

WashU ICPC Notebook

Contents

1	Techniques	
1.1	Various algorithm techniques	2
2	Dynamic Programming	
2.1	Divide and Conquer Optimization	3
2.2	Knuth Optimization	4
2.3	Longest Increasing Subsequence	4
2.4	Sum Over Subsets DP	4
3	Geometry	
3.1	Convex Hull Algorithm	4
3.2	Delaunay Triangulation	4
3.3	Various Geometry Functions	5
4	Graphs	
4.1	BFS	6
4.2	DFS	6
4.3	Dijkstra's Algorithm	6
4.4	Shortest Path (Dijkstra)	7
4.5	Shortest Path (Bellman-Ford)	7
4.6	Topological Sort (Kahn)	7
4.7	Minimum Spanning Tree (Kruskal)	7
4.8	Minimum Spanning Tree (Prim)	7
4.9	Max Flow (Dinic's Algorithm)	7
4.10	Max Flow (Edmonds-Karp Algorithm)	8
4.11	Min Cost Max Flow	9
4.12	Eulerian Path	9
4.13	Hopcroft-Karp Algorithm	9
4.14	2-SAT Kosaraju	10
4.15	Articulation points and bridges	10
4.16	Hungarian	10
4.17	Hungarian Navarro	10
4.18	Max Bipartite Cardinality Matching (Kuhn)	11
4.19	Lowest Common Ancestor	11
4.20	Strongly Connected Components	12
4.21	Strongly Connected Components (Kosaraju)	12
4.22	Tarjan	12
4.23	Zero One BFS	13
4.24	Union-Find Set	13
5	Tree	
5.1	Cartesian Tree	13
5.2	Segment Tree	14
6	Math with Numbers	
6.1	Extended Euclid's Algorithm	14
6.2	Fast Prime Number Sieve	14
6.3	GCD and LCM	15
6.4	Euler Phi/Totient	15
7	Math with Matrices?	
7.1	Modular Linear Equation Solver	15
7.2	Fast Fourier Transform	15
7.3	Gauss-Jordan Elimination	16
8	Strings	
8.1	Aho-Corasick Algorithm	16
8.2	Knuth-Morris-Pratt Algorithm	17
8.3	Suffix Array	17
8.4	Manacher (Longest Palindromic Substring)	18
8.5	Z Function	18
9	Python Stuff...	
10	Geometry	
10.1	Convex Hull	18
11	Dynamic Programming	
11.1	Max Sum Subarray (Kadane's Algorithm)	18
11.2	Longest Common Subsequence	18
11.3	Levenshtein Distance	19
11.4	Longest Increasing Subsequence	19
12	Graphs/Trees	
12.1	Graph structure example for our DFS and BFS algorithms	19
12.2	Breadth-First Search	19
12.3	Breadth-First Search Paths	19
12.4	Breadth-First Search Shortest Path	19
12.5	Depth-First Search	19
12.6	Depth-First Search Paths	19
12.7	Dijkstra's Algorithm	19
12.8	Kruskal's Algorithm (including Merge-Find set)	20
12.9	Bellman-Ford Algorithm	20
12.10	Floyd-Warshall Algorithm	20
12.11	Max Flow (Ford-Fulkerson Algorithm)	20
12.12	Segment Tree	21
13	Math	
13.1	Prime Number Sieve (generator)	22
13.2	GCD and Euler's Totient Function	22
13.3	Miller-Rabin Primality Test	22
13.4	Gauss-Jordan Elimination (Matrix inversion and linear system solving)	22
14	Strings	
14.1	Knuth-Morris-Pratt Algorithm (fast pattern matching)	22
14.2	Rabin-Karp Algorithm (multiple pattern matching)	23
1	Techniques	
1.1	Various algorithm techniques	
	Recursion	
	Divide and conquer	
	Finding interesting points in $N \log N$	
	Greedy algorithm	
	Scheduling	
	Max contiguous subvector sum	
	Invariants	
	Huffman encoding	
	Graph theory	
	Dynamic graphs (extra book-keeping)	
	Breadth first search	
	Depth first search	
	* Normal trees / DFS trees	
	Dijkstra's algorithm	
	MST: Prim's algorithm	
	Bellman-Ford	
	Konig's theorem and vertex cover	
	Min-cost max flow	
	Lovasz toggle	
	Matrix tree theorem	
	Maximal matching, general graphs	
	Hopcroft-Karp	
	Hall's marriage theorem	
	Graphical sequences	
	Floyd-Warshall	
	Eulercykler	
	Flow networks	
	* Augmenting paths	
	* Edmonds-Karp	
	Bipartite matching	
	Min. path cover	
	Topological sorting	
	Strongly connected components	
	2-SAT	
	Cutvertices, cutedges och biconnected components	

