

International Institute Of Information Technology - Hyderabad

ladaiLadai

Aryan Gupta, Aarush jain, Keyur Chaudhari

ICPC Regionals 2021 Ganpati Bappa Morya Mangal Murti Morya

Combinatorial (1)

1.1 Permutations

1.1.1 Factorial

n	123	4	5 6	7	8	9		10 28800 17 3.6e14 171 >DBL_MAX
n!	126	24 12	20 720	5040	4032	0 3628	380 362	28800
n	11	12	13	1	4	15	16	17
n!	4.0e7	⁷ 4.8e	8 6.26	9 8.7	e10 1.	3e12 2	2.1e13	3.6e14
n	20	25	30	40	50	100	150	171
$\overline{n!}$	2e18	2e25	3e32	8e47	3e64	9e157	6e262	>DBL_MAX

1.1.2 Derangements

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

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

1.1.3 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

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

1.2 Partitions and subsets

1.2.1 Partition function

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

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

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

1.2.2 Lucas' Theorem

Let n,m be non-negative integers and p a prime. Write $n=n_kp^k+\ldots+n_1p+n_0$ and $m=m_kp^k+\ldots+m_1p+m_0$. The m binomials \pmod{p} . multinomial.

1.3 General purpose numbers

1.3.1 Bernoulli numbers

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

Sums of powers:

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

Euler-Maclaurin formula for infinite sums:

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

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

1.3.2 "Stirling numbers of the first kind

Number of permutations on n items with k cycles.

$$c(n,k) = c(n-1,k-1) + (n-1)c(n-1,k), \ c(0,0) = 1$$

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

$$c(8, k) = 8, 0, 5040, 13068, 13132, 6769, 1960, 322, 28, 1$$

 $c(n, 2) = 0, 0, 1, 3, 11, 50, 274, 1764, 13068, 109584, \dots$

1.3.3 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. $\pi(j) > j$.

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

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

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

1.3.4 Bell numbers

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

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

1.3.5 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 : $(n-2)!/((d_1-1)!\cdots(d_n-1)!)$

1.3.6 Stirling numbers of the second kind

Partitions of n distinct elements into exactly k groups. S(n,k) = S(n-1,k-1) + kS(n-1,k) S(n,1) = S(n,n) = 1

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

1.3.7 Catalan numbers

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

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

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

- sub-diagonal monotone paths in an $n \times n$ grid.
- 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.
- permutations of [n] with no 3-term increasing subseq.

1.4 DP Optimizations

```
Quadrangle Inequality. f satisfies it if \forall a \leq b \leq c \leq d, f(a,d) - f(a,c) \geq f(b,d) - f(b,c).
```

```
1D-1D.h
```

```
Description: Applicable if dp_i = min_{j>i}(dp_j + cost(i, j) \text{ s.t. } opt_i \leq opt_j when i \leq j (which holds if quadrangle) 
Time: \mathcal{O}(n \log n)
```

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

```
#define until first
#define opt second
11 dp[100000];
ll cost(int i, int j) {
  return dp[j] /* + cost to jump from i to j*/;
void solve(int n) {
  dp[n] = 0;
  vector<PII> v;
  v.EB(n - 1, n);
  for (int i = n - 1, ipos = 0; i >= 1; i--) {
    while (ipos+1 < SZ(v) && i <= v[ipos+1].until) ipos
    dp[i] = cost(i, v[ipos].opt);
    while (v.back().until < i</pre>
      && cost(v.back().until, i)
         <= cost(v.back().until, v.back().opt))</pre>
      { v.pop back(); }
    int l = 1, r = min(i - 1, v.back().until);
    while (l <= r) {
      int mid = (l + r)/2;
      if (cost(mid, i) <= cost(mid, v.back().opt)) {</pre>
        l = mid + 1;
      } else { r = mid - 1; }
    if (l - 1 >= 1) v.EB(l - 1, i);
```

Divide-and-Conquer.h

```
Description: Works when dp_{k,i} = min_{j < i} (dp_{k-1,j} + cost(j,i)) and opt_k(i) \leq opt_k(i+1). (This holds when quadrangle) Usage: find dp[1], then: for(i = 2 to n) solve(i, 1, n, 1, n) Time: \mathcal{O}\left(kn\log n\right)
```

```
Il dp[100][100]; // set correctly
Il cost(int i, int j); // cost to go from i to j, 1-
   indexed.
void solve(int i, int l, int r, int optl, int optr) {
   const ll inf = 1e18; // set correctly
   if (l > r || optl > optr) return;
   int mid = (l + r)/2; pair<ll, int> best = {inf, -1};
```

```
for (int j = optl; j <= min(mid, optr); j++) {</pre>
    pair<ll,int> cand(dp[i - 1][j] + cost(j, mid), j);
    if (best.second == -1) best = cand;
    else best = min(best, cand);
 dp[i][mid] = best.first;
 solve(i, l, mid - 1, optl, best.second);
 solve(i, mid + 1, r, best.second, optr);
Dvnamic-CHT.h
Description: Add lines y = ax + b and query for min at given x
Time: \mathcal{O}(logn) per update/query.
const dbl INF = 1e16;
struct HullDynamic {
#define CLREF const Line&
 struct Line {
    int a; ll b, val=0; dbl xLeft = -INF; bool type=0;
    Line(int a = 0, ll b = 0): a(a), b(b) {}
    ll eval(int x) const{ return a * 1ll * x + b; }
    bool operator< (CLREF l2) const {</pre>
      return l2.type ? (xLeft > l2.val) : (a < l2.a);
 };
 using Iter = set<Line>::iterator;
 bool parallel(CLREF l1, CLREF l2) { return l1.a == l2
 dbl meetX(CLREF l1, CLREF l2) {
    return parallel(l1, l2) ? INF :
      (l2.b-l1.b) / (dbl(l1.a-l2.a));
 set<Line> hull;
 bool hasPrev(Iter it) { return it != hull.begin(); }
 bool hasNext(Iter it) {
    return it != hull.end() && next(it) != hull.end();
 bool bad(CLREF l1, CLREF l2, CLREF l3){
    return meetX(l1,l3) <= meetX(l1,l2);</pre>
 bool bad(Iter it) {
    return hasPrev(it) && hasNext(it)
        && (bad(*next(it), *it, *prev(it)));
 Iter upd left border(Iter it) {
   if(!hasNext(it)) return it;
    dbl val = meetX(*it, *next(it));
    Line buf(*it); it = hull.erase(it);
    buf.xLeft = val; return hull.insert(it, buf);
 void insert line(int a, ll b) {
    Line l3 = Line(a, b); auto it = hull.lower bound(l3
    if (it != hull.end() && parallel(*it , l3)) {
      if (it->b <= b) return;</pre>
      it = hull.erase(it);
    it = hull.insert(it, l3);
```

KnuthDP.h

Description: 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 $f(b,c) \le f(a,d)$ and $f(a,c)+f(b,d) \ge f(a,d)+f(b,c)$ for all $a \le b \le c \le d$. Consider also: LineContainer (ch. Data structures), monotone queues, ternary search. **Time:** $\mathcal{O}\left(N^2\right)$

Numerical (2)

2.1 Polynomials and recurrences

Polynomial.h

2ad8d0, 17 line:

```
PolyRoots.h
```

```
Description: Finds the real roots to a polynomial.
Usage: polyRoots(\{\{2,-3,1\}\},-1e9,1e9\}) // solve x^2-3x+2=
Time: \mathcal{O}\left(n^2\log(1/\epsilon)\right)
"Polynomial.h"
                                                     7118fe, 23 lines
vector<double> polyRoots(Poly p, double xmin, double
    xmax) {
 if (SZ(p.a) == 2) { return {-p.a[0]/p.a[1]}; }
  vector<double> ret;
  Poly der = p;
  der.diff();
  auto dr = polyRoots(der, xmin, xmax);
  dr.push back(xmin-1);
  dr.push back(xmax+1);
  sort(ALL(dr));
  REP(i,0,SZ(dr)-1) {
    double l = dr[i], h = dr[i+1];
    bool sign = p(l) > 0;
    if (sign ^{(p(h) > 0)}) {
      REP(it,0,60) { // while (h - l > 1e-8)
         double m = (l + h) / 2, f = p(m);
        if ((f \le 0) \land sign) l = m;
        else h = m:
      ret.push back((l + h) / 2);
  return ret;
```

PolyInterpolate.h

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

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

Matrices

IntDeterminant.h

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

```
Time: \mathcal{O}(N^3)
                                                                                                 669167, 18 lines
```

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

SolveLinear.h

x.assign(m, 0);

b[i] /= A[i][i];

x[col[i]] = b[i];

for (int i = rank; i--;) {

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

```
typedef vector<double> vd;
const double eps = 1e-12;
int solveLinear(vector<vd>& A, vd& b, vd& x) {
 int n = SZ(A), m = SZ(x), rank = 0, br, bc;
 if (n) assert(SZ(A[0]) == m);
 VI col(m); iota(ALL(col), 0);
 REP(i,0,n) {
   double v, bv = 0;
   REP(r,i,n) REP(c,i,m)
     if ((v = fabs(A[r][c])) > bv)
       br = r, bc = c, bv = v;
   if (bv <= eps) {
      REP(j,i,n) if (fabs(b[j]) > eps) return -1;
      break;
   swap(A[i], A[br]);
   swap(b[i], b[br]);
   swap(col[i], col[bc]);
   REP(j,0,n) swap(A[j][i], A[j][bc]);
   bv = 1/A[i][i];
   REP(j,i+1,n) {
      double fac = A[j][i] * bv;
      b[i] -= fac * b[i];
      REP(k,i+1,m) A[j][k] -= fac*A[i][k];
   rank++;
```

```
REP(j,0,i) b[j] -= A[j][i] * b[i];
return rank; // (multiple solutions if rank < m)</pre>
```

SolveLinear2.h

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

```
"SolveLinear.h"
                                                    5cad07, 7 lines
REP(j,0,n) if (j != i) // instead of REP(j,i+1,n)
// ... then at the end:
x.assign(m, undefined);
REP(i,0,rank) {
  REP(j,rank,m) if (fabs(A[i][j]) > eps) goto fail;
  x[col[i]] = b[i] / A[i][i];
fail:: }
```

SolveLinearBinary.h

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

```
typedef bitset<1000> bs;
int solveLinear(vector<bs>& A, VI& b, bs& x, int m) {
 int n = SZ(A), rank = 0, br;
 assert(m \le SZ(x));
 VI col(m); iota(ALL(col), 0);
 REP(i,0,n) {
    for (br=i; br<n; ++br) if (A[br].any()) break;</pre>
   if (br == n) {
      REP(j,i,n) if(b[j]) return -1;
      break:
    int bc = (int)A[br]. Find next(i-1);
    swap(A[i], A[br]);
    swap(b[i], b[br]);
    swap(col[i], col[bc]);
   REP(j,0,n) if (A[j][i] != A[j][bc]) {
      A[j].flip(i); A[j].flip(bc);
   REP(j,i+1,n) if (A[j][i]) {
      b[j] ^= b[i];
      A[i] ^= A[i];
    rank++:
 x = bs();
 for (int i = rank; i--;) {
   if (!b[i]) continue;
   x[col[i]] = 1;
   REP(j,0,i) b[j] ^= A[j][i];
 return rank; // (multiple solutions if rank < m)</pre>
```

