

Codebook- Team Far Behind

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1 Syntax

1.1 Template

```
1 #include <bits/stdc++.h>
2 using namespace std;
2 #include <ext/pb_ds/assoc_container.hpp>
2 using namespace __gnu_pbds;
2 template<class T> ostream& operator<<(ostream &os, const
2 vector<T> &V) {
2 os << "["; for(auto v : V) os << v << " "; return os;
2 os << "];"
3 template<class L, class R> ostream& operator<<(ostream
4 ostream &os, pair<L,R> P) {
4 return os << "(" << P.first << "," << P.second << ")";
5 #define TRACE
5 #ifdef TRACE
6 #define trace(...) _f(#__VA_ARGS__, __VA_ARGS__)
6 template <typename Arg1> _f(const char* name, Arg1&& arg1){
7 void _f(const char* name, Arg1&& arg1){
7 cout << name << " : " << arg1 << endl;
8 }
8 template <typename Arg1, typename... Args>
8 void _f(const char* names, Arg1&& arg1, Args&&... args){
9 args){
9 const char* comma = strchr(names + 1, ',');cout <<
10 write(names, comma - names)
10 << " : " << arg1 << " | ";_f(comma+1, args...);
10 }
10 #else
11 #define trace(...) 1
11 #endif
12 #define ll long long
12 #define ld long double
13 #define vll vector<ll>
13 #define pll pair<ll,ll>
14 #define vpll vector<pll>
14 #define I insert
14 #define pb push_back
14 #define F first
14 #define S second
14 #define endl "\n"
15 // int mod=1e9+7;
16 inline int mul(int a,int b){return (a*1ll*b)%mod;}
16 inline int add(int a,int b){a+=b;if(a>=mod)a-=mod;return
16 a;}
```



```

ll snk, cnt; // cnt for vis, no need to initialize ←
vis
vector<ll> par, vis;
ll dfs(ll u, ll curr_flow){
    vis[u] = cnt; if(u == snk) return curr_flow;
    if(adj[u].size() == 0) return 0;
    for(ll j=0; j<5; j++){ // random for good ←
        augmentation(**sometimes take time**)
        ll a = rand()%(adj[u].size());
        ll v = adj[u][a];
        if(vis[v] == cnt || capacity[u][v] == 0) ←
            continue;
        par[v] = u;
        ll f = dfs(v, min(curr_flow, capacity[u][v])) ←
            ; if(vis[snk] == cnt) return f;
    }
    for(auto v : adj[u]){
        if(vis[v] == cnt || capacity[u][v] == 0) ←
            continue;
        par[v] = u;
        ll f = dfs(v, min(curr_flow, capacity[u][v])); ←
        if(vis[snk] == cnt) return f;
    }
    return 0;
}
ll maxflow(ll s, ll t) {
    snk = t; ll flow = 0; cnt++;
    par = vll(n, -1); vis = vll(n, 0);
    while(ll new_flow = dfs(s, INF)){
        flow += new_flow; cnt++;
        ll cur = t;
        while(cur != s){
            ll prev = par[cur];
            capacity[prev][cur] -= new_flow;
            capacity[cur][prev] += new_flow;
            cur = prev;
        }
    }
    return flow;
}

```

3.2 Dinic

```

// Time:  $O(m \cdot n^2)$  and for any unit capacity network ←
//  $O(m \cdot n^{1/2})$ 
// Time:  $O(\min(fm, mn^2))$  (f: flow routed)
// (so for bipartite matching as well)
// In practice it is pretty fast for any bipartite ←
// network
// I/O:      n -> vertice; DinicFlow net(n);
//           for(z : edges) net.addEdge(z.F, z.S, cap);
//           max flow = maxFlow(s, t);
// e=(u,v), e.flow represents the effective flow ←
// from u to v
// (i.e f(u->v) - f(v->u)), vertices are 1-indexed
struct edge {

```

```

ll x, y, cap, flow; };
struct DinicFlow {
    // *** change inf accordingly *****
    const ll inf = (1e18);
    vector<edge> e;
    vector<ll> cur, d;
    vector<vector<ll>> adj;
    ll n, source, sink;
    DinicFlow() {}
    DinicFlow(ll v) {
        n = v;
        cur = vector<ll>(n + 1);
        d = vector<ll>(n + 1);
        adj = vector<vector<ll>>(n + 1);
    }
    void addEdge(ll from, ll to, ll cap) {
        edge e1 = {from, to, cap, 0};
        edge e2 = {to, from, 0, 0};
        adj[from].push_back(e.size()); e.push_back(←
            e1);
        adj[to].push_back(e.size()); e.push_back(←
            e2);
    }
    ll bfs() {
        queue<ll> q;
        for(ll i = 0; i <= n; ++i) d[i] = -1;
        q.push(source); d[source] = 0;
        while(!q.empty() and d[sink] < 0) {
            ll x = q.front(); q.pop();
            for(ll i = 0; i < (ll)adj[x].size(); ++i ←
                ) {
                ll id = adj[x][i], y = e[id].y;
                if(d[y] < 0 and e[id].flow < e[id].←
                    cap) {
                    q.push(y); d[y] = d[x] + 1;
                }
            }
        }
        return d[sink] >= 0;
    }
    ll dfs(ll x, ll flow) {
        if(!flow) return 0;
        if(x == sink) return flow;
        for(; cur[x] < (ll)adj[x].size(); ++cur[x]) {
            ll id = adj[x][cur[x]], y = e[id].y;
            if(d[y] != d[x] + 1) continue;
            ll pushed = dfs(y, min(flow, e[id].cap - ←
                e[id].flow));
            if(pushed) {
                e[id].flow += pushed;
                e[id ^ 1].flow -= pushed;
                return pushed;
            }
        }
        return 0;
    }
}

```

```

ll maxFlow(ll src, ll snk) {
    this->source = src; this->sink = snk;
    ll flow = 0;
    while(bfs()) {
        for(ll i = 0; i <= n; ++i) cur[i] = 0;
        while(ll pushed = dfs(source, inf)) {
            flow += pushed;
        }
    }
    return flow;
};

