Team notebook

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1.7.1 bit formula 8	1.16.2 Tonelli-Shanks(for solving quadratic	<pre>/// Original Recurrence: dp[i][j] = min(dp[i-1][k] +</pre>
1.8 bitset 8	congruence equation) 19	C[k][j]) for $k < j$ (*** $C[k+1][j]$ instead of $C[k][j]$ is also possible)
1.9 combinatorics 8	1.16.3 bigmod 20	/// Sufficient condition: A[i][j] <= A[i][j+1] where
1.9.1 combinatorics 8	1.16.4 factorial factorize 20	A[i][j] = smallest k that gives optimal answer
1.10 flow	1.16.5 mod inverse from 1 to N 20	/// 1 <= i <= n, 1 <= j <= m
1.10.1 Dinic with and without scaling 8	1.16.6 pollard rho 20	/// Original complexity: O(n*m^2), Optimized
1.10.2 Hopcroft karp 9	1.16.7 sieve	<pre>complexity: O(n*m*lg(m)) /// ** 1-based indexing.</pre>
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1.11 geometry	1.17.1 aho corasick	/// to eliminate possibility of overflow reduce the
1.11.1 comparison of doubles	1.17.2 manachers	value of infLL

```
long long C(int i, int j); /// precompute C function
     beforehand:
11 dp[MAXN][MAXM]; ///See if you can optimize memory
    by declaring dp[2][MAXM];
void compute(int idx, int 1, int r, int opt1, int optr)
 if (1 > r) return:
 int mid = (1 + r) >> 1:
 pair<long long, int> best = {infLL, -1}: /// take
       -infLL for max answer.
 for (int k = optl: k <= min(mid-1, optr): k++) {</pre>
       /// take < for max answer; take C(k+1, mid) if
            required:
       if (best.F > dp[idx-1][k] + C(k, mid) ) best =
            MP(dp[idx-1][k] + C(k, mid), k): /// watch
            out for possible dp[2][MAXM] memory
            optimization!
 dp[idx][mid] = best.first:
 int opt = best.second;
 compute(idx, 1, mid - 1, optl, opt);
 compute(idx, mid + 1, r, opt, optr);
int main() {
 int n, m;
 dp[0][0] = 0;
 for( int i = 1; i <= m; ++i ) dp[0][i] = infLL;</pre>
       ///take -infLL for max answer;
 for( int i = 1; i <= n; ++i ) compute( i, i, m, i-1,</pre>
      m); /// apparently compute(i, 1, m, 0, m)
       also works, i don't know why.
 cout << dp[n][m] << endl:</pre>
```

1.1.2 Dynamic Convex Hull trick

```
//Insertion - O(1gN), query - O(1gN)
//Querying in empty cht container gives runtime error!
//y = mx+b;
const ll IS_QUERY = -(1LL<<62); //-INF
struct Line {
    ll m, b;
    mutable function<const Line*()> succ;
    bool operator<(const Line& rhs) const {
        if (rhs.b != IS_QUERY) return m < rhs.m;
        const Line* s = succ();
        if (!s) return 0;
        ll x = rhs.m;</pre>
```

```
return (long double)b - s->b < (long double)(s->m -
        m) * x; //add precision only if wa;
};
// Keeps upper hull for maximums.
// add lines with -m and -b and return -ans to make
    this code working for minimums.
struct HullDynamic : public multiset<Line> {
 bool bad(iterator v) {
   auto z = next(v):
   if (v == begin()) {
     if (z == end()) return 0:
       return v->m == z->m && v->b <= z->b:
   auto x = prev(v):
   if (z == end()) return v > m == x - m && v - b <= x - b:
   return (long double)(x->b - v->b)*(z->m - v->m) >=
        (long double)(v->b-z->b)*(v->m-x->m):
        //add precision only if wa:
 7
 void insert line(ll m. ll b) {
   auto y = insert({ m, b }); // insert({-m, -b)}
        instead to work for minimums.
   v->succ = [=] { return next(v) == end() ? 0 :
        &*next(v); };
   if (bad(y)) { erase(y); return; }
   while (next(y) != end() && bad(next(y)))
        erase(next(v)):
   while (y != begin() && bad(prev(y))) erase(prev(y));
 11 query(11 x) {
   auto 1 = *lower_bound((Line) { x, IS_QUERY });
   return 1.m * x + 1.b; // return (-1LL)*(1.m * x +
       1.b) instead for getting minimums.
 }
}:
int main() {
 HullDvnamic cht:
 int T:
 cin >> T:
 while(T--) {
   cht.clear():
```

1.1.3 SOS dp

```
///0(N*2^N)
int dp[1<<N];
for(int i = 0; i<(1<<N); ++i) dp[i] = A[i];
for(int i = 0; i < N; ++i) {
```

1.1.4 Semi-offline convex hull trick

```
///Covers all types of conditions :)
struct ConvexHullTrick{
 vector<ll> m. b:
 int ptr = 0;
 bool bad(int 11, int 12, int 13) { ///add precision
      if only WA.
   return (long double)(b[13] - b[11])*(m[11] - m[12])
        \leq (long double)(b[12] - b[11])*(m[11] -
        m[13]); ///(slope dec+query min),(slope
        inc+query max)
   return (long double)(b[13] - b[11])*(m[11] - m[12])
        > (long double)(b[12] - b[11])*(m[11] -
        m[13]); ///(slope dec+query max), (slope
        inc+querv min)
 void insert_line(ll _m, ll _b) {
   m.push back( m):
   b.push back(b):
   int s = m.size();
   while(s >= 3 && bad(s-3, s-2, s-1)) {
     m.erase(m.end()-2):
     b.erase(b.end()-2);
 11 f(int i, ll x) { return m[i]*x + b[i]; }
 ///(slope dec+query min), (slope inc+query max) -> x
      increasing
 ///(slope dec+query max), (slope inc+query min) -> x
      decreasing
 11 query(11 x) {
```

```
3
```

1.2 Data Structures

1.2.1 DSU (path compression + union by size)

```
int par[10005];
int sz[10005];
int Find(int r) { //Path compression
 if( par[r] == r ) return r;
 par[r] = Find(par[r]);
 return par[r];
void Union(int a, int b) { //Union by size of subtrees.
 a = Find(a);
 b = Find(b):
 if (a != b) {
   if (sz[a] < sz[b]) swap(a, b);</pre>
   par[b] = a;
   sz[a] += sz[b]:
int main() {
 for( int i = 0: i < 10005: ++i ) {
   par[i] = i;
   sz[i] = 1:
```

1.2.2 mos

```
//Complexity: O(sqrt(N) * (N+M))
const int mx = 3e5+5;
const int block_sz = 550; // N ~ 3e5
int freq[mx], mo_cnt = 0;
int ret[mx];
inline void add(int idx) {
    ++freq[a[idx]];
    if(freq[a[idx]] == 1) ++mo_cnt;
```

```
inline void del(int idx) {
  --freq[a[idx]];
 if(freq[a[idx]] == 0) --mo_cnt;
inline int get_ans() {
 return mo_cnt;
struct queries {
 int l. r. idx:
  queries() { }
  queries(int _1, int _r, int _i) : 1(_1), r(_r),
       idx( i) { }
  bool operator < (const queries &p) const {</pre>
   if(1/block_sz != p.1/block_sz) return 1 < p.1;</pre>
   return ((1/block sz) & 1) ? r > p.r : r < p.r:
}:
void mo(vector<queries> &q) {
  sort(q.begin(), q.end());
  memset(ret, -1, sizeof ret):
 // 1 = 1, r = 0 if 1-indexed array
  int 1 = 0, r = -1;
  for(auto &gg : g) {
   while(qq.1 < 1) add(--1);</pre>
    while (qq.r > r) add (++r);
    while(qq.1 > 1) del(1++);
    while(qq.r < r) del(r--);</pre>
    ret[qq.idx] = get_ans();
```

1.2.3 sparse table (0-based indexing)

```
//1. range sum query (copied from cp-algorithms)
const int MAXN = 1e5+5:
long long arr[MAXN]:
const int LOG = log2(MAXN) + 1;
long long sp[MAXN][LOG+1]:
int N: //size of arr
void init() { //O(NlogN)
 for( int i = 0; i < N; ++i ) sp[i][0] = arr[i];</pre>
   for( int j = 1; j <= LOG; ++j )</pre>
     for( int i = 0; i + (1 << j) <= N; ++i)
       sp[i][j] = sp[i][j-1] + sp[i + (1 << (j -
            1))][i - 1];
long long query( int L, int R ) { //O(logN)
 long long sum = 0;
 for( int j = LOG; j >= 0; --j ) {
   if((1 << j) <= R - L + 1) {
```

```
sum += sp[L][i];
     L += 1 << i;
 return sum;
//2. rmq (copied from cp-algorithms)
const int MAXN = 1e5+5;
long long arr[MAXN]:
const int LOG = log2(MAXN) + 1;
long long sp[MAXN][LOG+1]:
int prelog[MAXN];
int N: //size of arr
void init_log() { //O(MAXN)
 prelog[1] = 0:
 for( int i = 2: i < MAXN: ++i ) prelog[i] =</pre>
      prelog[i/2] + 1;
void init() { //O(NlogN)
 for( int i = 0; i < N; ++i ) sp[i][0] = arr[i];</pre>
   for( int j = 1; j <= LOG; ++j )</pre>
     for( int i = 0; i + (1 << j) <= N; ++i)
       sp[i][j] = min(sp[i][j-1], sp[i + (1 << (j -
            1))][i - 1]);
long long query(int L, int R) \{ //0(1) \}
 int j = prelog[R - L + 1];
 long long minimum = min(sp[L][j], sp[R - (1 << j) +
       1][i]):
 return minimum;
//3. 2D rmg
const int MAXN = 1005;
const int MAXM = 1005;
long long arr[MAXN][MAXM]:
const int LOGN = log2(MAXN) + 1:
const int LOGM = log2(MAXM) + 1;
long long sp[MAXN][MAXM][LOGN + 1][LOGM + 1];
int prelog[max(MAXN, MAXM)];
int N. M: //size of arr is N*M
void init_log() { //O(MAXN)
 prelog[1] = 0:
 for( int i = 2: i < max(MAXN, MAXM): ++i ) prelog[i]</pre>
       = prelog[i/2] + 1:
void init() { //O(N*M*logN*logM)
 for( int i = 0; i < N; ++i )
   for( int j = 0; j < M; ++j ) sp[i][j][0][0] =</pre>
        arr[i][j];
 for( int k = 1; k <= LOGN; ++k ) {</pre>
   for( int i = 0; i + (1 << k) <= N; ++i ) {</pre>
     for( int j = 0; j < M; ++j ) {</pre>
```

```
4
```

1.3 FFT

1.3.1 2D FFT

Rule of thumb for precision issues- (by Errichto) Well written FFT has an error that depends mainly on the magnitude of numbers in the answer, so it works if the resulting polynomial doesn't have numbers bigger than 1015 for doubles and 1018 for long doubles. You can measure the error by finding max of abs(xllround(x)) over all x in the answer. If it's smaller than 0.2, you're good and rounding will be fine. ///No need to clear anything at the beginning of a test /// 2D FFT of a 2D polynomial ~ O(N*M*(lgN + lgM)) /// product of two 2D polynomials ~ O(N*M*(lgN + lgM)) struct FFT 2D { struct node { double x,v; node() {} node(double a, double b): x(a), y(b) {}

```
node operator + (const node &a) const {return
       node(this->x+a.x,this->y+a.y);}
  node operator - (const node a) const {return
       node(this->x-a.x,this->y-a.y);}
  node operator * (const node a) const {return
       node(this->x*a.x-this->y*a.y,this->x*a.y+a.x*this->y);}
};
int M[2]:
vector<vector<node>> A. B:
vector<node> w[2][2]:
vector<int>rev[2]:
long double pi;
FFT 2D() {
 pi = 3.1415926535897932384;
void init(int n, int m) {
 M[O] = 1, M[1] = 1:
  while(M[0] < n) M[0] <<= 1;</pre>
 M[0] <<= 1:
  while(M[1] < m) M[1] <<= 1:</pre>
 M[1] <<= 1:
  A.resize(M[0], vector<node>(M[1]));
  B.resize(M[0], vector<node>(M[1]));
  for( int z = 0; z < 2; ++z ) {
   w[z][0].resize(M[z]);
   w[z][1].resize(M[z]):
   rev[z].resize(M[z]);
   for (int i=0; i<M[z]; i++) {</pre>
     int j=i,y=0;
     for (int x=1; x<M[z]; x<<=1, j>>=1) (y<<=1)+=j\&1;
     rev[z][i]=v;
   for (int i=0; i<M[z]; i++) {</pre>
     w[z][0][i] = node(
          cos(2*pi*i/M[z]).sin(2*pi*i/M[z])):
     w[z][1][i] = node(
          cos(2*pi*i/M[z]), -sin(2*pi*i/M[z]));
 }
void ftransform 2D( vector<vector<node>> &A. int p ) {
  for( int z = 0: z < M[0]: z++ ) {
   for (int i=0: i<M[1]: i++)</pre>
     if (i<rev[1][i])</pre>
       swap(A[z][i].A[z][rev[1][i]]);
   for (int i=1; i<M[1]; i<<=1)</pre>
     for (int j=0,t=M[1]/(i<<1); j<M[1]; j+=i<<1)</pre>
       for (int k=0,1=0; k<i; k++,1+=t) {
         node x=A[z][i+j+k]*w[1][p][1];
         node v=A[z][j+k];
         A[z][j+k]=y+x;
         A[z][j+k+i]=v-x;
```

```
if (p)
       for (int i=0; i<M[1]; i++)</pre>
         A[z][i].x/=M[1], A[z][i].y/=M[1];
   for( int z = 0; z < M[1]; z++ ) {
     for (int i=0; i<M[0]; i++)</pre>
       if (i<rev[0][i])</pre>
         swap(A[i][z].A[rev[0][i]][z]):
     for (int i=1; i<M[0]; i<<=1)</pre>
       for (int j=0,t=M[0]/(i<<1); j<M[0]; j+=i<<1)</pre>
         for (int k=0,1=0; k<i; k++,1+=t) {</pre>
           node x=w[0][p][1]*A[i+i+k][z]:
           node y=A[j+k][z];
           A[i+k][z]=v+x:
           A[i+k+i][z]=v-x:
     if (p)
       for (int i=0: i<M[0]: i++)</pre>
         A[i][z].x/=M[0]:
  ///2degree polynomial P * 2degree polynomial Q =
       2degree polynomial res
 ///Degree of P is x1,x2; Degree of Q is y1, y2;
      Degree of res is x1+y1, x2+y2;
 void multiply_2D( vector<vector<long long>> &P,
       vector<vector<long long>> &Q, vector<vector<long</pre>
       long>> &res ) {
    init( max(P.size(), Q.size()), max(P[0].size(),
         Q[0].size()));
    for( int i = 0; i < M[0]; i++ )</pre>
     for( int j = 0; j < M[1]; j++ )</pre>
       A[i][j].x = A[i][j].y = B[i][j].x = B[i][j].y =
            0:
    for( int i = 0: i < P.size(): i++ )</pre>
     for( int j = 0; j < P[i].size(); j++ )</pre>
       A[i][i].x = P[i][i]:
    for( int i = 0; i < Q.size(); i++ )</pre>
     for( int j = 0; j < Q[i].size(); j++ )</pre>
       B[i][j].x = Q[i][j];
   ftransform 2D(A, 0):
   ftransform 2D(B, 0):
   for( int i = 0; i < M[0]; i++ )</pre>
     for( int j = 0; j < M[1]; j++ )</pre>
       A[i][i] = A[i][i]*B[i][i]:
   ftransform_2D(A, 1);
   res.resize(M[0], vector<ll>(M[1]));
   for( int i = 0; i < M[0]; i++ )</pre>
     for( int j = 0; j < M[1]; j++ )</pre>
       res[i][j] = round(A[i][j].x);
}fft_2D;
```

```
int main() {
 /// a*b = c
 /// Degree of a = (n1-1, n2-1), Degree of b = (m1-1, n2-1)
      m2-1), Degree of c = (n1+m1-2, n2+m2-2);
 vector<vector<ll>>> a, b, c;
 int n1, n2, m1, m2;
 cin >> n1 >> n2 >> m1 >> m2;
 a.resize(n1, vector<ll>(n2)), b.resize(m1,
      vector<11>(m2)), c.resize(n1+m1-1,
      vector<ll>(n2+m2-1)):
 for( int i = 0; i < n1; ++i )</pre>
   for( int j = 0; j < n2; ++j ) cin >> a[i][j];
 for( int i = 0; i < m1; ++i )</pre>
   for( int j = 0; j < m2; ++j ) cin >> b[i][j];
 fft 2D.multiply(a, b, c): /// product of a and b is
      kept in c;
```