- Edge coloring
 - * Trees
- Vertex coloring
 - * Bipartite graphs (\Rightarrow trees)
 - * 3^n (special case of set cover)
- Diameter and centroid
- K'th shortest path
- Shortest cycle
- Dynamic programming
 - Knapsack
 - Coin change
 - Longest common subsequence
 - Longest increasing subsequence
 - Number of paths in a dag
 - Shortest path in a dag
 - Dynprog over intervals
 - Dynprog over subsets
 - Dynprog over probabilities
 - Dynprog over trees
 - 3^n set cover
 - Divide and conquer
 - Knuth optimization
 - Convex hull optimizations
 - RMQ (sparse table a.k.a 2^k -jumps)
 - Bitonic cycle
 - Log partitioning (loop over most restricted)
- Combinatorics
 - Computation of binomial coefficients
 - Pigeon-hole principle
 - Inclusion/exclusion
 - Catalan number
 - Pick's theorem (Area = interior + boundary/2 - 1)
- Number theory
 - Integer parts
 - Divisibility
 - Euclidean algorithm
 - Modular arithmetic
 - * Modular multiplication
 - * Modular inverses
 - * Modular exponentiation by squaring
 - Chinese remainder theorem
 - Fermat's small theorem
 - Euler's theorem
 - Phi function
 - Frobenius number
 - Quadratic reciprocity
 - Pollard-Rho
 - Miller-Rabin
 - Hensel lifting
 - Vieta root jumping
- Game theory
 - Combinatorial games
 - Game trees
 - Mini-max
 - Nim
 - Games on graphs
 - Games on graphs with loops
 - Grundy numbers
 - Bipartite games without repetition
 - General games without repetition
 - Alpha-beta pruning
- Probability theory
- Optimization
 - Binary search
 - Ternary search
 - Unimodality and convex functions
 - Binary search on derivative
- Numerical methods
 - Numeric integration
 - Newton's method
 - Root-finding with binary/ternary search
 - Golden section search
- Matrices
 - Gaussian elimination
 - Exponentiation by squaring
- Sorting
 - Radix sort
- Geometry
 - Coordinates and vectors
 - * Cross product
 - * Scalar product
 - Convex hull
 - Polygon cut
 - Closest pair
 - Coordinate-compression
 - Quadrees
 - KD-trees
 - All segment-segment intersection
- Sweeping
 - Discretization (convert to events and sweep)
 - Angle sweeping
 - Line sweeping

- Discrete second derivatives
- Strings
 - Longest common substring
 - Palindrome subsequences
 - Knuth-Morris-Pratt
 - Tries
 - Rolling polynom hashes
 - Suffix array
 - Suffix tree
 - Aho-Corasick
 - Manacher's algorithm
 - Letter position lists
- Combinatorial search
 - Meet in the middle
 - Brute-force with pruning
 - Best-first (A^*)
 - Bidirectional search
 - Iterative deepening DFS / A^*
- Data structures
 - LCA (2^k -jumps in trees in general)
 - Pull/push-technique on trees
 - Heavy-light decomposition
 - Centroid decomposition
 - Lazy propagation
 - Self-balancing trees
 - Convex hull trick (wcipeg.com/wiki/Convex_hull_trick)
 - Monotone queues / monotone stacks / sliding queues
 - Sliding queue using 2 stacks
 - Persistent segment tree

2 Dynamic Programming

2.1 Divide and Conquer Optimization

```

/*****
 * DIVIDE AND CONQUER OPTIMIZATION ( dp[i][k] = min_{j<k} {dp[j][k-1] + C(j,i)} )
 * Description: searches for bounds to optimal point using the monotonicity condition
 * Condition: L[i][k] <= L[i+1][k]
 * Time Complexity: O(K*N^2) becomes O(K*N*logN)
 * Notation: dp[i][k]: optimal solution using k positions, until position i
 *           L[i][k]: optimal point, smallest j which minimizes dp[i][k]
 *           C(i,j): cost for splitting range [j,i] to j and i
 *****/

const int N = 1e3+5;

ll dp[N][N];

//Cost for using i and j
ll C(ll i, ll j);

void compute(ll l, ll r, ll k, ll optl, ll optr){
    // stop condition
    if(l > r) return;

    ll mid = (l+r)/2;
    //best : cost, pos
    pair<ll,ll> best = {LINF,-1};

    //searchs best: lower bound to right, upper bound to left
    for(ll i = optl; i <= min(mid, optr); i++){
        best = min(best, {dp[i][k-1] + C(i,mid), i});
    }
    dp[mid][k] = best.first;
    ll opt = best.second;

    compute(l, mid-1, k, optl, opt);
    compute(mid + 1, r, k, opt, optr);
}

//Iterate over k to calculate
ll solve(){
    //dimensions of dp[N][K]
    int n, k;

    //Initialize DP
    for(ll i = 1; i <= n; i++){
        //dp[i,1] = cost from 0 to i
        dp[i][1] = C(0, i);
    }

    for(ll l = 2; l <= k; l++){
        compute(1, n, l, 1, n);
    }
}