MatrixInverse.h

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

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

```
43464f, 35 lines
int matInv(vector<vector<double>>& A) {
 int n = SZ(A): VI col(n):
 vector<vector<double>> tmp(n, vector<double>(n));
 REP(i,0,n) tmp[i][i] = 1, col[i] = i;
 REP(i,0,n) {
    int r = i, c = i;
    REP(j,i,n) REP(k,i,n)
     if (fabs(A[j][k]) > fabs(A[r][c]))
        r = j, c = k;
    if (fabs(A[r][c]) < 1e-12) return i;
    A[i].swap(A[r]); tmp[i].swap(tmp[r]);
    REP(j,0,n)
      swap(A[j][i], A[j][c]), swap(tmp[j][i], tmp[j][c
          1);
    swap(col[i], col[c]);
    double v = A[i][i];
    REP(j,i+1,n) {
      double f = A[j][i] / v;
      A[i][i] = 0;
     REP(k,i+1,n) A[j][k] -= f*A[i][k];
     REP(k,0,n) tmp[j][k] -= f*tmp[i][k];
    REP(j, i+1, n) A[i][j] /= v;
    REP(j,0,n) tmp[i][j] /= v;
    A[i][i] = 1;
 for (int i = n-1; i > 0; --i) REP(j,0,i) {
    double v = A[i][i];
    REP(k,0,n) tmp[j][k] -= v*tmp[i][k];
 REP(i,0,n) REP(j,0,n) A[col[i]][col[j]] = tmp[i][j];
```

Tridiagonal.h

return n:

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

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

This is useful for solving problems on the type

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

```
where a_0, a_{n+1}, b_i, c_i and d_i are known. a can then be obtained from
   \{a_i\} = \text{tridiagonal}(\{1, -1, -1, ..., -1, 1\}, \{0, c_1, c_2, ..., c_n\},\
                     \{b_1, b_2, \ldots, b_n, 0\}, \{a_0, d_1, d_2, \ldots, d_n, a_{n+1}\}\}.
Fails if the solution is not unique.
If |d_i| > |p_i| + |q_{i-1}| for all i, or |d_i| > |p_{i-1}| + |q_i|, or the matrix is positive
definite, the algorithm is numerically stable and neither tr nor the check for
diag[i] == 0 is needed.
Time: \mathcal{O}(N)
typedef double T;
vector<T> tridiagonal(vector<T> diag, const vector<T>&
    const vector<T>& sub, vector<T> b) {
  int n = SZ(b); VI tr(n);
  REP(i,0,n-1) {
    if (abs(diag[i]) < 1e-9 * abs(super[i])) { // diag[
         i1 == 0
       b[i+1] -= b[i] * diag[i+1] / super[i];
       if (i+2 < n) b[i+2] -= b[i] * sub[i+1] / super[i</pre>
       diaq[i+1] = sub[i]; tr[++i] = 1;
    } else {
       diaq[i+1] -= super[i]*sub[i]/diag[i];
       b[i+1] -= b[i]*sub[i]/diag[i];
  for (int i = n; i--;) {
    if (tr[i]) {
       swap(b[i], b[i-1]);
       diaq[i-1] = diaq[i];
       b[i] /= super[i-1];
     } else {
       b[i] /= diag[i];
       if (i) b[i-1] -= b[i]*super[i-1];
  return b:
```

2.3 Fourier transforms

FastFourierTransform.h

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

```
a333b0, 35 lines
typedef complex<double> C;
typedef vector<double> vd:
void fft(vector<C>& a) {
 int n = SZ(a), L = 31 - builtin clz(n);
 static vector<complex<long double>> R(2, 1);
 static vector<C> rt(2, 1); // (^ 10% faster if
 for (static int k = 2; k < n; k *= 2) {
```

```
R.resize(n); rt.resize(n);
   auto x = polar(1.0L, acos(-1.0L) / k);
   REP(i,k,2*k) rt[i] = R[i] = i&1 ? R[i/2] * x : R[i
  VI rev(n);
  REP(i,0,n) rev[i] = (rev[i / 2] | (i \& 1) << L) / 2;
  REP(i,0,n) if (i < rev[i]) swap(a[i], a[rev[i]]);
  for (int k = 1; k < n; k *= 2)
   for (int i = 0; i < n; i += 2 * k) REP(j,0,k) {
      Cz = rt[j+k] * a[i+j+k]; // (25% faster if hand-
          rolled)
      a[i + j + k] = a[i + j] - z;
      a[i + j] += z;
vd conv(const vd& a, const vd& b) {
  if (a.empty() || b.empty()) return {};
 vd res(SZ(a) + SZ(b) - 1);
  int L = 32 - builtin clz(SZ(res)), n = 1 \ll L;
  vector<C> in(n), out(n);
  copy(ALL(a), begin(in));
  REP(i,0,SZ(b)) in[i].imag(b[i]);
  fft(in):
  for (C& x : in) x *= x;
  REP(i,0,n) out[i] = in[-i & (n - 1)] - conj(in[i]);
  fft(out):
  REP(i,0,SZ(res)) res[i] = imag(out[i]) / (4 * n);
  return res:
```

FastFourierTransformMod.h

REP(i,0,SZ(res)) {

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

```
Time: \mathcal{O}(N \log N), where N = |A| + |B| (twice as slow as NTT or FFT)
```

```
"FastFourierTransform.h"
typedef vector<ll> vl;
template<int M> vl convMod(const vl &a, const vl &b) {
 if (a.empty() || b.empty()) return {};
 vl res(SZ(a) + SZ(b) - 1);
  int B=32- builtin clz(SZ(res)), n=1<<B, cut=int(sqrt</pre>
      (M));
  vector<C> L(n), R(n), outs(n), outl(n);
  REP(i,0,SZ(a)) L[i] = C((int)a[i] / cut, (int)a[i] %
  REP(i,0,SZ(b)) R[i] = C((int)b[i] / cut, (int)b[i] %
      cut);
  fft(L), fft(R);
 REP(i,0,n) {
   int j = -i \& (n - 1);
   outl[j] = (L[i] + conj(L[j])) * R[i] / (2.0 * n);
   outs[j] = (L[i] - conj(L[j])) * R[i] / (2.0 * n) /
        1i:
 fft(outl), fft(outs);
```

Description: Transform to a basis with fast convolutions of the form c[z] =

```
ll av = ll(real(outl[i])+.5), cv = ll(imag(outs[i])
      +.5):
 II bv = II(imag(outl[i])+.5) + II(real(outs[i])+.5)
  res[i] = ((av % M * cut + bv) % M * cut + cv) % M;
return res;
```

NumberTheoreticTransform.h

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

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

```
FastSubsetTransform.h
```

```
\sum_{z=x \oplus y} a[x] \cdot b[y], where \oplus is one of AND, OR, XOR. The size of a must
be a power of two.
Time: \mathcal{O}(N \log N)
void FST(VI& a, bool inv) {
 for (int n = SZ(a), step = 1; step < n; step *= 2) {
    for (int i = 0; i < n; i += 2 * step) REP(j,i,i+
         step) {
      int \&u = a[j], \&v = a[j + step]; tie(u, v) =
        inv ? PII(v - u, u) : PII(v, u + v); // AND
        inv ? PII(v, u - v) : PII(u + v, u); // OR
        PII(u + v, u - v);
 if (inv) for (int& x : a) x /= SZ(a); // XOR only
VI conv(VI a, VI b) {
  FST(a, 0); FST(b, 0);
  REP(i,0,SZ(a)) a[i] *= b[i];
  FST(a, 1); return a;
```

WalshHadamard.h

```
Description: C_k = \sum_{i \otimes j = k} A_i B_j
Usage: Apply the transform, point multiply and invert
Time: \mathcal{O}(N \log N)
void WalshHadamard(poly &P, bool invert) {
```

```
for (int len = 1; 2 * len <= SZ(P); len <<= 1) {</pre>
  for (int i = 0; i < SZ(P); i += 2 * len) {
    REP(i, 0, len) {
      auto u = P[i + j], v = P[i + len + j];
      P[i + j] = u + v, P[i + len + j] = u - v; //
  }
if (invert) for (auto \&x : P) \times /= SZ(P);
```

```
Description: Given B_1, \ldots B_m, compute A_i = \sum_{j=1}^{i-1} A_j * B_{i-j}
Usage: 1-indexed, pad B[i] = 0 for i > m
Time: \mathcal{O}\left(N\log^2 N\right)
                                                        211338, 18 lines
void online(const Poly &B, CD a1, int n, Poly &A) {
  const int m = SZ(B) - 1;
  A.assign(n + 1, 0); A[1] = a1;
  auto bst = B.begin(), ast = A.begin();
  REP(i, 1, n) {
    A[i + 1] += A[i] * B[1];
    if (i + 2 \le n) A[i + 2] += A[i] * B[2];
    for (int pw = 2; i % pw == 0 \&\& pw + 1 <= m; pw <<=
```

Poly blockA(ast + i - pw, ast + i);

```
Poly blockB(bst + pw + 1, bst + min(pw * 2, m) +
   1);
Poly prod = conv(blockA, blockB);
REP(i, 0, SZ(prod)) {
 if (i + 1 + j \le n)
   A[i + 1 + j] += prod[j];
```

Number theory (3)

Modular arithmetic

ModInverse.h

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

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

ModPow.h

```
int power(long long x, unsigned int y, int p){
 int res = 1:
 x = x % p;
 if (x == 0)
   return 0;
 while (y > 0)
   if (y & 1)
      res = (res * x) % p;
   y = y >> 1;
   x = (x * x) % p:
 return res;
```

NCR.h

```
Description: Calculates ncr for large N and prime Mod
```

#define MAX N FACT (int)3e5 + 5

```
vector<long long> factorial(MAX N FACT, 1),
   inverse factorial(MAX_N_FACT, 1), inv(MAX_N_FACT,
```

```
1);
void prec factorials() {
    for (int i = 2; i < MAX N FACT; i++) {</pre>
        factorial[i] = factorial[i - 1] * i % mod;
        inv[i] = (mod - (mod / i) * inv[mod % i] % mod)
        inverse factorial[i] = (inv[i] *
            inverse factorial[i - 1]) % mod;
long long ncr(int n, int k) {
```

```
return factorial[n] * inverse_factorial[k] % mod *
   inverse_factorial[n - k] % mod;
```

ModLog.h

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

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

ModSum.h

return -1;

Description: Sums of mod'ed arithmetic progressions.