```

3.3 Edmond Karp

```

// running time -  $O(n*m^2)$ 
// The matrix capacity stores the capacity for every pair of vertices. adj
// is the adjacency list of the undirected graph, since we have also to use
// the reversed of directed edges when we are looking for augmenting paths.
// The function maxflow will return the value of the maximal flow. During
// the algorithm the matrix capacity will actually store the residual capacity
// of the network. The value of the flow in each edge will actually no stored,
// but it is easy to extent the implementation - by using an additional matrix
// - to also store the flow and return it.
const ll N = 3e3;
ll n; // number of vertices
ll capacity[N][N]; // adj matrix for capacity
vll adj[N]; // adj list of the corresponding undirected graph(**imp**)
// E = {1-2,2->3,3->2}, adj list should be => {1->2,2->1,2->3,3->2}
// *** vertices are 0-indexed ***
ll INF = (1e18);
ll bfs(ll s, ll t, vector<ll>& parent) {
    fill(parent.begin(), parent.end(), -1);
    parent[s] = -2;
    queue<pair<ll, ll>> q;
    q.push({s, INF});
    while (!q.empty()) {
        ll cur = q.front().first;
        ll flow = q.front().second;
        q.pop();
        for (ll next : adj[cur]) {
            if (parent[next] == -1 && capacity[cur][next]) {
                parent[next] = cur;
                ll new_flow = min(flow, capacity[cur][next]);
            }
        }
    }
    return parent[t] != -1 ? flow : 0;
}

```

```

        if (next == t)
            return new_flow;
        q.push({next, new_flow});
    }
}
return 0;
}
ll maxflow(ll s, ll t) {
    ll flow = 0; ll new_flow;
    vector<ll> parent(n);
    while (new_flow = bfs(s, t, parent)) {
        flow += new_flow;
        ll cur = t;
        while (cur != s) {
            ll prev = parent[cur];
            capacity[prev][cur] -= new_flow;
            capacity[cur][prev] += new_flow;
            cur = prev;
        }
    }
    return flow;
}

```

3.4 Push Relabel

```

// Adjacency list implementation of FIFO push relabel maximum flow
// with the gap relabeling heuristic. This implementation is
// significantly faster than straight Ford-Fulkerson. It solves
// random problems with 10000 vertices and 1000000 edges in a few
// seconds, though it is possible to construct test cases that
// achieve the worst-case.
// Time:  $O(V^3)$ 
// I/O:- addEdge(),src,snk ** vertices are 0-indexed
// - To obtain the actual flow values, look at all edges with
// capacity > 0 (zero capacity edges are residual edges).
struct edge {
    ll from, to, cap, flow, index;
    edge(ll from, ll to, ll cap, ll flow, ll index) :
        from(from), to(to), cap(cap), flow(flow), index(index) {}
};
struct PushRelabel {
    ll n;
    vector<vector<edge>> G;
    vector<ll> excess;
    vector<ll> dist, active, count;
    queue<ll> Q;
    PushRelabel(ll n) : n(n), G(n), excess(n), dist(n), active(n), count(2*n) {}
}

```

```

void addEdge(ll from, ll to, ll cap) {
    G[from].push_back(edge(from, to, cap, 0, G[to].size()));
    if (from == to) G[from].back().index++;
    G[to].push_back(edge(to, from, 0, 0, (ll)G[from].size() - 1));
}

void enqueue(ll v) {
    if (!active[v] && excess[v] > 0) { active[v] = true; Q.push(v); }
}

void push(edge &e) {
    ll amt = min(excess[e.from], e.cap - e.flow);
    if (dist[e.from] <= dist[e.to] || amt == 0) return;
    e.flow += amt;
    G[e.to][e.index].flow -= amt;
    excess[e.to] += amt;
    excess[e.from] -= amt;
    enqueue(e.to);
}

void gap(ll k) {
    for (ll v = 0; v < n; v++) {
        if (dist[v] < k) continue;
        count[dist[v]]--; dist[v] = max(dist[v], n+1);
        count[dist[v]]++; enqueue(v);
    }
}

void relabel(ll v) {
    count[dist[v]]--;
    dist[v] = 2*n;
    for (ll i = 0; i < G[v].size(); i++)
        if (G[v][i].cap - G[v][i].flow > 0)
            dist[v] = min(dist[v], dist[G[v][i].to] + 1);
    count[dist[v]]++;
    enqueue(v);
}

void discharge(ll v) {
    for (ll i = 0; excess[v] > 0 && i < G[v].size(); i++) push(G[v][i]);
    if (excess[v] > 0) {
        if (count[dist[v]] == 1) gap(dist[v]);
        else relabel(v);
    }
}

ll getMaxFlow(ll s, ll t) {
    count[0] = n-1;
    count[n] = 1;
    dist[s] = n;
    active[s] = active[t] = true;
    for (ll i = 0; i < G[s].size(); i++) {
        excess[s] += G[s][i].cap;
        push(G[s][i]);
    }
}

```

```

while (!Q.empty()) {
    ll v = Q.front(); Q.pop();
    active[v] = false; discharge(v);
}

ll totflow = 0;
for (ll i = 0; i < G[s].size(); i++) totflow += G[s][i].flow;
return totflow;
};

```

3.5 MCMF

```

// MCMF Theory:
// 1. If a network with negative costs had no negative cycle it is possible to transform it into one with nonnegative costs. Using  $C_{ij\_new}(\pi) = C_{ij\_old} + \pi(i) - \pi(j)$ , where  $\pi(x)$  is shortest path from s to x in network with an added vertex s. The objective value remains the same ( $z_{new} = z + \text{constant}$ ).  $z(x) = \sum(c_{ij} \cdot x_{ij})$ 
// (x->flow, c->cost, u->cap, r->residual cap).
// 2. Residual Network:  $c_{ji} = -c_{ij}$ ,  $r_{ij} = u_{ij} - x_{ij}$ ,  $r_{ji} = x_{ij}$ .
// 3. Note: If edge (i,j),(j,i) both are there then residual graph will have four edges b/w i,j (pairs of parallel edges).
// 4. let  $x^*$  be a feasible soln, its optimal iff residual network  $G_{x^*}$  contains no negative cost cycle.
// 5. Cycle Cancelling algo => Complexity  $O(n \cdot m^2 \cdot U \cdot C)$  (C->max abs value of cost, U->max cap) (m*U*C iterations).
// 6. Successive shortest path algo => Complexity  $O(n^3 \cdot B)$  /  $O(nm \log n)$  (using heap in Dijkstra) (B -> largest supply node).

// Works for negative costs, but does not work for negative cycles
// Complexity:  $O(\min(E^2 \cdot V \log V, E \log V \cdot \text{flow}))$ 
// to use -> graph G(n), G.add_edge(u,v,cap,cost), G.min_cost_max_flow(s,t)
// ***** INF is used in both flow_type and cost_type so change accordingly
const ll INF = 99999999;
// vertices are 0-indexed
struct graph {
    typedef ll flow_type; // **** flow type ****
    typedef ll cost_type; // **** cost type ****
    struct edge {
        int src, dst;
        flow_type capacity, flow;
        cost_type cost;
        size_t rev;
    };
};