```

```

    //++ Iterate over i to get min(dp[i][k]), don't forget cost from n to i
    for (ll i=1; i<=n; i++) {
        ll rest = ;
        ans = min(ans, dp[i][k] + rest);
    }
}
*/

```

2.2 Knuth Optimization

```

// Knuth DP Optimization - O(n^3) -> O(n^2)
//
// 1) dp[i][j] = min i<k<j { dp[i][k] + dp[k][j] } + C[i][j]
// 2) dp[i][j] = min k<i { dp[k][j-1] + C[k][i] }
//
// Condition: A[i][j-1] <= A[i][j] <= A[i+1][j]
// A[i][j] is the smallest k that gives an optimal answer to dp[i][j]
//
// reference (pt-br): https://algorithmmarch.wordpress.com/2016/08/12/a-otimizacao-de-pds-e-o-garcom-
// da-maratona/
//
// 1) dp[i][j] = min i<k<j { dp[i][k] + dp[k][j] } + C[i][j]
int n;
int dp[N][N], a[N][N];

// declare the cost function
int cost(int i, int j) {
    // ...
}

void knuth() {
    // calculate base cases
    memset(dp, 63, sizeof(dp));
    for (int i = 1; i <= n; i++) dp[i][i] = 0;

    // set initial a[i][j]
    for (int i = 1; i <= n; i++) a[i][i] = i;

    for (int j = 2; j <= n; j++) {
        for (int i = j; i >= 1; i--) {
            for (int k = a[i][j-1]; k <= a[i+1][j]; k++) {
                ll v = dp[i][k] + dp[k][j] + cost(i, j);

                // store the minimum answer for d[i][k]
                // in case of maximum, use v > dp[i][k]
                if (v < dp[i][j])
                    a[i][j] = k, dp[i][j] = v;
            }
        }
    }

    //++ Iterate over i to get min(dp[i][j]) for each j, don't forget cost from n to
}

// 2) dp[i][j] = min k<i { dp[k][j-1] + C[k][i] }
int n, maxj;
int dp[N][J], a[N][J];

// declare the cost function
int cost(int i, int j) {
    // ...
}

void knuth() {
    // calculate base cases
    memset(dp, 63, sizeof(dp));
    for (int i = 1; i <= n; i++) dp[i][1] = // ...

    // set initial a[i][j]
    for (int i = 1; i <= n; i++) a[i][1] = 1, a[n+1][i] = n;

    for (int j = 2; j <= maxj; j++) {
        for (int i = n; i >= 1; i--) {
            for (int k = a[i][j-1]; k <= a[i+1][j]; k++) {
                ll v = dp[k][j-1] + cost(k, i);

                // store the minimum answer for d[i][k]
                // in case of maximum, use v > dp[i][k]
                if (v < dp[i][j])
                    a[i][j] = k, dp[i][j] = v;
            }
        }
    }

    //++ Iterate over i to get min(dp[i][j]) for each j, don't forget cost from n to
}

```

2.3 Longest Increasing Subsequence

```

// Longest Increasing Subsequence - O(nlogn)
//
// dp(i) = max j<i { dp(j) | a[j] < a[i] } + 1
//
// int dp[N], v[N], n, lis;

memset(dp, 63, sizeof dp);
for (int i = 0; i < n; ++i) {
    // increasing: lower_bound
    // non-decreasing: upper_bound
    int j = lower_bound(dp, dp + lis, v[i]) - dp;
    dp[j] = min(dp[j], v[i]);
    lis = max(lis, j + 1);
}

```

2.4 Sum Over Subsets DP

```

// O(N * 2^N)
// A[i] = initial values
// Calculate F[i] = Sum of A[j] for j subset of i
for (int i = 0; i < (1 << N); i++)
    F[i] = A[i];
for (int i = 0; i < N; i++)
    for (int j = 0; j < (1 << N); j++)
        if (j & (1 << i))
            F[j] += F[j ^ (1 << i)];

