if path is
     ⇒single vertex, done
  p->splay_dir(d()); if (p->r()) return; // if parent is
     ⇒already at top, done
  // otherwise vertex is in between two edges
  assert(p->d() != d());
  if (d() == 1) p->rot();
  assert(d() == 0);
  p->splay_2(); // splay parent of parent to the top
  assert(d() == 0 \&\& p > d() == 1 \&\& p > p > r());
void splay() { // normal splay of an edge
  assert(!is vert);
  while (!r()) {
    if (!p->r()) {
      if (p->d() == d()) p->rot();
      else rot();
    rot();
// either brings root to top of splay tree, or to right child
   \hookrightarrow of top of splay tree
// cuts part of compress tree to the right of this
pt cut_right() {
  assert(is_vert && is_path);
  splay_vert();
  if (r() \mid \mid d() == 1) { // if last vertex on path, do
     \hookrightarrownothing
    assert(r() || (d() == 1 \&\& p->r()));
    assert(c[0] == nullptr); // why?
    return nullptr; // don't need to cut anything
  // goal: cut pa and everything to the right of it
  // make it a rake
  // before:
  // pa->p
  //
        pa
  // this (vertex, children are rake trees)
  // after:
  // pa->p
        this (pa as foster child)
  // or -> single vertex
  pt pa = p;
  assert (pa->r() | (pa->d() == 1 && pa->p->r()));
  assert(!pa->is_vert && pa->is_path);
```

```
assert (pa->c[0] == this && pa->c[2] == nullptr); // pa is
       ⇒path edge
    setLink(pa->p,this,pa->d()); // now this or this->p is root
    pa->is_path = false; pa->c[2] = pa->c[1]; // pa is now a
    FOR(i,2) setLink(pa,c[i],i); // pa inherits rake children
       \hookrightarrowof this
    c[0] = nullptr; setLink(this,pa,1); // set pa to be a rake
    assert(c[2] == nullptr);
    pa->update(); return pa;
 pt splice_non_path() { // cut some previous path, replace
    ⇒with new path
    assert(!is_path && !is_vert); splay(); // bring to top of
       ⇒rake tree
    assert (p && p->is_vert && p->is_path); // parent is a
    p->cut right(); // cut part to right of vertex
    if (!p->is_path) rot(); // rotate this to top of rake tree
    assert (p && p->is_vert && p->is_path); // same parent
    assert (p->r() | | (p->d() == 1 && p->p->r()));
    assert(p \rightarrow c[d()] == this && p \rightarrow c[!d()] == nullptr); //
       ⇒parent vertex only has one rake child? why?
    pt pa = p;
    setLink(pa->p,this,pa->d());
    F0R(i,2) setLink(pa,c[i],i);
    assert(c[2] && c[2]->is_path);
    c[1] = c[2]; // don't need to change parent
    setLink(this,pa,0);
    c[2] = nullptr;
    is_path = true;
    assert (d() != 0);
    pa->update(); return pa;
  // Return the topmost vertex which was spliced into
 pt splice_all() { // make this part of topmost path
    pt res = nullptr;
    for (pt cur = this; cur; cur = cur->p) {
      if (!cur->is_path) res = cur->splice_non_path();
      assert(cur->is_path);
    return res;
public:
 pt getRoot() {
    expose();
    pt v = this; while (v->p) v = v->p;
    assert (v);
    v->downdate();
    while (!v->is_vert) {
      assert (v->c[0]);
      v = v - c[0]; v - downdate();
    v->expose();
    return v:
  friend bool connected(pt a, pt b) {
    return a->getRoot() == b->getRoot();
  // Return the topmost vertex which was spliced into
  pt expose() { // yay makes sense!
    assert(is_vert);
    downdate_all(); // make sure to propagate everything above
    pt res = splice_all(); cut_right(); update_all();
```

```
// Brings edge to the top
// Return the topmost vertex which was spliced into
pt expose_edge() {
 assert(!is_vert); downdate_all();
 pt v = is_path ? c[1] : c[2]; v->downdate();
  // if is_path: path to right of edge
  // otherwise: compress tree under edge
  while (!v->is\_vert) { v = v->c[0]; v->downdate(); }
 pt res = v->splice_all(); v->cut_right(); v->update_all();
    \hookrightarrow // same as expose
  assert(!p && v == c[1]);
 return res;
// make sure path end only has one child
// Return the new root
pt meld path end() {
 assert(!p);
 pt rt = this;
  while (true) {
   rt->downdate();
   if (rt->is vert) break;
   rt = rt - c[1];
  assert(rt->is vert);
  rt->splay_vert();
  if (rt->c[0] && rt->c[1]) { // make sure path end has only
     ⇒one rake child ...
   pt ch = rt -> c[1];
   while (true) {
     ch->downdate();
     if (!ch->c[0]) break;
      ch = ch -> c[0];
   ch->splay();
   assert(ch->c[0] == nullptr);
   setLink(ch,rt->c[0],0); rt->c[0] = nullptr;
   ch->update();
  } else if (rt->c[0]) {
   rt->c[1] = rt->c[0];
   rt->c[0] = nullptr;
  assert(rt->c[0] == nullptr);
  return rt->update_all();
void make_root() {
  expose();
  pt rt = this;
  while (rt->p) {
   assert (rt->d() == 1);
   rt = rt - p;
  rt->do_flip_path(); rt->meld_path_end();
  expose(); assert(!p); // root path is now single node
// link v2 as a child of v1 with edge e
friend void link(pt e, pt v1, pt v2) {
 assert (e && v1 && v2); F0R(i,3) assert(!e->c[i]);
 e->is_path = true, e->is_vert = false;
 v1\rightarrow expose(); while (v1\rightarrow p) v1 = v1\rightarrow p;
 v2->make_root();
 assert(!v1-p \&\& !v2-p); // should both be at top
  setLink(e, v1, 0); setLink(e, v2, 1);
 e->update();
// Cuts the edge e
// Returns the top-tree-root of the two halves; they are not
  \hookrightarrownecessarily the split vertices.
friend pair<pt, pt> cut(pt e) {
 assert(!e->is_vert); e->expose_edge();
```

```
assert(!e->p \&\& e->is_path \&\& e->c[2] == nullptr);
   pt 1 = e - c[0], r = e - c[1]; assert(1 && r);
    e\rightarrow c[0] = e\rightarrow c[1] = nullptr; 1\rightarrow p = r\rightarrow p = nullptr; //
      \hookrightarrowdisconnect
   1 = 1->meld_path_end();
    return {1, r}:
 // bring path to the top
 friend pt get_path(pt a, pt b) {
    assert(a->is vert && b->is vert);
    a->make_root(); b->expose();
   if (a == b) { assert(!b->p); return b; } // top path is
       \hookrightarrowsingle node
    assert(!b->p->p); return b->p; // b is at end of path
 // root at rt, get subtree at n
 friend pt get_subtree(pt rt, pt n) {
   assert(rt->is vert && n->is vert);
    rt->make_root(); n->expose(); return n;
 // easy: just expose one and then the other
 friend pt lca(pt a, pt b) {
   assert(a->is vert && b->is vert);
   assert (connected(a,b));
    a->expose(); return b->expose() ?: b;
};
int main() {
 cin.tie(0)->sync_with_stdio(0);
 ints(N):
 V<top_tree_node> nodes(N);
 F0R(i.N) {
   pt n = &nodes[i];
   n->is_path = n->is_vert = true;
   n->update();
 V<top_tree_node> edges(N-1);
 F0R(i,N-1) {
   ints(u,v,t); --u,--v;
   edges[i].own_parity = t;
   link(&edges[i],&nodes[u],&nodes[v]);
 ints(M);
 rep(M) {
   ints(id); --id;
   edges[id].expose_edge();
   edges[id].own_parity ^= 1; edges[id].update();
   ps(edges[id].best_path);
```

## Geometry (8)

## 8.1 Primitives

## ComplexComp.h

Description: Allows you to sort complex numbers.

6f828b, 5 lines

```
#define x real()
#define y imag()
using P = complex<db>;
namespace std {
  bool operator<(P 1,P r) { return mp(1.x,1.y)<mp(r.x,r.y); } }</pre>
```

#### | PointShort.h

**Description:** Use in place of complex<T>.

cefef8, 36 lines

18