1.3.2 FWHT

```
///O(NlogN)
///Doubles are never used as in FFT. Always long long
    is used(modulo or not).
#define MOD 1000000007
#define OR O
#define AND 1
#define XOR 2
struct FWHT {
 void walsh transform(vector<11> &p. bool inv. int
   int n = p.size(); assert((n&(n-1))==0);
   for (int len=1; 2*len<=n; len <<= 1) {</pre>
     for (int i = 0; i < n; i += len+len){</pre>
       for (int i = 0: i < len: <math>i++) { ///Don't forget
            to include p[i+j] = (p[i+j]+MOD)%MOD,
            p[i+len+i] = (p[i+len+i]+MOD)%MOD at the
            end of this for loop(inside of course) when
            required.
         ll u = p[i+j], v = p[i+len+j];
         if( (flag == XOR) ) p[i+j]=u+v, p[i+len+j]=u-v;
         if( (flag == AND) && !inv ) p[i+j]=v,
              p[i+len+j]=u+v;
         if( (flag == OR) && !inv ) p[i+j]=u+v,
              p[i+len+j]=u;
         if( (flag == AND) && inv ) p[i+j] = -u+v,
              p[i+len+j]=u;
         if( (flag == OR) && inv ) p[i+j]=v,
              p[i+len+j]=u-v;
```

```
if(inv && (flag == XOR)) for(int i=0;i<n;i++)</pre>
        p[i]/=n; ///Don't forget modular inverse n
        when required
 /// For i = 0 to n - 1, j = 0 to n - 1
  /// v[i flag j] += a[i] * b[j]
  vector<ll> convo(vector<ll> &a. vector<ll> &b. int
      flag = XOR) {
   int n = 1, sz = max(a.size(), b.size());
    while(n<sz) n*=2:</pre>
    a.resize(n); b.resize(n); vector<11>res(n, 0);
    walsh_transform(a, 0, flag); walsh_transform(b, 0,
        flag):
   for(int i=0:i<n:i++) res[i] = a[i] * b[i]: ///</pre>
        Don't forget modular multiplication when
        required.
    walsh_transform(res, 1, flag);
   return res:
  ///compute A^k where A?A == A convolution A
  vector<ll> pow(vector<ll> &A, ll k, int flag = XOR) {
   int n = 1, sz = A.size();
   while(n \le z) n *= 2;
   A.resize(n):
   walsh_transform(A, 0, flag);
   for(int i=0; i<n; i++) A[i]=expo(A[i], k, MOD);</pre>
        ///MOD = infLL if there is no modulo
        required.(trix, modern problems require modern
        solutions.)
    walsh_transform(A, 1, flag);
   return A:
}fwht:
int main() {
 vector<ll> vec(6);
  vector<ll> temp = vec;
  vector<ll> store1 = fwht.convo(temp, temp, XOR);
      ///size of store1 will always be a power of 2.
  vector<11> store2 = fwht.pow(temp, 4, XOR); ///size
      of store2 will always be a power of 2.
```

1.3.3 fft by ashigul

```
/**
Rule of thumb for precision issues- (by Errichto)
```

```
Well written FFT has an error that depends mainly on
    the magnitude of numbers in the answer, so it
    works if the resulting polynomial doesn't have
    numbers bigger than 1015 for doubles and 1018 for
    long doubles. You can measure the error by finding
    max of abs(xllround(x)) over all x in the answer.
    If it's smaller than 0.2, you're good and rounding
    will be fine.
///No need to clear anything at the beginning of a test
/// FFT
struct FFT {
 struct node {
   double x.v:
   node() {}
   node(double a, double b): x(a), y(b) {}
   node operator + (const node &a) const {return
        node(this->x+a.x,this->y+a.y);}
   node operator - (const node a) const {return
        node(this->x-a.x,this->y-a.y);}
   node operator * (const node a) const {return
        node(this->x*a.x-this->y*a.y,this->x*a.y+a.x*this->y);
 int M:
 vector<node> A, B, w[2];
 vector<int>rev;
 long double pi;
 FFT() {
   pi = 3.1415926535897932384;
 void init(int n) {
   M = 1:
   while(M < n) M <<= 1:</pre>
   M <<= 1:
   A.resize(M):
   B.resize(M):
   w[0].resize(M):
   w[1].resize(M);
   rev.resize(M):
   for (int i=0: i<M: i++) {</pre>
     int j=i,y=0;
     for (int x=1: x<M: x<<=1, i>>=1) (v<<=1)+=i&1:
   for (int i=0; i<M; i++) {</pre>
     w[0][i] = node(cos(2*pi*i/M),sin(2*pi*i/M));
     w[1][i] = node(cos(2*pi*i/M), -sin(2*pi*i/M));
 void ftransform( vector<node> &A, int p ) {
```

5

```
for (int i=0; i<M; i++)</pre>
      if (i<rev[i])</pre>
       swap(A[i],A[rev[i]]);
    for (int i=1; i<M; i<<=1)</pre>
     for (int j=0,t=M/(i<<1); j<M; j+=i<<1)</pre>
       for (int k=0,1=0; k<i; k++,1+=t) {</pre>
         node x=w[p][1]*A[i+j+k];
         node y=A[j+k];
         A[i+k]=v+x:
         A[j+k+i]=y-x;
    if (p)
      for (int i=0: i<M: i++)</pre>
       A[i].x/=M:
  /// multiply P*Q and keeps the result in res
  ///degree of P is n and degree of Q is m
  ///P, Q is given in standard power form, in increasing
  void multiply( vector<long long> &P. vector<long</pre>
       long> &Q, vector<long long> &res) {
    init( max(P.size(),Q.size()) );
    for( int i = 0; i < M; i++ )</pre>
     A[i].x = A[i].y = B[i].x = B[i].y = 0;
    for( int i = 0; i < P.size(); i++ )</pre>
     A[i].x = P[i]:
    for( int i = 0; i < Q.size(); i++ )</pre>
     B[i].x = Q[i]:
    ftransform(A,0);
    ftransform(B,0);
    for (int k=0; k<M; k++)</pre>
     A[k] = A[k] * B[k];
    ftransform(A,1);
    res.resize(M);
    for( int i = 0; i < M; i++ )</pre>
     res[i] = round(A[i].x):
}fft:
int main() {
 /// a*b = c
 /// Degree of a = n-1, Degree of b = m-1, Degree of c
      = n+m-2:
 vector<ll> a. b. c:
 int n. m:
  cin >> n >> m:
 a.resize(n), b.resize(m), c.resize(n+m-1):
 for( int i = 0; i < n; ++i ) cin >> a[i];
 for( int i = 0; i < m; ++i ) cin >> b[i];
 fft.multiply(a, b, c); /// product of a and b is kept
       in c:
```

1.4 Math

return 1:

1.4.1 Gaussian elimination modulo 2 (32 times faster)

```
// calculates 32 times faster for modulo 2
  n rows/equations, m+1 columns, m variables
  calculates determinant, rank and ans[] -> value for
       variables
  returns {0, 1, INF} -> number of solutions
  gives truely row echelon form. (sort of Identity
       matrixish, but the independent column/variables
       may have non-zero values at the upper rows.)
const double EPS = 1e-9:
#define INF 100000000
#define MAX 105
int Rank: //Rank means number of non-zero rows in row
    echelon form
int Det:
int gauss(vector < bitset<MAX> > a, bitset<MAX> &ans,
    int n. int m) { // number of variables must be
    defined!
 ans.reset(); //reset all bits;
 vector<int> where(m, -1);
 Det = 1. Rank = 0:
 for(int col = 0, row = 0; col < m && row < n; ++col) {</pre>
   int sel = row:
   for(int i = row: i < n: ++i) if(a[i][col]) { sel =</pre>
        i: break: }
   if(!a[sel][col]) { Det = 0; continue; }
   swap(a[sel], a[row]):
   if(row != sel) Det = -Det; ///add mod? (idk, i made
        this up. munim vai kept this line as it is. So
        change only if WA)
   Det &= a[row][col]:
   where[col] = row:
   for(int i = 0: i < n: ++i) if (i != row &&
        a[i][col] > 0) a[i] ^= a[row];
   ++row, ++Rank;
 for(int i = 0; i < m; ++i) ans[i] = (where[i] == -1)</pre>
      ? 0 : a[where[i]][m]:
 for(int i = Rank; i < n; ++i) if(a[i][m]) return 0;</pre>
 for(int i = 0; i < m; ++i) if(where[i] == -1) return</pre>
```

```
}
```

1.4.2 Gaussian elimination

```
n rows/equations, m+1 columns, m variables
  calculates determinant, rank and ans [] -> value for
       variables
  returns {0, 1, INF} -> number of solutions
  gives truely row echelon form. (sort of Identity
       matrixish. but the independent column/variables
       may have non-zero values at the upper rows.)
///For "gaussian elimination modulo version"-
 1. add big mod function to the top.
 2. just do all the calculations in mod(determinant
      too, but not rank).
 3. change fabs() to abs() maybe??
 4. I have add comments to the code to help with
      "steps- 2 and 3".
 5. change all doubles to long long as all the
      calculations are done in long long. omit EPS.
const double EPS = 1e-9;
#define INF 100000000
int Rank; //Rank means number of non-zero rows in
    row-echelon form
double Det:
int gauss( vector<vector<double>> a, vector<double>
    &ans ) {
 int n = (int)a.size(), m = (int)a[0].size()-1;
 ans.clear(): ans.resize(m):
 vector<int> where(m, -1):
 Det = 1.0. Rank = 0:
 for(int col = 0. row = 0: col < m && row < n: ++col) {</pre>
   int sel = row:
   for(int i = row+1: i < n: ++i)</pre>
     if(fabs(a[i][col]) > fabs(a[sel][col])) sel = i;
         ///For modulo version: if(abs(a[i][col]) >
          abs(a[sel][col]))
   if(fabs(a[sel][col]) < EPS) { Det = 0.0; continue;</pre>
        } ///For modulo version: if(!a[sel][col])
   for(int j = 0; j <= m; ++j) swap(a[sel][j],</pre>
        a[row][i]);
   if(row != sel) Det = -Det; ///For modulo version:
        add mod? (idk, i made this up. munim vai kept
        this line as it is. So change only if WA)
   Det *= a[row][col]; ///For modulo version: mod mul
   where[col] = row:
```

```
double s = (1.0 / a[row][col]): ///For modulo
       version: mod inverse
  for(int j = 0; j <= m; ++j) a[row][j] *= s; ///For</pre>
       modulo version: mod mul
  for(int i = 0; i < n; ++i) if (i != row &&</pre>
       fabs(a[i][col]) > EPS) { ///For modulo
       version: if (i != row && a[i][col] > 0)
    double t = a[i][col]:
   for(int j = 0; j <= m; ++j) a[i][j] -= a[row][j]</pre>
        * t: ///For modulo version: mod mul and mod
         sub
  ++row, ++Rank;
for(int i = 0; i < m; ++i) ans[i] = (where[i] == -1)
     ? 0.0 : a[where[i]][m]:
for(int i = Rank: i < n: ++i) if(fabs(a[i][m]) > EPS)
     return 0: ///For modulo version: if(a[i][m])
for(int i = 0: i < m: ++i) if(where[i] == -1) return</pre>
     INF:
return 1;
```

1.5 Policy based Data structures

```
#include<bits/stdc++.h>
#include<ext/pb_ds/assoc_container.hpp>
#include<ext/pb ds/tree policy.hpp>
using namespace __gnu_pbds;
using namespace std;
template <typename T>
using ordered set = tree<T. null type, less<T>.
    rb tree tag, tree order statistics node update>:
    ///set in ascending order (the typical one)
template <tvpename T>
using ordered_set_of_pairs = tree<pair<T, size_t>,
    null type, less<pair<T, size t>>, rb tree tag.
    tree_order_statistics_node_update>; ///set of
    pairs in ascending order (the typical one)
template <typename T>
using ordered_set_desc = tree<T, null_type, greater<T>,
    rb_tree_tag, tree_order_statistics_node_update>;
    ///set in descending order
template <typename T>
using ordered_set_of_pairs_desc = tree<pair<T, size_t>,
    null_type, greater<pair<T, size_t>>, rb_tree_tag,
    tree_order_statistics_node_update>; //set of
```

```
pairs in descending order
#define optimize()
    ios_base::sync_with_stdio(0);cin.tie(0);cout.tie(0);
#define endl '\n'
#define MP make_pair
#define F first
#define S second
#define PB push back
ordered set<int> st1:
ordered set of pairs<int> st2:
ordered_set_desc<int> st3;
ordered set of pairs desc<int> st4:
int main() {
 optimize():
 for( int i = 0: i < 10: ++i ) st1.insert(i):</pre>
 cout << st1.order_of_key(2) << endl; ///how many</pre>
      elements in st1 less than 2? (output: 2)
 cout << *st1.find_by_order(5) << endl << endl;</pre>
      ///what is the 5th minimum element in st1?(0th
      based indexing) (output: 5)
 for( int i = 0; i < 10; ++i ) st2.insert(MP(i, i));</pre>
 cout << st2.order_of_key(MP(2, 3)) << endl;</pre>
      ///output: 3
 cout << st2.order_of_key(MP(2, 2)) << endl;</pre>
      ///output: 2
 cout << st2.order_of_key(MP(3, 2)) << endl;</pre>
      ///output: 3
 cout << st2.order_of_key(MP(3, -1)) << endl;</pre>
      ///output: 4 (i know, you were expecting 3. but
      giving negative numbers as second element gives
      unexpected results.)
 cout << (*st2.find by order(5)).F << " " <<
      (*st2.find by order(5)).S << endl << endl:
      ///output: 5 5
```