```

3 Geometry

3.1 Convex Hull Algorithm

```

bool compare(PT a, PT b) { return a.y < b.y || (a.y == b.y && a.x < b.x); }
double cross(PT o, PT a, PT b)
{
    return (a.x - o.x) * (b.y - o.y) - (a.y - o.y) * (b.x - o.x);
}

vector<PT> ConvexHull(vector<PT> p) {
    int n = p.size();
    vector<PT> h(2 * n);
    sort(p.begin(), p.end(), compare);
    // build lower hull
    for (int i = 0; i < n; i++) {
        while (k >= 2 && cross(h[k-2], h[k-1], p[i]) <= 0) k--;
        h[k++] = p[i];
    }
    // build top hull
    for (int i = n-2; i >= 0; i--) {
        while (k >= t && cross(h[k-2], h[k-1], p[i]) <= 0) k--;
        h[k++] = p[i];
    }
    h.resize(k);
    return h;
}

```

3.2 Delaunay Triangulation

```

/*
Stanford notebook
-----
Delaunay Algorithm Does not handle degenerate cases
Running time: O(n^4)
INPUT: x[] = x-coordinates
       y[] = y-coordinates
OUTPUT: triples = a vector containing m triples
        (indices corresponding to triangle vertices)
-----
*/

typedef double T;

struct triple {

```

```

    int i, j, k;
    triple() {}
    triple(int i, int j, int k) : i(i), j(j), k(k) {}
};

vector<triple> delaunayTriangulation(vector<T>& x, vector<T>& y)
{
    int n = x.size();
    vector<T> z(n);
    vector<triple> ret;

    for (int i = 0; i < n; i++)
        z[i] = x[i] * x[i] + y[i] * y[i];
    for (int i = 0; i < n-2; i++)
    {
        for (int j = i+1; j < n; j++)
        {
            for (int k = i+1; k < n; k++)
            {
                if (j == k) continue;
                double xn = (y[j]-y[i])*(z[k]-z[i]) - (y[k]-y[i])*(z[j]-z[i]);
                double yn = (x[k]-x[i])*(z[j]-z[i]) - (x[j]-x[i])*(z[k]-z[i]);
                double zn = (x[j]-x[i])*(y[k]-y[i]) - (x[k]-x[i])*(y[j]-y[i]);
                bool flag = zn < 0;
                for (int m = 0; flag && m < n; m++)
                    flag = flag && ((x[m]-x[i])*xn + (y[m]-y[i])*yn + (z[m]-z[i])*zn <= 0);
                if (flag) ret.push_back(triple(i, j, k));
            }
        }
    }
    return ret;
}

int main() {
    T xs[]={0, 0, 1, 0.9};
    T ys[]={0, 1, 0, 0.9};
    vector<T> x{xs[0], xs[4]}, y{ys[0], ys[4]};
    vector<triple> tri = delaunayTriangulation(x, y);
    //expected: 0 1 3
    //           0 3 2
    int i;
    for (i = 0; i < tri.size(); i++)
        printf("%d %d %d\n", tri[i].i, tri[i].j, tri[i].k);
    return 0;
}

```

3.3 Various Geometry Functions

```

// Stanford Notebook
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); }
};

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 cross(PT p, PT q) { return p.x*q.y-p.y*q.x; }

// rotate a point CCW or CW around the origin
PT RotateCCW90(PT p) { return PT(-p.y,p.x); }
PT RotateCW90(PT p) { return PT(p.y,-p.x); }
PT RotateCCW(PT p, double t) { return PT(p.x*cos(t)-p.y*sin(t), p.x*sin(t)+p.y*cos(t)); }

// 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;
    r = dot(c-a, b-a)/r;
    if (r < 0) return a;
    if (r > 1) 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,y,z) and plane ax+by+cz=d
double DistancePointPlane(double x, double y, 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; }

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 ||
            dist2(b, c) < EPS || dist2(b, d) < EPS) return true;
        if (dot(c-a, c-b) > 0 && dot(d-a, d-b) > 0 && dot(c-b, d-b) > 0)
            return false;
        return true;
    }
    if (cross(d-a, b-a) * cross(c-a, b-a) > 0)
        return false;
    if (cross(a-c, d-c) * cross(b-c, d-c) > 0)
        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 if
// 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);
}

// compute center of circle given three points
PT ComputeCircleCenter(PT a, PT b, PT c)
{
    b=(a+b)/2;
    c=(a+c)/2;
    return ComputeLineIntersection(b, b+RotateCW90(a-b), c, c+RotateCW90(a-c));
}

// determine if point is in a possibly non-convex polygon (by William
// Randolph Franklin); returns 1 for strictly interior points, 0 for
// strictly exterior points, and 0 or 1 for the remaining points.
// Note that it is possible to convert this into an *exact* test using
// integer arithmetic by taking care of the division appropriately
// (making sure to deal with signs properly) and then by writing exact
// tests for checking point on polygon boundary
bool PointInPolygon(const vector<PT> &p, PT q)
{
    bool c = 0;
    for (int i = 0; i < p.size(); i++)
    {
        int j = (i+1)%p.size();
        if ((p[i].y <= q.y && q.y < p[j].y ||
            p[j].y <= q.y && q.y < p[i].y) &&
            q.x * p[i].x + (p[j].x - p[i].x) * (q.y - p[i].y) / (p[j].y - p[i].y))
            c = !c;
    }
    return c;
}

// determine if point is on the boundary of a polygon
bool PointOnPolygon(const vector<PT> &p, PT q)
{
    for (int i = 0; i < p.size(); i++)
        if (dist2(ProjectPointSegment(p[i], p[(i+1)%p.size()], q), q) < EPS)
            return true;
    return false;
}