```

```

};
vector<edge> edges;
void add_edge(int src, int dst, flow_type cap, ←
    cost_type cost) {
    adj[src].push_back({src, dst, cap, 0, cost, adj[←
        dst].size()});
    adj[dst].push_back({dst, src, 0, 0, -cost, adj[←
        src].size()-1});
}
int n;
vector<vector<edge>> adj;
graph(int n) : n(n), adj(n) { }
pair<flow_type, cost_type> min_cost_max_flow(int s←
    int t) {
    flow_type flow = 0;
    cost_type cost = 0;
    for (int u = 0; u < n; ++u) // initialize
        for (auto &e: adj[u]) e.flow = 0;
    vector<cost_type> p(n, 0);
    auto rcost = [&](edge e) { return e.cost + p[e.←
        src] - p[e.dst]; };
    for (int iter = 0; ; ++iter) {
        vector<int> prev(n, -1); prev[s] = 0;
        vector<cost_type> dist(n, INF); dist[s] = 0;
        if (iter == 0) { // use Bellman-Ford to remove←
            negative cost edges
        }
        vector<int> count(n); count[s] = 1;
        queue<int> que;
        for (que.push(s); !que.empty(); ) {
            int u = que.front(); que.pop();
            count[u] = -count[u];
            for (auto &e: adj[u]) {
                if (e.capacity > e.flow && dist[e.dst] >←
                    dist[e.src] + rcost(e)) {
                    dist[e.dst] = dist[e.src] + rcost(e);
                    prev[e.dst] = e.rev;
                    if (count[e.dst] <= 0) {
                        count[e.dst] = -count[e.dst] + 1;
                        que.push(e.dst);
                    }
                }
            }
        }
        for(int i=0;i<n;i++) p[i] = dist[i]; // ←
        added it
        continue;
    } else { // use Dijkstra
        typedef pair<cost_type, int> node;
        priority_queue<node, vector<node>, greater<←
            node>> que;
        que.push({0, s});
        while (!que.empty()) {
            node a = que.top(); que.pop();
            if (a.S == t) break;
            if (dist[a.S] > a.F) continue;
            for (auto e: adj[a.S]) {

```

```

                if (e.capacity > e.flow && dist[e.dst] >←
                    a.F + rcost(e)) {
                    dist[e.dst] = dist[e.src] + rcost(e);
                    prev[e.dst] = e.rev;
                    que.push({dist[e.dst], e.dst});
                }
            }
        }
        if (prev[t] == -1) break;
        for (int u = 0; u < n; ++u)
            if (dist[u] < dist[t]) p[u] += dist[u] - ←
                dist[t];
        function<flow_type(int, flow_type)> augment = ←
            [&](int u, flow_type cur) {
                if (u == s) return cur;
                edge &r = adj[u][prev[u]], &e = adj[r.dst][r←
                    .rev];
                flow_type f = augment(e.src, min(e.capacity ←
                    - e.flow, cur));
                e.flow += f; r.flow -= f;
                return f;
            };
        flow_type f = augment(t, INF);
        flow += f;
        cost += f * (p[t] - p[s]);
    }
    return {flow, cost};
};

```

4 Geometry

4.1 Convex Hull

```

// code credits(PT struct) -->> https://github.com/jaehyunp/stanfordacm/blob/master/code/Geometry.cc
double INF = 1e100;
double EPS = 1e-9;
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 ); }
};

```



```

double dot(PT p, PT q)      { return p.x*q.x+p.y*q.y;←
}
double dist2(PT p, PT q)    { return dot(p-q,p-q); }
double dist(PT p, PT q)     { return sqrt(dist2(p,q));←
; }
double cross(PT p, PT q)    { return p.x*q.y-p.y*q.x;←
}
//print a point
ostream &operator<<(ostream &os, const PT &p) {
    return os << "(" << p.x << "," << p.y << ")";
}
//point of reference for making hull (leftmost and ←
    bottommost)
PT firstpoint;
//Returns 0 is x,y,z lie on a line, 1 is x->y->z is ←
    ccw direction and 2 if x->y->z is cw
ll orient(PT x,PT y,PT z){
    PT p,q;
    p=y-x;q=z-y;ld cr=cross(p,q);
    if(abs(cr)<EPS){return 0;}
    else if(cr>0) return 1;
    return 2;
}
//for sorting points in ccw(counter clockwise) ←
    direction w.r.t firstpoint (leftmost and ←
    bottommost)
bool compare(PT x,PT y){
    if(orient(firstpoint,x,y)!=2) return true;return ←
        false;
}
/*takes as input a vector of points containing input←
    points and an empty vector for making hull
the points forming convex hull are pushed in vector ←
    hull
returns hull containing minimum number of points in ←
    ccw order
***remove EPS for making integer hull
*/
void make_hull(vector<PT>& poi,vector<PT>& hull)
{
    pair<ld,ld> bl={INF,INF};
    ll n=poi.size();ll ind;
    for(ll i=0;i<n;i++){
        pair<ld,ld> pp={poi[i].y,poi[i].x};
        if(pp<bl){
            ind=i;bl={poi[i].y,poi[i].x};
        }
    }
    swap(bl.F,bl.S);firstpoint=PT(bl.F,bl.S);
    vector<PT> cons;
    for(ll i=0;i<n;i++){
        if(i==ind) continue;cons.pb(poi[i]);
    }
    sort(cons.begin(),cons.end(),compare);
    hull.pb(firstpoint);ll m;
    for(auto z:cons){
        if(hull.size()<=1){hull.pb(z);continue;}
        PT pr,ppr;bool fl=true;

```

```

        while((m=hull.size())>=2){
            pr=hull[m-1];ppr=hull[m-2];
            ll ch=orient(ppr,pr,z);
            if(ch==1){break;}
            else if(ch==2){hull.pop_back();continue;}
            else {
                ll d1,d2;
                d1=dist2(ppr,pr);d2=dist2(ppr,z);
                if(d1>d2){fl=false;break;}else {hull.pop_back()←
                    ;}
            }
        }
        if(fl){hull.push_back(z);}
    }
    return;
}

```

4.2 Convex Hull Trick

```

/*
maintains upper convex hull of lines ax+b and gives ←
    minimum value at a given x
to add line ax+b: sameoldcht.addline(a,b), to get ←
    min value at x: sameoldcht.getbest(x)
to get maximum value at x add -ax-b as lines instead←
    of ax+b and use -sameoldcht.getbest(x)
*/
const int N = 1e5 + 5;
int n;
int a[N];
int b[N];
long long dp[N];
struct line{
    long long a , b;
    double xleft;
    bool type;
    line(long long _a , long long _b){
        a = _a;
        b = _b;
        type = 0;
    }
    bool operator < (const line &other) const{
        if(other.type){
            return xleft < other.xleft;
        }
        return a > other.a;
    }
};
double meet(line x , line y){
    return 1.0 * (y.b - x.b) / (x.a - y.a);
}
struct cht{
    set < line > hull;
    cht(){
        hull.clear();
    }
    typedef set < line > :: iterator ite;
    bool hasleft(ite node){
        return node != hull.begin();
    }
}