1.6 bigint

```
#include <bits/stdc++.h>
using namespace std;

#define optimize()
    ios_base::sync_with_stdio(0);cin.tie(0);cout.tie(0);
#define endl '\n'

struct Bigint {
    // representations and structures
```

```
string a; // to store the digits
int sign; // sign = -1 for negative numbers, sign = 1
    otherwise
// constructors
Bigint() {} // default constructor
Bigint( string b ) { (*this) = b; } // constructor
    for string
// some helpful methods
int size() { // returns number of digits
 return a.size():
Bigint inverseSign() { // changes the sign
 sign *= -1:
 return (*this):
Bigint normalize( int newSign ) { // removes leading
    0. fixes sign
 for( int i = a.size() - 1: i > 0 && a[i] == '0':
      i-- )
   a.erase(a.begin() + i);
  sign = (a.size() == 1 && a[0] == '0') ? 1:
      newSign;
 return (*this):
// assignment operator
void operator = ( string b ) { // assigns a string to
 a = b[0] == '-' ? b.substr(1) : b;
 reverse( a.begin(), a.end() );
 this->normalize( b[0] == '-' ? -1 : 1 );
// conditional operators
bool operator < ( const Bigint &b ) const { // less
     than operator
 if( sign != b.sign ) return sign < b.sign;</pre>
 if( a.size() != b.a.size() )
   return sign == 1 ? a.size() < b.a.size() :</pre>
        a.size() > b.a.size():
  for( int i = a.size() - 1; i >= 0; i-- ) if( a[i]
      != b.a[i] )
   return sign == 1 ? a[i] < b.a[i] : a[i] > b.a[i];
 return false:
bool operator == ( const Bigint &b ) const { //
     operator for equality
 return a == b.a && sign == b.sign:
// mathematical operators
Bigint operator + ( Bigint b ) { // addition operator
    overloading
 if( sign != b.sign ) return (*this) -
      b.inverseSign();
 Bigint c;
```

```
for(int i = 0, carry = 0; i<a.size() || i<b.size()</pre>
      || carrv: i++ ) {
   carry+=(i<a.size() ? a[i]-48 : 0)+(i<b.a.size() ?</pre>
        b.a[i]-48:0);
   c.a += (carry % 10 + 48);
   carry /= 10;
 return c.normalize(sign);
Bigint operator - ( Bigint b ) { // subtraction
     operator overloading
 if( sign != b.sign ) return (*this) +
      b.inverseSign():
 int s = sign; sign = b.sign = 1;
 if((*this) < b) return ((b -
      (*this)).inverseSign()).normalize(-s):
 for( int i = 0, borrow = 0; i < a.size(); i++ ) {</pre>
   borrow = a[i] - borrow - (i < b.size() ? b.a[i] :</pre>
   c.a += borrow >= 0 ? borrow + 48 : borrow + 58:
   borrow = borrow \ge 0 ? 0 : 1:
 return c.normalize(s);
Bigint operator * ( Bigint b ) { // multiplication
     operator overloading
 Bigint c("0");
 for( int i = 0, k = a[i] - 48; i < a.size(); i++, k</pre>
      = a[i] - 48) {
   while(k--) c = c + b; // ith digit is k, so, we
        add k times
   b.a.insert(b.a.begin(), '0'); // multiplied by 10
 return c.normalize(sign * b.sign);
Bigint operator / ( Bigint b ) { // division operator
    overloading
 if( b.size() == 1 && b.a[0] == '0' ) b.a[0] /= (
      b.a[0] - 48):
 Bigint c("0"), d;
 for( int i = 0: i < a.size(): i++ ) d.a += "0":</pre>
 int dSign = sign * b.sign: b.sign = 1:
 for( int i = a.size() - 1: i >= 0: i-- ) {
   c.a.insert( c.a.begin(), '0');
   c = c + a.substr(i. 1):
   while(!(c < b)) c = c - b, d.a[i]++;
 return d.normalize(dSign);
Bigint operator % ( Bigint b ) { // modulo operator
     overloading
```

```
if( b.size() == 1 && b.a[0] == '0' ) b.a[0] /= (
       b.a[0] - 48):
   Bigint c("0");
   b.sign = 1;
   for( int i = a.size() - 1; i >= 0; i-- ) {
    c.a.insert( c.a.begin(), '0');
    c = c + a.substr(i, 1);
    while( !(c < b) ) c = c - b;
   return c.normalize(sign):
 // output method
 void print() {
   if( sign == -1 ) cout << '-';
   for( int i = a.size() - 1: i >= 0: i-- ) cout <<
}:
int main() {
 optimize():
 Bigint a, b, c; // declared some Bigint variables
 // taking Bigint input //
 string input; // string to take input
 cin >> input; // take the Big integer as string
 a = input; // assign the string to Bigint a
 cin >> input; // take the Big integer as string
 b = input; // assign the string to Bigint b
 // Using mathematical operators //
 c = a + b; // adding a and b
 c.print(): // printing the Bigint
 cout << endl: // newline</pre>
 c = a - b: // subtracting b from a
 c.print(): // printing the Bigint
 cout << endl: // newline</pre>
 c = a * b: // multiplying a and b
 c.print(); // printing the Bigint
 cout << endl: // newline</pre>
 c = a / b: // dividing a by b
 c.print(); // printing the Bigint
 cout << endl: // newline</pre>
 c = a \% b: // a modulo b
 c.print(); // printing the Bigint
 cout << endl; // newline</pre>
 // Using conditional operators //
 if( a == b ) puts("equal"); // checking equality
 else puts("not equal");
```

```
if( a < b ) puts("a is smaller than b"); // checking
    less than operator
  return 0;
}</pre>
```

1.7 bit math

1.7.1 bit formula

```
//& = bitwise and
//| = bitwise or
//^ = bitwise xor
1. a+b = (a^b) + 2*(a&b)
2. a+b = (a|b) + (a&b)
```

1.8 bitset

```
int main() {
 bitset<5005> bss[12] . now :
 bss[i].set(j); // sets jth bit of i'th bitset
 now = bss[i]&bss[j]; // can do logical operation
      betwn two bitset
 int here = now.count(); // counts number of 1s in the
 now[i] = 1; // sets i-th bit from right ( 0 based) to
 now.all() : // check if all the bits are set
 now.anv() : // checks if any bits are set
 now.none(): //checks if none of the bits are set
 bitset<5005> bs1, bs2, bs3; //Complexity of declaring
      bitsets is O(N/32)
 bs3 = bs1 ^bs2 (Complexity is O(N/32))
 bs3 = bs1 & bs2 (Complexity is O(N/32))
 bs3 = bs1 | bs2 (Complexity is O(N/32))
 bs3 = bs2 (Complexity is O(N/32))
 bs3 = "bs3 // inverse all bits.
 bs3.Find_next(idx); (Complexity is O(N/32)) //Finds
      the position of set bit after idx;
 bs3.Find_First(); (Complexity is O(N/32)) //Finds the
      first set bit; return bs3.size() if there is no
      set bit:
 bs3.reset(); //rests all bits
 bs3.set(): //set all bits
 bs3.flip(); //flip all bits
 bs3 = "bs3; //flip all bits
```

```
cout<<bb.to_ullong()<<"\n"; // to unsigned long long
   int
  return 0;</pre>
```

1.9 combinatorics

1.9.1 combinatorics

```
    nCr = nC(r-1) * ((n-r+1)/r)
    nCr = (n-1)Cr + (n-1)C(r-1)
    nC0+nC2+nC4+.... = nC1+nC3+nC5.... = 2^(n-1)
    summation of x*nCx such that (1 <= x <= n or 0 <= x <= n) = n*2^(n-1). (easy to prove).</li>
    Catalan Number(1, 1, 2, 5, 14, 42, 132, 429, 1430, 4862, 16796, 58786, 208012, 742900, 2674440, 9694845, 35357670, 129644790, 477638700, 1767263190)
    1 Catn = (1/(n+1))*((2n)Cn)
    2 Catn = summation of Cat(k-1)*Catk(n-k) such that 1 <= k <= n</li>
    3 Cat(n+1) = ((2*(2n+1))/(n+2)) * Catn
```

1.10 flow

1.10.1 Dinic with and without scaling

```
Dinic algorithm with scaling and not scaling.
Complexity with scaling : O(VE * lg(U)). here, U is the
    maximum capacity
Complexity without scaling : O((V^2)E)
Complexity in bipartite graph matching: O(E*sqrt(V)).
Complexity in unit graph: O(E*sqrt(V)). (but. Chilli
    says it is wrong and actual complexity is
    O(E*min(V^{(2/3)}, E^{(1/2)})).
A unit network is a network in which all the edges have
    unit capacity, and for any vertex except s and t
    either incoming or outgoing edge is unique. That's
    exactly the case with the network we build to
    solve the maximum matching problem with flows. So,
    Dinic gives similar performance to hopcroft karp
    algorithm incase of maximum bipartite matching as
    it has the same time complexity.
//////
1. Works on directed graph
```

```
2. Works on undirected graph
3. Works on multi-edge(directed/undirected) graph
4. Works on self-loop(directed/undirected) graph
Can find the actual flow.
Can find the non-maxflow upto a certain value
Implement it on your own. very easy. just a simple O(n)
Can find minimum cut sets(A and B).
A contains source itself and the nodes that are
    reachable from source using non-saturated edges.
B contains sink and the nodes that are not reachable
    from source using non-saturated edges.
Value of minimum cut capacity is summation of the
    per_cap of all edges u->v such that u belongs to
    set A. v belongs to set B.
Value of minimum cut flow is = (summation of the flow
    of all edges u->v such that u belongs to set A. v
    belongs to set B) - (summation of the flow of all
    edges v->u such that v belongs to set B, u belongs
    to set A)
Value of minimum cut capacity == maxflow.
To find the minimum cut sets(A and B) first find the
    maxflow. Then, we apply dfs from the source. This
    dfs will be such that- only
the nodes of set A will eventually be marked visited.
    Suppose we have reached in node u, that means u
    belongs to set A. Now,
If (flow of edge u->v) < (per_cap of edge u->v) then v
    belongs to set A.
else if(flow of edge u \rightarrow v) == (per cap of edge u \rightarrow v)
    then v belongs to set B.
To find the minimum cut sets with minimum cardinality.
    first multiply (E+1) with all the edges. Then add
    1 to all the edges. Then run the maxflow algorithm.
111111
**/
#include <bits/stdc++.h>
using namespace std:
#define optimize()
    ios base::svnc with stdio(0):cin.tie(0):cout.tie(0):
#define endl '\n'
template <class flow_t> struct Dinic { //int/long long;
 const static bool SCALING = true; // non-scaling =
      V^2E, Scaling=VElog(U) with higher constant
 long long lim = 1;
 const flow_t INF = numeric_limits<flow_t>::max();
```

```
bool K2 = false;
struct edge {
 int to, rev;
 flow_t cap, flow, per_cap;
int s, t;
vector<int> level, ptr;
vector< vector<edge> > adj;
Dinic(int NN) : s(NN-2), t(NN-1), level(NN), ptr(NN),
     adi(NN) {}
void flow limit( flow t val ) { //non-maxflow upto K.
 K2 = true:
 K = val;
void addEdge(int a, int b, flow_t cap, bool
     isDirected = true) {
 adi[a].push back({b, (int)adi[b].size(), cap. 0.
       cap}):
  adi[b].push back({a, (int)adi[a].size() - 1,
      isDirected ? 0 : cap, 0, isDirected ? 0 :
      cap}):
bool bfs() {
 queue<int> q({s});
 fill( level.begin(), level.end(), -1 );
 level[s] = 0;
 while (!q.empty() && level[t] == -1) {
   int v = q.front();
   q.pop();
   for (auto e : adj[v]) {
     if (level[e.to] == -1 && e.flow < e.cap &&</pre>
          (!SCALING || e.cap - e.flow >= lim)) {
       q.push(e.to);
       level[e.to] = level[v] + 1;
 return level[t] != -1:
flow t dfs(int v. flow t flow) {
 if (v == t || !flow) return flow;
   for (: ptr[v] < adi[v].size(): ptr[v]++) {</pre>
     edge &e = adi[v][ptr[v]]:
     if (level[e.to] != level[v] + 1) continue:
     if (flow t pushed = dfs(e.to, min(flow, e.cap -
          e.flow))) {
       e.flow += pushed;
       adj[e.to][e.rev].flow -= pushed;
       return pushed;
 return 0;
```

```
flow_t max_flow(int source, int sink) {
   s = source, t = sink;
   long long flow = 0;
   for (lim = SCALING ? (1LL << 30) : 1; lim > 0; lim
        >>= 1) { //Here, lim = SCALING?(U):1 ; Here U
        is an int/long long strictly greater than the
        max capacity ;
     while (bfs()) {
      fill( ptr.begin(), ptr.end(), 0 );
      while (flow_t pushed = dfs(s,
           ((K2==true)?K:INF))) {
        flow += pushed:
        if(K2) {
          K -= pushed;
          if( K == 0 ) break:
       if( K2 && (K == 0) ) break:
   return flow:
 vector<pair<int,int>,long long>> getActualFlow()
   vector<pair<int,int>, long long>> vec;
   for( int i = 0; i < adj.size(); ++i ) {</pre>
    for( int j = 0; j < adj[i].size(); ++j ) {</pre>
      if( adj[i][j].flow > 0 ) {
        vec.push_back( make_pair(make_pair(i,
             adj[i][j].to), adj[i][j].flow));
     }
   return vec;
int main() {
 optimize():
 int T:
 cin >> T:
 for( int test = 1; test <= T; ++test ) {</pre>
   int N. M. s. t: // no. of nodes: no. of edges:
        source: sink:
   cin >> N >> M >> s >> t:
   Dinic<int> fl(N+1); //for long long change int to
        long long: no. of nodes+1:
   for( int i = 1; i <= M; ++i ) {</pre>
     int u, v, w;
     cin >> u >> v >> w;
    fl.addEdge(u, v, w); //Directed graph;
     //fl.addEdge(u, v, w, false); //Undirected graph;
```

1.10.2 Hopcroft karp

```
//Complexity: O(\operatorname{sqrt}(V) * E), constant may be a bit
    high.
//Works on self loops.
#include <bits/stdc++.h>
using namespace std;
struct Hopcroft_karp {
 int n;
 vector< vector<int> > edge;
 vector<int> dis, parent, L, R;
 vector<int> Q;
 Hopcroft_karp(int n_) : n(n_), edge(n_+1), dis(n_+1),
      parent(n_{+1}), L(n_{+1}), R(n_{+1}), Q(n_{+1}) {};
 void add edge( int u, int v ) {
   edge[u].push back(v):
 bool dfs(int i) {
   int len = edge[i].size();
   for (int j = 0; j < len; j++) {
     int x = edge[i][i];
     if (L[x] == -1 \mid | (parent[L[x]] == i))  {
       if (L[x] == -1 \mid | dfs(L[x])) {
         L[x] = i:
         R[i] = x;
         return (true):
     }
   return false;
 bool bfs() {
   int x, f = 0, 1 = 0;
   fill(dis.begin(), dis.end(), -1);
   for (int i = 1; i <= n; i++) {</pre>
     if (R[i] == -1) {
```

```
Q[1++] = i;
       dis[i] = 0;
   }
   while (f < 1) {</pre>
     int i = Q[f++];
     int len = edge[i].size();
     for (int j = 0; j < len; j++) {</pre>
       x = edge[i][i]:
       if (L[x] == -1) return true;
       else if (dis[L[x]] == -1) {
         parent[L[x]] = i:
         dis[L[x]] = dis[i] + 1:
         Q[1++] = L[x];
     }
   return false:
  int matching() {
   int counter = 0; //How many nodes are part of the
        maximum matching?
   fill( L.begin(), L.end(), -1 );
   fill( R.begin(), R.end(), -1 );
   while (bfs()) {
     for (int i = 1; i <= n; i++) {
       if (R[i] == -1 \&\& dfs(i)) counter++:
   return counter;
int main() {
 int n, m; //no. of nodes; no. of edges;
 cin >> n >> m:
 Hopcroft karp h(n):
 for( int i = 0: i < m: ++i ) {
   int u. v:
   cin >> u >> v:
   /** //Undirected: //(hopcroft karp in undirected
        graph is impossible, i dont know why i added
        this. :|)
   h.add edge(u, v):
   h.add edge(v. u):
   //Directed:
   h.add_edge(u, v);
 int ans = h.matching(); //How many nodes are part of
      the maximum matching?
 cout << ans << endl;</pre>
 //print the actual maximum matching
 for( int i = 1; i <= n; ++i ) {</pre>
```

```
if( h.L[i] != -1 ) cout << i << " " << h.L[i] <<
    endl;
}
</pre>
```