```
using T = db; // or 11
const T EPS = 1e-9; // adjust as needed
using P = pair<T,T>; using vP = V<P>; using Line = pair<P,P>;
int sgn(T a) { return (a>EPS)-(a<-EPS); }</pre>
T sq(T a) { return a*a; }
T norm(P p) { return sq(p.f)+sq(p.s); }
T abs(P p) { return sqrt(norm(p)); }
T arg(P p) { return atan2(p.s,p.f); }
P conj(P p) { return P(p.f,-p.s); }
P perp(P p) { return P(-p.s,p.f); }
P dir(T ang) { return P(cos(ang), sin(ang)); }
P operator+(P 1, P r) { return P(1.f+r.f,1.s+r.s); }
P operator-(P 1, P r) { return P(1.f-r.f,1.s-r.s); }
P operator*(P 1, T r) { return P(1.f*r,1.s*r); }
P operator/(P 1, T r) { return P(1.f/r,1.s/r); }
P operator*(P 1, P r) { // complex # multiplication
 return P(1.f*r.f-1.s*r.s,1.s*r.f+1.f*r.s); }
P operator/(P 1, P r) { return 1*conj(r)/norm(r); }
P unit(const P& p) { return p/abs(p); }
T dot(const P& a, const P& b) { return a.f*b.f+a.s*b.s; }
T dot(const P& p, const P& a, const P& b) { return dot(a-p,b-p)
T cross(const P& a, const P& b) { return a.f*b.s-a.s*b.f; }
T cross(const P& p, const P& a, const P& b) {
 return cross(a-p,b-p); }
P reflect (const P& p, const Line& 1) {
 P a = 1.f, d = 1.s-1.f;
 return a+conj((p-a)/d)*d; }
P foot (const P& p, const Line& 1) {
 return (p+reflect(p,1))/(T)2; }
bool onSeg(const P& p, const Line& 1) {
 return sgn(cross(1.f,1.s,p)) == 0 && sgn(dot(p,1.f,1.s)) <= 0
ostream& operator << (ostream& os, P p) {
 return os << "(" << p.f << "," << p.s << ")"; }
```

#### AngleCmp.h

**Description:** Sorts points in ccw order about origin in the same way as atan2, which returns real in  $(-\pi, \pi]$  so points on negative x-axis come last.

"Point.N" 2df5fc, 12 lines

// WARNING: you will get unexpected results if you mistype this

→ as bool instead of int

// -1 if lower half, 0 if origin, 1 if upper half
int half(P x) { return x.s != 0 ? sgn(x.s) : -sgn(x.f); }
bool angleCmp(P a, P b) { int A = half(a), B = half(b);
 return A == B ? cross(a,b) > 0 : A < B; }

/\* Usage:

\* vP v;

\* sort(all(v),[](P a, P b) { return
 atan2(a.s,a.f) < atan2(b.s,b.f); });

\* sort(all(v),angleCmp); // should give same result

\*/

## SegDist.h

**Description:** computes distance between P and line (segment) AB

```
return lineDist(p,1); }
```

## SegIsect.h

**Description:** computes the intersection point(s) of line (segments) a and b7f6ba0, 26 lines

```
// {unique intersection point} if it exists
// {b.f,b.s} if input lines are the same
// empty if lines do not intersect
vP lineIsect(const Line& a, const Line& b) {
 T = a\theta = cross(a.f,a.s,b.f), al = cross(a.f,a.s,b.s);
 if (a0 == a1) return a0 == 0 ? vP{b.f,b.s} : vP{};
  return { (b.s*a0-b.f*a1) / (a0-a1) };
// point in interior of both segments a and b, if it exists
vP strictIsect(const Line& a, const Line& b) {
 T = a0 = cross(a.f,a.s,b.f), a1 = cross(a.f,a.s,b.s);
 T b\theta = cross(b.f,b.s,a.f), b1 = cross(b.f,b.s,a.s);
  if (sgn(a0) * sgn(a1) < 0 && sgn(b0) * sgn(b1) < 0)
   return { (b.s*a0-b.f*a1) / (a0-a1) };
  return {};
// intersection of segments, a and b may be degenerate
vP segIsect(const Line& a, const Line& b) {
 vP v = strictIsect(a,b); if (sz(v)) return v;
  set<P> s;
  \#define i(x,y) if (onSeg(x,y)) s.ins(x)
 i(a.f,b); i(a.s,b); i(b.f,a); i(b.s,a);
  return {all(s)};
```

## Polygons

## PolygonCenArea.h

Description: centroid (center of mass) of a polygon with constant mass per unit area and SIGNED area

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

```
4ca221, 7 lines
"Point.h"
pair<P,T> cenArea(const vP& v) { assert(sz(v) >= 3);
 P cen{}; T area{};
  F0R(i,sz(v)) {
   int j = (i+1) %sz(v); T a = cross(v[i],v[j]);
    cen += a*(v[i]+v[j]); area += a; }
  return {cen/area/(T)3,area/2}; // area is SIGNED
```

### InPolygon.h

Description: Tests whether point is inside, on, or outside of a polygon (returns -1, 0, or 1). Both CW and CCW polygons are ok.

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

```
"Point.h"
                                                      2ed683, 9 lines
int inPolygon(const P& p, const vP& poly) {
 int n = sz(poly), ans = 0;
  F0R(i,n) {
   P x = poly[i], y = poly[(i+1)%n]; if (x.s > y.s) swap(x,y);
   if (onSeg(p, {x,y})) return 0;
    ans ^= (x.s <= p.s && p.s < y.s && cross(p,x,y) > 0);
  return ans ? -1 : 1;
```

#### ConvexHull.h

Description: top-bottom convex hull

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

```
868655, 18 lines
"Point.h"
pair<vi, vi> ulHull(const vP& v) {
 vi p(sz(v)), u, l; iota(all(p), 0);
  sort(all(p), [&v](int a, int b) { return v[a] < v[b]; });</pre>
 each(i,p) {
    \#define ADDP(C, cmp) while (sz(C) > 1 && cross(\
      v[C[sz(C)-2]], v[C.bk], v[i]) cmp 0) C.pop_back(); C.pb(i);
    ADDP(u, >=); ADDP(1, <=);
 return {u,1};
vi hullInd(const vP& v) {
 vi u,1; tie(u,1) = ulHull(v); if (sz(1) \le 1) return 1;
 if (v[1[0]] == v[1[1]]) return {0};
 1.insert (end(1),1+rall(u)-1); return 1;
vP hull(const vP& v) {
 vi w = hullInd(v); vP res; each(t,w) res.pb(v[t]);
 return res; }
```

#### ConvexHull2.h

Description: Graham Scan

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

```
"Point.h"
                                                     5758fd, 13 lines
vi hullInd(const vP& v) {
 int ind = int(min_element(all(v))-begin(v));
 vi cand, C\{ind\}; FOR(i,sz(v)) if (v[i] != v[ind]) cand.pb(i);
 sort(all(cand),[&](int a, int b) {
    // sort by angle, tiebreak by distance
   P x = v[a]-v[ind], y = v[b]-v[ind]; T t = cross(x,y);
   return t != 0 ? t > 0 : norm(x) < norm(y); });
 each(c,cand) {
   while (sz(C) > 1 \&\& cross(v[end(C)[-2]],v[C.bk],v[c]) \le 0)
     C.pop_back();
   C.pb(c); }
 return C;
```

#### Diameter.h

**Description:** rotating caliphers, gives greatest distance between two points

### **Time:** $\mathcal{O}(N)$ given convex hull

```
"ConvexHull.h"
                                                       ee92de, 9 lines
db diameter(vP P) {
 P = hull(P);
 int n = sz(P), ind = 1; db ans = 0;
 if (n > 1) F0R(i,n) for (int j = (i+1)%n;;ind = (ind+1)%n) {
   ckmax(ans,abs(P[i]-P[ind]));
    if (cross(P[j]-P[i],P[(ind+1)%n]-P[ind]) <= 0) break;
 return ans;
```

#### LineHull.h

Description: lineHull accepts line and ccw convex polygon. If all vertices in poly lie to one side of the line, returns a vector of closest vertices to line as well as orientation of poly with respect to line (±1 for above/below). Otherwise, returns the range of vertices that lie on or below the line. extrVertex returns the point of a hull with the max projection onto a line.

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

```
"Point.h"
                                                      40e5a6, 41 lines
using Line = AR<P,2>;
#define cmp(i,j) sgn(-dot(dir,poly[(i)%n]-poly[(j)%n]))
#define extr(i) cmp(i+1,i) >= 0 && cmp(i,i-1+n) < 0
int extrVertex(const vP& poly, P dir) {
```