```

```
// compute intersection of line through points a and b with
// circle centered at c with radius r > 0
vector<PT> CircleLineIntersection(PT a, PT b, PT c, double r)
{
    vector<PT> ret;
    b = b-a; a = a-c;
    double A = dot(b, b);
    double B = dot(a, b);
    double C = dot(a, a) - r*r;
    double D = B*B - A*C;
    if (D < -EPS) return ret;
    ret.push_back(c+a+b*(-B+sqrt(D+EPS))/A);
    if (D > EPS)
        ret.push_back(c+a+b*(-B-sqrt(D))/A);
    return ret;
}

// compute intersection of circle centered at a with radius r
// with circle centered at b with radius R
vector<PT> CircleCircleIntersection(PT a, PT b, double r, double R)
{
    vector<PT> ret;
    double d = sqrt(dist2(a, b));
    if (d > r+R || d+min(r, R) < max(r, R)) return ret;
    double x = (d*d-R*R+r*r)/(2*d);
    double y = sqrt(r*r-x*x);
    PT v = (b-a)/d;
    ret.push_back(a+v*x + RotateCCW90(v)*y);
    if (y > 0) ret.push_back(a+v*x - RotateCCW90(v)*y);
    return ret;
}

// 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++)
    {
        int j = (i+1) % p.size();
        area += p[i].x*p[j].y - p[j].x*p[i].y;
    }
    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++)
    {
        int j = (i+1) % p.size();
        c = c + (p[i]+p[j])*(p[i].x*p[j].y - p[j].x*p[i].y);
    }
    return c / scale;
}

// tests whether or not a given polygon (in CW or CCW order) is simple
bool IsSimple(const vector<PT> &p)
{
    for (int i = 0; i < p.size(); i++)
    {
        for (int k = i+1; k < p.size(); k++)
        {
            int j = (i+1) % p.size();
            int l = (k+1) % p.size();
            if (i == 1 || j == k) continue;
            if (SegmentsIntersect(p[i], p[j], p[k], p[l]))
                return false;
        }
    }
    return true;
}
}
```

4 Graphs

4.1 BFS

```
/* *****
 * BFS (BREADTH-FIRST SEARCH)
 * ***** */
```

```
* Time complexity: O(V+E)
* Usage: bfs(node)
* Notation: s: starting node
* adj[i]: adjacency list for node i
* vis[i]: visited state for node i (0 or 1)
* *****/

const int N = 1e5+10; // Maximum number of nodes
int dist[N], par[N];
vector<int> adj[N];
queue<int> q;

void bfs (int s) {
    memset(dist, 63, sizeof(dist));
    dist[s] = 0;
    q.push(s);

    while (!q.empty()) {
        int u = q.front(); q.pop();
        for (auto v : adj[u]) if (dist[v] > dist[u] + 1) {
            par[v] = u;
            dist[v] = dist[u] + 1;
            q.push(v);
        }
    }
}
```

4.2 DFS

```
/* *****
 * DFS (DEPTH-FIRST SEARCH)
 * Time complexity: O(V+E)
 * ***** */

// Notation: adj[x]: adjacency list for node x
// vis[i]: visited state for node i (0 or 1)
// *****/

const int N = 1e5+10;
int vis[N];
vector<int> adj[N];

void dfs(int u) {
    vis[u] = 1;
    for (int v : adj[u]) {
        if (!vis[v]) {
            dfs(v);
        }
    }
    // vis[u] = 0;
    // Uncomment the line above if you need to
    // traverse only one path at a time (backtracking)
}
```