1.10.3 minCostflow by dacin21(1)

```
/**
// Push-Relabel implementation of the cost-scaling
    algorithm
// Runs in O( <max flow> * log(V * max edge cost)) = O(
    V^3 * log(V * C)
// Operates on integers
// Works on regular directed graphs.
// Can't operate on doubles. For this, use bellman
    ford/diikstra method.
//Whether or not it works on undirected or multi-edge
    or self-loop graphs is yet to be verified.
// To get the actual flow collect all the edges which
    have (f > 0) - this works just like the actual
    flow finding in normal max flow algorithms;
// Can't find nonMaxflow value upto K. The code is too
    complex for me to add this function. Maybe its not
     possible at all and you can only do it with the
    bellmanFord/Djikstra method.
#include<bits/stdc++.h>
using namespace std;
#define optimize()
    ios base::svnc with stdio(0):cin.tie(0):cout.tie(0):
#define endl '\n'
template<typename flow_t = int, typename cost_t = int>
    ///int/long long; int/long long;
struct mcSFlow{
 struct Edge{
   cost t c:
   flow t f:
   int to, rev:
   Edge(int _to, cost_t _c, flow_t _f, int
        _rev):c(_c), f(_f), to(_to), rev(_rev){}
 }:
 const cost t INFCOST =
      numeric_limits<cost_t>::max()/2; ///divide by
      slightly bigger number if overflow;
 const cost_t INFFLOW =
      numeric_limits<flow_t>::max()/2; ///divide by
      slightly bigger number if overflow:
 cost_t epsilon;
 int N. S. T:
 vector<vector<Edge> > G;
 vector<unsigned int> isEnqueued, state;
```

```
mcSFlow(int _N):epsilon(0), N(_N), G(_N){}
void add_edge(int a, int b, cost_t cost, flow_t cap){
 if(a==b){assert(cost>=0); return;}
 cost*=N;// to preserve integer-values
 epsilon = max(epsilon, abs(cost));
 assert(a>=0&&a<N&&b>=0&&b<N);
 G[a].emplace_back(b, cost, cap, G[b].size());
 G[b].emplace_back(a, -cost, 0, G[a].size()-1);
flow t calc max flow(){ // Dinic max-flow
 vector<flow t> dist(N). state(N):
 vector<Edge*> path(N):
 auto cmp = [](Edge*a, Edge*b){return a->f < b->f:}:
 flow t addFlow. retflow=0::
   fill(dist.begin(), dist.end(), -1);
   dist[S]=0:
   auto head = state.begin(). tail = state.begin():
   for(*tail++ = S:head!=tail:++head){
     for(Edge const&e:G[*head]){
      if(e.f && dist[e.to]==-1){
        dist[e.to] = dist[*head]+1;
         *tail++=e.to:
   addFlow = 0:
   fill(state.begin(), state.end(), 0);
   auto top = path.begin();
   Edge dummy(S, 0, INFFLOW, -1);
   *top++ = &dummy;
   while(top != path.begin()){
     int n = (*prev(top))->to;
       auto next top = min element(path.begin(), top.
       flow t flow = (*next top)->f:
       while(--top!=path.begin()){
        Edge &e=**top, &f=G[e.to][e.rev];
        e.f-=flow:
        f.f+=flow:
       addFlow=1:
       retflow+=flow:
       top = next_top;
       continue:
     for(int &i=state[n], i_max = G[n].size(), need
         = dist[n]+1;;++i){
       if(i==i max){
        dist[n]=-1;
        --top;
        break:
```

```
if(dist[G[n][i].to] == need && G[n][i].f){
         *top++ = &G[n][i];
         break;
  }while(addFlow);
 return retflow:
vector<flow t> excess:
vector<cost t> h:
void push(Edge &e. flow t amt){
 //cerr << "push: " << G[e.to][e.rev].to << " -> "
      << e.to << " (" << e.f << "/" << e.c << ") : "
      << amt. << "\n":
 if(e.f < amt) amt=e.f:</pre>
 e.f-=amt:
 excess[e.to]+=amt:
 G[e.to][e.rev].f+=amt:
 excess[G[e.to][e.rev].to]-=amt:
void relabel(int vertex){
 cost_t newHeight = -INFCOST;
 for(unsigned int i=0;i<G[vertex].size();++i){</pre>
   Edge const&e = G[vertex][i];
   if(e.f && newHeight < h[e.to]-e.c){</pre>
     newHeight = h[e.to] - e.c;
     state[vertex] = i:
 h[vertex] = newHeight - epsilon;
const int scale=2:
pair<flow_t, cost_t> minCostFlow(int _S, int _T){
 S = S, T = T:
 cost t retCost = 0:
 for(int i=0:i<N:++i){</pre>
   for(Edge &e:G[i]){
     retCost += e.c*(e.f):
 //find feasible flow
 flow t retFlow = calc max flow():
  excess.resize(N):h.resize(N):
 aueue<int> a:
 isEnqueued.assign(N, 0); state.assign(N,0);
 for(;epsilon;epsilon>>=scale){
   //refine
   fill(state.begin(), state.end(), 0);
   for(int i=0;i<N;++i)</pre>
     for(auto &e:G[i])
```

```
if(h[i] + e.c - h[e.to] < 0 && e.f) push(e,
             e.f):
     for(int i=0;i<N;++i){</pre>
      if(excess[i]>0){
        q.push(i);
        isEnqueued[i]=1;
     }
     while(!a.emptv()){
      int cur=q.front();q.pop();
      isEngueued[cur]=0:
      // discharge
      while(excess[cur]>0){
        if(state[cur] == G[cur].size()){
          relabel(cur):
         for(unsigned int &i=state[cur]. max i =
             G[curl.size():i<max i:++i){</pre>
          Edge &e=G[curl[i]:
          if(h[cur] + e.c - h[e.to] < 0)
            push(e, excess[cur]);
            if(excess[e.to]>0 && isEnqueued[e.to]==0){
              q.push(e.to);
              isEnqueued[e.to]=1;
            if(excess[cur]==0) break;
     if(epsilon>1 && epsilon>>scale==0){
       epsilon = 1<<scale;
   for(int i=0:i<N:++i){</pre>
     for(Edge &e:G[i]){
      retCost -= e.c*(e.f):
   //cerr << " -> " << retFlow << " / " << retCost <<
        " bzw. " << retCost/2/N << "\n":
   return make pair(retFlow. retCost/2/N):
 flow t getFlow(Edge const &e){
   return G[e.to][e.rev].f:
 }
int main() {
 optimize();
 int T;
 cin >> T;
 for( int test = 1; test <= T; ++test ) {</pre>
```

}:

```
int N, M, S, T; /// Number of nodes; number of
      edges: source: sink:
  cin >> N >> M >> S >> T;
 mcSFlow<> fl(N+1); ///Add long long, long if
      necessary:
 for( int i = 1; i <= M; ++i ) {</pre>
   int u, v, c, w; /// node; node; cost; capacity;
   cin >> u >> v >> c >> w;
   fl.add_edge(u, v, c, w); //Directed graph;
  cout << fl.minCostFlow(S, T).first << " " <<</pre>
      fl.minCostFlow(S, T).second: ///flow: cost:
}
```

1.11 geometry

1.11.1 comparison of doubles

```
1. bool equalTo ( double a, double b ){ if ( fabs ( a -
    b) <= eps ) return true; else return false; }</pre>
2. bool notEqual ( double a, double b ){if ( fabs ( a -
    b ) > eps ) return true; else return false; }
3. bool lessThan (double a, double b){ if (a + eps <
    b) return true: else return false: }
4. bool lessThanEqual ( double a, double b ){if ( a < b
    + eps ) return true; else return false;}
5. bool greaterThan (double a, double b){if (a > b +
    eps ) return true;else return false;}
6. bool greaterThanEqual ( double a, double b ) if ( a
    + eps > b ) return true:else return false:}
```

1.11.2 convex hull 2D

```
//Complexity : O(NlogN)
// returns points of convex hull in CW direction.
double EPS = 1e-12:
struct PT {
 double x, y;
 PT() {}
 PT(double x, double y) : x(x), y(y) {}
 PT(const PT &p) : x(p.x), y(p.y) {}
 PT operator + (const PT &p) const { return PT(x+p.x,
      v+p.v): }
 PT operator - (const PT &p) const { return PT(x-p.x,
      v-p.v); }
 PT operator * (double c) const { return PT(x*c, y*c
```

```
PT operator / (double c) const { return PT(x/c, y/c
      ): }
 bool operator <(const PT &p) const {</pre>
   return x < p.x | | (x == p.x && y < p.y);
};
ostream & operator << (ostream & os, const PT & p) {
 os << "(" << p.x << "," << p.y << ")";
// checks if a-b-c is CW or not.
bool isPointsCW(PT a. PT b. PT c) {
 return a.x*(b.y-c.y)+b.x*(c.y-a.y)+c.x*(a.y-b.y)+EPS
}
// checks if a-b-c is CCW or not.
bool isPointsCCW(PT a, PT b, PT c) {
 return a.x*(b.v-c.v)+b.x*(c.v-a.v)+c.x*(a.v-b.v) >
// checks if a-b-c is collinear or not.
bool isPointsCollinear(PT a, PT b, PT c) {
 return abs(a.x*(b.y-c.y)+b.x*(c.y-a.y)+c.x*(a.y-b.y))
void convex_hull(vector<PT>& a) { ///^ == upper hull, U
     == lower hull
 int a sz = a.size():
 if (a_sz == 1) return;
  sort(a.begin(), a.end());
 PT p1 = a[0], p2 = a.back();
 vector<PT> up, down; // up will store upper
      hull/lower envelope, down will store lower
      hull/upper envelope
 up.push_back(p1);
 down.push back(p1):
 int up sz = 1, down sz = 1:
  for (int i = 1: i < a sz: i++) {
   if (i == a sz - 1 || isPointsCW(p1, a[i], p2)) {
///include all the collinear points of the boundary
     instead of only the outer two points:-
///while (up_sz >= 2 && (!isPointsCW(up[up_sz-2],
    up[up sz-1], a[i]) &&
     !isPointsCollinear(up[up sz-2], up[up sz-1].
     a[i])))
     while (up_sz >= 2 && !isPointsCW(up[up_sz-2],
          up[up sz-1], a[i])) {
       up.pop_back();
       --up_sz;
     up.push_back(a[i]);
     ++up_sz;
   if (i == a_sz - 1 || isPointsCCW(p1, a[i], p2)) {
```

```
///include all the collinear points of the boundary
    instead of only the outer two points:-
///while(down_sz >= 2 && (!isPointsCCW(down[down_sz-2],
    down[down sz-1], a[i]) &&
    !isPointsCollinear(down[down_sz-2],
    down[down_sz-1], a[i])) )
     while(down_sz >= 2 &&
          !isPointsCCW(down[down_sz-2],
          down[down sz-1], a[i])) {
      down.pop_back();
       --down sz:
     down.push back(a[i]):
     ++down_sz;
 // up has upper hull/lower envelope
 // down has lower hull/upper envelope
 // a has the overall convex hull
 a.clear():
 for (int i = 0; i < up_sz; i++) a.push_back(up[i]);</pre>
 for (int i = down_sz - 2; i > 0; i--)
      a.push back(down[i]):
```