```

```

}
bool hasright(ite node){
    return node != prev(hull.end());
}
void updateborder(ite node){
    if(hasright(node)){
        line temp = *next(node);
        hull.erase(temp);
        temp.xleft = meet(*node , temp);
        hull.insert(temp);
    }
    if(hasleft(node)){
        line temp = *node;
        temp.xleft = meet(*prev(node) , temp);
        hull.erase(node);
        hull.insert(temp);
    }
    else{
        line temp = *node;
        hull.erase(node);
        temp.xleft = -1e18;
        hull.insert(temp);
    }
}
bool useless(line left , line middle , line ←
right){
    double x = meet(left , right);
    double y = x * middle.a + middle.b;
    double ly = left.a * x + left.b;
    return y > ly;
}
bool useless(ite node){
    if(hasleft(node) && hasright(node)){
        return useless(*prev(node) , *node , *←
next(node));
    }
    return 0;
}
void addline(long long a , long long b){
    line temp = line(a , b);
    auto it = hull.lower_bound(temp);
    if(it != hull.end() && it -> a == a){
        if(it -> b > b){
            hull.erase(it);
        }
        else{
            return;
        }
    }
    hull.insert(temp);
    it = hull.find(temp);
    if(useless(it)){
        hull.erase(it);
        return;
    }
    while(hasleft(it) && useless(prev(it))){
        hull.erase(prev(it));
    }
    while(hasright(it) && useless(next(it))){

```

```

        hull.erase(next(it));
    }
    updateborder(it);
}
long long getbest(long long x){
    if(hull.empty()){
        return 1e18;
    }
    line query(0 , 0);
    query.xleft = x;
    query.type = 1;
    auto it = hull.lower_bound(query);
    it = prev(it);
    return it -> a * x + it -> b;
}
};
cht sameoldcht;
int main()
{
    scanf("%d" , &n);
    for(int i = 1 ; i <= n ; ++i){
        scanf("%d" , a + i);
    }
    for(int i = 1 ; i <= n ; ++i){
        scanf("%d" , b + i);
    }
    sameoldcht.addline(b[1] , 0);
    for(int i = 2 ; i <= n ; ++i){
        dp[i] = sameoldcht.getbest(a[i]);
        sameoldcht.addline(b[i] , dp[i]);
    }
    printf("%lld\n" , dp[n]);
}

```

5 Trees

5.1 LCA

```

const int N = int(1e5)+10;
const int LOGN = 20;
set<int> g[N];
int level[N];
int DP[LOGN][N];
int n,m;
/*----- Pre-Processing -----*/
/* Code Cridits : Tanuj Khattar codeforces ←
submission */
void dfs0(int u)
{
    for(auto it=g[u].begin();it!=g[u].end();it++)
        if(*it!=DP[0][u])
        {
            DP[0][*it]=u;
            level[*it]=level[u]+1;
            dfs0(*it);
        }
}

```



```

void preprocess()
{
    level[0]=0;
    DP[0][0]=0;
    dfs0(0);
    for(int i=1;i<LOGN;i++)
        for(int j=0;j<n;j++)
            DP[i][j] = DP[i-1][DP[i-1][j]];
}

int lca(int a,int b)
{
    if(level[a]>level[b])swap(a,b);
    int d = level[b]-level[a];
    for(int i=0;i<LOGN;i++)
        if(d&(1<<i))
            b=DP[i][b];
    if(a==b)return a;
    for(int i=LOGN-1;i>=0;i--)
        if(DP[i][a]!=DP[i][b])
            a=DP[i][a],b=DP[i][b];
    return DP[0][a];
}

int dist(int u,int v)
{
    return level[u] + level[v] - 2*level[lca(u,v)];
}

```

5.2 Centroid Decomposition

```

/*
nx:maximum number of nodes
adj:adjacency list of tree,adj1: adjacency list of ←
    centroid tree
par:parents of nodes in centroid tree,timstamp: ←
    timestamps of nodes when they became centroids (←
    helpful in comparing which of the two nodes ←
    became centroid first)
ssize,vis:utility arrays for storing subtree size ←
    and visit times in dfs
tim: utility for doing dfs (for deciding which nodes←
    to visit)
cntrorder: centroids stored in order in which they ←
    were formed
dist[nx]: vector of vectors with dist[i][0][j]=←
    number of nodes at distance of k in subtree of i ←
    in centroid tree and dist[i][j][k]=number of ←
    nodes at distance k in jth child of i in centroid←
    tree *** (use adj while doing dfs instead of adj1←
    )***
dfs: find subtree sizes visiting nodes starting from←
    root without visiting already formed centroids
dfs1: root- starting node, n- subtree size remaining←
    after removing centroids -> returns centroid in ←
    subtree of root
preprocess: stores all values in dist array
*/
const int nx=1e5;

```

```

vector<int> adj[nx],adj1[nx]; //adj is adjacency ←
    list of tree and adj1 is adjacency list for ←
    centroid tree
int par[nx],timstamp[nx],ssize[nx],vis[nx]; //par is ←
    parent of each node in centroid tree,ssize is ←
    subtree size of each node in centroid tree,vis ←
    and timstamp are auxillary arrays for visit times ←
    in dfs- timstamp contains nonzero values only for ←
    centroids
int tim=1;
vector<int> cntrorder; //contains list of centroids ←
    generated (in order)
vector<vector<int>> > dist[nx];
int dfs(int root)
{
    vis[root]=tim;
    int t=0;
    for(auto i:adj[root])
    {
        if(!timstamp[i]&&vis[i]<tim)
            t+=dfs(i);
    }
    ssize[root]=t+1;return t+1;
}

int dfs1(int root,int n)
{
    vis[root]=tim;pair<int,int> mxc={0,-1};bool poss=←
    true;
    for(auto i:adj[root])
    {
        if(!timstamp[i]&&vis[i]<tim)
            poss&=(ssize[i]<=n/2),mxc=max(mxc,{ssize[i],i});
    }
    if(poss&&(n-ssize[root])<=n/2)return root;
    return dfs1(mxc.second,n);
}

int findc(int root)
{
    dfs(root);
    int n=ssize[root];tim++;
    return dfs1(root,n);
}

void cntrdecom(int root,int p)
{
    int cntr=findc(root);
    cntrorder.push_back(cntr);
    timstamp[cntr]=tim++;
    par[cntr]=p;
    if(p>=0)adj1[p].push_back(cntr);
    for(auto i:adj[cntr])
        if(!timstamp[i])
            cntrdecom(i,cntr);
}

void dfs2(int root,int nod,int j,int dst)
{
    if(dist[root][j].size()==dst)dist[root][j].←
    push_back(0);
    vis[nod]=tim;
    dist[root][j][dst]+=1;
}