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

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

```
c5bc5, 16 lines
```

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

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

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

ModSqrt.h

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

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

```
// a^{(n+3)/8} or 2^{(n+3)/8} * 2^{(n-1)/4} works if p % 8
ll s = p - 1, n = 2;
int r = 0, m;
while (s % 2 == 0)
  ++r. s /= 2:
while (modpow(n, (p - 1) / 2, p) != p - 1) ++n;
II x = modpow(a, (s + 1) / 2, p);
11 b = modpow(a, s, p), q = modpow(n, s, p);
for (;; r = m) {
  ll t = b:
  for (m = 0; m < r \&\& t != 1; ++m)
   t = t * t % p;
  if (m == 0) return x;
  II gs = modpow(g, 1LL \ll (r - m - 1), p);
  q = qs * qs % p;
 x = x * gs % p;
  b = b * q % p;
```

3.2 Primality

FastEratosthenes.h

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

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

MillerRabin.h

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

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

```
"ModMulLL.h" 60dcd1, 12 lines

bool isPrime(ull n) {
```

Factor.h

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

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

```
"ModMulLL.h", "MillerRabin.h"
                                                  65f857, 18 lines
ull pollard(ull n) {
  auto f = [n](ull x) \{ return modmul(x, x, n) + 1; \};
  ull x = 0, y = 0, t = 30, prd = 2, i = 1, q;
  while (t++ % 40 | | gcd(prd, n) == 1) {
    if (x == y) x = ++i, y = f(x);
    if ((q = modmul(prd, max(x,y) - min(x,y), n))) prd
    x = f(x), y = f(f(y));
  return gcd(prd, n);
vector<ull> factor(ull n) {
 if (n == 1) return {};
  if (isPrime(n)) return {n};
  ull x = pollard(n);
  auto l = factor(x), r = factor(n / x);
 l.insert(l.end(), ALL(r));
  return l;
```

3.3 Divisibility

euclid.h

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

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

CRT.h

Description: Chinese Remainder Theorem. crt(a, m, b, n) computes x such that

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

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

04d93a, 7 lines

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

3.3.1 Bézout's identity

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

$$ax + by = d$$

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

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

phiFunction.h

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

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

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

Fractions

ContinuedFractions.h

Description: Given N and a real number $x \geq 0$, finds the closest rational approximation p/q with $p, q \leq N$. It will obey $|p/q - x| \leq 1/qN$. For consecutive convergents, $p_{k+1}q_k - q_{k+1}p_k = (-1)^k$. $(p_k/q_k$ alternates between > x and < x.) If x is rational, y eventually becomes ∞ ; if x is the root of a degree 2 polynomial the a's eventually become cyclic. Time: $\mathcal{O}(\log N)$

```
typedef double d; // for N \sim 1e7; long double for N \sim 1
pair<ll, ll> approximate(d x, ll N) {
 ll LP = 0, LQ = 1, P = 1, Q = 0, inf = LLONG MAX; d y
   ll lim = min(P ? (N-LP) / P : inf, Q ? (N-LQ) / Q :
```

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

FracBinarySearch.h

Description: Given f and N, finds the smallest fraction $p/q \in [0,1]$ such that f(p/q) is true, and $p, q \leq N$. You may want to throw an exception from f if it finds an exact solution, in which case N can be removed. Usage: fracBS([](Frac f) { return f.p>=3*f.q; }, 10); // {1,3}

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

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

3.5 Pythagorean Triples

The Pythagorean triples are uniquely generated by

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

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

3.6 Primes

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

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

3.7 **Estimates**

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

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

Mobius Function

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

Mobius Inversion:

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

Other useful formulas/forms:

```
\sum_{d|n} \mu(d) = [n=1] (very useful)
g(n) = \sum_{n|d} f(d) \Leftrightarrow f(n) = \sum_{n|d} \mu(d/n)g(d)
g(n) = \sum_{1 \le m \le n} f(\left|\frac{n}{m}\right|) \Leftrightarrow f(n) = \sum_{1 \le m \le n} \mu(m)g(\left|\frac{n}{m}\right|)
```

Mobius.h

```
Description: Dirichlet - H(n) = \sum_{xy=n} a_x b_y, 1 \le n \le N
                                                                  306cf1, 21 lines
VI mobius(int N) {
  VI mu(N + 1, 1);
```

```
vector<bool> ispr(N + 1, 1);
for (int i = 2; i \le N; ++i) {
  if (!ispr[i]) continue;
```

OrderStatisticTree HashMap Trie Treap LineContainer

Data structures (4)

4.1 Set and Map like

OrderStatisticTree.h

Description: A set (not multiset!) with support for finding the n'th element, and finding the index of an element. To get a map, change null_type.

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

782797, 16 lines

HashMap.h

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

d77092, 7 lines

```
gnu pbds::gp hash table<ll,int,chash> h({},{},{},{},{}),{
    1<<16}):
Trie.h
Description: krishna
Time: \mathcal{O}(\log N)
                                                  1cba01, 26 lines
const int NX = int(1e6) + int(5e5);
int arr[NX][26];
int root:
int lastocc;
void Trie() {
    root = 0, lastocc = 0;
    memset(arr, 0, sizeof(int) * NX * 26);
void insert(const string &x) {
    int curptr = root:
    for (auto ch : x) {
        if (arr[curptr][ch - 'a'] == 0)
            arr[curptr][ch - 'a'] = ++lastocc;
        curptr = arr[curptr][ch - 'a'];
int search(const string &x) {
    int curptr = root:
    for (auto ch : x) {
        if (arr[curptr][ch - 'a'] == 0)
            return 0:
        else
            curptr = arr[curptr][ch - 'a'];
    return 1;
Treap.h
```

Description: A short self-balancing tree. It acts as a sequential container with log-time splits/joins, and is easy to augment with additional data.

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

```
struct Node {
   Node *l = 0, *r = 0;
   int val, y, c = 1;
   Node(int val) : val(val), y(rand()) {}
   void recalc();
};

int cnt(Node* n) { return n ? n->c : 0; }
void Node::recalc() { c = cnt(l) + cnt(r) + 1; }

template<class F> void each(Node* n, F f) {
   if (n) { each(n->l, f); f(n->val); each(n->r, f); }
}

pair<Node*, Node*> split(Node* n, int k) {
   if (!n) return {};
   if (cnt(n->l) >= k) { // "n->val >= k" for
        lower_bound(k)
        auto pa = split(n->l, k);
```

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

LineContainer.h

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

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

8ec1c7, 30 lines

```
struct Line {
  mutable ll k, m, p;
  bool operator<(const Line& o) const { return k < o.k;
  }
  bool operator<(ll x) const { return p < x; }
};

struct LineContainer : multiset<Line, less<>> {
  // (for doubles, use inf = 1/.0, div(a,b) = a/b)
  static const ll inf = LLONG_MAX;
  ll div(ll a, ll b) { // floored division
   return a / b - ((a ^ b) < 0 && a % b); }</pre>
```

```
bool isect(iterator x, iterator y) {
    if (y == end()) return x -> p = inf, 0;
    if (x->k == y->k) x->p = x->m > y->m ? inf : -inf;
    else x -> p = div(y -> m - x -> m, x -> k - y -> k);
    return x - p >= y - p;
  void add(ll k, ll m) {
    auto z = insert(\{k, m, 0\}), y = z++, x = y;
    while (isect(y, z)) z = erase(z);
    if (x != begin() \&\& isect(--x, y)) isect(x, y =
        erase(y));
    while ((y = x) != begin() \&\& (--x)->p >= y->p)
      isect(x, erase(y));
  ll query(ll x) {
    assert(!empty());
    auto l = *lower bound(x);
    return l.k * x + l.m;
};
```

UnionFindRollback.h

 $\begin{tabular}{ll} \textbf{Description:} & Disjoint-set data structure with undo. If undo is not needed, skip st, time() and rollback(). \end{tabular}$

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

f79afc, 18 lines

d8e33f, 26 lines

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

4.2 Matrix

Matrix.h

Description: Basic operations on square matrices.

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

```
template<class T, int N> struct Matrix {
  typedef Matrix M;
  array<array<T, N>, N> d{};
  M operator*(const M& m) const {
```

```
REP(i,0,N) REP(j,0,N)
      REP(k,0,N) a.d[i][j] += d[i][k]*m.d[k][j];
 vector<T> operator*(const vector<T>& vec) const {
   vector<T> ret(N);
   REP(i,0,N) REP(j,0,N) ret[i] += d[i][j] * vec[j];
    return ret:
 M operator^(ll p) const {
   assert(p >= 0);
   M a, b(*this);
   REP(i,0,N) \text{ a.d}[i][i] = 1;
   while (p) {
     if (p&1) a = a*b;
      b = b*b:
      p >>= 1;
   return a;
};
```

4.3 Range DS

SeamentTree.h

Description: Zero-indexed max-tree. Bounds are inclusive to the left and exclusive to the right. Can be changed by modifying T, f and unit.

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

```
710629, 19 lines
template <typename T>
struct segTree {
 T unit;
 T (*f) (T obj1, T obj2);
 vector<T> s;
 seqTree(int n, T (*c)(T obj1, T obj2), T def) : s(2 *
       n, def), n(n), f(c), unit(def) {}
 void update(int pos, T val) {
   for (s[pos += n] = val; pos /= 2;)
      s[pos] = f(s[pos * 2], s[pos * 2 + 1]);
 T query(int b, int e) { // query [b, e]
   T ra = unit, rb = unit;
    for (b += n, e += n; b < e; b /= 2, e /= 2) {
     if (b % 2) ra = f(ra, s[b++]);
      if (e \% 2) rb = f(s[--e], rb);
    return f(ra, rb);}};
```

LazySegmentTree.h

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

```
Usage: Node* tr = new Node(v, \theta, SZ(v)); Time: \mathcal{O}(\log N).
```

faa5ec, 58 lines

```
struct node {
   int sum, width; // width = # of leaves for node
    node operator+(const node &n) {
        return {max(sum, n.sum), width + n.width);}};
struct update {
    int value; bool type;// 0 = add, 1 = reset
   node operator()(const node &n) {
        if (type) return {n.width * value, n.width};
        else return {n.sum + value, n.width};}
   update operator+(const update &u) {
        if (u.tvpe) return u:
        return {type, value + u.value};}};
template <typename T, typename U>
struct seg tree lazy {
   int S, H; T zero;
    vector<T> value; U noop;
    vector<bool> dirty; vector<U> prop;
    seg tree lazy<T, U>(int S, T zero = T(), U noop
        = U())
        zero = zero, noop = _noop;
        for (S = 1, H = 1; S < S;) S *= 2, H++;
        value.resize(2 * S, zero), dirty.resize(2 * S,
            false):
        prop.resize(2 * S, noop);}
    void set leaves(vector<T> &leaves) {
        copy(leaves.begin(), leaves.end(), value.begin
            () + S):
        for (int i = S - 1; i > 0; i--)
            value[i] = value[2 * i] + value[2 * i + 1];
    void apply(int i, U &update) {
        value[i] = update(value[i]);
        if (i < S) {
            prop[i] = prop[i] + update;
            dirty[i] = true;}}
    void rebuild(int i) {
        for (int l = i / 2; l; l /= 2) {
            T combined = value[2 * l] + value[2 * l +
            value[l] = prop[l](combined):}}
    void propagate(int i) {
        for (int h = H; h > 0; h--) {
            int l = i >> h:
            if (dirty[l]) {
                apply(2 * 1, prop[l]);
                apply(2 * l + 1, prop[l]);
                prop[l] = noop;
                dirty[l] = false;}}
   void upd(int i, int j, U update) {
        i += S, j += S;
        propagate(i), propagate(j);
        for (int l = i, r = j; l <= r; l /= 2, r /= 2)
            if ((l & 1) == 1) apply(l++, update);
            if ((r & 1) == 0) apply(r--, update);}
        rebuild(i), rebuild(j);}
   T query(int i, int j) {
```

```
i += S, j += S;
propagate(i), propagate(j);
T res left = zero, res right = zero;
for (; i \le j; i \ne 2, j \ne 2) {
    if ((i & 1) == 1) res left = res left +
        value[i++];
    if ((j \& 1) == 0) res right = value[j--] +
        res right;}
return res left + res right;}};
```

FenwickTree.h

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

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

b06af0, 22 lines

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

FenwickTree2d.h

Description: Computes sums a[i,i] for all i<1, i<J, and increases single elements a[i,i]. Requires that the elements to be updated are known in advance (call fakeUpdate() before init()).