1.11.3 geo routine

```
#include <bits/stdc++.h>
using namespace std;
#define optimize()
    ios base::svnc with stdio(0):cin.tie(0):cout.tie(0):
#define endl '\n'
typedef long long 11;
typedef vector<int> vi;
typedef vector<ll> v1;
typedef pair<int.int> pii:
typedef pair<11.11> pll:
typedef vector<pii> vii:
typedef vector<pll> vll:
typedef long double ld;
#define F first
#define S second
#define MP make_pair
#define PB push_back
double INF = 1e100;
double EPS = 1e-12;
struct PT {
 double x, v;
 PT() {}
 PT(double x, double y) : x(x), y(y) {}
 PT(const PT \&p) : x(p.x), y(p.y) {}
```

```
PT operator + (const PT &p) const { return PT(x+p.x,
      v+p.v); }
 PT operator - (const PT &p) const { return PT(x-p.x,
      v-p.v); }
 PT operator * (double c) const { return PT(x*c, y*c
 PT operator / (double c) const { return PT(x/c, y/c
      ): }
 bool operator <(const PT &p) const {
   return x < p.x || (x == p.x \&\& y < p.y);
 bool operator ==(const PT &p) const {
   return (fabs(x-p.x) <= EPS && fabs(y-p.y) <= EPS):
double dot(PT p, PT q) { return p.x*q.x+p.v*q.v: }
double dist2(PT p. PT a) { return dot(p-a.p-a): }
double cross(PT p, PT q) { return p.x*q.y-p.y*q.x; }
ostream &operator << (ostream &os. const PT &p) {
 os << "(" << p.x << "," << p.v << ")":
// checks if a-b-c is CW or not.
bool isPointsCW(PT a, PT b, PT c) {
        a.x*(b.y-c.y)+b.x*(c.y-a.y)+c.x*(a.y-b.y)+EPS
// checks if a-b-c is CCW or not.
bool isPointsCCW(PT a, PT b, PT c) {
   return a.x*(b.y-c.y)+b.x*(c.y-a.y)+c.x*(a.y-b.y) >
// checks if a-b-c is collinear or not.
bool isPointsCollinear(PT a, PT b, PT c) {
   return
        abs(a.x*(b.v-c.v)+b.x*(c.v-a.v)+c.x*(a.v-b.v))
// rotate a point CCW or CW around the origin
PT RotateCCW90(PT p) { return PT(-p.v.p.x); }
PT RotateCW90(PT p) { return PT(p.y,-p.x); }
PT RotateCCW(PT p. double t) { // rotate a point CCW t
    degrees around the origin
 return PT(p.x*cos(t)-p.v*sin(t).
      p.x*sin(t)+p.v*cos(t)):
// rotate a point p CCW around another point q by angle
    t(radian)
PT RotateCCW_around_point(PT p, PT q, double t) {
 return PT((p.x-q.x)*cos(t)-(p.y-q.y)*sin(t)+q.x,
      (p.x-q.x)*sin(t)+(p.y-q.y)*cos(t)+q.y);
// project point c onto line through a and b
```

```
// assuming a != b
PT ProjectPointLine(PT a, PT b, PT c) {
 return a + (b-a)*dot(c-a, b-a)/dot(b-a, b-a);
// project point c onto line segment through a and b
PT ProjectPointSegment(PT a, PT b, PT c) {
 double r = dot(b-a,b-a);
 if (fabs(r) < EPS) return a;</pre>
 r = dot(c-a, b-a)/r:
 if (r+EPS < 0) return a:
 if (r > 1+EPS) return b:
 return a + (b-a)*r:
// compute distance from c to segment between a and b
double DistancePointSegment(PT a, PT b, PT c) {
 return sqrt(dist2(c, ProjectPointSegment(a, b, c)));
// compute distance between point (x.v.z) and plane
    ax+bv+cz=d
double DistancePointPlane(double x, double v, double z,
                       double a. double b. double c.
                            double d) {
 return fabs(a*x+b*y+c*z-d)/sqrt(a*a+b*b+c*c);
// determine if lines from a to b and c to d are
     parallel or collinear
bool LinesParallel(PT a, PT b, PT c, PT d) {
 return fabs(cross(b-a, c-d)) < EPS;</pre>
bool LinesCollinear(PT a, PT b, PT c, PT d) {
 return LinesParallel(a, b, c, d)
   && fabs(cross(a-b, a-c)) < EPS
   && fabs(cross(c-d, c-a)) < EPS;
// determine if line segment from a to b intersects with
// line segment from c to d
bool SegmentsIntersect(PT a. PT b. PT c. PT d) {
 if (LinesCollinear(a, b, c, d)) {
   if (dist2(a, c) < EPS || dist2(a, d) < EPS ||</pre>
     dist2(b, c) < EPS || dist2(b, d) < EPS) return
          true:
   if (dot(c-a, c-b) > EPS \&\& dot(d-a, d-b) > EPS \&\&
        dot(c-b, d-b) > EPS)
     return false:
   return true:
 if (cross(d-a, b-a) * cross(c-a, b-a) > EPS) return
 if (cross(a-c, d-c) * cross(b-c, d-c) > EPS) return
      false:
 return true;
// compute intersection of line passing through a and b
```

```
// with line passing through c and d, assuming that
     unique
// intersection exists; for segment intersection, check
// segments intersect first
PT ComputeLineIntersection(PT a, PT b, PT c, PT d) {
 b=b-a; d=c-d; c=c-a;
 assert(dot(b, b) > EPS && dot(d, d) > EPS);
 return a + b*cross(c, d)/cross(b, d):
// shift the straight line passing through points a and
// by distance Dist.
// If Dist is negative the line is shifted rightwards
     or upwards.
// If Dist is positive the line is shifted leftwards or
    downwards.
// The new line passes through points c and d
pair<PT.PT> ShiftLineBvDist(PT a, PT b, double Dist) {
 double r = sqrt( dist2(a, b) );
 double delx = (Dist*(a.y-b.y))/r;
 double dely = (Dist*(b.x-a.x))/r;
 PT c = PT(a.x+delx, a.y+dely);
 PT d = PT(b.x+delx, b.y+dely);
 return MP(c, d);
// This code computes the area or centroid of a
     (possibly nonconvex)
// polygon, assuming that the coordinates are listed in
    a clockwise or
// counterclockwise fashion. Note that the centroid is
    often known as
// the "center of gravity" or "center of mass".
double ComputeSignedArea(const vector<PT> &p) {
 double area = 0:
 for(int i = 0: i < p.size(): i++) {</pre>
   int j = (i+1) % p.size();
   area += p[i].x*p[i].v - p[i].x*p[i].v:
 return area / 2.0:
double ComputeArea(const vector<PT> &p) {
 return fabs(ComputeSignedArea(p)):
PT ComputeCentroid(const vector<PT> &p) {
 PT c(0.0):
 double scale = 6.0 * ComputeSignedArea(p);
 for (int i = 0; i < p.size(); i++){</pre>
   int j = (i+1) % p.size();
   c = c + (p[i]+p[j])*(p[i].x*p[j].v - p[j].x*p[i].v);
 return c / scale;
```

```
// angle from p2->p1 to p2->p3, returns -PI to PI (but
    you may choose to return the absolute value only
     which is "more practical").
double angle(PT p1, PT p2, PT p3) {
 PT va = p1-p2, vb=p3-p2;
  double x,v;
 x=dot(va,vb);
 v=cross(va,vb);
  return(atan2(y,x)); /// return abs( atan2(y,x) ) to
       get the absolute value of angle. (tested)
double angle_to_radian( double theta ) {
 return ((M PI/180.0)*theta):
double radian_to_angle( double x ) {
 return ((180.0/M PI)*x):
int main() {
 // expected: (-5.2)
  cerr << RotateCCW90(PT(2.5)) << endl:</pre>
  // expected: (5,-2)
  cerr << RotateCW90(PT(2,5)) << endl;</pre>
 // expected: (-5,2)
  cerr << RotateCCW(PT(2,5),M_PI/2) << endl;</pre>
  // expected: (5,2)
  cerr << ProjectPointLine(PT(-5,-2), PT(10,4),</pre>
       PT(3.7)) << end1:
  // expected: (5,2) (7.5,3) (2.5,1)
  cerr << ProjectPointSegment(PT(-5,-2), PT(10,4),</pre>
       PT(3.7)) << " "
    << ProjectPointSegment(PT(7.5,3), PT(10,4),</pre>
        PT(3.7)) << " '
    << ProjectPointSegment(PT(-5,-2), PT(2.5,1),</pre>
        PT(3,7)) << end1;
  // expected: 6.78903
  cerr << DistancePointPlane(4.-4.3.2.-2.5.-8) << endl:
  // expected: 1 0 1
  cerr << LinesParallel(PT(1.1), PT(3.5), PT(2.1),
      PT(4,5)) << " "
    << LinesParallel(PT(1,1), PT(3,5), PT(2,0),</pre>
        PT(4.5)) << " '
    << LinesParallel(PT(1.1), PT(3.5), PT(5.9).</pre>
        PT(7.13)) << endl:
  // expected: 0 0 1
  cerr << LinesCollinear(PT(1.1), PT(3.5), PT(2.1),
       PT(4.5)) << " "
    << LinesCollinear(PT(1,1), PT(3,5), PT(2,0),</pre>
        PT(4.5)) << " '
    << LinesCollinear(PT(1,1), PT(3,5), PT(5,9),</pre>
        PT(7,13)) << endl;
  // expected: (1,2)
  cerr << ComputeLineIntersection(PT(0,0), PT(2,4),</pre>
       PT(3,1), PT(-1,3)) << endl;
```

```
// area should be 5.0
// centroid should be (1.1666666, 1.166666)
PT pa[] = \{ PT(0,0), PT(5,0), PT(1,1), PT(0,5) \};
vector<PT> p(pa, pa+4);
PT c = ComputeCentroid(p);
cerr << "Area: " << ComputeArea(p) << endl;</pre>
cerr << "Centroid: " << c << endl;</pre>
// expected: 0
cerr << isPointsCCW( PT(5, 6), PT(10, 10), PT(11, 5)
    ) << endl:
// expected: 1
cerr << isPointsCCW( PT(5, 6), PT(10, 2), PT(11, 5) )
// expected: 1
cerr << isPointsCW( PT(5, 6), PT(10, 10), PT(11, 5) )</pre>
// expected: 0
cerr << isPointsCW( PT(5, 6), PT(10, 2), PT(11, 5) )
     << endl:
// expected: 0
cerr << isPointsCollinear( PT(5, 6), PT(10, 2),
    PT(11, 5) ) << endl;
// expected: 1
cerr << isPointsCollinear( PT(5, 6), PT(10, 6),
    PT(11, 6) ) << endl;
// expected: (-0.437602,12.6564) (2.5624,14.6564)
cerr << ShiftLineByDist( PT(4, 6), PT(7, 8), 8 ).F <<</pre>
     " " << ShiftLineByDist( PT(4, 6), PT(7, 8), 8
    ).S << endl:
// expected: (8.4376,-0.656402) (11.4376,1.3436)
cerr << ShiftLineByDist( PT(4, 6), PT(7, 8), -8 ).F
    << " " << ShiftLineByDist( PT(4, 6), PT(7, 8),</pre>
     -8 ).S << endl;
```

1.11.4 half plane intersection Radewoosh style (OFFLINE)

```
// OFFLINE
// Complexity: O(NlgN)
// very easy concept and implementation
double INF = 1e100;
double EPS = 1e-12;
struct PT {
   double x, y;
   PT() {}
   PT(double x, double y) : x(x), y(y) {}
   PT(const PT &p) : x(p.x), y(p.y) {}
   PT operator + (const PT &p) const { return PT(x+p.x, y+p.y); }
```

```
PT operator - (const PT &p) const { return PT(x-p.x,
      y-p.y); }
 PT operator * (double c) const { return PT(x*c, y*c
      ); }
 PT operator / (double c) const { return PT(x/c, y/c
 bool operator <(const PT &p) const {</pre>
   return x < p.x || (x == p.x \&\& y < p.y);
ostream &operator << (ostream &os. const PT &p) {
 os << "(" << p.x << "," << p.v << ")";
int steps = 600;
vector<PT> lower_hull, upper_hull;
int lower hull sz. upper hull sz:
bool leBorder = 0. riBorder = 0:
double func( double xx. double val ) {
 double ans1 = INF, ans2 = -INF, ans:
 for( int i = 0: i < lower hull sz-1: ++i ) {</pre>
   if( leBorder && (i == 0) ) continue;
   PT a = lower_hull[i], b = lower_hull[i+1]; //
        straight line passes through points a and b
   double m = (a.y-b.y)/(a.x-b.x); // slope of the
        straight line; if the TL is strict, then
        better precalculate all the slopes and store
        them beforehand
   double c = a.y - a.x*(m); // intercept of the
        straight line; if the TL is strict, then
        better precalculate all the intercepts and
        store them beforehand
   double aa = m*xx;
   double bb = c:
   double cc = aa+bb:
   ans1 = min(ans1, cc):
 for( int i = 0: i < upper hull sz-1: ++i ) {</pre>
   if( riBorder && (i == upper_hull_sz-2) ) continue;
   PT a = upper_hull[i], b = upper_hull[i+1]; //
        straight line passes through points a and b
   double m = (a.v-b.v)/(a.x-b.x): // slope of the
        straight line: if the TL is strict, then
        better precalculate all the slopes and store
        them beforehand
   double c = a.v - a.x*(m): // intercept of the
        straight line; if the TL is strict, then
        better precalculate all the intercepts and
        store them beforehand
   double aa = m*xx:
   double bb = c;
   double cc = aa+bb;
   ans2 = max(ans2, cc);
```

```
ans = ans1-ans2:
 return ans:
bool Ternary_Search(double val) {
 double lo = -INF, hi = INF, mid1, mid2;
 leBorder = 0, riBorder = 0;
 if( lower_hull[0].x == lower_hull[1].x ) lo =
      lower hull[0].x+val. leBorder = 1:
 if( upper_hull[upper_hull_sz-2].x ==
      upper hull[upper hull sz-1].x ) hi =
      upper_hull[upper_hull_sz-1].x-val, riBorder = 1;
 if( lo > hi ) return 0:
 for( int i = 0; i < steps; ++i ) {</pre>
   mid1 = (10*2.0 + hi)/3.0:
   mid2 = (lo + 2.0*hi)/3.0:
   double ff1 = func(mid1, val):
   double ff2 = func(mid2, val):
   if( ff1 >= 0 || ff2 >= 0 ) return 1:
   if( ff1 > ff2 ) hi = mid2:
   else lo = mid1:
 if( func(lo, val) >= 0 ) return 1;
 return 0;
```

1.12 graph

1.12.1 FloydWarshall

```
// order of nodes: k->i->j
```

1.12.2 articulation bridges

```
//complexity: O(E+V)
//nodes are 1-based indexed.
int n; // number of nodes
vector<vector<int>> adjlist; // adjacency list of graph
vector<bool> visited;
vector<int>> tin, low;
int timer;
void dfs_art(int u, int par = -1) {
  visited[u] = true;
  tin[u] = low[u] = timer++;
  for (auto to : adjlist[u]) {
    if (to == par) continue;
    if (visited[to]) {
        low[u] = min(low[u], tin[to]);
    }
}
```

```
}
else {
    dfs_art(to, u);
    low[u] = min(low[u], low[to]);
    if (low[to] > tin[u]) {
        IS_BRIDGE(u, to);
    }
}

void find_bridges() {
    timer = 0;
    visited.clear(); tin.clear(); low.clear();
    visited.assign(n+1, false);
    tin.assign(n+1, -1);
    low.assign(n+1, -1);
    for (int i = 1; i <= n; ++i) {
        if (!visited[i]) dfs_art(i);
    }
}</pre>
```

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1.12.3 centroid decomposition

```
const int MAXN = 1e5+5: /// Maximum no. of nodes:
int n; /// Number of nodes;
vector<int> adjlist[MAXN];
int sub_cd[MAXN]; /// sub_cd[u] = subtree size of node
    u after each decomposing:
bool done_cd[MAXN]; /// done_cd[u] = node u have been
    decomposed:
void dfs sub cd(int u. int par) {
 \sup cd[u] = 1:
  for( auto v: adjlist[u] ) {
   if ( done cd[v] || v == par ) continue:
   dfs sub cd(v, u):
   sub cd[u] += sub cd[v]:
}
int dfs find centroid( int u, int par, int sz ) {
 for( auto v: adilist[u] ) {
   if( !done cd[v] && v != par && sub cd[v] > sz ) {
     return dfs_find_centroid( v, u, sz );
 return u;
void dfs_decompose( int u ) {
 dfs_sub_cd(u, -1);
 int centroid = dfs_find_centroid(u, -1, sub_cd[u]/2);
 done_cd[centroid] = 1;
 for( auto v : adjlist[centroid] ) {
```

```
if( !done_cd[v] ) dfs_decompose(v);
}
```

1.12.4 djikstra

```
///O((V+E)logV)
///Works on negative edges(a bit slow) but not negative
const int MAXN = 1e5+5;
vi adjlist[MAXN];
int dist[MAXN];
void djikstra( int source ) {
 for( int i = 0; i < MAXN; ++i ) dist[i] = inf;</pre>
 dist[source] = 0:
 priority_queue< pii, vector<pii>, greater<pii> > pq;
 pq.push(MP(0, source));
 while (!pq.empty()) {
   pii Front = pq.top(); pq.pop();
   int d = Front.first. u = Front.second:
   if (d > dist[u]) continue:
   for (int j = 0; j < adjlist[u].size(); j++) {</pre>
     pii v = adjlist[u][j];
     if (dist[u] + v.second < dist[v.first]) {</pre>
       dist[v.first] = dist[u] + v.second:
       pq.push(MP(dist[v.first], v.first));
 }
```

1.12.5 dsu on trees (contest version)

```
int cnt[MAXN];
bool big[MAXN];
void add(int u, int par, int x){
  cnt[ col[u] ] += x;
  for(auto v: adjlist[u])
    if(v != par && !big[v]) add(v, u, x);
}
void dfs(int u, int par, bool keep){
  int mx_val = -1, bigChild = -1;
  for(auto v: adjlist[u])
    if(v != par && sz[v] > mx_val) mx_val = sz[v],
        bigChild = v;
  for(auto v: adjlist[u])
    if(v != par && v != bigChild)
```

```
dfs(v, u, 0); // run a dfs on small childs and
          clear them from cnt
 if(bigChild != -1)
   dfs(bigChild, u, 1), big[bigChild] = 1; // bigChild
        marked as big and not cleared from cnt
 //now cnt[c] is the number of vertices in subtree of
      vertex u that has color c. You can answer the
      queries easily.
 if(bigChild != -1) big[bigChild] = 0;
 if(keep == 0) add(u, par, -1);
But why it is O(nlogn)? You know that why dsu has
    O(qlogn) time (for q queries); the code uses the
    same method. Merge smaller to greater.
If you have heard heavy-light decomposition you will
    see that function add will go light edges only.
    because of this, code works in O(nlogn) time. Any
    problems of this type can be solved with same dfs
    function and just differs in add function.
```