```
int n = sz(poly), lo = 0, hi = n;
 if (extr(0)) return 0;
  while (lo+1 < hi) {
   int m = (lo+hi)/2;
    if (extr(m)) return m;
    int 1s = cmp(1o+1, 1o), ms = cmp(m+1, m);
    (ls < ms \mid | (ls == ms \&\& ls == cmp(lo, m)) ? hi : lo) = m;
 return lo;
vi same (Line line, const vP& poly, int a) {
 // points on same parallel as a
 int n = sz(poly); P dir = perp(line[0]-line[1]);
 if (cmp(a+n-1,a) == 0) return \{(a+n-1) n,a\};
 if (cmp(a,a+1) == 0) return \{a,(a+1)\%n\};
 return {a};
#define cmpL(i) sqn(cross(line[0],line[1],poly[i]))
pair<int, vi> lineHull(Line line, const vP& poly) {
 int n = sz(poly); assert(n>1);
 int endA = extrVertex(poly,perp(line[0]-line[1])); // lowest
 if (cmpL(endA) >= 0) return {1, same(line, poly, endA)};
 int endB = extrVertex(poly,perp(line[1]-line[0])); // highest
 if (cmpL(endB) <= 0) return {-1, same(line, poly, endB)};</pre>
 AR<int,2> res:
 F0R(i,2) {
    int lo = endA, hi = endB; if (hi < lo) hi += n;
    while (lo < hi) {
     int m = (lo+hi+1)/2;
      if (cmpL(m%n) == cmpL(endA)) lo = m;
      else hi = m-1;
    res[i] = lo%n; swap(endA,endB);
 if (cmpL((res[0]+1)%n) == 0) res[0] = (res[0]+1)%n;
 return {0, {(res[1]+1)%n, res[0]}};
```

## HalfPlaneIsect2.h

**Description:** Returns vertices of half-plane intersection. A half-plane is the area to the left of a ray, which is defined by a point p and a direction dp. Area of intersection should be sufficiently precise when all inputs are integers with magnitude  $< 10^5$ . Assumes intersection is bounded (easiest way to ensure this is to uncomment the code below).

```
Time: \mathcal{O}(N \log N)
"AngleCmp.h"
                                                       d9e261, 46 lines
struct Ray {
  P p, dp; // origin, direction
  P isect(const Ray& L) const {
    return p+dp*(cross(L.dp,L.p-p)/cross(L.dp,dp)); }
  bool operator<(const Ray& L) const {
    return angleCmp(dp,L.dp); }
};
vP halfPlaneIsect(V<Ray> _segs) {
  // int DX = 1e9, DY = 1e9; // bound input by rectangle [0,DX]
     \hookrightarrow x [0,DY]
  // _segs.pb({P{0,0},P{1,0}});
  // _segs.pb({P{DX,0},P{0,1}});
  // _segs.pb({P{DX,DY},P{-1,0}});
  // _segs.pb({P{0,DY},P{0,-1}});
  sor(_segs); // sort planes by angle
  V<Ray> segs; // remove parallel planes
  each(t,_segs) {
    if (!sz(segs) || segs.bk < t) { segs.pb(t); continue; }</pre>
    if (cross(t.dp,t.p-segs.bk.p) > 0) segs.bk = t;
  auto bad = [&] (const Ray& a, const Ray& b, const Ray& c) {
```

```
P p1 = a.isect(b), p2 = b.isect(c);
 if (dot(p2-p1,b.dp) <= EPS) { // this EPS is required ...
   if (cross(a.dp,c.dp) \le 0) return 2; // isect(a,b,c) =
   return 1;
 return 0; // isect(a,c) == isect(a,b,c)
};
#define reduce(t) \
 while (sz(polv) > 1) \{ \
   int b = bad(poly.at(sz(poly)-2),poly.bk,t); \
   if (b == 2) return {}; \
   if (b == 1) poly.pop_back(); \
   else break; \
deque<Ray> poly;
each(t, segs) { reduce(t); poly.pb(t); }
for(;;poly.pop_front()) {
 reduce(poly[0]);
 if (!bad(poly.bk,poly[0],poly[1])) break;
assert(sz(poly) >= 3); // if you reach this point, area
  ⇒should be nonzero
vP poly points; F0R(i,sz(poly))
 poly_points.pb(poly[i].isect(poly[(i+1)%sz(poly)]));
return poly_points;
```

## 8.3 Circles

## Circle.h

**Description:** represent circle as {center,radius}

```
"Point.h" 91f3fc, 6 lines
using Circ = pair<P,T>;
int in(const Circ& x, const P& y) { // -1 if inside, 0, 1
   return sgn(abs(y-x.f)-x.s); }
T arcLength(const Circ& x, P a, P b) {
   // precondition: a and b on x
   P d = (a-x.f)/(b-x.f); return x.s*acos(d.f); }
```

#### CircleIsect.h

 $\bf Description:$  Circle intersection points and intersection area. Tangents will be returned twice.

```
a0b0f8, 22 lines
"Circle.h"
vP isect(const Circ& x, const Circ& y) { // precondition: x!=y
 T d = abs(x.f-y.f), a = x.s, b = y.s;
  if (sgn(d) == 0) { assert(a != b); return {}; }
  T C = (a*a+d*d-b*b)/(2*a*d);
  if (abs(C) > 1+EPS) return {};
  T S = sqrt(max(1-C*C, (T)0)); P tmp = (y.f-x.f)/d*x.s;
  return \{x.f+tmp*P(C,S),x.f+tmp*P(C,-S)\};
vP isect(const Circ& x, const Line& y) {
  P c = foot(x.f,y); T sq_dist = sq(x.s) - norm(x.f-c);
  if (sgn(sq_dist) < 0) return {};</pre>
  P offset = unit(y.s-y.f) *sqrt(max(sq_dist,T(0)));
  return {c+offset,c-offset};
T isect_area(Circ x, Circ y) { // not thoroughly tested
  T d = abs(x.f-y.f), a = x.s, b = y.s; if (a < b) swap(a,b);
  if (d >= a+b) return 0;
  if (d <= a-b) return PI*b*b;
  T ca = (a*a+d*d-b*b)/(2*a*d), cb = (b*b+d*d-a*a)/(2*b*d);
  T s = (a+b+d)/2, h = 2*sqrt(s*(s-a)*(s-b)*(s-d))/d;
  return a*a*acos(ca)+b*b*acos(cb)-d*h;
```

## CircleTangents.h

```
Description: internal and external tangents between two circles
```

```
d9a76f, 22 lines
P tangent (P x, Circ y, int t = 0) {
 y.s = abs(y.s); // abs needed because internal calls y.s < 0
 if (y.s == 0) return y.f;
 T d = abs(x-y.f);
 P = pow(y.s/d, 2) * (x-y.f) + y.f;
 P b = \operatorname{sqrt} (d*d-y.s*y.s)/d*y.s*unit(x-y.f)*dir(PI/2);
 return t == 0 ? a+b : a-b;
V<pair<P,P>> external(Circ x, Circ y) {
 V<pair<P,P>> v;
 if (x.s == v.s) {
   P \text{ tmp} = \text{unit}(x.f-y.f)*x.s*dir(PI/2);
   v.eb(x.f+tmp,y.f+tmp);
   v.eb(x.f-tmp, v.f-tmp);
 } else {
   P p = (y.s*x.f-x.s*y.f)/(y.s-x.s);
    FOR(i,2) v.eb(tangent(p,x,i),tangent(p,y,i));
 return v;
V<pair<P,P>> internal(Circ x, Circ y) {
 return external({x.f,-x.s},y); }
```

## Circumcenter.h

#### **Description:** returns {circumcenter,circumradius}

#### MinEnclosingCirc.h

#### Description: minimum enclosing circle

### 

## 8.4 Misc

## ClosestPair.h

## **Description:** line sweep to find two closest points **Time:** $\mathcal{O}(N \log N)$

for (; v[i].f-v[ind].f >= bes.f; ++ind)

```
S.erase({v[ind].s,v[ind].f});
for (auto it = S.ub({v[i].s-bes.f,INF});
   it != end(S) && it->f < v[i].s+bes.f; ++it) {
   P t = {it->s,it->f};
   ckmin(bes,{abs(t-v[i]),{t,v[i]}});
}
S.insert({v[i].s,v[i].f});
}
return bes.s;
```

## DelaunayIncremental.h

**Description:** Bowyer-Watson where not all points collinear. Works for  $|x|, |y| \le 10^4$ , assuming that all circumradii in final triangulation are  $\ll 10^9$ . **Time:**  $\mathcal{O}\left(N^2 \log N\right)$ 