4.3 Dijkstra's Algorithm

```
//Dijkstra Algorithm
int t,n,m,s,e;
vector<ii> edges[N]; //pair<NodeEnd,dist>
int distances[N]; // =INF=0x3f3f3f3f
int parent[N]; // =-1

int Dijkstra()
{
    vector<ii> :: iterator it;
    priority_queue<ii, vector<ii>, greater<ii> > pq;
    distances[s]=0;
    pq.push(ii(distances[s],s));
    while(!pq.empty())
    {
        ii p = pq.top();
        pq.pop();
        int d=p.first;
        int a=p.second;
        for(it=edges[a].begin();it!=edges[a].end();++it)
        {
            if(distances[it->first]>distances[a]+it->second)
            {
                distances[it->first]=distances[a]+it->second;
                parent[it->first]=a;
                pq.push(ii(distances[it->first],it->first));
            }
        }
    }
}
```

```

    return distances[e];
}

```

4.4 Shortest Path (Dijkstra)

```

/*****
 * DIJKSTRA'S ALGORITHM (SHORTEST PATH TO A VERTEX)
 * Time complexity:  $O((V+E)\log E)$ 
 * Usage: dist[node]
 * Notation: m:          number of edges
 *           (a, b, w):  edge between a and b with weight w
 *           s:          starting node
 *           par[v]:     parent node of u, used to rebuild the shortest path
 *****/

vector<int> adj[N], adjw[N];
int dist[N];

memset(dist, 63, sizeof(dist));
priority_queue<pii> pq;
pq.push(mp(0,0));

while (!pq.empty()) {
    int u = pq.top().nd;
    int d = -pq.top().st;
    pq.pop();

    if (d > dist[u]) continue;
    for (int i = 0; i < adj[u].size(); ++i) {
        int v = adj[u][i];
        int w = adjw[u][i];
        if (dist[u] + w < dist[v])
            dist[v] = dist[u] + w, pq.push(mp(-dist[v], v));
    }
}

```

4.5 Shortest Path (Bellman-Ford)

```

/*****
 * BELLMAN-FORD ALGORITHM (SHORTEST PATH TO A VERTEX - WITH NEGATIVE COST)
 * Time complexity:  $O(VE)$ 
 * Usage: dist[node]
 * Notation: m:          number of edges
 *           n:          number of vertices
 *           (a, b, w):  edge between a and b with weight w
 *           s:          starting node
 *****/

const int N = 1e4+10; // Maximum number of nodes
vector<int> adj[N], adjw[N];
int dist[N], v, w;

memset(dist, 63, sizeof(dist));
dist[0] = 0;
for (int i = 0; i < n-1; ++i)
    for (int u = 0; u < n; ++u)
        for (int j = 0; j < adj[u].size(); ++j)
            v = adj[u][j], w = adjw[u][j],
            dist[v] = min(dist[v], dist[u] + w);

```

4.6 Topological Sort (Kahn)

```

/*****
 * KAHN'S ALGORITHM (TOPOLOGICAL SORTING)
 *****/

*
 * Time complexity:  $O(V+E)$ 
 * Notation: adj[i]:  adjacency matrix for node i
 *           n:       number of vertices
 *           e:       number of edges
 *           a, b:    edge between a and b
 *           inc:     number of incoming arcs/edges
 *           q:       queue with the independent vertices
 *           tsort:   final topo sort, i.e. possible order to traverse graph
 *****/

vector<int> adj[N];
int inc[N]; // number of incoming arcs/edges

```

```

// undirected graph: inc[v] <= 1
// directed graph:  inc[v] == 0

queue<int> q;
for (int i = 1; i <= n; ++i) if (inc[i] <= 1) q.push(i);

while (!q.empty()) {
    int u = q.front(); q.pop();
    for (int v : adj[u])
        if (inc[v] > 1 and --inc[v] <= 1)
            q.push(v);
}

```

4.7 Minimum Spanning Tree (Kruskal)

```

/*****
 * KRUSKAL'S ALGORITHM (MINIMAL SPANNING TREE - INCREASING EDGE SIZE)
 * Time complexity:  $O(E\log E)$ 
 * Usage: cost, sz[find(node)]
 * Notation: cost: sum of all edges which belong to such MST
 *           sz:   vector of subsets sizes, i.e. size of the subset a node is in
 *****/

// + Union-find

int cost = 0;
vector<pair<int, pair<int, int>>> edges; //mp(dist, mp(node1, node2))

int main () {
    // ...
    sort(edges.begin(), edges.end());
    for (auto e : edges)
        if (find(e.nd.st) != find(e.nd.nd))
            unite(e.nd.st, e.nd.nd), cost += e.st;

    return 0;
}

```

4.8 Minimum Spanning Tree (Prim)

```

// Prim - MST  $O(E\log E)$ 
vi adj[N], adjw[N];
int vis[N];

priority_queue<pii> pq;
pq.push(mp(0, 0));

while (!pq.empty()) {
    int u = pq.top().nd;
    pq.pop();
    if (vis[u]) continue;
    vis[u] = 1;
    for (int i = 0; i < adj[u].size(); ++i) {
        int v = adj[u][i];
        int w = adjw[u][i];
        if (!vis[v]) pq.push(mp(-w, v));
    }
}