```

```

for(auto i:adj[nod])
{
    if((timestamp[i]<=timestamp[root])||(vis[i]==vis[nod]))↵
        continue;
    vis[i]=tim;dfs2(root,i,j,dst+1);
}
}
void preprocess()
{
    for(int i=0;i<cntorder.size();i++)
    {
        int root=cntorder[i];
        vector<int> temp;
        dist[root].push_back(temp);
        temp.push_back(0);
        ++tim;
        dfs2(root,root,0,0);
        int cnt=0;
        for(int j=0;j<adj[root].size();j++)
        {
            int nod=adj[root][j];
            if(timestamp[nod]<timestamp[root])
                continue;
            dist[root].push_back(temp);
            ++tim;
            dfs2(root,nod,++cnt,1);
        }
    }
}

```

6 Maths

6.1 Chinese Remainder Theorem

```

#include<bits/stdc++.h>
using namespace std;
#define ll long long
/*solves system of equations x=rem[i]%mods[i] for ↵
any mod (need not be coprime)
input:vector of remainders and moduli
output: pair of answer(x%lcm of modulo) and lcm of ↵
all the modulo (returns -1 if it is inconsistent)↵
*/
ll GCD(ll a, ll b) { return (b == 0) ? a : GCD(b, a ↵
% b); }
inline ll LCM(ll a, ll b) { return a / GCD(a, b) * b↵
; }
inline ll normalize(ll x, ll mod) { x %= mod; if (x ↵
< 0) x += mod; return x; }
struct GCD_type { ll x, y, d; };
GCD_type ex_GCD(ll a, ll b)
{
    if (b == 0) return {1, 0, a};
    GCD_type pom = ex_GCD(b, a % b);
    return {pom.y, pom.x - a / b * pom.y, pom.d};
}

```

```

pair<ll,ll> CRT(vector<ll> &rem,vector<ll> &mods)
{
    ll n=rem.size();
    ll ans=rem[0];
    ll lcm=mods[0];
    for(ll i=1;i<n;i++)
    {
        auto pom=ex_GCD(lcm,mods[i]);
        ll x1=pom.x;
        ll d=pom.d;
        if((rem[i]-ans)%d!=0) return {-1,0};
        ans=normalize(ans+x1*(rem[i]-ans)/d%(mods[i]↵
)/d)*lcm,lcm*mods[i]/d);
        lcm=LCM(lcm,mods[i]); // you can save time ↵
        by replacing above lcm * n[i] /d by lcm =↵
        lcm * n[i] / d
    }
    return {ans,lcm};
}

```

6.2 Discrete Log

```

// Discrete Log , Baby-Step Giant-Step , e-maxx
// The idea is to make two functions,
// f1(p) , f2(q) and find p,q s.t.
// f1(p) = f2(q) by storing all possible values of ↵
f1,
// and checking for q. In this case  $a^x = b \pmod m$  ↵
) is
// solved by substituting x by p.n-q , where
// is choosen optimally , usually sqrt(m).

// credits : https://cp-algorithms.com/algebra/↵
discrete-log.html
// returns a soln. for  $a^x = b \pmod m$ 
// for given a,b,m . -1 if no. soln.
// complexity :  $O(\sqrt{m} \cdot \log(m))$ 
// use unordered_map to remove log factor.
// IMP : works only if a,m are co-prime. But can be ↵
modified.

int solve (int a, int b, int m) {
    int n = (int) sqrt (m + .0) + 1;
    int an = 1;
    for (int i=0; i<n; ++i)
        an = (an * a) % m;
    map<int,int> vals;
    for (int i=1, cur=an; i<=n; ++i) {
        if (!vals.count(cur))
            vals[cur] = i;
        cur = (cur * an) % m;
    }
    for (int i=0, cur=b; i<=n; ++i) {
        if (vals.count(cur)) {
            int ans = vals[cur] * n - i;
            if (ans < m)

```

```

        return ans;
    }
    cur = (cur * a) % m;
}
return -1;
}

```

6.3 NTT

```

/*
Kevin's different Code: https://s3.amazonaws.com/codechef\_shared/download/Solutions/JUNE15/tester/MOREFB.cpp
****There is no problem that FFT can solve while this NTT cannot
Case1: If the answer would be small choose a small enough NTT prime modulus
Case2: If the answer is large(> ~1e9) FFT would not work anyway due to precision issues
In Case2 use NTT. If max_answer_size=n*(largest_coefficient^2)
So use two or three modulus to solve it
****Compute a*b%mod if a%mod*b%mod would result in overflow in O(log(a)) time:
ll mulmod(ll a, ll b, ll mod) {
    ll res = 0;
    while (a != 0) {
        if (a & 1) res = (res + b) % m;
        a >>= 1;
        b = (b << 1) % m;
    }
    return res;
}
Fastest NTT (can also do polynomial multiplication if max coefficients are upto 1e18 using 2 modulus and CRT)
How to use:
P=A*B
Polynomial1 = A[0]+A[1]*x^1+A[2]*x^2+...+A[n-1]*x^n-1
Polynomial2 = B[0]+B[1]*x^1+B[2]*x^2+...+B[n-1]*x^n-1
P=multiply(A,B)
A and B are not passed by reference because they are changed in multiply function
For CRT after obtaining answer modulo two primes p1 and p2:
x = a1 mod p1, x = a2 mod p2 => x=((a1*(m2^-1)%m1)*m2+(a2*(m1^-1)%m2)*m1)%m1m2
*** Before each call to multiply:
set base=1, roots={0,1}, rev={0,1}, max_base=x (such that if mod=c*(2^k)+1 then x<=k and 2^x is greater than equal to nearest power of 2 of 2*n)
root=primitive_root^((mod-1)/(2^max_base))
For P=A*A use square function
Some useful modulo and examples
mod1=463470593 = 1768*2^18+1 primitive root = 3 => max_base=18, root=3^1768
mod2=469762049 = 1792*2^18+1 primitive root = 3 => max_base=18, root=3^1792
(mod1^-1)%mod2=313174774 (mod2^-1)%mod1=154490124