```
"DelaunayFast.h"
                                                      57c54d, 23 lines
// include inCircle from DelaunayFast
const T BIG = 1e9; // >> (10^4)^2
V<AR<int,3>> triIncrement(vP v) {
 v.pb({-BIG,-BIG}); v.pb({BIG,0}); v.pb({0,BIG});
 V<AR<int,3>> ret, tmp;
  ret.pb(\{sz(v)-3, sz(v)-2, sz(v)-1\});
 FOR(i,sz(v)-3) {
    map<pi,int> m;
    each(a,ret) {
      if (inCircle(v[i], v[a[0]], v[a[1]], v[a[2]]))
        m[{a[0],a[1]}]++, m[{a[1],a[2]}]++, m[{a[0],a[2]}]++;
      else tmp.pb(a);
    each(a, m) if (a.s == 1) {
      AR < int, 3 > x{a.f.f,a.f.s,i};
      sor(x); tmp.pb(x);
    swap(ret,tmp); tmp.clear();
 each(a, ret) if (a[2] < sz(v)-3) tmp.pb(a);
  return tmp;
```

## DelaunayFast.h

53963d, 13 lines

**Description:** Fast Delaunay triangulation assuming no duplicates and not all points collinear (in latter case, result will be empty). Should work for doubles as well, though there may be precision issues in 'circ'. Returns triangles in ccw order. Each circumcircle will contain none of the input points. If coordinates are ints at most B then  $\mathbb T$  should be large enough to support ints on the order of  $B^4$ .

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

```
"Point.h"
                                                        0e7085, 82 lines
// using T = 11; (if coords are < 2e4)
using 111 = __int128;
// return true if p strictly within circumcircle(a,b,c)
bool inCircle(P p, P a, P b, P c) {
  a \rightarrow p, b \rightarrow p, c \rightarrow p; // assert(cross(a,b,c)>0);
 lll x = (lll) norm(a) *cross(b,c) + (lll) norm(b) *cross(c,a)
      +(111) norm(c) *cross(a,b);
  return x*(cross(a,b,c)>0?1:-1) > 0;
P arb(LLONG_MAX, LLONG_MAX); // not equal to any other point
using Q = struct Quad*;
struct Ouad {
  bool mark; Q o, rot; P p;
  P F() { return r()->p; }
  Q r() { return rot->rot; }
  Q prev() { return rot->o->rot; }
  Q next() { return r()->prev(); }
```

## ManhattanMST Point3D Hull3D

```
Q makeEdge(P orig, P dest) {
  Q q[]{new Quad{0,0,0,orig}, new Quad{0,0,0,arb},
     new Quad{0,0,0,dest}, new Quad{0,0,0,arb}};
  FOR(i, 4) q[i] \rightarrow o = q[-i \& 3], q[i] \rightarrow rot = q[(i+1) \& 3];
  return *a;
void splice(Q a, Q b) { swap(a->o->rot->o, b->o->rot->o); swap(
  \hookrightarrow a->0, b->0); }
0 connect(0 a, 0 b) {
  Q = makeEdge(a->F(), b->p);
  splice(q, a->next()); splice(q->r(), b);
  return q;
pair<0,0> rec(const vP& s) {
  if (sz(s) \le 3) {
    Q = makeEdge(s[0], s[1]), b = makeEdge(s[1], s.bk);
   if (sz(s) == 2) return { a, a->r() };
    splice(a->r(), b);
   auto side = cross(s[0], s[1], s[2]);
   0 c = side ? connect(b, a) : 0;
    return {side < 0 ? c -> r() : a, side < 0 ? c : b -> r() };
#define H(e) e \rightarrow F(), e \rightarrow p
#define valid(e) (cross(e->F(),H(base)) > 0)
  O A, B, ra, rb;
  int half = sz(s) / 2;
  tie(ra, A) = rec({all(s)-half});
  tie(B, rb) = rec(\{sz(s)-half+all(s)\});
  while ((cross(B->p,H(A)) < 0 && (A = A->next())) ||
       (cross(A->p,H(B)) > 0 && (B = B->r()->o)));
  Q base = connect(B->r(), A);
  if (A->p == ra->p) ra = base->r();
  if (B->p == rb->p) rb = base;
#define DEL(e, init, dir) Q e = init->dir; if (valid(e)) \
    while (inCircle(e->dir->F(), H(base), e->F())) {
      0 t = e -> dir; \setminus
      splice(e, e->prev()); \
     splice(e->r(), e->r()->prev()); \
     e = t; \
  while (1) {
   DEL(LC, base->r(), o); DEL(RC, base, prev());
    if (!valid(LC) && !valid(RC)) break;
   if (!valid(LC) || (valid(RC) && inCircle(H(RC), H(LC))))
     base = connect(RC, base->r());
    else base = connect(base->r(), LC->r());
  return {ra, rb};
V<AR<P,3>> triangulate(vP pts) {
  sor(pts); assert(unique(all(pts)) == end(pts)); // no
     →duplicates
  if (sz(pts) < 2) return {};
  Q = rec(pts).f; V<Q>q = {e};
  while (cross(e->o->F(), e->F(), e->p) < 0) e = e->o;
#define ADD { Q c = e; do { c->mark = 1; pts.pb(c->p); \
  q.pb(c->r()); c = c->next(); } while (c != e); }
  ADD; pts.clear();
  int qi = 0; while (qi < sz(q)) if (!(e = q[qi++]) -> mark) ADD;
  V<AR<P,3>> ret(sz(pts)/3);
  FOR(i, sz(pts)) ret[i/3][i%3] = pts[i];
  return ret;
```

## ManhattanMST.h

**Description:** Given N points, returns up to 4N edges which are guaranteed to contain a MST for graph with edge weights w(p,q) = |p.x-q.x| + |p.y-q.y|. Edges are in the form {dist, {src, dst}}.

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

```
"DSU.h"
                                                    b7a3bd, 24 lines
// use standard MST algorithm on result to find final MST
V<pair<int,pi>> manhattanMst(vpi v) {
 vi id(sz(v)); iota(all(id), 0);
 V<pair<int,pi>> ed;
 F0R(k, 4) {
   sort(all(id),[&](int i, int j) {
     return v[i].f+v[i].s < v[j].f+v[j].s; });
    map<int,int> sweep; // find first octant neighbors
    each(i,id) { // those in sweep haven't found neighbor yet
      for (auto it = sweep.lb(-v[i].s);
       it != end(sweep); sweep.erase(it++)) {
       int j = it -> s;
       pi d{v[i].f-v[j].f,v[i].s-v[j].s};if (d.s>d.f)break;
       ed.pb({d.f+d.s,{i,j}});
     sweep[-v[i].s] = i;
    each(p,v) {
     if (k\&1) p.f *=-1;
     else swap(p.f,p.s);
 return ed:
```

## 8.5 3D

## Point3D.h

Description: Basic 3D geometry.

```
10a63a, 82 lines
using P3 = AR<T,3>; using Tri = AR<P3,3>; using vP3 = V<P3>;
T norm(const P3% x) {
 T sum = 0; F0R(i,3) sum += sq(x[i]);
 return sum; }
T abs(const P3& x) { return sqrt(norm(x)); }
P3& operator+=(P3& 1, const P3& r) { F0R(i,3) 1[i] += r[i];
 return 1: }
P3& operator = (P3& 1, const P3& r) { F0R(i,3) 1[i] -= r[i];
 return 1; }
P3& operator*=(P3& 1, const T& r) { F0R(i,3) 1[i] *= r;
 return 1; }
P3& operator/=(P3& 1, const T& r) { F0R(i,3) 1[i] /= r;
 return 1; }
P3 operator-(P3 1) { 1 *= -1; return 1; }
P3 operator+(P3 1, const P3& r) { return 1 += r; }
P3 operator-(P3 1, const P3& r) { return 1 -= r; }
P3 operator*(P3 1, const T& r) { return 1 *= r; }
P3 operator* (const T& r, const P3& 1) { return 1*r; }
P3 operator/(P3 1, const T& r) { return 1 /= r; }
P3 unit(const P3& x) { return x/abs(x); }
T dot(const P3& a, const P3& b) {
 T sum = 0; FOR(i,3) sum += a[i]*b[i];
 return sum; }
P3 cross(const P3& a, const P3& b) {
 return {a[1]*b[2]-a[2]*b[1],a[2]*b[0]-a[0]*b[2],
      a[0]*b[1]-a[1]*b[0]; }
P3 cross(const P3& a, const P3& b, const P3& c) {
 return cross(b-a,c-a); }
P3 perp(const P3& a, const P3& b, const P3& c) {
```