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

```
d5c1b7, 22 lines
struct FT2 {
 vector<VI> ys; vector<FT> ft;
 FT2(int limx) : ys(limx) {}
 void fakeUpdate(int x, int y) {
   for (; x < SZ(ys); x = x + 1) ys[x].push back(y);
 void init() {
    for (VI& v : ys) sort(ALL(v)), ft.emplace back(SZ(v
       ));
 int ind(int x, int y) {
```

```
return (int)(lower bound(ALL(ys[x]), y) - ys[x].
         begin()); }
  void update(int x, int y, ll dif) {
    for (; x < SZ(ys); x = x + 1)
      ft[x].update(ind(x, y), dif);
 ll query(int x, int y) {
    ll sum = 0;
    for (; x; x \&= x - 1)
      sum += ft[x-1].query(ind(x-1, y));
    return sum:
RMQ.h
Description: Range Minimum Queries on an array. Returns min(V[a], V[a
+ 1], ... V[b - 1]) in constant time.
Usage: RMQ rmq(values);
rmq.query(inclusive, exclusive);
Time: \mathcal{O}(|V|\log|V|+Q)
                                                    9a1bbf, 16 lines
template<class T>
struct RMQ {
```

```
vector<vector<T>> jmp;
  RMQ(const \ vector< T>\&\ V) : jmp(1,\ V)  {
    for (int pw = 1, k = 1; pw * 2 <= SZ(V); pw *= 2.
        ++k) {
      jmp.emplace back(SZ(V) - pw * 2 + 1);
      REP(j,0,SZ(jmp[k]))
        jmp[k][j] = min(jmp[k - 1][j], jmp[k - 1][j +
            ; ([wq
 T query(int a, int b) {
    assert(a < b): // or return inf if a == b
    int dep = 31 - builtin clz(b - a);
    return min(jmp[dep][a], jmp[dep][b - (1 << dep)]);</pre>
 }
};
```

MoQueries.h

Description: Answer interval or tree path queries by finding an approximate TSP through the queries, and moving from one query to the next by adding/removing points at the ends. If values are on tree edges, change step to add/remove the edge (a, c) and remove the initial add call (but keep in).

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

```
void add(int ind, int end) { ... } // add a[ind] (end =
     0 or 1)
void del(int ind, int end) { ... } // remove a[ind]
int calc() { ... } // compute current answer
VI mo(vector<PII> 0) {
 int L = 0, R = 0, blk = 350; // \sim N/sqrt(Q)
 VI s(SZ(Q)), res = s;
#define K(x) PII(x.first/blk, x.second ^ -(x.first/blk
    & 1))
  iota(ALL(s), 0);
```

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

MoWithUpdates.h

Description: Supports point updates at position

```
Time: \mathcal{O}(n^{5/3}) when block = n^{2/3}
```

```
struct Query { int l, r, id, t; }
struct Update { int pos, pre, now; };
void MoWithUpdates(vector<Query> qs, vector<Update> upd
    ) {
  int BLK; // set block size
  sort(qs.begin(), qs.end(), [&](Query a, Query b) {
    return {a.l/BLK, a.r/BLK, a.t} < {b.l/BLK, b.r/BLK,
         b.t;;);
  for (auto g : gs) {
    while (t < q.t) ++t, apply(upd[t].pos, upd[t].now);</pre>
    while (t > q.t) apply(upd[t].pos, upd[t].pre), --t;
```

```
while (l > q.l) add(--l);
while (l < q.l) remove(l++);
while (r < q.r) add(++r);
while (r > q.r) remove(r--);
ans[q.id] = qet();}
```

Strings (5)

String Matching

KMP.h

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

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

```
VI pi(const string& s) {
  VI p(SZ(s));
  REP(i,1,SZ(s)) {
    int q = p[i-1];
    while (g \&\& s[i] != s[g]) g = p[g-1];
    p[i] = g + (s[i] == s[g]);
  return p;
VI match(const string& s, const string& pat) {
  VI p = pi(pat + '\0' + s), res;
  REP(i,SZ(p)-SZ(s),SZ(p))
    if (p[i] == SZ(pat)) res.push back(i - 2 * SZ(pat))
  return res;
```

Zfunc.h

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

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

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

AhoCorasick.h

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

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

f7553c, 66 lines

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

```
int n = 0;
    VI res; // ll count = 0;
    for (char c : word) {
      n = N[n].next[c - first];
      res.push back(N[n].end);
      // count += N[n].nmatches;
    return res;
  vector<VI> findAll(vector<string>& pat, string word)
    VI r = find(word);
    vector<VI> res(SZ(word));
    REP(i,0,SZ(word)) {
      int ind = r[i];
      while (ind != -1) {
        res[i - SZ(pat[ind]) + 1].push back(ind);
        ind = backp[ind];
    return res;
};
```

Hashing.h

```
Description: Self-explanatory methods for string hashing.
// Arithmetic mod 2^64-1. 2x slower than mod 2^64 and
// code, but works on evil test data (e.g. Thue-Morse,
    where
// ABBA... and BAAB... of length 2^10 hash the same mod
// "typedef ull H;" instead if you think test data is
// or work mod 10^9+7 if the Birthday paradox is not a
    problem.
struct H {
 typedef uint64 t ull;
  ull x; H(ull x=0) : x(x) {}
#define OP(0,A,B) H operator O(H \circ) { ull r = x; asm \
  (A "addq %%rdx, %0\n adcq $0,\%0" : "+a"(r) : B);
      return r: }
  OP(+,,"d"(o.x)) OP(*,"mul %1\n", "r"(o.x) : "rdx")
  H operator-(H o) { return *this + \sim0.X; }
  ull get() const { return x + !\sim x; }
  bool operator==(H o) const { return get() == o.get();
  bool operator<(H o) const { return get() < o.get(); }</pre>
static const H C = (ll)1e11+3; // (order ~ 3e9; random
    also ok)
struct HashInterval {
 vector<H> ha, pw;
 HashInterval(string& str) : ha(SZ(str)+1), pw(ha) {
    pw[0] = 1;
    REP(i,0,SZ(str))
```

```
ha[i+1] = ha[i] * C + str[i],
      pw[i+1] = pw[i] * C;
  H hashInterval(int a, int b) { // hash [a, b) and 0
      indexed
    return ha[b] - ha[a] * pw[b - a];
};
MinRotation.h
Description: Finds the lexicographically smallest rotation of a string.
             rotate(v.begin(), v.begin()+minRotation(v),
v.end());
Time: \mathcal{O}(N)
int minRotation(string s) {
  int a=0, N=SZ(s); s += s;
  REP(b,0,N) REP(k,0,N) {
    if (a+k == b || s[a+k] < s[b+k]) {b += max(0, k-1);}
          break;}
    if (s[a+k] > s[b+k]) { a = b; break; }
```

5.2 Palindromes

Manacher.h

return a;

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

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

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

5.3 Suffix DS

SuffixArray.h

Description: Builds suffix array for a string. sa[i] is the starting index of the suffix which is i'th in the sorted suffix array. The returned vector is of size n+1, and sa[0]=n. The lcp array contains longest common prefixes for neighbouring strings in the suffix array: lcp[i]=lcp(sa[i],sa[i-1]),lcp[0]=0. The input string must not contain any zero bytes. Time: $\mathcal{O}(n\log n)$

```
struct SuffixArray {
  VI sa, lcp;
```

```
SuffixArray(string& s, int lim=256) { // or
      basic string<int>
   int n = SZ(s) + 1, k = 0, a, b;
   VI x(ALL(s)+1), y(n), ws(max(n, lim)), rank(n);
    sa = lcp = v, iota(ALL(sa), 0);
   for (int j = 0, p = 0; p < n; j = max(1, j * 2),
        \lim = p) {
      p = j, iota(ALL(y), n - j);
     REP(i,0,n) if (sa[i] >= j) y[p++] = sa[i] - j;
      fill(ALL(ws), 0);
      REP(i,0,n) ws[x[i]]++;
      REP(i,1,lim) ws[i] += ws[i - 1];
      for (int i = n; i--;) sa[--ws[x[y[i]]]] = y[i];
      swap(x, y), p = 1, x[sa[0]] = 0;
      REP(i,1,n) a = sa[i - 1], b = sa[i], x[b] =
        (y[a] == y[b] \&\& y[a + j] == y[b + j]) ? p - 1
   REP(i,1,n) rank[sa[i]] = i;
   for (int i = 0, j; i < n - 1; lcp[rank[i++]] = k)
     for (k \&\& k--, j = sa[rank[i] - 1];
         s[i + k] == s[j + k]; k++);
};
```

SuffixAutomaton.h

Description: Each path in the automaton is a substring (if it ends in a terminal node, it is a suffix) And no. of occurences = no. of ways to reach a terminal node. Or keep reverse edges of suffix links(all prefixes for that substring), then no. of ways to reach a root.

 $\mbox{\bf Time:} \ \mathcal{O} \, (len) \ \mbox{map accesses, map can be at most of size alphabet, can also use unordered_map}$

```
37fe84, 34 lines
struct SuffixAutomaton {
 vector<map<char, int>> edges;
 VI link, length; // length[i]: longest string in i-th
       class
 int last;
                   // index of equivalence class of
      whole string
 SuffixAutomaton(string s) : edges{}, link{-1}, length
      {0}, last(0) {
    edges.emplace back();
    REP(i, 0, SZ(s)) {
      edges.emplace back();
      length.push back(i + 1);
      link.push back(0);
      int r = SZ(edges) - 1, p = last;
      while (p \ge 0 \&\& edges[p].find(s[i]) == edges[p].
          end()) {
        edges[p][s[i]] = r, p = link[p];
      if (p != -1) {
        const int q = edges[p][s[i]];
        if (length[p] + 1 == length[q]) link[r] = q;
          edges.push back(edges[q]);
          length.push back(length[p] + 1);
          link.push back(link[q]);
```

Graph (6)

BellmanFord.h

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

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

```
const ll inf = LLONG MAX:
struct Ed { int a, b, w, s() { return a < b ? a : -a; }
struct Node { ll dist = inf; int prev = -1; };
void bellmanFord(vector<Node>& nodes, vector<Ed>& eds,
    int s) {
  nodes[s].dist = 0;
  sort(ALL(eds), [](Ed a, Ed b) { return a.s() < b.s();</pre>
  int lim = SZ(nodes) / 2 + 2; // /3+100 with shuffled
      vertices
  REP(i,0,lim) for (Ed ed : eds) {
    Node cur = nodes[ed.a], &dest = nodes[ed.b];
    if (abs(cur.dist) == inf) continue;
    ll d = cur.dist + ed.w;
    if (d < dest.dist) {</pre>
      dest.prev = ed.a;
      dest.dist = (i < lim-1 ? d : -inf);}}</pre>
  REP(i,0,lim) for (Ed e : eds) {
    if (nodes[e.a].dist == -inf)
      nodes[e.b].dist = -inf;}}
```

FloydWarshall.h

Description: Calculates all-pairs shortest path in a directed graph that might have negative edge weights. Input is an distance matrix m, where $m[i][j] = \inf$ if i and j are not adjacent. As output, m[i][j] is set to the shortest distance between i and j, \inf if no path, or $-\inf$ if the path goes through a negative-weight cycle.

```
REP(i,0,n) m[i][i] = min(m[i][i], OLL);
REP(k,0,n) REP(i,0,n) REP(j,0,n)
 if (m[i][k] != inf && m[k][j] != inf) {
    auto newDist = max(m[i][k] + m[k][j], -inf);
    m[i][j] = min(m[i][j], newDist);
REP(k,0,n) if (m[k][k] < 0) REP(i,0,n) REP(j,0,n)
  if (m[i][k] != inf && m[k][j] != inf) m[i][j] = -
      inf;}
```