```

Some prime modulus and primitive root

635437057	11
639631361	6
645922817	3
648019969	17
666894337	5
683671553	3
710934529	17
715128833	3
740294657	3
754974721	11
786432001	7
799014913	13
824180737	5
880803841	26
897581057	3
899678209	7
918552577	5
924844033	5
935329793	3
943718401	7
950009857	7
962592769	7
975175681	17
985661441	3
998244353	3

```

*/
//x = a1 mod m1, x = a2 mod m2, invm2m1 = (m2^-1)%m1, invm1m2 = (m1^-1)%m2, gives x%m1*m2
#define chinese(a1,m1, invm2m1,a2,m2, invm1m2) ((a1 *1ll* invm2m1 % m1 * 1ll*m2 + a2 *1ll* invm1m2 % m2 * 1ll*m1) % (m1 *1ll* m2))
int mod; //reset mod everytime with required modulus
inline int mul(int a,int b){return (a*1ll*b)%mod;}
inline int add(int a,int b){a+=b;if(a>=mod)a-=mod;return a;}
inline int sub(int a,int b){a-=b;if(a<0)a+=mod;return a;}
inline int power(int a,int b){int rt=1;while(b>0){if(b&1)rt=mul(rt,a);a=mul(a,a);b>>=1;}return rt;}
inline int inv(int a){return power(a,mod-2);}
inline void modadd(int &a,int &b){a+=b;if(a>=mod)a-=mod;}

int base = 1;
vector<int> roots = {0, 1};
vector<int> rev = {0, 1};
int max_base=18; //x such that 2^x|(mod-1) and 2^x>max answer size(=2*n)
int root=202376916; //primitive root^((mod-1)/(2^max_base))
void ensure_base(int nbase) {
    if (nbase <= base) {
        return;
    }
    assert(nbase <= max_base);
    rev.resize(1 << nbase);
    for (int i = 0; i < (1 << nbase); i++) {
        rev[i] = (rev[i >> 1] >> 1) + ((i & 1) << (nbase - 1));
    }
    roots.resize(1 << nbase);
    while (base < nbase) {

```

```

int z = power(root, 1 << (max_base - 1 - base));
for (int i = 1 << (base - 1); i < (1 << base); i++) {
    roots[i << 1] = roots[i];
    roots[(i << 1) + 1] = mul(roots[i], z);
}
base++;
}
}

void fft(vector<int> &a) {
    int n = (int) a.size();
    assert((n & (n - 1)) == 0);
    int zeros = __builtin_ctz(n);
    ensure_base(zeros);
    int shift = base - zeros;
    for (int i = 0; i < n; i++) {
        if (i < (rev[i] >> shift)) {
            swap(a[i], a[rev[i] >> shift]);
        }
    }
    for (int k = 1; k < n; k <= 1) {
        for (int i = 0; i < n; i += 2 * k) {
            for (int j = 0; j < k; j++) {
                int x = a[i + j];
                int y = mul(a[i + j + k], roots[j + k]);
                a[i + j] = x + y - mod;
                if (a[i + j] < 0) a[i + j] += mod;
                a[i + j + k] = x - y + mod;
                if (a[i + j + k] >= mod) a[i + j + k] -= mod;
            }
        }
    }
}

vector<int> multiply(vector<int> a, vector<int> b, ←
    int eq = 0) {
    int need = (int) (a.size() + b.size() - 1);
    int nbase = 0;
    while ((1 << nbase) < need) nbase++;
    ensure_base(nbase);
    int sz = 1 << nbase;
    a.resize(sz);
    b.resize(sz);
    fft(a);
    if (eq) b = a; else fft(b);
    int inv_sz = inv(sz);
    for (int i = 0; i < sz; i++) {
        a[i] = mul(mul(a[i], b[i]), inv_sz);
    }
    reverse(a.begin() + 1, a.end());
    fft(a);
    a.resize(need);
    return a;
}

vector<int> square(vector<int> a) {
    return multiply(a, a, 1);
}

```

6.4 Lagrange Interpolation

```

/* Input :
Degree of polynomial: k
Polynomial values at x=0,1,2,3,...,k
Output :
Polynomial value at x
Complexity: O(degree of polynomial)
Works only if the points are equally spaced
*/

ll lagrange(vll& v , int k, ll x, int mod)
{
    if(x <= k)
        return v[x];
    ll inn = 1;
    ll den = 1;
    for(int i = 1; i <= k; i++)
    {
        inn = (inn*(x - i))%mod;
        den = (den*(mod - i))%mod;
    }
    inn = (inn*inv(den % mod))%mod;
    ll ret = 0;
    for(int i = 0; i <= k; i++){
        ret = (ret + v[i]*inn)%mod;
        ll md1 = mod - ((x-i)*(k-i))%mod;
        ll md2 = ((i+1)*(x-i-1))%mod;
        if(i!=k)
            inn = (((inn*md1)%mod)*inv(md2 % mod))%←
                mod;
    }
    return ret;
}

```

6.5 Matrix Struct

```

struct matrix{
    ld B[N][N], n;
    matrix(){n = N; memset(B, 0, sizeof B);}
    matrix(int _n){
        n = _n; memset(B, 0, sizeof B);
    }
    void iden(){
        for(int i = 0; i < n; i++)
            B[i][i] = 1;
    }
    void operator += (matrix M){
        for(int i = 0; i < n; i++)
            for(int j = 0; j < n; j++)
                B[i][j] = add(B[i][j], M.B[i][j]);
    }
    void operator -= (matrix M){
        for(int i = 0; i < n; i++)
            for(int j = 0; j < n; j++)
                B[i][j] = sub(B[i][j], M.B[i][j]);
    }
    void operator *= (ld b){

```

```

        for(int i = 0; i < n; i++)
            for(int j = 0; j < n; j++)
                B[i][j] = mul(b, B[i][j]);
    }
    matrix operator - (matrix M){
        matrix ret = (*this);
        ret -= M; return ret;
    }
    matrix operator + (matrix M){
        matrix ret = (*this);
        ret += M; return ret;
    }
    matrix operator * (matrix M){
        matrix ret = matrix(n); memset(ret.B, 0, ←
        sizeof ret.B);
        for(int i = 0; i < n; i++)
            for(int j = 0; j < n; j++)
                for(int k = 0; k < n; k++){
                    ret.B[i][j] = add(ret.B[i][j], ←
                    mul(B[i][k], M.B[k][j]));
                }
        return ret;
    }
    matrix operator *= (matrix M){ *this = ((*this) ←
    * M); }
    matrix operator * (int b){
        matrix ret = (*this); ret *= b; return ret;
    }
    vector<double> multiply(const vector<double> & v←
    ) const{
        vector<double> ret(n);
        for(int i = 0; i < n; i++)
            for(int j = 0; j < n; j++){
                ret[i] += B[i][j] * v[j];
            }
        return ret;
    }
};

```

6.6 nCr(Non Prime Modulo)

```

// sandwich, jtnydv25 video
// calculates nCr, for small
// non-prime modulo, and (very) big n,r.
ll phimod;
vll pr, prn; vll fact;
ll power(ll a, ll x, ll mod){
    ll ans=1;
    while(x){
        if((1LL)&(x)) ans=(ans*a)%mod;
        a=(a*a)%mod; x>>=1LL;
    }
    return ans;
}
// prime factorization of x.
// pr-> prime ; prn -> it's exponent
void getprime(ll x){