```
return unit(cross(a,b,c)); }
bool isMult(const P3& a, const P3& b) { // for long longs
 P3 c = cross(a,b); FOR(i,sz(c)) if (c[i] != 0) return 0;
bool collinear(const P3& a, const P3& b, const P3& c) {
 return isMult(b-a,c-a); }
T DC(const P3&a,const P3&b,const P3&c,const P3&p) {
 return dot(cross(a,b,c),p-a); }
bool coplanar(const P3&a,const P3&b,const P3&c,const P3&p) {
 return DC(a,b,c,p) == 0; }
bool op(const P3& a, const P3& b) {
 int ind = 0; // going in opposite directions?
  FOR(i,1,3) if (std::abs(a[i]*b[i])>std::abs(a[ind]*b[ind]))
 return a[ind] *b[ind] < 0;</pre>
// coplanar points, b0 and b1 on opposite sides of a0-a1?
bool opSide(const P3&a,const P3&b,const P3&c,const P3&d) {
 return op(cross(a,b,c),cross(a,b,d)); }
// coplanar points, is a in Triangle b
bool inTri(const P3& a, const Tri& b) {
 FOR(i,3) if (opSide(b[i],b[(i+1)%3],b[(i+2)%3],a)) return 0;
 return 1; }
// point-sea dist
T psDist(const P3&p,const P3&a,const P3&b) {
 if (dot(a-p,a-b) \le 0) return abs(a-p);
 if (dot(b-p,b-a) <= 0) return abs(b-p);
 return abs(cross(p,a,b))/abs(a-b);
// projection onto line
P3 foot(const P3& p, const P3& a, const P3& b) {
 P3 d = unit(b-a); return a+dot(p-a,d)*d; }
// rotate p about axis
P3 rotAxis(const P3& p, const P3& a, const P3& b, T theta) {
 P3 dz = unit(b-a), f = foot(p,a,b);
  P3 dx = p-f, dy = cross(dz, dx);
 return f+cos(theta)*dx+sin(theta)*dy;
// projection onto plane
P3 foot(const P3& a, const Tri& b) {
 P3 c = perp(b[0],b[1],b[2]);
 return a-c*(dot(a,c)-dot(b[0],c)); }
// line-plane intersection
P3 lpIntersect (const P3&a0, const P3&a1, const Tri&b) {
 P3 c = unit(cross(b[2]-b[0],b[1]-b[0]));
 T x = dot(a0,c) - dot(b[0],c), y = dot(a1,c) - dot(b[0],c);
 return (y*a0-x*a1)/(y-x);
```

#### Hull3D.h

**Description:** Incremental 3D convex hull where not all points are coplanar. Normals to returned faces point outwards. If coordinates are into at most B then  $\mathbb{T}$  should be large enough to support into on the order of  $B^3$ . Changes order of points.

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

292bc6, 92 lines

// using T=11;

T above (const P3&a, const P3&b, const P3&c, const P3&p) {
    return DC(a,b,c,p) > 0; } // is p strictly above plane

void prep(vP3& p) { // rearrange points such that
    shuffle(all(p),rng); // first four are not coplanar
    int dim = 1;

FOR(i,1,sz(p))
    if (dim == 1) {
        if (p[0] != p[i]) swap(p[1],p[i]), ++dim;
    } else if (dim == 2) {
```

566170, 15 lines

fcc3f7, 13 lines

```
if (!collinear(p[0],p[1],p[i]))
       swap(p[2],p[i]), ++dim;
   } else if (dim == 3) {
     if (!coplanar(p[0],p[1],p[2],p[i]))
        swap(p[3],p[i]), ++dim;
  assert (dim == 4);
using F = AR<int,3>; // face
V<F> hull3d(vP3& p) {
  // s.t. first four points form tetra
  prep(p); int N = sz(p); V<F> hull; // triangle for each face
  auto ad = [\&] (int a, int b, int c) { hull.pb(\{a,b,c\}); };
  // +new face to hull
  ad(0,1,2), ad(0,2,1); // initialize hull as first 3 points
  V<vb> in(N,vb(N)); // is zero before each iteration
  FOR(i,3,N) { // incremental construction
    V<F> def, HULL; swap(hull, HULL);
    // HULL now contains old hull
    auto ins = [&](int a, int b, int c) {
     if (in[b][a]) in[b][a] = 0; // kill reverse face
     else in[a][b] = 1, ad(a,b,c);
    };
    each(f, HULL) {
     if (above(p[f[0]],p[f[1]],p[f[2]],p[i]))
       FOR(j,3) ins(f[j],f[(j+1)%3],i);
        // recalc all faces s.t. point is above face
     else def.pb(f);
    each(t,hull) if (in[t[0]][t[1]]) // edge exposed,
     in[t[0]][t[1]] = 0, def.pb(t); // add a new face
    swap(hull,def);
  return hull;
V<F> hull3dFast(vP3& p) {
 prep(p); int N = sz(p); V < F > hull;
  vb active; V<vi> rvis; V<AR<pi,3>> other;
  // whether face is active
  // points visible from each face
  // other face adjacent to each edge of face
  V<vi> vis(N); // faces visible from each point
  auto ad = [&](int a, int b, int c) {
   hull.pb({a,b,c}); active.pb(1); rvis.eb(); other.eb(); };
  auto ae = [&](int a, int b) { vis[b].pb(a), rvis[a].pb(b); };
  auto abv = [&] (int a, int b) {
   f f=hull[a]; return above(p[f[0]],p[f[1]],p[f[2]],p[b]);};
  auto edge = [&](pi e) -> pi {
   return {hull[e.f][e.s],hull[e.f][(e.s+1)%3]}; };
  auto glue = [&] (pi a, pi b) { // link two faces by an edge
   pi x = edge(a); assert(edge(b) == mp(x.s, x.f));
   other[a.f][a.s] = b, other[b.f][b.s] = a;
  }; // ensure face 0 is removed when i=3
  ad(0,1,2), ad(0,2,1); if (abv(1,3)) swap(p[1],p[2]);
  FOR(i,3) glue(\{0,i\},\{1,2-i\});
  FOR(i,3,N) ae(abv(1,i),i); // coplanar points go in rvis[0]
  vi label (N, -1);
  FOR(i,3,N) { // incremental construction
    vi rem; each(t,vis[i]) if (active[t]) active[t]=0, rem.pb(t
    if (!sz(rem)) continue; // hull unchanged
    int st = -1;
    each(r, rem) FOR(j, 3) {
      int o = other[r][j].f;
      if (active[o]) { // create new face!
        int a,b; tie(a,b) = edge(\{r,j\}); ad(a,b,i); st = a;
        int cur = sz(rvis)-1; label[a] = cur;
       vi tmp; set_union(all(rvis[r]),all(rvis[o]),
```

```
back_inserter(tmp));
      // merge sorted vectors ignoring duplicates
      each(x,tmp) if (abv(cur,x)) ae(cur,x);
      glue({cur,0},other[r][j]); // glue old w/ new face
  for (int x = st, y; x = y) { // glue new faces together
   int X = label[x]; glue({X,1},{label[y=hull[X][1]],2});
    if (y == st) break;
V<F> ans; F0R(i,sz(hull)) if (active[i]) ans.pb(hull[i]);
return ans;
```

## PolvSaVol.h

Description: surface area and volume of polyhedron, normals to faces must point outwards

```
52fc2b, 8 lines
"Hull3D.h"
pair<T,T> SaVol(vP3 p, V<F> faces) {
 T s = 0, v = 0;
 each(i,faces) {
   P3 a = p[i[0]], b = p[i[1]], c = p[i[2]];
   s += abs(cross(a,b,c)); v += dot(cross(a,b),c);
 return {s/2, v/6};
```

## Delaunav3.h

Description: Delaunay triangulation with 3D hull. Fails when all points collinear. If coordinates are into at most B, T should be large enough to support ints on the order of  $B^4$ .

```
"Point.h", "Hull3D.h", "AngleCmp.h"
                                                      14907e, 15 lines
V<AR<P,3>> triHull(vP p) {
 V<P3> p3; V<AR<P,3>> res; each(x,p) p3.pb({x.f,x.s,norm(x)});
 bool ok = 0; each(t,p3) ok = !coplanar(p3[0],p3[1],p3[2],t);
 if (!ok) { // all points concyclic
   sort(1+all(p),[&p](P a, P b) {
      return cross(a-p[0],b-p[0])>0; });
    FOR(i,1,sz(p)-1) res.pb({p[0],p[i],p[i+1]});
    #define nor(x) P(p3[x][0], p3[x][1])
    each(t,hull3dFast(p3))
      if (dot(cross(p3[t[0]],p3[t[1]],p3[t[2]]), \{0,0,1\}) < 0)
        res.pb(\{nor(t[0]), nor(t[2]), nor(t[1])\});
 return res;
```