Network flow 6.1

Dinic.h

Description: Flow algorithm with complexity $O(VE \log U)$ where U =max |cap|. $O(\min(E^{1/2}, V^{2/3})E)$ if U = 1; $O(\sqrt{V}E)$ for bipartite matchabfd54, 42 lines

```
struct Dinic {
 struct Edge {
   int to, rev;
   ll c. oc:
   Il flow() { return max(oc - c, OLL); } // if you
        need flows
 };
 VI lvl, ptr, q;
 vector<vector<Edge>> adi:
 Dinic(int n) : lvl(n), ptr(n), q(n), adj(n) {}
 void addEdge(int a, int b, ll c, ll rcap = 0) {
    adj[a].push back({b, SZ(adj[b]), c, c});
    adj[b].push back({a, SZ(adj[a]) - 1, rcap, rcap});
 ll dfs(int v, int t, ll f) {
   if (v == t || !f) return f;
    for (int& i = ptr[v]; i < SZ(adj[v]); i++) {</pre>
      Edge& e = adi[v][i];
     if (lvl[e.to] == lvl[v] + 1)
       if (ll p = dfs(e.to, t, min(f, e.c))) {
          e.c -= p, adj[e.to][e.rev].c += p;
          return p;
       }
    return 0;
 ll calc(int s, int t) {
   II flow = 0; q[0] = s;
    REP(L,0,31) do { // 'int L=30' maybe faster for
        random data
      lvl = ptr = VI(SZ(q));
      int gi = 0, ge = lvl[s] = 1;
      while (qi < qe && !lvl[t]) {</pre>
       int v = q[qi++];
        for (Edge e : adj[v])
         if (!lvl[e.to] && e.c >> (30 - L))
            q[qe++] = e.to, lvl[e.to] = lvl[v] + 1;
      while (ll p = dfs(s, t, LLONG MAX)) flow += p;
    } while (lvl[t]);
    return flow;
```

```
bool leftOfMinCut(int a) { return lvl[a] != 0; }
};
MCMF-SPFA.h
Description: Multiedges and negative costs allowed.
Time: Approximately \mathcal{O}(V^2E^2)
template <typename FLOW, typename COST> struct MCMF {
 const COST INFC = 1e9, EPSC = 0;
  const FLOW INFF = 1e9, EPSF = 0;
 struct Edge {
    int from. to:
    FLOW flow, cap;
    COST cost;
  int nodes, src, dest, m = 0;
  vector<vector<int>> adi:
  vector<Edge> edges;
  void add(int u, int v, FLOW cap, COST cost) {
    edges.EB(u, v, 0, cap, cost);
    adi[u].PB(m++);
    edges.EB(v, u, 0, 0, -cost);
    adi[v].PB(m++);
  vector<COST> dis:
  vector<bool> in0;
  VI par;
  pair<FLOW, COST> SPFA() {
    fill(ALL(dis), INFC);
    fill(ALL(inQ), false);
    queue<int> Q;
    Q.push(src), dis[src] = 0, inQ[src] = true;
    while (!Q.empty()) {
      int u = Q.front(); Q.pop();
      inQ[u] = false;
      for (int i : adj[u]) {
        auto &e = edges[i];
        if (e.cap - e.flow > EPSF
            && dis[e.to] - (dis[u] + e.cost) > EPSC) {
          dis[e.to] = dis[u] + e.cost;
          par[e.to] = i;
          if (!inQ[e.to]) { Q.push(e.to), inQ[e.to] =
              true: }
      }
    if (dis[dest] + EPSC >= INFC) return {0, 0};
    FLOW aug = INFF;
    for (int u = dest; u != src; u = edges[par[u]].from
      aug = min(aug, edges[par[u]].cap - edges[par[u]].
          flow):
    for (int u = dest; u != src; u = edges[par[u]].from
      edges[par[u]].flow += aug;
      edges[par[u] ^ 1].flow -= aug;
```

```
return {aug, aug * dis[dest]};
  MCMF(int n, int s, int t)
  : nodes(n), src(s), dest(t), adj(n), dis(n), inQ(n),
      par(n) {}
  pair<FLOW, COST> mincostmaxflow() {
    pair<FLOW, COST> ans(0, 0);
    while (true) {
      auto cur = SPFA();
      if (cur.first <= EPSF) break;</pre>
      ans.first += cur.first;
      ans.second += cur.second;
    return ans;
};
```

MinCostMaxFlow.h

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

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

q.push({0, s});

```
92261c, 81 lines
#include <bits/extc++.h>
const ll INF = numeric limits<ll>::max() / 4;
tvpedef vector<ll> VL:
struct MCMF {
 int N;
  vector<VI> ed, red;
  vector<VL> cap, flow, cost;
  VI seen;
  VL dist, pi;
  vector<PII> par:
  MCMF(int N) :
    N(N), ed(N), red(N), cap(N, VL(N)), flow(cap), cost
    seen(N), dist(N), pi(N), par(N) {}
  void addEdge(int from, int to, ll cap, ll cost) {
    this->cap[from][to] = cap;
    this->cost[from][to] = cost;
    ed[from].push back(to);
    red[to].push back(from);
  void path(int s) {
    fill(ALL(seen), 0);
    fill(ALL(dist), INF);
    dist[s] = 0; ll di;
    gnu pbds::priority queue<pair<ll, int>> q;
```

vector<decltype(g)::point iterator> its(N);

```
auto relax = [&](int i, ll cap, ll cost, int dir) {
      ll val = di - pi[i] + cost;
      if (cap && val < dist[i]) {
        dist[i] = val;
        par[i] = \{s, dir\};
       if (its[i] == q.end()) its[i] = q.push({-dist[i
            l, i});
        else q.modify(its[i], {-dist[i], i});
   };
   while (!q.empty()) {
      s = q.top().second; q.pop();
      seen[s] = 1; di = dist[s] + pi[s];
      for (int i : ed[s]) if (!seen[i])
        relax(i, cap[s][i] - flow[s][i], cost[s][i], 1)
      for (int i : red[s]) if (!seen[i])
        relax(i, flow[i][s], -cost[i][s], 0);
    REP(i,0,N) pi[i] = min(pi[i] + dist[i], INF);
  pair<ll, ll> maxflow(int s, int t) {
   Il totflow = 0, totcost = 0;
    while (path(s), seen[t]) {
     II fl = INF;
      for (int p,r,x = t; tie(p,r) = par[x], x != s; x
        fl = min(fl, r ? cap[p][x] - flow[p][x] : flow[
            x][p]);
      totflow += fl;
      for (int p,r,x = t; tie(p,r) = par[x], x != s; x
        if (r) flow[p][x] += fl;
        else flow[x][p] -= fl;
    REP(i,0,N) REP(j,0,N) totcost += cost[i][j] * flow
    return {totflow, totcost};
  // If some costs can be negative, call this before
      maxflow:
  void setpi(int s) { // (otherwise, leave this out)
   fill(ALL(pi), INF); pi[s] = 0;
    int it = N, ch = 1; ll v;
    while (ch-- && it--)
      REP(i,0,N) if (pi[i] != INF)
       for (int to : ed[i]) if (cap[i][to])
          if ((v = pi[i] + cost[i][to]) < pi[to])
            pi[to] = v, ch = 1;
    assert(it >= 0); // negative cost cycle
};
```

GlobalMinCut.h

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

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

```
pair<int, VI> globalMinCut(vector<VI> mat) {
 pair<int, VI> best = {INT MAX, {}};
 int n = SZ(mat);
 vector<VI> co(n);
 REP(i,0,n) co[i] = \{i\};
 REP(ph,1,n) {
   VI w = mat[0];
   size t s = 0, t = 0;
   REP(it, 0, n-ph) { // O(V^2) -> O(E \log V) with prio.
         queue
     w[t] = INT MIN;
     s = t, t = max element(ALL(w)) - w.begin();
      REP(i,0,n) w[i] += mat[t][i];
   best = min(best, \{w[t] - mat[t][t], co[t]\});
   co[s].insert(co[s].end(), ALL(co[t]));
   REP(i,0,n) mat[s][i] += mat[t][i];
   REP(i,0,n) mat[i][s] = mat[s][i];
   mat[0][t] = INT MIN;
 return best:
```

6.2 Matching

MinimumVertexCover.h

Description: Finds a minimum vertex cover in a bipartite graph. The size is the same as the size of a maximum matching, and the complement is a maximum independent set.

```
"DFSMatching.h"
VI cover(vector<VI>& q, int n, int m) {
 VI match(m, -1);
 int res = dfsMatching(g, match);
 vector<bool> lfound(n, true), seen(m);
 for (int it : match) if (it != -1) lfound[it] = false
 VI a. cover:
 REP(i,0,n) if (lfound[i]) q.push back(i);
 while (!q.empty()) {
    int i = q.back(); q.pop back();
    lfound[i] = 1;
    for (int e : q[i]) if (!seen[e] && match[e] != -1)
      seen[e] = true;
      q.push back(match[e]);
 REP(i,0,n) if (!lfound[i]) cover.push back(i);
 REP(i,0,m) if (seen[i]) cover.push back(n+i);
 assert(SZ(cover) == res);
 return cover;
```

WeightedMatching.h

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

```
Time: \mathcal{O}\left(N^2M\right)
pair<int, VI> hungarian(const vector<VI> &a) {
  if (a.empty()) return {0, {}};
  int n = SZ(a) + 1, m = SZ(a[0]) + 1;
  VI u(n), v(m), p(m), ans(n - 1);
  REP(i,1,n) {
    p[0] = i;
    int j0 = 0; // add "dummy" worker 0
    VI dist(m, INT MAX), pre(m, -1);
    vector<bool> done(m + 1);
    do { // dijkstra
      done[i0] = true;
      int i0 = p[j0], j1, delta = INT MAX;
      REP(j,1,m) if (!done[j]) {
        auto cur = a[i0 - 1][j - 1] - u[i0] - v[j];
        if (cur < dist[j]) dist[j] = cur, pre[j] = j0;
        if (dist[j] < delta) delta = dist[j], j1 = j;</pre>
      REP(j,0,m) {
        if (done[j]) u[p[j]] += delta, v[j] -= delta;
        else dist[j] -= delta;
      j0 = j1;
    } while (p[j0]);
    while (j0) { // update alternating path
      int j1 = pre[j0];
      p[j0] = p[j1], j0 = j1;
  REP(j,1,m) if (p[j]) ans[p[j] - 1] = j - 1;
  return {-v[0], ans}; // min cost
```

6.3 DFS algorithms

SCC.Ł

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

Usage: $scc(graph, [\&](VI\& v) \{ \dots \})$ visits all components in reverse topological order. comp[i] holds the component index of a node (a component only has edges to components with lower index). ncomps will contain the number of components.