1.12.6 euler circuit and path(2 in one)

```
///The implementation for euler circuit and euler path
    is exactly same.
///The implementation for directed and undirected graph
    is exactly same.
 1. Path traversing all the edges just once starting
      from one node and ending at that node is euler
 2. If the path does not end at the source node then
      that path is euler path.
1. euler circuit/euler path for undirected graph
 1. The graph must be a single component. So, you have
      to check it.
 2. All the nodes should have even degree. So, check
      it. But if just 2 nodes have odd degree then
      there will an eulerian path only. Choose any of
      these 2 nodes as source node in that case.
 3. If node u has a self loop then deg[u] += 2; i.e.
      the degree of node u is increased by 2.
2. Euler circuit/euler path for directed graph
 1. The graph -if converted to undirected must be a
      single component. So, you have to check it.
 2. in-degree and out-degree of all nodes should be
      equal. So, check it. But if just 2 nodes have
```

```
unequal in-degree and out-degree then there will
      an eulerian path only if and only if one these
      node(source node to be exact) have
      out-degree==1+in-degree and the other have in
      degree==1+out-degree. Choose any of these 2
      nodes as source node in that case.
 3. If node u has a self loop then in-degree[u]++; and
      out-degree[u]++:
vector<pii> gvec[10005]; ///adjacency list; gvec[u]
    contains pairs (v. id) where v is node and id is
    the edge no.
bool bad[10005]: ///size of this array is the number of
vector<int> circuit: ///contains the path
void eularcircuit(int src) {
 stack<int> curr path:
 circuit.clear():
 if(gvec[src].emptv())return;
 int curr v = src:
 while (1) {
   if (!gvec[curr_v].empty()) {
     auto it=gvec[curr_v].back();
     int next_v = it.F, id = it.S;
     gvec[curr_v].pop_back();
     if( bad[id] ) continue;
     bad[id] = 1:
     curr_path.push(curr_v);
     curr_v = next_v;
   else {
     circuit.push_back(curr_v);
     if(curr_path.empty()) break;
     curr_v = curr_path.top();
     curr path.pop():
 reverse(circuit.begin().circuit.end()):
int main() {
 int n, m; /// no. of nodes, no. of edges;
 cin >> n >> m:
 for( int i = 0: i < 10005: ++i ) gvec[i].clear():</pre>
 memset( bad, 0, sizeof(bad) );
 circuit.clear():
 for( int i = 0: i < m: ++i ) {</pre>
   int u, v;
   cin >> u >> v:
   gvec[u].PB( MP(v, i) ); ///Directed graph
   /** Undirected graph
     gvec[u].PB( MP(v, i) );
     gvec[v].PB( MP(u, i) ):
```

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```
}
eularcircuit( source );
}
```

1.12.7 hld

```
///HLD path query + Euler tour(pre-order traversal)
    subtree query
const int MAXN = 1e5+5: /// max no. of nodes:
int n: ///number of nodes:
vector<int> adjlist[MAXN];
int sub[MAXN]: ///sub[u] = size of subtree u:
int head[MAXN]; ///head[u] = head of node u's chain;
int par[MAXN]: ///par[u] = parent of node u: parent of
    root node is -1:
int depth[MAXN]; ///depth[u] = depth of node u;
int tin[MAXN];
int tout[MAXN];
int timer;
No need to clear at the beginning of a test case: sub,
    head, par, depth, tin, tout;
**/
1. Position of node u in segment tree is: tin[u]:
2. Subtree query(u): [tin[u], tout[u]]
3. Path querv(u, v): [tin[u], tin[v]] (here u and v are
    of the same chain and depth[u] <= depth[v])
4. u--v is a heavy edge if and only if depth[u] ==
    depth[v]-1 and adjlist[u][0] = v; Thus u and v are
    of the same chain:
int arr[MAXN]: /// contains the values for the segment
    tree
int tree[MAXN * 4]:
void init_hld( int node, int 1, int r ) {
 if( 1 == r ) {
   tree[node] = arr[1];
   return;
 int mid = (1 + r)/2;
 init_hld( 2*node, 1, mid );
 init_hld( 2*node+1, mid+1, r );
 tree[node] = max( tree[2*node], tree[2*node+1] );
```

```
void update_hld( int node, int 1, int r, int b, int x )
    { ///point update
 if(1 == r) {
   tree[node] = x;
   return:
 int mid = (1 + r)/2;
 if( mid >= b ) update_hld( 2*node, 1, mid, b, x );
 else update_hld(2*node+1, mid+1, r, b, x );
 tree[node] = max( tree[2*node], tree[2*node+1] );
int querv hld( int node, int 1, int r, int b, int e ) {
 if (1 > e \mid | r < b) return -inf:
 if( 1 >= b && r <= e ) return tree[node]:</pre>
 int mid = (1 + r)/2:
 return max( query_hld( 2*node, 1, mid, b, e ),
      querv hld( 2*node+1, mid+1, r, b, e ) );
void dfs sub( int u, int p, int dep ) {
 sub[u] = 1;
 par[u] = p;
 depth[u] = dep;
 int cur = 0;
 for( auto &v: adjlist[u] ) {
   if( v == p ) continue:
   dfs_sub(v, u, dep+1);
   sub[u] += sub[v];
   if(sub[v] > cur) {
     cur = sub[v]:
     swap(v, adjlist[u][0]);
void dfs hld( int u, int p ) {
 tin[u] = timer++:
 for( auto v: adilist[u] ) {
   if( v == p ) continue:
   head[v] = (v == adjlist[u][0] ? head[u] : v);
   dfs hld(v, u):
 }
 tout[u] = timer-1:
void preprocess(int root) {
 timer = 1:
 head[root] = root;
 dfs_sub(root, -1, 0);
 dfs_hld(root, -1);
```

```
int ultimate_query_hld(int u, int v) {
 int ans = -inf:
 while(head[u] != head[v]) {
   if(depth[head[u]] > depth[head[v]]) swap(u, v);
   ans = max(ans, query_hld(1, 1, n, tin[head[v]],
        tin[v]));
   v = par[head[v]];
 if(depth[u] > depth[v]) swap(u, v);
 ans = \max(\text{ans. querv hld}(1, 1, n, tin[u], tin[v])):
 return ans:
void ultimate_update_hld(int u, int v, int x) {
 while(head[u] != head[v]) {
   if(depth[head[u]] > depth[head[v]]) swap(u, v);
   update hld(1, 1, n, tin[head[v]], tin[v], x):
   v = par[head[v]]:
 if(depth[u] > depth[v]) swap(u, v);
 update_hld(1, 1, n, tin[u], tin[v], x);
int lca(int u, int v) {
 while(head[u] != head[v]) {
   if(depth[head[u]] > depth[head[v]]) swap(u, v);
   v = par[head[v]]:
 if(depth[u] > depth[v]) swap(u, v);
 return u;
```

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1.12.8 lca

```
int n, 1;
vector<vector<int>> adjlist;
int timer;
vector<int>> tin, tout;
vector<vector<int>> up;
vector<int>> depth;
void dfs_lca(int u, int par, int dep) {
  depth[u] = dep;
  tin[u] = ++timer;
  up[u][0] = par;
  for (int i = 1; i <= 1; ++i) up[u][i] =
      up[up[u][i-1]][i-1];
  for (auto v : adjlist[u]) {
    if (v != par) dfs_lca(v, u, dep+1);
  }
  tout[u] = ++timer;
}</pre>
```

```
bool is_ancestor(int u, int v) { ///is u an ancestor of
 return tin[u] <= tin[v] && tout[u] >= tout[v];
int lca(int u, int v) {
 if (is_ancestor(u, v)) return u;
 if (is_ancestor(v, u)) return v;
 for (int i = 1; i >= 0; --i) {
   if (!is_ancestor(up[u][i], v)) u = up[u][i];
 return up[u][0]:
int kth ancestor(int u. int k) { ///kth ancestor of u
 for( int i = 1; i >= 0; --i )
   if ((1LL << i) & k) u = up[u][i]:</pre>
   return u:
int get dist( int u, int v ) {
 return depth[u]+depth[v]-2*depth[lca(u,v)]:
void preprocess(int root) {
 tin.resize(n+1);
 tout.resize(n+1):
 depth.resize(n+1);
 timer = 0:
 1 = ceil(log2(n+1));
 up.assign(n+1, vector<int>(1 + 1));
 dfs_lca(root, root, 0);
```

1.12.9 topological sort (tarian)

```
///complexity: 0(N)
vi adjlist[10005];
bool vis[10005];
vector<int> toposort;
void dfs_topo( int u ) {
    vis[u] = 1;
    for( auto v : adjlist[u] ) {
        if( !vis[v] ) dfs_topo(v);
    }
    toposort.PB(u);
}
int main() {
    for( int i = 1; i <= n; ++i ) {
        if( !vis[i] ) dfs_topo(i);
    }
    reverse( toposort.begin(), toposort.end() );
}</pre>
```

1.12.10 virtual tree

```
///Complexity for each query: O(3*klogk) ~ O(klogk); k
    is the number of special nodes in each query;
-Before approaching a virtual tree problem you must be
    able to understand this code!!!! This code is very
    easy. This code can be revised in 5 minutes(worst
    case maybe?)
-But no need to understand the logic of why the code is
    written that wav.
**Brief explanation of the code**
For each query:
1. sort the nodes according to their dfs starting time
2. find the lca of the adjacent nodes and insert them
    to the same container
3. sort the nodes again according to their dfs starting
4. edge banao
-if two nodes u and v belongs to an auxiliary tree then
    lca(u.v) also belongs to that auxiliary tree.
-for each query: for k special nodes there will be
    highest 2*k-1 nodes in the auxiliary tree.
const int mx = 1e5+5: ///max no. of nodes:
/**********LCA code needed here*******/
bool cmp( const int &lhs. const int &rhs ) {
 return tin[lhs] < tin[rhs]: /// tin is the starting
      dfs time of a node
vi aux[mx]; ///auxiliary tree
int main() {
 /**
   Problem:
   You will be given a tree.
   You will be given many queries. In each query you
        will be given k special nodes. Now build an
        auxiliary tree out of these k special nodes.
    Constraints:
```

```
Summation of k over all gueries does not exceed 10<sup>5</sup>
 int q; ///queries;
 cin >> q;
 while(q--) {
   int k; ///how many special nodes?
   vector<int> special; ///this contains the special
   for( int i = 0: i < k: ++i ) {</pre>
     int node:
     cin >> node:
     special.PB(node):
   /**Node selection for the auxiliary tree - start**/
   sort( special.begin(), special.end(), cmp );
        ///sort the nodes according to their starting
        dfs time:
   special.erase( unique(special.begin(),
        special.end()), special.end() );
   for( int i = special.size()-1; i > 0; --i ) {
     special.PB( lca( special[i-1], special[i] ) );
   sort( special.begin(), special.end(), cmp );
        ///sort the nodes according to their starting
        dfs time:
   special.erase( unique(special.begin(),
        special.end()), special.end() );
   /**Node selection for the auxiliary tree - end**/
   /**Edge selection for the auxiliary tree - start**/
   for( auto node : special ) {
     aux[node].clear(); ///clearing the auxiliary tree
          graph
   stack<int> st:
   st.push(special[0]):
   for( int i = 1; i < special.size(); ++i ) {</pre>
     while( !is ancestor(st.top(), special[i]) )
          st.pop();
     int u = st.top(), v = special[i]:
     aux[u].PB(v):
     aux[v].PB(u):
     st.push(special[i]):
   /**Edge selection for the auxiliary tree - end**/
   /**Thus aux contains the auxiliary tree graph**/
}
```

1.13 graph or integer sequence hashing

1.13.1 graph hashing using modulo and prime(Xellos approved)

```
///Graph hashing is not the same as string hashing;
///Be intuitive:
///No idea how this works. But is works:
///***Here double hashing was used. Single hashing will
     also do.
Q) How to hash the path u-v-w?
A) hash value of the path =
     ((po1[u]+po1[v]+po1[w])%MOD1,
     (po2[u]+po2[v]+po2[w])%MOD2)
const int MAXN = 1e5+5; //MAX value of nodes;
int n: ///no. of nodes:
int base1 = 131, base2 = 137; //fairly big prime
ll po1[MAXN], po2[MAXN]; /// po1[u] and po2[u] stores
    the double hash value of node u:
#define MOD1 1000000007
#define MOD2 1000000009
int main() {
 po1[0] = po2[0] = 1;
 for( int i = 1; i <= n; ++i ) {
   po1[i] = po1[i-1]*base1;
   po1[i] %= MOD1;
   po2[i] = po2[i-1]*base2;
   po2[i] %= MOD2;
 }
```

1.13.2 graph hashing using xor(Xellos approved)

```
///Graph hashing is not the same as string hashing;
///Be intuitive;
///No idea how this works. But is works;

///***Here double hashing was used. Single hashing will also do.
/**
Q) How to hash the path u-v-w?
A) There are two ways-
1. hash value of the path = ((xx1[u]+xx1[v]+xx1[w]), (xx2[u]+xx2[v]+xx2[w]))
```

1.14 important C++ features

```
/// 1. Unique vector
sort(vec.begin(), vec.end());
vec.erase( unique( vec.begin(), vec.end() ), vec.end()
    );
/// 2. bit manipulation
inline bool checkBit(ll n, int i) { return n&(1LL<<i); }</pre>
inline 11 setBit(11 n, int i) { return n|(1LL<<i);; }</pre>
inline ll resetBit(ll n. int i) { return n&(~(1LL<<i)):
    }
///3. dir arrav
int dx[] = \{0, 0, +1, -1\}:
int dy[] = \{+1, -1, 0, 0\};
//int dx[] = \{+1, 0, -1, 0, +1, +1, -1, -1\};
//int dv[] = \{0, +1, 0, -1, +1, -1, +1, -1\};
///4. constructor overwriting
struct Vertex {
 int next[K];
 bool leaf = false;
  int p = -1;
  char pch;
  int link = -1;
  int go[K]:
  Vertex(int p=-1, char ch='$') : p(p), pch(ch) {
   fill(begin(next), end(next), -1);
   fill(begin(go), end(go), -1);
```

```
};
///5. substring copy(Very fast. faster than hand-made
     custom, can avoid TLE)
string s;
s.substr(idx); //copy substring of range [idx,
     s.size()-1]
s.substr(idx, x); //copy substring of range [idx,
///6. priority queue in ascending order
priority_queue<int, vector<int>, greater<int> > pq;
///7. random number generator
     rng((unsigned)chrono::system_clock::now().time_since_epoch
    //mt199937 64 for 11 (**use this above main
    function**)
shuffle( vec.begin(), vec.end(), rng ); //shuffles a
int temp = rng(); ///generates a random number
///8. make map faster
ump.reserve(1<<10); ump.max_load_factor(0.25);</pre>
///9. Architectural improvement. (Use this above header
    file) (Noticeable incase of bitset)
#pragma GCC optimize("Ofast")
#pragma GCC
    target("sse,sse2,sse3,ssse3,sse4,popcnt,abm,mmx,avx,avx2,f
#pragma GCC optimize("unroll-loops")
///10. count the number of set bits in a number
__builtin_popcount(num); // for integer;
builtin popcountll(num): // for long long:
```

1.15 matrix exponentiation

```
struct matrix {
  matrix() {}
  vector<vector<ll>> mat;
  matrix(int n, int m) : mat(n, vector<ll>(m)) {}
};
matrix operator * (matrix a, matrix b) { /// pass by
    reference if TLE
  int n = a.mat.size();
  int m = a.mat[0].size();
  assert(b.mat.size() == m);
  int k = b.mat[0].size();
  matrix ans (n, k);
```