## Strings (9)

## 9.1 Light

#### KMP.h

**Description:** f[i] is length of the longest proper suffix of the *i*-th prefix of s that is a prefix of sTime:  $\mathcal{O}(N)$ 

```
4538e4, 13 lines
vi kmp(str s) {
 int N = sz(s); vi f(N+1); f[0] = -1;
  FOR(i, 1, N+1) {
    for (f[i]=f[i-1];f[i]!=-1&&s[f[i]]!=s[i-1];)f[i]=f[f[i]];
    ++f[i]; }
  return f;
vi getOc(str a, str b) { // find occurrences of a in b
```

```
vi f = kmp(a+"@"+b), ret;
FOR(i, sz(a), sz(b)+1) if (f[i+sz(a)+1] == sz(a))
  ret.pb(i-sz(a));
return ret;
```

## Z.h

**Description:** f[i] is the max len such that s.substr(0,len) == s.substr(i,len) Time:  $\mathcal{O}(N)$ 

```
vi z(str s) {
 int N = sz(s), L = 1, R = 0; s += '#';
 vi ans(N); ans[0] = N;
 FOR(i,1,N) {
   if (i \le R) ans[i] = min(R-i+1, ans[i-L]);
   while (s[i+ans[i]] == s[ans[i]]) ++ans[i];
   if (i+ans[i]-1 > R) L = i, R = i+ans[i]-1;
 return ans;
vi getPrefix(str a, str b) { // find prefixes of a in b
 vi t = z(a+b); t = vi(sz(a)+all(t));
 each(u,t) ckmin(u,sz(a));
 return t;
```

#### Manacher.h

Description: length of largest palindrome centered at each character of string and between every consecutive pair

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

```
vi manacher(str _S) {
  str S = "@"; each(c, S) S += c, S += "#";
  S.bk = '&';
  vi ans(sz(S)-1); int lo = 0, hi = 0;
  FOR(i, 1, sz(S)-1) {
    if (i != 1) ans[i] = min(hi-i, ans[hi-i+lo]);
    while (S[i-ans[i]-1] == S[i+ans[i]+1]) ++ans[i];
    if (i+ans[i] > hi) lo = i-ans[i], hi = i+ans[i];
  ans.erase(begin(ans));
  FOR(i, sz(ans)) if (i%2 == ans[i]%2) ++ans[i];
  return ans:
```

## LyndonFactor.h

**Description:** A string is "simple" if it is strictly smaller than any of its own nontrivial suffixes. The Lyndon factorization of the string s is a factorization  $s = w_1 w_2 \dots w_k$  where all strings  $w_i$  are simple and  $w_1 \geq w_2 \geq \dots \geq w_k$ . Min rotation gets min index i such that cyclic shift of s starting at i is mini-

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

40c2f1, 19 lines

```
vs duval(str s) {
 int N = sz(s); vs factors;
 for (int i = 0; i < N; ) {
    int j = i+1, k = i;
    for (; j < n \&\& s[k] <= s[j]; j++) {
     if (s[k] < s[j]) k = i;
      else ++k;
    for (; i \le k; i += j-k) factors.pb(s.substr(i, j-k));
 return factors;
int minRotation(str s) {
 int n = sz(s); s += s;
 auto d = duval(s); int ind = 0, ans = 0;
```

```
while (ans+sz(d[ind]) < n) ans += sz(d[ind++]);
while (ind && d[ind] == d[ind-1]) ans -= sz(d[ind--]);
return ans;</pre>
```

## HashRange.h

Description: Polynomial hash for substrings with two bases. fc0b90, 24 lines

```
using H = AR<int,2>; // bases not too close to ends
H makeH(char c) { return {c,c}; }
uniform_int_distribution<int> BDIST(0.1*MOD, 0.9*MOD);
const H base{BDIST(rng),BDIST(rng)};
H operator+(H l, H r) {
 FOR(i,2) if ((l[i] += r[i]) >= MOD) l[i] -= MOD;
  return 1; }
H operator-(H 1, H r) {
 FOR(i,2) if ((1[i] -= r[i]) < 0) 1[i] += MOD;
  return 1; }
H operator*(H l, H r) {
 FOR(i,2) 1[i] = (11)1[i] *r[i] %MOD;
  return 1; }
V<H> pows{{1,1}};
struct HashRange {
  str S; V<H> cum{{}};
  void add(char c) { S += c; cum.pb(base*cum.bk+makeH(c)); }
  void add(str s) { each(c,s) add(c); }
  void extend(int len) { while (sz(pows) <= len)</pre>
   pows.pb(base*pows.bk); }
  H hash(int 1, int r) { int len = r+1-1; extend(len);
    return cum[r+1]-pows[len]*cum[1]; }
```

### ReverseBW.h

**Description:** Used only once. Burrows-Wheeler Transform appends # to a string, sorts the rotations of the string in increasing order, and constructs a new string that contains the last character of each rotation. This function reverses the transform.

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

e400d8, 7 lines

```
str reverseBW(str t) {
  vi nex(sz(t)); iota(all(nex),0);
  stable_sort(all(nex),[&t](int a,int b){return t[a]<t[b];});
  str ret; for (int i = nex[0]; i; )
    ret += t[i = nex[i]];
  return ret;
}</pre>
```

## ACfixed.h

**Description:** Aho-Corasick for fixed alphabet. For each prefix, stores link to max length suffix which is also a prefix.

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

c0b364, 28 lines

```
struct ACfixed { // fixed alphabet
    static const int ASZ = 26;
    struct node { AR<int, ASZ> to; int link; };
    V<node> d = {{}};
    int add(str s) { // add word
        int v = 0;
        each(C,s) {
            int c = C-'a';
            if (!d[v].to[c]) d[v].to[c] = sz(d), d.eb();
            v = d[v].to[c];
        }
        return v;
    }
    void init() { // generate links
        d[0].link = -1;
        queue<int> q; q.push(0);
```

```
while (sz(q)) {
   int v = q.ft; q.pop();
   FOR(c,ASZ) {
     int u = d[v].to[c]; if (!u) continue;
     d[u].link = d[v].link == -1 ? 0 : d[d[v].link].to[c];
     q.push(u);
   }
   if (v) FOR(c,ASZ) if (!d[v].to[c])
     d[v].to[c] = d[d[v].link].to[c];
}
};
```

## SuffixArray.h

**Description:** Sort suffixes. First element of sa is sz(S), isa is the inverse of sa, and lcp stores the longest common prefix between every two consecutive elements of sa.

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

```
"RMQ.h"
                                                     27a566, 30 lines
struct SuffixArray {
 str S; int N; vi sa, isa, lcp;
 void init(str \_S) { N = sz(S = \_S)+1; genSa(); genLcp(); }
 void genSa() { // sa has size sz(S)+1, starts with sz(S)
   sa = isa = vi(N); sa[0] = N-1; iota(1+all(sa), 0);
    sort(1+all(sa),[&](int a, int b) { return S[a] < S[b]; });
   FOR(i, 1, N)  { int a = sa[i-1], b = sa[i];
     isa[b] = i > 1 \&\& S[a] == S[b] ? isa[a] : i; }
    for (int len = 1; len < N; len *= 2) { // currently sorted
     // by first len chars
     vi s(sa), is(isa), pos(N); iota(all(pos),\theta);
     each(t,s) {int T=t-len; if (T>=0) sa[pos[isa[T]]++] = T;}
     FOR(i, 1, N)  { int a = sa[i-1], b = sa[i];
       isa[b] = is[a] == is[b] \&\&is[a+len] == is[b+len]?isa[a]:i;
 void genLcp() { // Kasai's Algo
   lcp = vi(N-1); int h = 0;
   FOR(b, N-1) { int a = sa[isa[b]-1];
     while (a+h < sz(S) \&\& S[a+h] == S[b+h]) ++h;
     lcp[isa[b]-1] = h; if (h) h--; }
   R.init(lcp);
 RMO<int> R;
 int getLCP(int a, int b) { // lcp of suffixes starting at a,b
   if (a == b) return sz(S)-a;
   int l = isa[a], r = isa[b]; if (l > r) swap(l,r);
   return R.query(1,r-1);
```