BiconnectedComponents 2sat EulerWalk EdgeColoring

num.assign(SZ(ed), 0);

```
if (low == val[j]) {
    do √
     x = z.back(); z.pop back();
      comp[x] = ncomps;
      cont.push back(x);
    } while (x != j);
   f(cont); cont.clear();
   ncomps++;}
  return val[j] = low;}
template<class G, class F> void scc(G& g, F f) {
 int n = SZ(q);
 val.assign(n, 0); comp.assign(n, -1);
 Time = ncomps = 0;
 REP(i,0,n) if (comp[i] < 0) dfs(i, q, f);}
```

BiconnectedComponents.h

void bicomps(F f) {

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

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

69fe6f, 33 lines

```
VI num, st;
vector<vector<PII>>> ed:
int Time;
template<class F>
int dfs(int at, int par, F& f) {
  int me = num[at] = ++Time, e, y, top = me;
  for (auto pa : ed[at]) if (pa.second != par) {
    tie(y, e) = pa;
    if (num[y]) {
      top = min(top, num[y]);
      if (num[y] < me)
        st.push back(e);
    } else {
      int si = SZ(st);
      int up = dfs(y, e, f);
      top = min(top, up);
      if (up == me) {
        st.push back(e);
        f(VI(st.begin() + si, st.end()));
        st.resize(si);
      else if (up < me) st.push back(e);</pre>
      else { /* e is a bridge */ }
  return top;
template<class F>
```

```
REP(i,0,SZ(ed)) if (!num[i]) dfs(i, -1, f);
2sat.h
Description: Calculates a valid assignment to boolean variables a,
b, c,... to a 2-SAT problem, so that an expression of the type
(a|||b)\&\&(!a|||c)\&\&(d|||!b)\&\&... becomes true, or reports that it is unsat-
isfiable. Negated variables are represented by bit-inversions (\simx).
Usage: TwoSat ts(number of boolean variables);
ts.either(0, \sim3); // Var 0 is true or var 3 is false
ts.setValue(2); // Var 2 is true
ts.atMostOne({0,\sim1,2}); // <= 1 of vars 0, \sim1 and 2
are true
ts.solve(); // Returns true iff it is solvable
ts.values[0..N-1] holds the assigned values to the vars
Time: \mathcal{O}(N+E), where N is the number of boolean variables, and E is the
number of clauses.
struct TwoSat {
```

```
int N;
vector<VI> gr;
VI values; // 0 = false, 1 = true
TwoSat(int n = 0) : N(n), gr(2*n) {}
int addVar() { // (optional)
  gr.emplace back();
  gr.emplace back();
  return N++:
void either(int f, int j) {
  f = max(2*f, -1-2*f);
  j = max(2*j, -1-2*j);
  gr[f].push back(j^1);
  gr[j].push back(f^1);
void setValue(int x) { either(x, x); }
void atMostOne(const VI& li) { // (optional)
  if (SZ(li) <= 1) return;</pre>
  int cur = \simli[0];
  REP(i,2,SZ(li)) {
    int next = addVar();
    either(cur, ~li[i]);
    either(cur, next);
    either(~li[i], next);
    cur = \sim next;
  either(cur, ~li[1]);
VI val, comp, z; int time = 0;
int dfs(int i) {
  int low = val[i] = ++time, x; z.push_back(i);
  for(int e : gr[i]) if (!comp[e])
    low = min(low, val[e] ?: dfs(e));
```

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

EulerWalk.h

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

```
VI eulerWalk(vector<vector<PII>>& gr, int nedges, int
    src=0) {
  int n = SZ(ar):
  VI D(n), its(n), eu(nedges), ret, s = {src};
  D[src]++; // to allow Euler paths, not just cycles
  while (!s.empty()) {
    int x = s.back(), y, e, &it = its[x], end = SZ(gr[x
    if (it == end){ ret.push back(x); s.pop back();
        continue: }
    tie(y, e) = qr[x][it++];
    if (!eu[e]) {
      D[x] -- , D[y] ++;
      eu[e] = 1; s.push back(y);
  for (int x : D) if (x < 0 \mid \mid SZ(ret) != nedges+1)
      return {};
```

6.4 Coloring

return {ret.rbegin(), ret.rend()};

EdgeColoring.h

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

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

```
VI edgeColoring(int N, vector<PII> eds) {
 VI cc(N + 1), ret(SZ(eds)), fan(N), free(N), loc;
  for (PII e : eds) ++cc[e.first], ++cc[e.second];
```

BinaryLifting LCA CompressTree HLD LinkCutTree

```
int u, v, ncols = *max element(ALL(cc)) + 1;
vector<VI> adj(N, VI(ncols, -1));
for (PII e : eds) {
  tie(u, v) = e;
  fan[0] = v;
  loc.assign(ncols, 0);
  int at = u, end = u, d, c = free[u], ind = 0, i =
  while (d = free[v], !loc[d] && (v = adj[u][d]) !=
    loc[d] = ++ind, cc[ind] = d, fan[ind] = v;
  cc[loc[d]] = c;
  for (int cd = d; at != -1; cd ^= c ^ d, at = adj[at
     ][cd])
    swap(adj[at][cd], adj[end = at][cd ^ c ^ d]);
  while (adj[fan[i]][d] != -1) {
    int left = fan[i], right = fan[++i], e = cc[i];
    adj[u][e] = left;
    adj[left][e] = u;
    adj[right][e] = -1;
    free[right] = e;
  adi[u][d] = fan[i];
  adj[fan[i]][d] = u;
  for (int y : {fan[0], u, end})
   for (int& z = free[y] = 0; adj[y][z] != -1; z++);
REP(i,0,SZ(eds))
 for (tie(u, v) = eds[i]; adj[u][ret[i]] != v;) ++
      ret[i]:
return ret;
```

6.5 Trees

BinaryLifting.h

Description: Assumes the root node points to itself.

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

c4e44c, 19 lines

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

```
if (c != d) a = c, b = d;}
  return tbl[0][a];}
LCA.h
Description: Data structure for computing lowest common ancestors in a
tree (with 0 as root). C should be an adjacency list of the tree, either directed
or undirected.
Time: \mathcal{O}(N \log N + Q)
"../data-structures/RMQ.h"
                                                     cbd116, 16 lines
struct LCA {
 int T = 0;
 VI time, path, ret;
  RMQ<int> rmg;
 LCA(vector < VI > \& C) : time(SZ(C)), rmq((dfs(C,0,-1),
       ret)) {}
  void dfs(vector<VI>& C, int v, int par) {
    time[v] = T++;
    for (int y : C[v]) if (y != par) {
      path.push back(v), ret.push back(time[v]);
      dfs(C, y, v);}}
 int lca(int a, int b) {
    if (a == b) return a;
    tie(a, b) = minmax(time[a], time[b]);
    return path[rmq.query(a, b)];}
 //dist(a,b) {return depth[a] + depth[b] - 2*depth[lca(
      a.b)1:}
};
```

CompressTree.h

Description: Assumes the root node points to itself. **Time:** construction $\mathcal{O}(N \log N)$, queries $\mathcal{O}(\log N)$

0292cd, 13 lines

```
bool cmp(int u,int v){return arr[u]<arr[v];}</pre>
int create tree(){//return root of tree
set<int> S;//get distinct nodes
REP(i,k)S.insert(Q[i]);k=0;for(auto it : S)Q[k++]=it;
sort(Q,Q+k,cmp);int kk = k;//distinct initial nodes
//add lca of adjacent pairs
for(int i=0; i< kk-1; i++) \{ int x = lca(0[i], 0[i+1]) \}
if(S.count(x))continue;Q[k++]=x;S.insert(x);
}sort(Q,Q+k,cmp);stack<int> s;s.push(Q[0]);
for(int i=1;i<k;i++){</pre>
while(!anc(s.top(),Q[i]))s.pop();
tree[s.top()].PB(Q[i]);tree[Q[i]].PB(s.top());
s.push(Q[i]);}return Q[0];}
```

HLD.h

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

```
Time: \mathcal{O}\left((\log N)^2\right)
                                                                             c4840c, 21 lines
VI sz, sc, hd, en, ex, par, dep;
```

```
seg tree lazy<node, update> st(1, \{0, 0\}, \{0, 0\}); //
    for alter accordingly
int timer = -1:
void hld(int u, int p, int ch, int d) {
    hd[u] = ch; en[u] = ++timer; par[u] = p; dep[u] = d
    if (sc[u] != -1) hld(sc[u], u, ch, d + 1);
    for (auto e : q[u]) {
        int v = U[e] ^ V[e] ^ u;
        if (v == p || v == sc[u]) continue;
        hld(v, u, v, d + 1);
    ex[u] = timer;}
int path(int x, int y) {
    int ma = (int) - 1e9;
    while (hd[x] != hd[y]) {
        if (dep[hd[x]] < dep[hd[y]]) swap(x, y);
        ma = max(st.query(en[hd[x]], en[x]).sum, ma);
            // for hd[x] \rightarrow x
        x = par[hd[x]];
    if (dep[x] < dep[y]) swap(x, y);
    ma = max(ma, st.query(en[y], en[x]).sum); // for y
        -> X
    return ma:}
```

LinkCutTree.h

Description: Represents a forest of unrooted trees. You can add and remove edges (as long as the result is still a forest), and check whether two nodes are in the same tree.

```
Time: All operations take amortized \mathcal{O}(\log N).
```

```
struct Node { // Splay tree. Root's pp contains tree's
    parent.
 Node *p = 0, *pp = 0, *c[2];
 bool flip = 0;
 Node() { c[0] = c[1] = 0; fix(); }
  void fix() {
   if (c[0]) c[0]->p = this;
   if (c[1]) c[1] -> p = this;
   // (+ update sum of subtree elements etc. if wanted
 void pushFlip() {
   if (!flip) return;
   flip = 0; swap(c[0], c[1]);
   if (c[0]) c[0]->flip ^= 1;
   if (c[1]) c[1]->flip ^= 1;}
  int up() { return p ? p->c[1] == this : -1; }
  void rot(int i, int b) {
   int h = i ^ b;
   Node *x = c[i], *y = b == 2 ? x : x->c[h], *z = b ?
   if ((y->p = p)) p->c[up()] = y;
   c[i] = z - > c[i ^ 1]:
    if (b < 2) {
      x - c[h] = y - c[h ^ 1];
      z \rightarrow c[h \land 1] = b ? x : this;
   v - c[i ^1] = b ? this : x;
```

DirectedMST CentriodDecomposition

```
fix(); x->fix(); y->fix();
    if (p) p->fix();
    swap(pp, y->pp);}
 void splay() {
    for (pushFlip(); p; ) {
      if (p->p) p->p->pushFlip();
      p->pushFlip(); pushFlip();
      int c1 = up(), c2 = p->up();
      if (c2 == -1) p->rot(c1, 2);
      else p->p->rot(c2, c1 != c2);}}
 Node* first() {
    pushFlip();
    return c[0] ? c[0]->first() : (splay(), this);}};
struct LinkCut {
 vector<Node> node;
 LinkCut(int N) : node(N) {}
 void link(int u, int v) { // add an edge (u, v)
    assert(!connected(u, v));
    makeRoot(&node[u]);
    node[u].pp = &node[v];}
 void cut(int u, int v) { // remove an edge (u, v)
    Node *x = &node[u], *top = &node[v];
    makeRoot(top); x->splay();
    assert(top == (x-pp ?: x-c[0]));
    if (x->pp) x->pp = 0;
   else {
      x \rightarrow c[0] = top \rightarrow p = 0;
      x->fix();}}
 bool connected(int u, int v) { // are u, v in the
      same tree?
    Node* nu = access(&node[u])->first();
    return nu == access(&node[v])->first();}
 void makeRoot(Node* u) {
    access(u);
    u->splay();
    if(u->c[0]) {
      u - c[0] - p = 0;
      u - c[0] - flip ^= 1;
      u - c[0] - pp = u;
      u - c[0] = 0;
      u->fix();}}
 Node* access(Node* u) {
    u->splay();
    while (Node* pp = u->pp) {
      pp->splay(); u->pp = 0;
      if (pp->c[1]) {
        pp - c[1] - p = 0; pp - c[1] - pp = pp; 
      pp->c[1] = u; pp->fix(); u = pp;
    return u;}};
```