```

```

    pr.clear(); prn.clear();
    ll i,j,k;
    for(i=2; (i*i)<=x; i++){
        k=0; while((x%i)==0){k++; x/=i;}
        if(k>0){pr.pb(i); prn.pb(k);}
    }
    if(x!=1){pr.pb(x); prn.pb(1);}
    return;
}
// factorials are calculated ignoring
// multiples of p.
void primeproc(ll p, ll pe){ // p , p^e
    ll i,d;
    fact.clear(); fact.pb(1); d=1;
    for(i=1; i<pe; i++){
        if(i%p){fact.pb((fact[i-1]*i)%pe);}
        else {fact.pb(fact[i-1]);}
    }
    return;
}
// again note this has ignored multiples of p
ll Bigfact(ll n, ll mod){
    ll a,b,c,d,i,j,k;
    a=n/mod; a%=phimod; a=power(fact[mod-1], a, mod);
    b=n%mod; a=(a*fact[b])%mod;
    return a;
}
// Chinese Remainder Thm.
vll crtval, crtmod;
ll crt(vll &val, vll &mod){
    ll a,b,c,d,i,j,k; b=1;
    for(ll z:mod) b*=z;
    ll ans=0;
    for(i=0; i<mod.size(); i++){
        a=mod[i]; c=b/a;
        d=power(c, (((a/pr[i])*(pr[i]-1))-1), a);
        c=(c*d)%b; c=(c*val[i])%b; ans=(ans+c)%b;
    }
    return ans;
}
// calculate for prime powers and
// take crt. For each prime power,
// first ignore multiples of p,
// and then do recursively, calculating
// the powers of p separately.
ll Bigncr(ll n, ll r, ll mod){
    ll a,b,c,d,i,j,k; ll p, pe;
    getprime(mod); ll Fnum=1; ll Fden;
    crtval.clear(); crtmod.clear();
    for(i=0; i<pr.size(); i++){
        Fnum=1; Fden=1;
        p=pr[i]; pe=power(p, prn[i], 1e17);
        primeproc(p, pe);
        a=1; d=0;
        phimod=(pe*(p-1LL))/p;
        ll n1=n, r1=r, nr=n-r;
        while(n1){
            Fnum=(Fnum*(Bigfact(n1, pe)))%pe;
            Fden=(Fden*(Bigfact(r1, pe)))%pe;

```

```

Fden=(Fden*(Bigfact(nr,pe)))%pe;
d+=n1-(r1+nr);
n1/=p;r1/=p;nr/=p;
}
Fnum=(Fnum*(power(Fden,(phimod-1LL),pe)))%pe;
if(d>=prn[i])Fnum=0;
else Fnum=(Fnum*(power(p,d,pe)))%pe;
crtmod.pb(pe);crtval.pb(Fnum);
}
// you can just iterate instead of crt
// for(i=0;i<mod;i++){
//     bool cg=true;
//     for(j=0;j<crtmod.size();j++){
//         if(i%crtmod[j]!=crtval[j])cg=false;
//     }
//     if(cg)return i;
// }
return crt(crtval,crtmod);
}

```

6.7 Primitive Root Generator

```

/*To find generator of U(p),we check for all
g in [1,p]. But only for powers of the
form phi(p)/p_j, where p_j is a prime factor of
phi(p). Note that p is not prime here.
Existence, if one of these :
1. p = 1,2,4
2. p = q^k, where q -> odd prime.
3. p = 2.(q^k), where q-> odd prime
Note that a.g^(phi(p)) = 1 (mod p)
    b.there are phi(phi(p)) generators if exists.
*/
// Finds "a" generator of U(p),
// multiplicative group of integers mod p.
// here calc_phi returns the totient function for p
// Complexity : O(Ans.log(phi(p)).log(p)) + time for factorizing phi(p).
// By some theorem, Ans = O((log(p))^6). Should be fast generally.
int generator(int p) {
    vector<int> fact;
    int phi = calc_phi(p), n = phi;
    for (int i=2; i*i<=n; ++i)
        if (n % i == 0) {
            fact.push_back(i);
            while (n % i == 0)
                n /= i;
        }
    if (n > 1)
        fact.push_back(n);
    for (int res=2; res<=p; ++res) {
        if(gcd(res,p)!=1)continue;
        bool ok = true;

```

```

        for (size_t i=0; i<fact.size() && ok; ++i)
            ok &= powmod(res, phi / fact[i], p) != 1;
        if (ok) return res;
    }
    return -1;
}

```

7 Strings

7.1 Hashing Theory

If order not imp. and count/frequency imp. use this as hash fn:-
 $(a_1+h)^k + (a_2+h)^k + (a_3+h)^k + (a_4+h)^k \pmod{p}$
 p. Select : h,k,p
 Alternate:
 $((x)^{(a_1)}+(x)^{(a_2)}+\dots+(x)^{(a_k)})\%mod$ where x and mod are fixed and $a_1\dots a_k$ is an unordered set

7.2 Manacher

```

// Same idea as Z_algo, Time : O(n)
// [l,r] represents : boundaries of rightmost detected subpalindrom(with max r)
// takes string s and returns a vector of lengths of odd length palindrom
// centered around that char(e.g abac for 'b' returns 2(not 3))
vll manacher_odd(string s){
    ll n = s.length(); vll d1(n);
    for(ll i = 0, l = 0, r = -1; i<n; i++){
        d1[i] = 1;
        if(i <= r){
            d1[i] = min(r-i+1,d1[l+r-i]); // use prev val
        }
        while(i+d1[i] < n && i-d1[i] >= 0 && s[i+d1[i]] == s[i-d1[i]]) d1[i]++; // trivial matching
        if(r < i+d1[i]-1) l=i-d1[i]+1, r=i+d1[i]-1; // update r
    }
    return d1;
}
// takes string s and returns vector of lengths of even length ...
// (it's centered around the right middle char, bb is centered around the later 'b')
vll manacher_even(string s){
    ll n = s.length(); vll d2(n);
    for(ll i = 0, l = 0, r = -1; i<n; i++){
        d2[i] = 0;
        if(i <= r){
            d2[i] = min(r-i+1,d2[l+r+1-i]);

```



```

    }
    while(i+d2[i] < n && i-d2[i]-1 >= 0 && s[i+d2[i]] == s[i-d2[i]-1]) d2[i]++;
    if(d2[i] > 0 && r < i+d2[i]-1) l=i-d2[i], r=i+d2[i]-1;
}
return d2;
}
// Other mtd : To do both things in one pass, add special char e.g string "abc" => "$a$b$c$"