## SuffixArrayLinear.h

Description: Linear Time Suffix Array

```
vi sa_is(const vi& s, int upper) {
  int n = sz(s); if (!n) return {};
  vi sa(n); vb ls(n);
  R0F(i,n-1) ls[i] = s[i] == s[i+1] ? ls[i+1] : s[i] < s[i+1];
  vi sum_l(upper), sum_s(upper);
  F0R(i,n) (ls[i] ? sum_l[s[i]+1] : sum_s[s[i]])++;
  F0R(i,upper) {
    if (i) sum_l[i] += sum_s[i-1];
    sum_s[i] += sum_l[i];
  }
  auto induce = [&] (const vi& lms) {
    fill(all(sa),-1);
    vi buf = sum_s;
    for (int d: lms) if (d != n) sa[buf[s[d]]++] = d;
    buf = sum_l; sa[buf[s[n-1]]++] = n-1;</pre>
```

```
int v = sa[i]-1;
    if (v \ge 0 \&\& !ls[v]) sa[buf[s[v]]++] = v;
  R0F(i,n) {
   int v = sa[i]-1;
    if (v >= 0 \&\& ls[v]) sa[--buf[s[v]+1]] = v;
};
vi lms_map(n+1,-1), lms; int m = 0;
FOR(i,1,n) if (!ls[i-1] && ls[i]) lms_map[i]=m++, lms.pb(i);
induce(lms); // sorts LMS prefixes
vi sorted_lms; each(v,sa)if (lms_map[v]!=-1)sorted_lms.pb(v);
vi rec s(m); int rec upper = 0; // smaller subproblem
FOR(i,1,m) { // compare two lms substrings in sorted order
  int l = sorted_lms[i-1], r = sorted_lms[i];
  int end l = lms map[l]+1 < m ? lms[lms map[l]+1] : n;
  int end_r = lms_map[r]+1 < m ? lms[lms_map[r]+1] : n;
  bool same = 0; // whether 1ms substrings are same
  if (end 1-1 == end r-r) {
    for (:1 < end 1 && s[1] == s[r]; ++1,++r);
    if (1 != n \&\& s[1] == s[r]) same = 1;
  rec_s[lms_map[sorted_lms[i]]] = (rec_upper += !same);
vi rec sa = sa is(rec s.rec upper+1);
FOR(i,m) sorted_lms[i] = lms[rec_sa[i]];
induce(sorted lms); // sorts LMS suffixes
return sa:
```

## TandemRepeats.h

**Description:** Find all (i, p) such that s.substr(i,p) == s.substr(i+p,p). No two intervals with the same period intersect or touch.

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

## 9.2 Heavy

## PalTree.h

ed0bb4, 46 lines

**Description:** Used infrequently. Palindromic tree computes number of occurrences of each palindrome within string. ans [1] [0] stores min even x such that the prefix s[1.i] can be split into exactly x palindromes, ans [1] [1] does the same for odd x.

Time:  $\mathcal{O}(N \sum)$  for addChar,  $\mathcal{O}(N \log N)$  for updAns

8a7d31, 41 lines

```
struct PalTree {
  static const int ASZ = 26;
  struct node {
    AR<int,ASZ> to = AR<int,ASZ>();
    int len, link, oc = 0; // # occurrences of pal
    int slink = 0, diff = 0;
```

## SuffixAutomaton SuffixTree CircularLCS

```
AR<int,2> seriesAns;
 node(int _len, int _link) : len(_len), link(_link) {}
str s = "@"; V<AR<int, 2>> ans = \{\{0, MOD\}\};
V < node > d = \{\{0,1\}, \{-1,0\}\}; // dummy pals of len 0,-1
int last = 1:
int getLink(int v) {
 while (s[sz(s)-d[v].len-2] != s.bk) v = d[v].link;
void updAns() { // serial path has O(log n) vertices
  ans.pb({MOD,MOD});
  for (int v = last; d[v].len > 0; v = d[v].slink) {
   d[v].seriesAns=ans[sz(s)-1-d[d[v].slink].len-d[v].diff];
   if (d[v].diff == d[d[v].link].diff)
     FOR(i,2) ckmin(d[v].seriesAns[i],
           d[d[v].link].seriesAns[i]);
    // start of previous oc of link[v]=start of last oc of v
   FOR(i,2) ckmin(ans.bk[i],d[v].seriesAns[i^1]+1);
void addChar(char C) {
 s += C; int c = C-'a'; last = getLink(last);
 if (!d[last].to[c]) {
   d.eb(d[last].len+2,d[getLink(d[last].link)].to[c]);
   d[last].to[c] = sz(d)-1;
   auto& z = d.bk; z.diff = z.len-d[z.link].len;
   z.slink = z.diff == d[z.link].diff
     ? d[z.link].slink : z.link;
  } // max suf with different dif
  last = d[last].to[c]; ++d[last].oc;
  updAns();
void numOc() { ROF(i,2,sz(d)) d[d[i].link].oc += d[i].oc; }
```

## SuffixAutomaton.h

Description: Used infrequently. Constructs minimal deterministic finite automaton (DFA) that recognizes all suffixes of a string. len corresponds to the maximum length of a string in the equivalence class, pos corresponds to the first ending position of such a string, 1nk corresponds to the longest suffix that is in a different class. Suffix links correspond to suffix tree of the reversed string!

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

```
a99c6<u>d, 67 lines</u>
struct SuffixAutomaton {
  int N = 1; vi lnk\{-1\}, len\{0\}, pos\{-1\}; // suffix link,
  // max length of state, last pos of first occurrence of state
  V<map<char,int>> nex{1}; V<bool> isClone{0};
  // transitions, cloned -> not terminal state
  V<vi> iLnk; // inverse links
  int add(int p, char c) { // \sim p \text{ nonzero if } p != -1
    auto getNex = [&]() {
      if (p == -1) return 0;
      int q = nex[p][c]; if (len[p]+1 == len[q]) return q;
      int clone = N++; lnk.pb(lnk[q]); lnk[q] = clone;
     len.pb(len[p]+1), nex.pb(nex[q]),
     pos.pb(pos[q]), isClone.pb(1);
      for (; ~p && nex[p][c] == q; p = lnk[p]) nex[p][c]=clone;
      return clone;
    // if (nex[p].count(c)) return getNex();
    // ^ need if adding > 1 string
    int cur = N++; // make new state
    lnk.eb(), len.pb(len[p]+1), nex.eb(),
   pos.pb(pos[p]+1), isClone.pb(0);
    for (; \sim p \&\& !nex[p].count(c); p = lnk[p]) nex[p][c] = cur;
    int x = getNex(); lnk[cur] = x; return cur;
```

```
void init(str s) { int p = 0; each(x,s) p = add(p,x); }
 // inverse links
 void genIlnk() {iLnk.rsz(N); FOR(v,1,N) iLnk[lnk[v]].pb(v);}
 // APPLICATIONS
 void getAllOccur(vi& oc, int v) {
    if (!isClone[v]) oc.pb(pos[v]); // terminal position
    each(u,iLnk[v]) getAllOccur(oc,u); }
 vi allOccur(str s) { // get all occurrences of s in automaton
    each(x,s) {
     if (!nex[cur].count(x)) return {};
     cur = nex[cur][x]; }
    // convert end pos -> start pos
    vi oc; getAllOccur(oc,cur); each(t,oc) t += 1-sz(s);
    sort(all(oc)); return oc;
 vl distinct;
 11 getDistinct(int x) {
    // # distinct strings starting at state x
   if (distinct[x]) return distinct[x];
    distinct[x]=1;each(y,nex[x]) distinct[x]+=getDistinct(y.s);
    return distinct[x]; }
 11 numDistinct() { // # distinct substrings including empty
    distinct.rsz(N); return getDistinct(0); }
 11 numDistinct2() { // assert(numDistinct() == numDistinct2());
   ll ans = 1; FOR(i,1,N) ans += len[i]-len[lnk[i]];
    return ans: }
SuffixAutomaton S:
vi sa; str s;
void dfs(int x) {
 if (!S.isClone[x]) sa.pb(sz(s)-1-S.pos[x]);
 V<pair<char,int>> chr;
 each(t,S.iLnk[x]) chr.pb(\{s[S.pos[t]-S.len[x]],t\});
 sort(all(chr)); each(t,chr) dfs(t.s);
int main() {
 re(s); reverse(all(s));
 S.init(s); S.genIlnk();
 dfs(0); ps(sa); // generating suffix array for s
```