DirectedMST.h

Description: Finds a minimum spanning tree/arborescence of a directed graph, given a root node. If no MST exists, returns -1.

```
\begin{tabular}{lll} \textbf{Time: } \mathcal{O}\left(E\log V\right) \\ & & & & & & & & & \\ & & & & & & & \\ & & & & & & & \\ & & & & & & & \\ & & & & & & \\ & & & & & & \\ & & & & & & \\ & & & & & & \\ & & & & & & \\ & & & & & \\ & & & & & \\ & & & & & \\ & & & & & \\ & & & & & \\ & & & & & \\ & & & & & \\ & & & & & \\ & & & & & \\ & & & & & \\ & & & & & \\ & & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & & \\ & & & \\ & & & & \\ & & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & & \\ & & \\ & & & \\ & & \\ & & & \\ & & \\ & & \\ & & &
```

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

```
return {res, par};
```

CentriodDecomposition.h

Description: Assumes the root node points to itself. **Time:** construction $\mathcal{O}(N \log N)$, queries $\mathcal{O}(\log N)$

f8ca72, 14 lines

```
VI U, V, W, isDel;
int dp[n][log2(n) + 1];
// (avoid deleted edges) in all 3 DFS
void decompose(int root, int p) {
   dfs sz(root, -1); // calc sizes of subtrees
    int c = get centroid(root, -1, sz[root]); // if sz
        [v] * 2 > sz[root] return get centroid(v) else
        return u
   if (p == -1) p = root:
    // Add edge btwn p and c here
   dfs(c); // to compute functions
    for (auto e : q[root]) {
        if (isDel[e]) continue;
        isDel[e] = 1;
        int v = U[e] ^ V[e] ^ u;
        decompose(v, root);}}
```

6.6 Math

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

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

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

Mirsky's Theorem Max length chain is equal to min partitioning into antichains. Max chain is height of poset.

Dilworth's Theorem Min partition into chains is equal to max length antichain. From poset create bipartite graph. Any edge from v_i - v_j implies LV_i - RV_j . Let A be the set of vertices such that neither LV_i nor RV_i are in vertex cover. A is an antichain of size n-max matching. To get min partition into chains, take a vertex from left side, keep taking vertices till a matching exist. Consider this as a chain. Its size is n-max matching.

Geometry (7)

7.1 Geometric primitives

Point.h

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

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

lineDistance.h

Description:

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



SegmentDistance.h

Description:

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

```
 \begin{array}{lll} \textbf{Usage:} \; \text{Point} < \text{double} > \; \text{a, b(2,2), p(1,1);} \\ \text{bool onSegment = segDist(a,b,p)} \; & \; \text{1e-10;} \\ \end{array}
```

SegmentIntersection.h

Description:

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

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

```
"Point.h", "OnSegment.h"

template<class P> vector<P> segInter(P a, P b, P c, P d
    ) {
    auto oa = c.cross(d, a), ob = c.cross(d, b),
        oc = a.cross(b, c), od = a.cross(b, d);
    // Checks if intersection is single non-endpoint
        point.

if (sgn(oa) * sgn(ob) < 0 && sgn(oc) * sgn(od) < 0)
        return {(a * ob - b * oa) / (ob - oa)};
    set<P> s;
    if (onSegment(c, d, a)) s.insert(a);
    if (onSegment(a, b, c)) s.insert(c);
    if (onSegment(a, b, d)) s.insert(d);
    return {ALL(s)};
}
```

lineIntersection.h

Description:

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



sideOf.h

"Point.h"

res

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

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

```
template<class P>
int sideOf(P s, P e, P p) { return sgn(s.cross(e, p));
}
```

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

OnSegment.h

"Point.h"

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

```
template<class P> bool onSegment(P s, P e, P p) {
   return p.cross(s, e) == 0 && (s - p).dot(e - p) <= 0;
}</pre>
```

linearTransformation.h Description:

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

Angle.h

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

```
Usage:
                        vector<Angle> v = \{w[0], w[0].t360() ...\}; //
sorted
int j = 0; REP(i,0,n) { while (v[j] < v[i].t180())
++j; }
// sweeps j such that (j-i) represents the number of
positively oriented triangles with vertices at \theta and i of0602,35 lines
struct Angle {
     int x, y;
     int t;
     Angle(int x, int y, int t=0) : x(x), y(y), t(t) {}
     Angle operator-(Angle b) const { return {x-b.x, y-b.y}
                , t}; }
     int half() const {
          assert(x || v);
          return y < 0 \mid | (y == 0 \&\& x < 0);
     Angle t90() const { return \{-y, x, t + (half() \&\& x\}\}
    Angle t180() const { return {-x, -y, t + half()}; }
     Angle t360() const { return \{x, y, t + 1\}; \}
bool operator<(Angle a, Angle b) {</pre>
    // add a.dist2() and b.dist2() to also compare
                distances
     return make tuple(a.t, a.half(), a.y * (ll)b.x) <</pre>
                      make tuple(b.t, b.half(), a.x * (ll)b.y);
// Given two points, this calculates the smallest angle
// them, i.e., the angle that covers the defined line
          segment.
pair<Angle, Angle > segmentAngles(Angle a, Angle b) {
     if (b < a) swap(a, b);
     return (b < a.t180() ?
                         make pair(a, b) : make pair(b, a.t360()));
Angle operator+(Angle a, Angle b) { // point a + vector
     Angle r(a.x + b.x, a.y + b.y, a.t);
     if (a.t180() < r) r.t--;
     return r.t180() < a ? r.t360() : r;
Angle angleDiff(Angle a, Angle b) { // angle b - angle
     int tu = b.t - a.t; a.t = b.t;
     return \{a.x*b.x + a.y*b.y, a.x*b.y - a.y*b.x, tu - (b.x., tu -
                  < a);
```

7.2 Circles

CircleIntersection.h

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

```
84d6d3, 11 lines
typedef Point<double> P;
bool circleInter(P a,P b,double r1,double r2,pair<P, P</pre>
    >* out) {
  if (a == b) { assert(r1 != r2); return false; }
  P \text{ vec} = b - a;
  double d2 = vec.dist2(), sum = r1+r2, dif = r1-r2,
          p = (d2 + r1*r1 - r2*r2)/(d2*2), h2 = r1*r1 -
               p*p*d2;
  if (sum*sum < d2 || dif*dif > d2) return false;
  P \text{ mid} = a + \text{vec*p}, \text{ per} = \text{vec.perp()} * \text{sqrt(fmax(0, h2))}
       ) / d2);
  *out = {mid + per, mid - per};
  return true;
```

CircleTangents.h

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

```
"Point.h"
                                                 b0153d, 13 lines
template<class P>
vector<pair<P, P>> tangents(P c1, double r1, P c2,
    double r2) {
 P d = c2 - c1;
 double dr = r1 - r2, d2 = d.dist2(), h2 = d2 - dr *
 if (d2 == 0 || h2 < 0) return {};
 vector<pair<P, P>> out;
 for (double sign : {-1, 1}) {
   P v = (d * dr + d.perp() * sqrt(h2) * sign) / d2;
   out.push back(\{c1 + v * r1, c2 + v * r2\});
 if (h2 == 0) out.pop back();
 return out;
```

CircleLine.h

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

```
"Point.h"
                                                                                     eea4d5, 10 lines
```

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

```
if (h2 < 0) return {};
if (h2 == 0) return {p};
P h = ab.unit() * sqrt(h2);
return \{p - h, p + h\};
```

CirclePolygonIntersection.h

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

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

```
"../../content/geometry/Point.h"
                                                                                   f5c096, 19 lines
```

```
typedef Point<double> P;
#define arg(p, q) atan2(p.cross(q), p.dot(q))
double circlePoly(P c, double r, vector<P> ps) {
 auto tri = [&](P p, P q) {
   auto r2 = r * r / 2;
   Pd = q - p;
   auto a = d.dot(p)/d.dist2(), b = (p.dist2()-r*r)/d.
   auto det = a * a - b;
   if (det \le 0) return arg(p, q) * r2;
   auto s = max(0., -a-sqrt(det)), t = min(1., -a+sqrt)
        (det));
   if (t < 0 | | 1 \le s) return arg(p, q) * r2;
   Pu = p + d * s, v = p + d * t;
    return arg(p,u) * r2 + u.cross(v)/2 + arg(v,q) * r2
  auto sum = 0.0;
  REP(i,0,SZ(ps))
   sum += tri(ps[i] - c, ps[(i + 1) % SZ(ps)] - c);
  return sum;
```

circumcircle.h

Description:

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



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

MinimumEnclosingCircle.h

shuffle(ALL(ps), mt19937(time(0)));

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

```
<u>"circumcircle.h"</u>
pair<P, double> mec(vector<P> ps) {
```

```
P o = ps[0];
double r = 0, EPS = 1 + 1e-8;
REP(i,0,SZ(ps)) if ((o - ps[i]).dist() > r * EPS) {
  o = ps[i], r = 0;
  REP(j,0,i) if ((o - ps[j]).dist() > r * EPS) {
     o = (ps[i] + ps[j]) / 2;
     r = (o - ps[i]).dist();
  REP(k,0,j) if ((o - ps[k]).dist() > r * EPS) {
     o = ccCenter(ps[i], ps[j], ps[k]);
     r = (o - ps[i]).dist();
}}return {0, r};
```

7.3 Polygons

InsidePolygon.h

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

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

"Point.h", "OnSegment.h", "SegmentDistance.h" 2261c4,1

PolygonArea.h

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

PolygonCenter.h

Description: Returns the center of mass for a polygon.

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

```
return res / A / 3;
PolygonCut.h
Description:
Returns a vector with the vertices of a polygon with every-
thing to the left of the line going from s to e cut away.
Usage: vector<P> p = ...;
p = polygonCut(p, P(0,0), P(1,0));
"Point.h", "lineIntersection.h"
typedef Point<double> P;
vector<P> polygonCut(const vector<P>& poly, P s, P e) {
  if (SZ(poly) <= 2) return {};</pre>
  vector<P> res:
  REP(i,0,SZ(poly)) {
    P cur = poly[i], prev = i ? poly[i-1] : poly.back()
    if (zero(s.cross(e, cur))) {
       res.push back(cur):
       continue;
    bool side = s.cross(e, cur) < 0;
    if (side != (s.cross(e, prev) < 0))
       res.push back(lineInter(s, e, cur, prev).second);
    if (side)
       res.push back(cur);
  return res;
ConvexHull.h
Description:
Returns a vector of the points of the convex hull in counter-
clockwise order. Points on the edge of the hull between two
other points are not considered part of the hull.
Time: \mathcal{O}(n \log n)
"Point.h"
                                                       c5c490, 13 lines
typedef Point<ll> P;
vector<P> convexHull(vector<P> pts) {
  if (SZ(pts) <= 1) return pts;</pre>
```

HullDiameter.h

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

```
"Point.h" 261063, 12 lines
```

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

PointInsideHull.h

typedef Point<ll> P;

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

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

```
"Point.h", "sideOf.h", "OnSegment.h" efb6da, 14 line
```

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

LineHullIntersection.h

return sgn(l[a].cross(l[b], p)) < r;</pre>

int n = SZ(poly), lo = 0, hi = n;

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

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

ClosestPair PolyhedronVolume Point3D 3dHull

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

7.4 Misc. Point Set **Problems**

ClosestPair.h

Description: Finds the closest pair of points.