```

4.9 Max Flow (Dinic's Algorithm)

```

// Stanford Notebook
typedef long long LL;

struct Edge
{
    int from, to, cap, flow, index;
    Edge(int from, int to, int cap, int flow, int index) :
        from(from), to(to), cap(cap), flow(flow), index(index) {}
    LL rcap() { return cap - flow; }
};

struct Dinic
{
    int N;
    vector<vector<Edge>> > G;
    vector<vector<Edge>*> > Lf;
    vector<int> layer;

```

4.10 Max Flow (Edmonds-Karp Algorithm)

```

vector<int> Q;
Dinic(int N) : N(N), G(N), Q(N) {}
void AddEdge(int from, int to, int cap)
{
    if (from == to) return;
    G[from].push_back(Edge(from, to, cap, 0, G[to].size()));
    G[to].push_back(Edge(to, from, 0, 0, G[from].size() - 1));
}

LL BlockingFlow(int s, int t)
{
    layer.clear();
    layer.resize(N, -1);
    layer[s] = 0;
    Lf.clear(); Lf.resize(N);

    int head = 0, tail = 0;
    Q[tail++] = s;
    while (head < tail)
    {
        int x = Q[head++];
        for (int i = 0; i < G[x].size(); i++)
        {
            Edge &e = G[x][i]; if (e.rcap() <= 0) continue;
            if (layer[e.to] == -1)
            {
                layer[e.to] = layer[e.from] + 1;
                Q[tail++] = e.to;
            }
            if (layer[e.to] > layer[e.from])
            {
                Lf[e.from].push_back(&e);
            }
        }
    }
    if (layer[t] == -1) return 0;
    LL totflow = 0;
    vector<Edge*> P;
    while (!Lf[s].empty())
    {
        int curr = P.empty() ? s : P.back()->to;
        if (curr == t)
        {
            // Augment
            LL amt = P.front()->rcap();
            for (int i = 0; i < P.size(); ++i)
            {
                amt = min(amt, P[i]->rcap());
            }
            totflow += amt;
            for (int i = P.size() - 1; i >= 0; --i)
            {
                P[i]->flow += amt;
                G[P[i]->to][P[i]->index].flow -= amt;
                if (P[i]->rcap() <= 0)
                {
                    Lf[P[i]->from].pop_back();
                    P.resize(i);
                }
            }
        }
        else if (Lf[curr].empty())
        {
            // Retreat
            P.pop_back();
            for (int i = 0; i < N; ++i)
            {
                for (int j = 0; j < Lf[i].size(); ++j)
                {
                    if (Lf[i][j]->to == curr)
                        Lf[i].erase(Lf[i].begin() + j);
                }
            }
        }
        else
        {
            // Advance
            P.push_back(Lf[curr].back());
        }
    }
    return totflow;
}

LL GetMaxFlow(int s, int t)
{
    LL totflow = 0;
    while (LL flow = BlockingFlow(s, t))
        totflow += flow;
    return totflow;
}
};

```

```

/*
Stanford Notebook
MinCostMaxFlow (adjacency matrix, Edmonds and Karp 1972)
This implementation keeps track of forward and reverse edges separately
(so you can set cap[i][j] != cap[j][i]). For a regular max flow, set all edge costs to 0.
Running time, O(|V|^2) cost per augmentation
    max flow: O(|V|^3) augmentations
    min cost max flow: O(|V|^4 * MAX_EDGE_COST) augmentations
INPUT: - graph, constructed using AddEdge()
        - source
        - sink
OUTPUT: - (maximum flow value, minimum cost value)
        - To obtain the actual flow, look at positive values only.
*/

typedef vector<int> VI;
typedef vector<VI> VVI;
typedef long long L;
typedef vector<L> VL;
typedef vector<VL> VVL;
typedef pair<int, int> PII;
typedef vector<PII> VPII;

const L INF = numeric_limits<L>::max() / 4;

struct MinCostMaxFlow {
    int N;
    VVL cap, flow, cost;
    VI found;
    VL dist, pi, width;
    VPII dad;

    MinCostMaxFlow(int N) :
        N(N), cap(N, VL(N)), flow(N, VL(N)), cost(N, VL(N)),
        found(N), dist(N), pi(N), width(N), dad(N) {}

    void AddEdge(int from, int to, L cap, L cost)
    {
        this->cap[from][to] = cap;
        this->cost[from][to] = cost;
    }

    void Relax(int s, int k, L cap, L cost, int dir)
    {
        L val = dist[s] + pi[s] - pi[k] + cost;
        if (cap && val < dist[k])
        {
            dist[k] = val;
            dad[k] = make_pair(s, dir);
            width[k] = min(cap, width[s]);
        }
    }

    L Dijkstra(int s, int t)
    {
        fill(found.begin(), found.end(), false);
        fill(dist.begin(), dist.end(), INF);
        fill(width.begin(), width.end(), 0);
        dist[s] = 0;
        width[s] = INF;

        while (s != -1)
        {
            int best = -1;
            found[s] = true;
            for (int k = 0; k < N; k++)
            {
                if (found[k]) continue;
                Relax(s, k, cap[s][k] - flow[s][k], cost[s][k], 1);
                Relax(s, k, flow[k][s], -cost[k][s], -1);
                if (best == -1 || dist[k] < dist[best]) best = k;
            }
            s = best;
        }

        for (int k = 0; k < N; k++)
            pi[k] = min(pi[k] + dist[k], INF);
        return width[t];
    }

    pair<L, L> GetMaxFlow(int s, int t)
    {
        L totflow = 0, totcost = 0;
        while (L amt = Dijkstra(s, t))
        {
            totflow += amt;
            for (int x = t; x != s; x = dad[x].first)
            {