```
for (int i=0; i<n; i++)
    for (int j=0; j<k; j++)
        for (int l=0; l<m; l++)
            ans.mat[i][j] = (ans.mat[i][j] +
                 a.mat[i][l]*b.mat[l][j])%MOD;
    return ans;
}
matrix power( matrix b, ll p ) { /// pass by reference
        if TLE
matrix res = I, x = b;
while(p) {
        if (p&1LL ) res = res*x;
        x = x*x;
        p >>= 1LL;
    }
    return res;
}
```

1.16 number theory

1.16.1 Mobius inversion

1.16.2 Tonelli-Shanks(for solving quadratic congruence equation)

```
//solves TIMUS 1132
//can't figure out the time complexity, but its fast
//This function solves quadratic equation modulo prime
//returns -1 if there is no possible solution. i.e the
    congruence equation is not valid.
//returns an answer(if exists) as integer less than
    prime p.
//This algorithm may work for prime power moduli
    (according to wikipedia, but not tested)
//solves x*x == a \pmod{p}:
***Important notes***
**For even prime moduli(p = 2):
1. There is exactly 1 solution in range [1,p-1]. But
    when range is not concerned then there are
    infinite solutions. If a solution is x then the
    other solutions can be defined as k*p+x or k*p-x
    where k is any integer. you can prove it by
    showing that (k*p+x)^2 and x^2 are congruent
    modulo p or (k*p-x)^2 and x^2 are congruent modulo
    p.
```

```
**For odd prime moduli:
1. If the congruence equation is valid then there are
    always exactly 2 solutions in range [1, p-1]. if a
    solution is x(returned by the function) then the
    other solution is p-x. (Exception: if congruence
    equation is valid and x = 0 is returned by the
    function, then p-x = 0. So, even though the
    congruence equation is valid- there are no
    solutions in range [1,p-1]. However, if range is
    not concerned then there are infinitly many
    solutions.)
2.But, if range is not concerned then there exists
    infinite solution if the congruence equation is
    valid. If a solution is x then the other solutions
    can be defined as k*p+x or k*p-x where k is anv
    integer, you can prove it by showing that
    (k*p+x)^2 and x^2 are congruent modulo p or
     (k*p-x)^2 and x^2 are congruent modulo p.
11 expo( ll b, ll power, ll m ) {
 ll res = 1LL, x = b\%m;
 while(power) {
   if ( power&1LL ) res = ( res*x ) % m;
   x = (x*x) \% m;
   power >>= 1LL:
 return res;
//solves x*x == a \pmod{p}; [Here, p is prime. We find a
    possible value of x.]
int solvequadratic(int a, int p) {
 if(p == 2) {
   if(a&1) return 1:
   return 0:
 if(a > p) a = a\%p;
 while(a < 0) a = a+p:
 if(a == 0) return 0;
 if(expo(a,(p-1)/2,p) != 1) return -1:
 int n = 0, k = p-1:
 while(k \% 2 == 0) k/=2, n++:
 while (\exp(q, (p-1)/2, p) != (p-1)) q++;
 int t = \exp(a,(k+1)/2,p);
 int r = \exp(a,k,p);
 while(true) {
   int s = 1;
   int i = 0;
   while(expo(r,s,p) != 1) i+=1, s*=2;
```

```
if(i == 0) return t;
int e = 1<<(n-i-1);
int u = expo(q, k*e,p);
t = ((long long)t*u)%p;
r = ((long long)r*u*u)%p;
}</pre>
```

20

1.16.3 bigmod

```
11 expo( 11 b, 11 p, 11 m ) {
    11 res = 1LL, x = b%m;
    while(p) {
        if ( p&1LL ) res = ( res*x ) % m;
        x = ( x*x ) % m;
        p >>= 1LL;
    }
    return res;
}
```

1.16.4 factorial factorize

```
///How many times prime p occurs in n! (O(lgn))
long long factorialPrimePower (long long n. long long
    p ) {
 long long freq = 0;
 long long x = n;
 while ( x ) {
   freq += x / p;
   x = x / p;
 return freq;
///prime factorization of n! (O(nlgn))
void factFactorize ( long long n ) {
 for ( int i = 0; i < prime.size() && prime[i] <= n;</pre>
      i++ ) {
   11 x = n;
   11 frea = 0:
   while (x) {
     freq += x / prime[i];
     x = x / prime[i]:
   cout << prime[i] << '^' << freq << endl;</pre>
```

1.16.5 mod inverse from 1 to N

```
///mod inverse from 1 to N
///Complexity: O(N)
const int mx = 100005;
ll inv[mx];
void generate_modinv() {
  inv[1] = 1;
  for( int i = 2; i < mx; i++ ) {
    inv[i] = (-(MOD/i) * inv[MOD%i] ) % MOD;
    inv[i] = inv[i] + MOD;
}</pre>
```

1.16.6 pollard rho

```
Range: 10^18 (tested), should be okay up to 2^63-1
miller_rabin(n)
   returns 1 if prime, 0 otherwise
   Identifies 70000 18 digit primes in 1 second on Toph
pollard_rho(n):
   If n is prime, returns n
   Otherwise returns a proper divisor of n
   Able to factorize ~120 18 digit semiprimes in 1
        second on Toph
   Able to factorize ~700 15 digit semiprimes in 1
        second on Toph
At the very beginning, call rho::init(): this makes
    miller_rabin and pollard rho functions to work
(Copied from sh shahin. I slightly modified the
    primeFactorize function to make it a bit faster.
    Changed rand() to rng()).
namespace rhof
 mt.19937
      rng(chrono::steady_clock::now().time_since_epoch().count());
 const int MAXP = 1000010:
 11 sea[MAXP]:
 int primes[MAXP], spf[MAXP];
 inline 11 mod_add(11 x, 11 y, 11 m){
   return (x += y) < m ? x : x - m;
 inline 11 mod_mul(11 x, 11 y, 11 m){
   11 res = x * y - (11)((1d)x * y / m + 0.5) * m;
   return res < 0 ? res + m : res:
 inline 11 mod_pow(11 x, 11 n, 11 m){
   ll res = 1 % m;
   for (; n; n >>= 1){}
```

```
if (n & 1) res = mod_mul(res, x, m);
   x = mod_mul(x, x, m);
  return res;
/// O(k logn logn logn), k = number of rounds
    performed
inline bool miller_rabin(ll n){
 if (n <= 2 || (n & 1 ^ 1)) return (n == 2);</pre>
 if (n < MAXP) return spf[n] == n;</pre>
 11 c. d. s = 0, r = n - 1:
 for (: !(r & 1): r >>= 1, s++) {}
  /// each iteration is a round
  for (int i = 0; primes[i] < n && primes[i] < 32;</pre>
   c = mod pow(primes[i], r, n):
   for (int j = 0; j < s; j++){
     d = mod mul(c, c, n):
     if (d == 1 && c != 1 && c != (n - 1)) return
          false:
     c = d:
   if (c != 1) return false;
  return true;
inline void init(){
  int i, j, k, cnt = 0;
  for (i = 2; i < MAXP; i++){</pre>
   if (!spf[i]) primes[cnt++] = spf[i] = i;
   for (j = 0, k; (k = i * primes[j]) < MAXP; j++){</pre>
     spf[k] = primes[j];
            if(spf[i] == spf[k]) break;
/// Expected complexity O(n^{(1/4)})
ll pollard rho(ll n){
  while (1){
   11 x = rng() \% n, y = x, c = rng() \% n, u = 1, v,
        t = 0:
   11 *px = seq, *py = seq;
   while (1) {
     *pv++ = v = mod add(mod mul(v, v, n), c, n):
     *py++ = y = mod_add(mod_mul(y, y, n), c, n);
     if((x = *px++) == v) break:
     u = mod_mul(u, abs(y - x), n);
     if (!u) return __gcd(v, n);
     if (++t == 32){
      t = 0;
       if ((u = _- gcd(u, n)) > 1 \&\& u < n) return u;
```

```
if (t \&\& (u = \_gcd(u, n)) > 1 \&\& u < n) return u;
 vector <1l> primeFactorize(ll n){
   if (n == 1) return vector <11>();
   if (miller_rabin(n)) return vector<ll> {n};
   vector <11> v, w, vw;
   while (n > 1 && n < MAXP)
     vw.push_back(spf[n]);
     n = spf[n]:
   }
   if (n \ge MAXP) {
     11 x = pollard_rho(n);
         v = primeFactorize(x):
         w = primeFactorize(n / x):
         vw.resize( (int)v.size()+(int)w.size() );
         merge( v.begin(), v.end(), w.begin(), w.end().
             vw.begin() ):
   return vw:
}
```

1.16.7 sieve

1.17 strings

1.17.1 aho corasick

```
///Aho Corasick - O(N); N = sum of lengths of all
    strings in the set;
```

```
/// Vertex means state, state means vertex. They are
    used interchangeably:
1. Proper suffix of root vertex is the root itself. So
    its suffix link will point to itself;
2. Suffix link of vertex of depth 1 will point to the
3. root vertex is 0;
const int ALPHA = 26: ///max no. of alphabets
struct AhoCorasick{
 struct Vertex {
   int edge[ALPHA]; /// edges indicate letters of
        alphabet:
   bool leaf = false; /// leaf = true- means this
        state represents a string of the set:
   int p = -1: /// parent of this state:
   char pch: /// this state's parent ----pch---> this
        state:
   int link = -1: /// suffix link of this state: ///
        link = -1 means suffix link has not been
   int go[ALPHA]; /// go to next state through a char
    Vertex(int p=-1, char ch='$') : p(p), pch(ch) {
     memset( edge, -1, sizeof(edge) );
     memset( go, -1, sizeof(go) );
 };
 vector<Vertex> trie;
 AhoCorasick() : trie(1) {}
 void add string(string const& s) {
   int node = 0:
   for (char ch : s) {
     int ch idx = ch - 'a':
     if (trie[node].edge[ch_idx] == -1) {
       trie[node].edge[ch_idx] = trie.size();
       trie.push_back( Vertex(node, ch) );
     node = trie[node].edge[ch idx]:
   trie[node].leaf = true;
 int get_link(int node) { /// get suffix link;
   if (trie[node].link == -1) {
     if (node == 0 || trie[node].p == 0)
       trie[node].link = 0;
     else
```

1.17.2 manachers

```
void manachers(string s, vector<int> &d1, vector<int>
    &d2) {
 int n = s.size():
 /// odd length palindromes.
 d1.resize(n):
 for (int i = 0, l = 0, r = -1; i < n; i++) {
   int k = (i > r) ? 1 : min(d1[1 + r - i], r - i + 1);
   while (0 \le i - k \&\& i + k \le n \&\& s[i - k] == s[i + k]
        kl) k++:
   d1[i] = k--:
   if (i + k > r) {
    l = i - k:
    r = i + k:
 /// even length palindromes, suppose "abba" is a
      palindrome. Here, 2nd index(0-based indexing) is
      the center.
 d2.resize(n):
 for (int i = 0, l = 0, r = -1; i < n; i++) {
   int k = (i > r) ? 0 : min(d2[1 + r - i + 1], r - i
   while (0 \le i - k - 1 \&\& i + k \le n \&\& s[i - k - 1]
        == s[i + k]) k++;
   d2[i] = k--;
   if (i + k > r) {
    1 = i - k - 1;
     r = i + k:
```

}

1.17.3 suffix array(better version by Ashiquul)

```
/**
1. sa[i] = i'th suffix, i from 0 to n-1
2. everything is in 0'th base
3. lcp[i] = lcp of (i-1)th and ith suffix, i from 0 to
    n-1. (In the code, height[] is the lcp array)
4. adjust the alpha, usually for string ALPHA = 128
     (max ascii value)
5. notice if clearing may cause tle
6. remove range lcp init() if not required
7. rev sa[sa[i]] = i:
const int MAXN = 1010; /// always take 10 extra.
const int ALPHA = 256, LOG = 12; //LOG is log2(MAXN)+3
struct SuffixArray {
 int sa[MAXN].data[MAXN],rnk[MAXN],height[MAXN],n;
  int wa[MAXN],wb[MAXN],wws[MAXN],wv[MAXN];
  int lg[MAXN], rmq[MAXN][LOG], rev_sa[MAXN];
  int cmp(int *r,int a,int b,int 1){
   return (r[a]==r[b]) && (r[a+1]==r[b+1]):
  void DA(int *r,int *sa,int n,int m){
   int i,j,p,*x=wa,*y=wb,*t;
   for(i=0:i<m:i++) wws[i]=0:</pre>
   for(i=0;i<n;i++) wws[x[i]=r[i]]++;</pre>
    for(i=1:i<m:i++) wws[i]+=wws[i-1]:</pre>
    for(i=n-1:i>=0:i--) sa[--wws[x[i]]]=i:
    for(j=1.p=1:p<n:j*=2.m=p) {</pre>
     for(p=0,i=n-j;i<n;i++) y[p++]=i;
     for(i=0;i<n;i++) if(sa[i]>=j) y[p++]=sa[i]-j;
     for(i=0;i<n;i++) wv[i]=x[y[i]];</pre>
     for(i=0:i<m:i++) wws[i]=0:</pre>
     for(i=0:i<n:i++) wws[wv[i]]++:</pre>
     for(i=1:i<m:i++) wws[i]+=wws[i-1]:</pre>
     for(i=n-1:i>=0:i--) sa[--wws[wv[i]]]=v[i]:
     for(t=x,x=y,y=t,p=1,x[sa[0]]=0,i=1;i<n;i++)
       x[sa[i]] = cmp(y,sa[i-1],sa[i],j)?p-1:p++;
  void calheight(int *r,int *sa,int n){
   int i,j,k=0;
   for(i=1;i<=n;i++) rnk[sa[i]]=i;</pre>
   for(i=0:i<n:height[rnk[i++]]=k)</pre>
     for(k?k--:0, j=sa[rnk[i]-1];r[i+k]==r[j+k];k++);
  void suffix_array (string &A) {
   n = A.size():
```

```
for(int i=0;i<max(n+5,ALPHA);i++)</pre>
         sa[i]=data[i]=rnk[i]=height[i]=wa[i]=wb[i]=wws[i]=wv[i]=0;cin >> s;
    for (int i = 0; i < n; i++) data[i] = A[i];</pre>
    DA(data,sa,n+1,ALPHA);
    calheight(data,sa,n);
    for(int i = 0; i < n; i++) sa[i] = sa[i+1],
        height[i] = height[i+1], rev_sa[sa[i]] = i;
    range_lcp_init();
  /** LCP for range : build of rmg table **/
  void range lcp init() {
    for(int i = 0; i < n; i++) rmq[i][0] = height[i];</pre>
   for(int j = 1; j < LOG; j++) {</pre>
     for(int i = 0; i < n; i++) {</pre>
       if (i+(1<<j)-1 < n) rmq[i][i] =</pre>
            \min(\text{rma}[i][i-1],\text{rma}[i+(1<<(i-1))][i-1]):
       else break:
    lg[0] = lg[1] = 0;
    for(int i = 2: i <= n: i++) {
     lg[i] = lg[i/2] + 1;
  /** lcp between l'th to r'th suffix **/
  int query_lcp(int 1, int r) {
    assert(1 <= r):
    assert(1>=0 && 1<n && r>=0 && r<n);
    if(1 == r) return n-sa[1];
    int k = lg[r-l+1];
    return min(rmq[1][k],rmq[r-(1<<k)+1][k]);
}SA:
///substring sort comparator function. (it is used to
     sort all possible substrings of a string)
bool cmp(pair<int.int> &lhs. pair<int.int> &rhs) {
  int 11 = lhs.first, r1 = lhs.second, 12 = rhs.first,
       r2 = rhs.second:
  bool f = 0:
  if(SA.rev sa[12] < SA.rev sa[11]) {</pre>
   swap(11, 12):
   swap(r1, r2):
   f ^= 1:
  int len1 = r1-11+1, len2 = r2-12+1;
  int com = SA.query_lcp(SA.rev_sa[11], SA.rev_sa[12]);
 if(com < min(len1, len2)) return f ^ 1;</pre>
 return (len1 < len2) ^ f:</pre>
int main() {
```