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

```
"Point.h"
                                                   ac393c, 17 lines
typedef Point<ll> P;
pair<P, P> closest(vector<P> v) {
  assert(SZ(v) > 1);
  set<P> S:
  sort(ALL(v), [](P a, P b) { return a.v < b.v; });
  pair<ll, pair<P, P>> ret{LLONG MAX, {P(), P()}};
  int j = 0;
  for (P p : v) {
    P d\{1 + (ll) sqrt(ret.first), 0\};
    while (v[j].y \le p.y - d.x) S.erase(v[j++]);
```

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

7.5 3D

PolyhedronVolume.h

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

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

Point3D.h

Description: Class to handle points in 3D space. T can be e.g. double or long long.

```
template<class T> struct Point3D {
 typedef Point3D P:
 typedef const P& R;
 T x. v. z:
 explicit Point3D(T x=0, T y=0, T z=0) : x(x), y(y), z
 bool operator<(R p) const {</pre>
    return tie(x, y, z) < tie(p.x, p.y, p.z); }</pre>
 bool operator==(R p) const {
    return tie(x, y, z) == tie(p.x, p.y, p.z); }
 P operator+(R p) const { return P(x+p.x, y+p.y, z+p.z
```

```
); }
P operator-(R p) const { return P(x-p.x, y-p.y, z-p.z
P operator*(T d) const { return P(x*d, y*d, z*d); }
P operator/(T d) const { return P(x/d, y/d, z/d); }
T dot(R p) const { return x*p.x + y*p.y + z*p.z; }
P cross(R p) const {
  return P(y*p.z - z*p.y, z*p.x - x*p.z, x*p.y - y*p.
      x);
T dist2() const { return x*x + y*y + z*z; }
double dist() const { return sqrt((double)dist2()); }
//Azimuthal angle (longitude) to x-axis in interval
    [-pi, pi]
double phi() const { return atan2(y, x); }
//Zenith angle (latitude) to the z-axis in interval
double theta() const { return atan2(sqrt(x*x+y*y),z);
```

P unit() const { return *this/(T)dist(); } //makes

dist()=1

```
//returns unit vector normal to *this and p
  P normal(P p) const { return cross(p).unit(); }
  //returns point rotated 'angle' radians ccw around
      axis
  P rotate(double angle, P axis) const {
    double s = sin(angle), c = cos(angle); Pu = axis.
    return u*dot(u)*(1-c) + (*this)*c - cross(u)*s;
};
```

3dHull.h

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

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

0754b0 49 lines

```
"Point3D.h"
typedef Point3D<double> P3;
struct PR {
 void ins(int x) { (a == -1 ? a : b) = x; }
  void rem(int x) { (a == x ? a : b) = -1; }
  int cnt() { return (a != -1) + (b != -1); }
 int a, b;
struct F { P3 q; int a, b, c; };
vector<F> hull3d(const vector<P3>& A) {
  assert(SZ(A) >= 4):
  vector<vector<PR>> E(SZ(A), vector<PR>(SZ(A), {-1, -1
#define E(x,y) E[f.x][f.y]
  vector<F> FS;
  auto mf = [\&](int i, int j, int k, int l) {
    P3 q = (A[i] - A[i]).cross((A[k] - A[i]));
   if (q.dot(A[l]) > q.dot(A[i]))
      q = q * -1;
    F f{q, i, j, k};
    E(a,b).ins(k); E(a,c).ins(j); E(b,c).ins(i);
    FS.push back(f);
  REP(i,0,4) REP(j,i+1,4) REP(k,j+1,4)
   mf(i, j, k, 6 - i - j - k);
  REP(i,4,SZ(A)) {
   REP(j,0,SZ(FS)) {
      F f = FS[i];
      if(f.q.dot(A[i]) > f.q.dot(A[f.a])) {
        E(a,b).rem(f.c);
        E(a,c).rem(f.b);
        E(b,c).rem(f.a);
        swap(FS[i--], FS.back());
        FS.pop back();
   int nw = SZ(FS);
   REP(j,0,nw) {
```

Mathematics (8)

Equations

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

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

$$ax + by = e cx + dy = f \Rightarrow x = \frac{ed - bf}{ad - bc} y = \frac{af - ec}{ad - bc}$$

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

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

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

Recurrences

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

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

Non-distinct roots r become polynomial factors, e.g. $a_n = (d_1 n + d_2)r^n.$

8.3 Trigonometry

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

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

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

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

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

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

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

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

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

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

8.4 Geometry

8.4.1 Triangles

Side lengths: a, b, c

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

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

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

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

Length of median (divides triangle into two equal-area

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

Length of bisector (divides angles in two):

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

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

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

8.4.2 Quadrilaterals

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

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

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

8.5 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:

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

8.6 Sums

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

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

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

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

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

8.7 Series

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

TernarySearch IntervalCover ConstantIntervals

$\sin x = x - \frac{x^3}{3!} + \frac{x^5}{5!} - \frac{x^7}{7!} + \dots, (-\infty < x < \infty)$ $\cos x = 1 - \frac{x^2}{2!} + \frac{x^4}{4!} - \frac{x^6}{6!} + \dots, (-\infty < x < \infty)$

8.8 Probability theory

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

Expectation is linear:

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

For independent X and Y,

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

8.8.1 Discrete distributions Binomial distribution

First success distribution

The number of trials needed to get the first success in independent yes/no experiments, each wich yields success with probability p is Fs(p), $0 \le p \le 1$.

$$p(k) = p(1-p)^{k-1}, k = 1, 2, \dots$$

$$\mu = \frac{1}{n}, \sigma^2 = \frac{1-p}{n^2}$$

Poisson distribution

The number of events occurring in a fixed period of time t if these events occur with a known average rate κ and independently of the time since the last event is $Po(\lambda)$, $\lambda = t\kappa$.

$$p(k) = e^{-\lambda} \frac{\lambda^k}{k!}, k = 0, 1, 2, \dots$$
$$\mu = \lambda, \sigma^2 = \lambda$$

8.8.2 Continuous distributions Uniform distribution

Exponential distribution

The time between events in a Poisson process is $\operatorname{Exp}(\lambda), \ \lambda > 0.$

$$f(x) = \begin{cases} \lambda e^{-\lambda x} & x \ge 0\\ 0 & x < 0 \end{cases}$$
$$\mu = \frac{1}{\lambda}, \, \sigma^2 = \frac{1}{\lambda^2}$$

Normal distribution

Most real random values with mean μ and variance σ^2 are well described by $\mathcal{N}(\mu, \sigma^2)$, $\sigma > 0$.

$$f(x) = \frac{1}{\sqrt{2\pi\sigma^2}} e^{-\frac{(x-\mu)^2}{2\sigma^2}}$$

If $X_1 \sim \mathcal{N}(\mu_1, \sigma_1^2)$ and $X_2 \sim \mathcal{N}(\mu_2, \sigma_2^2)$ then

$$aX_1 + bX_2 + c \sim \mathcal{N}(\mu_1 + \mu_2 + c, a^2\sigma_1^2 + b^2\sigma_2^2)$$

Miscellaneous (9)

9.1 RNG, Intervals, T.S

TernarySearch.h

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

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

Time: $\mathcal{O}\left(\log(b-a)
ight)$

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

IntervalCover.h

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

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

ConstantIntervals.h

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

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

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

FastKnapsack.h

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

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

IntervalContainer.h

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

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

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

RNGs.h

SEED = chrono::steady clock::now().time since epoch(). count(); // or use 'high resolution clock' random device rd; auto SEED = rd();

```
mt19937 rng(SEED);
uniform int distribution<> dis(MIN, MAX); // usage: dis
// others: uniform real distribution,
```

DebuggingTricks.cpp

```
Description: Debug
```

Time: $\mathcal{O}\left(k\log\frac{n}{t}\right)$

```
1. signal(SIGSEGV, [](int) { Exit(0); });
converts segfaults into Wrong Answers. Similarly one
    can catch SIGABRT (assertion failures) and SIGFPE (
    zero divisions). GLIBCXX DEBUG failures generate.
    SIGABRT (or SIGSEGV on gcc 5.4.0 apparently).
2. feenableexcept(29);
```

kills the program on NaNs(1), 0-divs (4), infinities (8) **and** denormals (16).

9.2 **Optimization tricks**

builtin ia32 ldmxcsr(40896); disables denormals (which make floats 20x slower near their minimum value).

9.2.1 Bit hacks

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

9.2.2 Pragmas

- #pragma GCC optimize ("Ofast") will make GCC auto-vectorize loops and optimizes floating points better.
- #pragma GCC target ("avx2") can double performance of vectorized code, but causes crashes on old machines.
- #pragma GCC optimize ("trapv") kills the program on integer overflows (but is really slow).
- **#pragma** GCC optimize("unroll-loops")
- target("sse,sse2,sse3,sse4,popcnt,abm,mmx, avx")

FastMod.h

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

```
typedef unsigned long long ull;
struct FastMod {
  ull b, m;
  FastMod(ull b) : b(b), m(-1ULL / b) {}
  ull reduce(ull a) \{ // a \% b + (0 \text{ or } b) \}
    return a - (ull)(( uint128 t(m) * a) >> 64) * b;
};
```

FastInput.h

Description: Usage requires your program to pipe in input from file.

Usage: ./a.out < input.txt</pre>

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

```
7b3c70, 16 lines
inline char qc() { // like getchar()
  static char buf[1 << 16];</pre>
  static size t bc, be;
  if (bc >= be) {
    buf[0] = 0, bc = 0;
    be = fread(buf, 1, sizeof(buf), stdin);
  return buf[bc++]; // returns 0 on EOF
int readInt() {
  int a, c;
  while ((a = qc()) < 40);
  if (a == '-') return -readInt();
  while ((c = gc()) >= 48) a = a * 10 + c - 480;
  return a - 48:
```

Contest (10)

template.cpp

```
// #pragma GCC optimize("03,unroll-loops")
// #pragma GCC target("avx2,bmi,bmi2,lzcnt,popcnt")
#include <bits/stdc++.h>
using namespace std;
#include <ext/pb ds/assoc container.hpp>
#include <ext/pb ds/tree policy.hpp>
template <class T>
using o set = tree<T, null type, less<T>, rb tree tag,
   tree order statistics node update>;
// order of key (val): no. of values less than val
// find by order (k): kth largest element.(0-based)
#define int long long
#define FOR(i, a, b) for (int i = (a); i < (b); ++i)
#define REP(i, a, b) for (int i = (a); i < (b); ++i)
#define ALL(x) begin(x), end(x)
#define SZ(x) ((int)(x).size())
#define SET(a, v) memset((a), (v), sizeof(a))
```

```
#define PB push back
#define EB emplace back
#define MP make pair
#define F first
#define S second
using LL = long long;
using dbl = double;
using II = pair<int, int>;
using VI = vector<int>;
using VII = vector<II>;
using VVI = vector<VI>;
#define endl "\n"
const long long mod = 1e9 + 7;
signed main() {
 // freopen("sample.in", "r", stdin);
// freopen("sample.out", "w", stdout);
  cin.tie(0)->sync with stdio(0);
  cin.exceptions(cin.failbit); // RTE if out of bound
  return 0;
```