```

7.3 Suffix Array

```

//code credits - https://cp-algorithms.com/string/suffix-array.html
/*Theory :-
Sorted array of suffixes = sorted array of cyclic shifts of string$.
We consider a prefix of len.  $2^k$  of the cyclic, in the kth iteration.
And find the sorted order, using values for (k-1)th iteration and
kind of radix sort. Could be thought as some kind of binary lifting.
String of len.  $2^k$  -> combination of 2 strings of len.  $2^{k-1}$ , whose
order we know. Just radix sort on pair for next iteration.

Applications :-
Finding the smallest cyclic shift; Finding a substring in a string;
Comparing two substrings of a string; Longest common prefix of two substrings;
Number of different substrings.
*/
vector<ll> sort_cyclic_shifts(string const& s) {
    ll n = s.size();
    const ll alphabet = 256;
    //***** change the alphabet size accordingly and indexing *****
    vector<ll> p(n), c(n), cnt(max(alphabet, n), 0);
    // p -> sorted order of 1-len prefix of each cyclic shift index.
    // c -> class of a index
    // pn -> same as p for kth iteration. ||ly cn.
    for (ll i = 0; i < n; i++)
        cnt[s[i]]++;
    for (ll i = 1; i < alphabet; i++)
        cnt[i] += cnt[i-1];
    for (ll i = 0; i < n; i++)
        p[--cnt[s[i]]] = i;
    c[p[0]] = 0;
    ll classes = 1;
    for (ll i = 1; i < n; i++) {
        if (s[p[i]] != s[p[i-1]])
            classes++;
    }
}

```

```

    c[p[i]] = classes - 1;
}
vector<ll> pn(n), cn(n);
for (ll h = 0; (1 << h) < n; ++h) {
    for (ll i = 0; i < n; i++) { // sorting w.r.t second part.
        pn[i] = p[i] - (1 << h);
        if (pn[i] < 0)
            pn[i] += n;
    }
    fill(cnt.begin(), cnt.begin() + classes, 0);
    for (ll i = 0; i < n; i++)
        cnt[c[pn[i]]]++;
    for (ll i = 1; i < classes; i++)
        cnt[i] += cnt[i-1];
    for (ll i = n-1; i >= 0; i--)
        p[--cnt[c[pn[i]]]] = pn[i]; // sorting w.r.t first (more significant) part.
    cn[p[0]] = 0;
    classes = 1;
    for (ll i = 1; i < n; i++) { // determining new classes in sorted array.
        pair<ll, ll> cur = {c[p[i]], c[(p[i] + (1 << h)) % n]};
        pair<ll, ll> prev = {c[p[i-1]], c[(p[i-1] + (1 << h)) % n]};
        if (cur != prev)
            ++classes;
        cn[p[i]] = classes - 1;
    }
    c.swap(cn);
}
return p;
}

vector<ll> suffix_array_construction(string s) {
    s += "$";
    vector<ll> sorted_shifts = sort_cyclic_shifts(s);
    sorted_shifts.erase(sorted_shifts.begin());
    return sorted_shifts;
}

// For comparing two substring of length l starting at i, j.
// k -  $2^k > l/2$ . check the first  $2^k$  part, if equal
// check last  $2^k$  part. c[k] is the c in kth iter of S.A construction.
int compare(int i, int j, int l, int k) {
    pair<int, int> a = {c[k][i], c[k][(i+1-(1 << k)) % n]};
    pair<int, int> b = {c[k][j], c[k][(j+1-(1 << k)) % n]};
    return a == b ? 0 : a < b ? -1 : 1;
}

/*
Kasai's Algo for LCP construction :
Longest Common Prefix for consecutive suffixes in suffix array.
*/

```

lcp[i]=length of lcp of ith and (i+1)th suffix in the suffix array.

1. Consider suffixes in decreasing order of length.
2. Let $p = s[i...n]$. It will be somewhere in the S. A. We determine its $\text{lcp} = k$.
3. Then lcp of $q = s[(i+1)...n]$ will be atleast $k-1$. Why?
4. Remove the first char of p and its successor in the S.A. These are suffixes with lcp $k-1$.
5. But note that these 2 may not be consecutive in S.A. But however lcp of strings in b/w have to be also atleast $k-1$.

```

*/
vector<ll> lcp_construction(string const& s, vector<ll>
ll> const& p) {
    ll n = s.size();
    vector<ll> rank(n, 0);
    for (ll i = 0; i < n; i++)
        rank[p[i]] = i;

    ll k = 0;
    vector<ll> lcp(n-1, 0);
    for (ll i = 0; i < n; i++) {
        if (rank[i] == n - 1) {
            k = 0;
            continue;
        }
        ll j = p[rank[i] + 1];
        while (i + k < n && j + k < n && s[i+k] == s[j+k])
            k++;
        lcp[rank[i]] = k;
        if (k)
            k--;
    }
    return lcp;
}

```

7.4 Trie

```

const ll AS = 26; // alphabet size
ll go[MAX][AS];
ll cnt[MAX]; ll cn=0;
// cn -> index of next new node
// convert all strings to vll
ll newNode() {
    for (ll i=0; i<AS; i++)
        go[cn][i]=0;
    return cn++;
}
// call newNode once *****
// before adding anything **
void addTrie(vll &x) {
    ll v = 0;
    cnt[v]++;
    for (ll i=0; i<x.size(); i++) {
        ll y=x[i];
        if (go[v][y]==-1)

```

```

        go[v][y]=newNode();
        v=go[v][y];
        cnt[v]++;
    }
}
// returns count of substrings with prefix x
ll getcount(vll &x){
    ll v=0;
    for (i=0; i<x.size(); i++){
        ll y=x[i];
        if (go[v][y]==-1)
            go[v][y]=newNode();
        v=go[v][y];
    }
    return cnt[v];
}

```

7.5 Z-algorithm

```

// [l,r] -> indices of the rightmost segment match
// (the detected segment that ends rightmost(with max r))
// 2 cases -> 1st.  $i \leq r$  :  $z[i]$  is atleast  $\min(r-i+1, z[i-1])$ , then match trivially
// 2nd. o.w compute  $z[i]$  with trivial matching
// update l,r
// Time :  $O(n)$  (asy. behavior), Proof : each iteration of inner while loop make r pointer advance to right,
// Applications: 1) Search substring(text t, pattern p)  $s = p + '$' + t$ .
// 3) String compression( $s = t+t+...+t$ , then find  $|t|$ )
// 2) Number of distinct substrings (in  $O(n^2)$ )
// (useful when appending or deleting characters online from the end or beginning)

vector<ll> z_function(string s) {
    ll n = (ll) s.length();
    vector<ll> z(n);
    for (ll i = 1, l = 0, r = 0; i < n; ++i) {
        if (i <= r)
            z[i] = min(r - i + 1, z[i - l]); // use previous z val
        while (i + z[i] < n && s[z[i]] == s[i + z[i]])
            ++z[i];
        if (i + z[i] - 1 > r)
            l = i, r = i + z[i] - 1; // update rightmost segment matched
    }
    return z;
}

```