## SuffixTree.h

Description: Used infrequently. Ukkonen's algorithm for suffix tree. Longest non-unique suffix of s has length len[p]+lef after each call to add terminates. Each iteration of loop within add decreases this quantity by one. Time:  $\mathcal{O}(N \log \Sigma)$ 39751c, 51 lines

```
struct SuffixTree {
 str s; int N = 0;
 vi pos, len, lnk; V<map<char,int>> to;
 int make (int POS, int LEN) { // lnk[x] is meaningful when
   // x!=0 and len[x] != MOD
   pos.pb(POS);len.pb(LEN);lnk.pb(-1);to.eb();return N++; }
 void add(int& p, int& lef, char c) { // longest
   // non-unique suffix is at node p with lef extra chars
   s += c; ++lef; int lst = 0;
   for (; lef; p?p=lnk[p]: lef--) { // if p != root then lnk[p]
     // must be defined
     while (lef>1 && lef>len[to[p][s[sz(s)-lef]]])
       p = to[p][s[sz(s)-lef]], lef -= len[p];
     // traverse edges of suffix tree while you can
     char e = s[sz(s)-lef]; int& q = to[p][e];
     // next edge of suffix tree
     if (!q) q = make(sz(s)-lef,MOD), lnk[lst] = p, lst = 0;
     // make new edge
     else {
```

```
char t = s[pos[q]+lef-1];
       if (t == c) { lnk[lst] = p; return; } // suffix not
       int u = make(pos[q],lef-1);
        // new node for current suffix-1, define its link
       to[u][c] = make(sz(s)-1,MOD); to[u][t] = q;
        // new, old nodes
       pos[q] += lef-1; if (len[q] != MOD) len[q] -= lef-1;
        q = u, lnk[lst] = u, lst = u;
 void init(str s) {
   make (-1, 0); int p = 0, lef = 0;
    each(c, s) add(p,lef,c);
    add(p,lef,'$'); s.pop_back(); // terminal char
 int maxPre(str x) { // max prefix of x which is substring
    for (int p = 0, ind = 0;;) {
     if (ind == sz(x) || !to[p].count(x[ind])) return ind;
     p = to[p][x[ind]];
     FOR(i,len[p]) {
       if (ind == sz(x) \mid \mid x[ind] != s[pos[p]+i]) return ind;
 vi sa; // generate suffix array
 void genSa(int x = 0, int Len = 0) {
   if (!sz(to[x])) sa.pb(pos[x]-Len); // found terminal node
    else each(t,to[x]) genSa(t.s,Len+len[x]);
};
```

## Various (10)

## 10.1 Dynamic programming

When doing DP on intervals:  $a[i][j] = \min_{i < k < j} (a[i][k] + a[k][j]) + f(i, j),$  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) < 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.

#### CircularLCS.h

**Description:** Used only twice. For strs A, B calculates longest common subsequence of A with all rotations of BTime:  $\mathcal{O}(|A| \cdot |B|)$ 

```
int circular_lcs(str A, str B) {
 B += B;
 int max_lcs = 0;
 V < vb > dif_left(sz(A)+1, vb(sz(B)+1)), dif_up(sz(A)+1, vb(sz(B))
```

```
auto recalc = [&](int x, int y) { assert(x && y);
 int res = (A.at(x-1) == B.at(y-1))
   dif_up[x][y-1] | dif_left[x-1][y];
 dif_left[x][y] = res-dif_up[x][y-1];
 dif_up[x][y] = res-dif_left[x-1][y];
FOR(i,1,sz(A)+1) FOR(j,1,sz(B)+1) recalc(i,j);
FOR(j, sz(B)/2) {
  // 1. zero out dp[.][j], update dif_left and dif_right
  if (i) for (int x = 1, y = i; x \le sz(A) \&\& y \le sz(B); ) {
   int pre_up = dif_up[x][y];
   if (y == j) dif_up[x][y] = 0;
   else recalc(x,y);
    (pre_up == dif_up[x][y]) ? ++x : ++y;
  // 2. calculate LCS(A[0:sz(A)),B[j:j+sz(B)/2))
  int cur_lcs = 0;
 FOR(x, 1, sz(A) + 1) cur_lcs += dif_up[x][j+sz(B)/2];
  ckmax(max_lcs,cur_lcs);
return max_lcs;
```

## SMAWK.h

**Description:** Given negation of totally monotone matrix with entries of type  $\mathbb{D}$ , find indices of row maxima (their indices increase for every submatrix). If tie, take lesser index. f returns matrix entry at (r,c) in O(1). Use in place of divide & conquer to remove a log factor.

Time:  $\mathcal{O}(R+C)$ , can be reduced to  $\mathcal{O}(C(1+\log R/C))$  evaluations of f

```
template < class F, class D=11> vi smawk (F f, vi x, vi y) {
  vi ans(sz(x),-1); // x = rows, y = cols
  \#define upd() if (ans[i] == -1 || w > mx) ans[i] = c, mx = w
  if (\min(sz(x), sz(y)) \le 8) {
   FOR(i,sz(x)) { int r = x[i]; D mx;
      each(c, y) \{ D w = f(r, c); upd(); \} \}
    return ans;
  if (sz(x) < sz(y)) { // reduce subset of cols to consider
   vi Y; each(c,y) {
     for (; sz(Y); Y.pop\_back()) \{ int X = x[sz(Y)-1]; \}
       if (f(X,Y.bk) >= f(X,c)) break; }
     if (sz(Y) < sz(x)) Y.pb(c);
   y = Y;
  } // recurse on half the rows
  vi X; for (int i = 1; i < sz(x); i += 2) X.pb(x[i]);
  vi ANS = smawk(f,X,y); FOR(i,sz(ANS)) ans[2*i+1] = ANS[i];
  for (int i = 0, k = 0; i < sz(x); i += 2) {
    int to = i+1 < sz(ans) ? ans[i+1] : y.bk; D mx;
    for(int r = x[i];;++k) {
     int c = y[k]; D w = f(r,c); upd();
     if (c == to) break; }
 return ans;
};
```

## 10.2 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 violations 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).

## 10.3 Optimization tricks

## 10.3.1 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.
- FOR(b,k) FOR(i,1<<K) if (i&1<<b) D[i] += D[i^(1<<b)]; computes all sums of subsets.

## 10.3.2 Pragmas

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

  Also consider older #pragma GCC target ("sse4").
- #pragma GCC optimize ("trapv") kills the program on integer overflows (but is really slow).

## FastIO.h

Description: Fast input and output.

**Time:** input is  $\sim 300 \text{ms}$  faster for  $10^6$  long longs on CF  $_{6a2f14, \ 37 \text{ lines}}$ 

```
namespace FastIO {
 const int BSZ = 1<<15; ///// INPUT</pre>
 char ibuf[BSZ]; int ipos, ilen;
 char nc() { // next char
   if (ipos == ilen) {
     ipos = 0; ilen = fread(ibuf, 1, BSZ, stdin);
     if (!ilen) return EOF;
   return ibuf[ipos++];
 void rs(str& x) { // read str
   char ch; while (isspace(ch = nc()));
   do { x += ch; } while (!isspace(ch = nc()) && ch != EOF);
 tcT> void ri(T& x) { // read int or 11
   char ch; int sgn = 1;
   while (!isdigit(ch = nc())) if (ch == '-') sqn \star= -1;
   x = ch'' \theta'; while (isdigit(ch = nc())) x = x*10+(ch'' \theta');
   x \star = sqn;
 tcT, class... Ts> void ri(T& t, Ts&... ts) {
   ri(t); ri(ts...); } // read ints
  ///// OUTPUT (call initO() at start)
 char obuf[BSZ], numBuf[100]; int opos;
```

```
void flushOut() { fwrite(obuf,1,opos,stdout); opos = 0; }
void wc(char c) { // write char
   if (opos == BSZ) flushOut();
   obuf[opos++] = c; }
void ws(str s) { each(c,s) wc(c); } // write str
tcT> void wi(T x, char after = '\0') {
   if (x < 0) wc('-'), x *= -1;
   int len = 0; for (;x>=10;x/=10) numBuf[len++] = '0'+(x%10);
   wc('0'+x); R0F(i,len) wc(numBuf[i]);
   if (after) wc(after);
}
void initO() { assert(atexit(flushOut) == 0); }
```

## 10.4 Other languages

## Python3.py

Description: not PyPy3, solves CF Factorisation Collaboration

```
47 lines
```

25

```
from math import *
import sys, random
def nextInt():
 return int(input())
def nextStrs():
 return input().split()
def nextInts():
 return list(map(int,nextStrs()))
n = nextInt()
def process(x):
 global v
  x = abs(x)
  for t in v: # print(type(t)) -> <class 'int'>
    g = gcd(t, x)
    if g != 1:
      V.append(g)
    if g != t:
      V.append(t//g)
for i in range (50):
 x = random.randint(0, n-1)
 if gcd(x,n) != 1:
    process(x)
    sx = x * x * n \# assert(qcd(sx,n) == 1)
    print(f"sqrt {sx}") # print value of var
    sys.stdout.flush()
    X = nextInt()
    process(x+X)
    process(x-X)
print(f'! {len(v)}',end='')
for i in v:
 print(f' {i}',end='')
print()
sys.stdout.flush() # sys.exit(0) -> exit
# sys.setrecursionlimit(int(1e9)) -> stack size
# print(f'{ans:=.6f}') -> print ans to 6 decimal places
from decimal import * # arbitrary precision decimals
ctx = getcontext()
ctx.prec = 28
print(Decimal(1) / Decimal(7)) # 0.1428571428571428571428571429
print(ctx.power(Decimal(10),-30)) # 1E-30
```