```
string s;
SA.suffix_array(s);
SA.range_lcp_init();
vector<pair<int,int>> vec;
for( int i = 0; i < s.size(); ++i ) {</pre>
 for( int j = i; j < s.size(); ++j ) {</pre>
   sub_string.PB(MP(i, j));
sort(sub string.begin(), sub string.end(), cmp);
     ///substring sorted.
```

1.17.4 z-algorithm

```
vector<int> z_function(string s) {
 int n = (int) s.length();
 vector<int> z(n);
 for( int i = 1, l = 0, r = 0; i < n; ++i ) {
   if (i \le r) z[i] = min (r - i + 1, z[i - 1]);
   while (i + z[i] < n \&\& s[z[i]] == s[i + z[i]])
   if (i + z[i] - 1 > r) 1 = i, r = i + z[i] - 1:
 return z:
```

1.18 ternary search

```
// complexity: O(lgN)
// Be careful about choosing the boundary
What is Unimodal function?
-We are given a function f(x) which is unimodal on an
    interval [1.r]. By unimodal function, we mean one
    of two behaviors of the function:
1. The function strictly increases first, reaches a
     maximum (at a single point or over an interval),
     and then strictly decreases.
 2. The function strictly decreases first, reaches a
     minimum, and then strictly increases.
**/
/// Ternary search on function f
```

```
This part is an unproven concept. Invented by me. Use
      it at your own risk.
 If the extrema of the unimodal function is a segment
      instead of a point then to get the leftmost or
      rightmost extremum return value of the function
      we may consider adding >= instead of > in the
      "if()" part of the code only if necessary.
//1. Finding the minimum function (doubles precision)
int steps = 200:
void ternarv search() {
 double lo = -inf, hi = inf, mid1, mid2;
 for(int i = 0: i < steps: i++) {</pre>
   mid1 = (lo*2+hi) / 3.0:
   mid2 = (1o+2*hi) / 3.0:
   if(f(mid1) > f(mid2)) lo = mid1:
   else hi = mid2:
 double ans = f(lo):
 return ans:
//2. Finding the maximum function (doubles precision)
void ternary_search() {
 double lo = -inf, hi = inf, mid1, mid2;
 for(int i = 0; i < steps; i++) {</pre>
   mid1 = (lo*2+hi) / 3.0;
   mid2 = (1o+2*hi) / 3.0;
   if(f(mid1) > f(mid2)) hi = mid2;
   else lo = mid1:
 double ans = f(lo);
 return ans:
//3. Finding the minimum function (search on integer
    range)
void ternarv search() {
 int lo = -inf, hi = inf, mid1, mid2;
 while(hi - lo > 4) {
   mid1 = (lo + hi) / 2:
   mid2 = (lo + hi) / 2 + 1:
   if(f(mid1) > f(mid2)) lo = mid1:
   else hi = mid2:
 for(int i = lo; i <= hi; i++) ans = min(ans , f(i));</pre>
//4. Finding the maximum function (search on integer
    range)
void ternary_search() {
 int lo = -inf, hi = inf, mid1, mid2;
 while(hi - lo > 4) {
```

```
mid1 = (lo + hi) / 2;
mid2 = (lo + hi) / 2 + 1;
if(f(mid1) > f(mid2)) hi = mid2;
else lo = mid1;
}
ans = inf;
for(int i = lo; i <= hi; i++) ans = max(ans , f(i));</pre>
```

2 omar bhai

2.1 Data Structure

2.1.1 2D Fenwick tree

```
int sum[10][10], row,col;
void update(int x,int y,int val){
   while(x<=row){
      while(y<=col){
            sum[x][y]+=val;
            y+=y&(-y);
      }
      x+=x&(-x);
   }
int getSum(int x,int y){
   int res=0;
   while(x){
      while(y){
      res+=sum[x][y];
      y-=y&(-y);
   }
   x-=x&(-x);
}
return res;
}</pre>
```

2.2 Hashing

```
///***Double Hashing***
///***Complexity: O(N)
struct simpleHash{
  vector<long long>p;
  vector<long long>h;
  long long base, len, mod = (long long)1e9+7;
  simpleHash(){}
  simpleHash(string &str, long long b){
```

```
base = b; len = str.size(); //0 base index str
   p.resize(len.1):
   h.resize(len+1,0);
   for(int i=1; i<len; i++) p[i] = ( p[i-1]*base )%mod;</pre>
   for(int i=1; i<=len; i++) h[i] = (h[i-1]*base +
        (str[i-1]-'a' + 3)) \mbox{mod}:
  long long rangeHash(int 1,int r){ //1 and r inclusive
      and also 0 based
   return ( h[r+1]-( (h[l] * p[r-l+1]) % mod )+ mod) %
 }
}:
struct doubleHashing{
 simpleHash h1.h2:
  doubleHashing(string &str){
   h1 = simpleHash(str. 149):
   h2 = simpleHash(str. 223):
  long long rangeHash(int l,int r){ //l and r inclusive
      and also 0 based
   return ( h1.rangeHash(1,r)<<32LL) ^</pre>
        h2.rangeHash(1,r);
};
```

2.3 Number theory

2.3.1 ONumber theory rules

```
1. If N = (p1^a1) * (p2^a2) * ... * (pN^aN) then
    NOD(N) = (a1+1) * (a2+1) * ... * (aN+1)
2. If N = (p1^a1) * (p2^a2) * ... * (pN^aN) then,
    SOD(N) = ((p1^(a1+1)) - 1) / (p1-1) *...* (
    (pN^{(aN+1)} - 1) / (pN-1)
1. If M & N are co-prime then the formula holds :: [
    Phi ( M ) * Phi ( N ) = Phi ( M*N ) ]
2. The Numbers(a) less than or Equal to N who all
    have [\gcd(a,N) = d] will be Phi(N/d)
3. (sum of all the [Phi(d)) = N]: where d
    represents all the divisors of N
4. For N>2 Phi(N) is always Even
5. Sum of all the Numbers less than or Equal to N that
    are Co Prime with N is [ N * Phi( N ) / 2 ]
6. [ Lcm(1, n) + Lcm(2, n) + . . . + Lcm(1, n)
    = (n/2) * ((sum of all Phi(d) * d) + 1)]:
    d is the divisors of n
1. Mod Inverse can be solved recursively with the
    following formula :: inv[ a ] = (-Floor( Mod / A )
```

```
* inv[ Mod%A ] + Mod) %Mod

1. logB(x) = ( logC(x) / logC(B) ) , [ Here , logB(x)
    means log(x) based B];

2. What does log10(X) means? 10^(fractional part of the
    result) means the leading digit!!!

3. lgamma(x) = log(1)+...log(x-1) // this is like
    magic!!
```

2.3.2 Chinese remainder theorem

```
///*** Finding the x that satisfies all the congruence
    equation***
///*** Complexity: O(N log(N) )***
 basically chinese remainder theorem is solved using
      extended euclid algorithm.
 the problems can be converted to a diophantine
      equation of ax+by=c format.
 x=a1 \pmod{n1}
 x=a2 \pmod{n2}
 we can rewrite these equations as follow:
 x=a1+n1*k1
 x=a2+n2*k2
 from the above we can write
 a1+n1*k1 = a2+n2*k2
 or, n1*(-k1) + n2*k2 = a1-a2 [which is a diophantine
      equation. so the solution will exists if and
      only if (a1-a2)\%gcd(n1,n2)==0
  [changing the equation we get :: x = a1 +
      x'*[(a2-a1)/d]*n1 where d is the gcd of n1 and
      n2. x' is the value we get from extended euclid
struct ChineseRemainderTheorem{
 long long ans,lcm;
 bool haveSolution:
 ChineseRemainderTheorem(){}
 ChineseRemainderTheorem(vector<long
      long>&a.vector<long long>&n){
   haveSolution=true:
   ans=a[0]:
   1cm=n[0]:
   for(int i=1;i<n.size();i++){</pre>
     long long x,v;
     long long gcd=extendedEuclid(lcm,n[i],x,y);
     if((a[i]-ans)%gcd){
       haveSolution=false:
       return;
     ans=(ans + ( ( x * ((a[i]-ans )/gcd) )
          %(n[i]/gcd) ) * lcm);
```

```
lcm=LCM(lcm,n[i],gcd);
     ans=((ans%lcm)+lcm)%lcm:
   //so ans variable holds the required result that
        will satisfy all the congruence equations.
 long long extendedEuclid(long long a,long long b,long
      long &x,long long &y){
   if(!b){
    x=1; y=0;
     return a:
   long long x0.v0:
   long long gcd=extendedEuclid(b,a%b,x0,y0);
   x=v0:
   y=x0-(a/b)*y0;
   return gcd:
 long long LCM(long long x,long long y,long long gcd){
   return (x/gcd)*y;
 }
};
```

2.3.3 Linear Diophantine equation

```
///*** Linear Diophantine[ax+by=c] one actual solution
     and solution count between range (x1,x2) &
     (v1, v2)***
///*** Complexity: O(log N)***
1. Linear Diophantine Equation [ ax+by=c ] will have
    solution only if [c % gcd(a,b) == 0]:
2. The other solutions of [ax+by=c] are for any integer
    value of k [x + k*(b/gcd(a,b)), y -
    k*(a/gcd(a,b))
3. The largest number that can't be expressed in ax+by
    form where ( [a.b.x.v all are >=0] and a.b are
    co-prime) is [ab-a-b].
4. All multiples of gcd(a, b) greater than [lcm(a,
    n)-a-n] are representable in the form [ax+by]
    where [a,b,x,v] all are >=0]
5. There are exactly [(a-1)*(b-1)/2] values that can't
    be expressed in ax+by where [a,b,x,y] all are >=0]
6. Compare [ax + by = gcd(a,b)] with [(a\%b)x + by =
    gcd(a,b)] or [(a - (Floor(a/b)*b))x + by =
    gcd( a,b )] For Recursion.
7. Extended Euclid Algorithm finds the x, y and gcd(
     a,b) of Bzouts lemma [ax + by = gcd(a,b)]
///***important :: this algo will not work if [a==0 ||
    b==0], we will have to compute the result
```

```
separately, in those case***
struct LinearDiophantineEquation{
 int extendedEuclid(int a,int b,long long &x,long long
   if(!b){
     x=1, y=0;
     return a;
   long long x0, v0;
   int gcd=extendedEuclid(b.a%b.x0.v0):
   y=x0-(a/b)*y0;
   return gcd;
 bool getOneSolution(int a,int b,int c,long long
      &x.long long &v.long long &gcd){
   gcd=extendedEuclid(abs(a).abs(b).x.v):
   if(c%gcd)return false:
   x*=(c/gcd):
   v*=(c/gcd):
   if(a<0)x*=-1;
   if(b<0)y*=-1;
   return true:
 void shiftResult(int a,int b,long long &x,long long
      &v,long long cnt){
   x+=b*cnt: //this can cause overflow in int
   y-=a*cnt; //this can cause overflow in int
 int getAllSolution(int a,int b,int c,int xLow,int
      xHigh, int yLow, int yHigh) {
   long long x,v,gcd;
   if(!getOneSolution(a,b,c,x,y,gcd))return 0;
   a/=gcd;
   b/=gcd:
   int sign a=(a>0)?+1:-1:
   int sign b=(b>0)?+1:-1:
   shiftResult(a,b,x,y,(xLow-x)/b);
   if(x<xLow)shiftResult(a,b,x,y,sign_b);</pre>
   if(x>xHigh)return 0:
   long long lx1=x;
   shiftResult(a,b,x,y,(xHigh-x)/b);
   long long rx1=x:
   shiftResult(a,b,x,y,-(yLow-y)/a);
   if(y<yLow)shiftResult(a,b,x,y,-sign_a);</pre>
   if(v>vHigh)return 0:
   long long lx2=x;
   shiftResult(a,b,x,y,-(yHigh-y)/a);
   long long rx2=x;
   if(lx2>rx2)swap(lx2,rx2);
   long long lx=max(lx1,lx2);
   long long rx=min(rx1,rx2);
   if(lx>rx)return 0:
```

```
return (rx-lx)/abs(b) +1;
};
```

2.3.4 Sum of NOD of numbers from 1 to N

```
///*** Complexity: 0( sqrt(N) ) ***
long long SNOD(long long n=10){
  long long res=0;
  long long len=sqrtl(n);
  for(int i=1;i<=len;i++)res+=(n/i)-i; //finding the
      ordered pair like a<b and a*b<n
  res*=2LL; //converting pair count to single value
  res+=len; //adding the equal values like (1,1),
      (2,2)..
  return res;
}</pre>
```

2.4 Rotating Callipers

```
///***Complexity: O(N)
struct RotatingCalipers{
 long double res=1e18;
 int len;
 /*this is just a variation of rotating calipers
      finding the minimum distance of two parallel
      line in which the polygon can pass through*/
 RotatingCalipers(vector<pair<int,int> >&hull){
      //convex hull.
   len=hull.size():
   for(int i=0, j=i+1; i<len; i++){</pre>
     pair<int,int>&p1=hull[i];
     pair<int.int>&p2=hull[(i+1)%len]:
     while(area(p1, p2, hull[i%len]) < area(p1, p2,
         hull[(i+1)%len]))i+=1:
     res=min(res. dist(p1. p2. hull[i%len])):
 //Cartesian Distance betweet point p1 & p2
 typedef pair<long double , long double > pld;
 long double dist(pld p1,pld p2){
   long double xDist=p1.first-p2.first;
   long double vDist=p1.second-p2.second;
   return sqrtl((xDist*xDist)+(vDist*vDist)):
 typedef pair<long long ,long long > pll;
 //returns area*2.0 enclosed by p1 ,p2 ,p3 points
 long long area(pll p1, pll p2, pll p3){
```

```
return llabs(p1.first*p2.second +
                                                               }
      p2.first*p3.second + p3.first*p1.second
                                                             };
    -p1.second*p2.first-p2.second*p3.first-p3.second*p1.first);
//distance from p3 to line going through p1 & p2
long double dist(pld p1, pld p2, pld p3){
                                                                 shaad bhai
 long double up=(p2.second-p1.second)*p3.first
          -(p2.first-p1.first)*p3.second
          +p2.first*p1.second
                                                             3.1 Intersection of Two Path
          -p1.first*p2.second;
  long double down=sqrtl(
       (p2.second-p1.second)*(p2.second-p1.second)
                                                             //Calculating common path between node (a,b) and (c,d)
       +(p2.first-p1.first)*(p2.first-p1.first));
                                                             int main() {
  return fabsl(up/down);
                                                               scanf("%d %d %d %d",&a,&b,&c,&d);
```

```
int t1=lca(a,b),t2=lca(c,d);
if (dep[t1]>dep[t2]) {
    swap(a,c); swap(b,d); swap(t1,t2);
}
int p1=lca(a,c),p2=lca(b,d);
int p3=lca(a,d),p4=lca(b,c);
int k1,k2;
if (p1==t1) k1=p4; else k1=p1;
if (p2==t1) k2=p3; else k2=p2;
if (dis(k1,k2)==1&&t2!=k1) // no intersection else // the path is (k1,k2)
}
```