```

```

        if (dad[x].second == 1)
        {
            flow[dad[x].first][x] += amt;
            totcost += amt * cost[dad[x].first][x];
        }
        else
        {
            flow[x][dad[x].first] -= amt;
            totcost -= amt * cost[x][dad[x].first];
        }
    }
    return make_pair(totflow, totcost);
}
};

```

4.11 Min Cost Max Flow

```

// USE INF = 1e9!

/*****
 * MIN COST MAX FLOW (MINIMUM COST TO ACHIEVE MAXIMUM FLOW)
 * Description: Given a graph which represents a flow network where every edge has
 * a capacity and a cost per unit, find the minimum cost to establish the maximum
 * possible flow from s to t.
 * Note: When adding edge (a, b), it is a directed edge!
 * Usage: min_cost_max_flow()
 * add_edge(from, to, cost, capacity)
 * Notation: flw: max flow
 *           cst: min cost to achieve flw
 * Testcase:
 * add_edge(src, 1, 0, 1); add_edge(1, snk, 0, 1); add_edge(2, 3, 1, INF);
 * add_edge(src, 2, 0, 1); add_edge(2, snk, 0, 1); add_edge(3, 4, 1, INF);
 * add_edge(src, 2, 0, 1); add_edge(3, snk, 0, 1);
 * add_edge(src, 2, 0, 1); add_edge(4, snk, 0, 1); => flw = 4, cst = 3
 *****/

// w: weight or cost, c : capacity
struct edge {int v, f, w, c; };

int n, flw_lmt=INF, src, snk, flw, cst, p[N], d[N], et[N];
vector<edge> e;
vector<int> g[N];

void add_edge(int u, int v, int w, int c) {
    int k = e.size();
    g[u].push_back(k);
    g[v].push_back(k+1);
    e.push_back({ v, 0, w, c });
    e.push_back({ u, 0, -w, 0 });
}

void clear() {
    flw_lmt = INF;
    for(int i=0; i<=n; ++i) g[i].clear();
    e.clear();
}

void min_cost_max_flow() {
    flw = 0, cst = 0;
    while (flw < flw_lmt) {
        memset(et, 0, (n+1) * sizeof(int));
        memset(d, 63, (n+1) * sizeof(int));
        deque<int> q;
        q.push_back(src); d[src] = 0;

        while (!q.empty()) {
            int u = q.front(); q.pop_front();
            et[u] = 2;

            for(int i : g[u]) {
                edge &dir = e[i];
                int v = dir.v;
                if (dir.f < dir.c and d[u] + dir.w < d[v]) {
                    d[v] = d[u] + dir.w;
                    if (et[v] == 0) q.push_back(v);
                    else if (et[v] == 2) q.push_front(v);
                    et[v] = 1;
                    p[v] = i;
                }
            }
        }

        if (d[snk] > INF) break;

        int inc = flw_lmt - flw;
        for (int u=snk; u != src; u = e[p[u]^1].v) {
            edge &dir = e[p[u]];

```

```

            inc = min(inc, dir.c - dir.f);
        }

        for (int u=snk; u != src; u = e[p[u]^1].v) {
            edge &dir = e[p[u]], &rev = e[p[u]^1];
            dir.f += inc;
            rev.f -= inc;
            cst += inc * dir.w;
        }

        if (!inc) break;
        flw += inc;
    }
}

```

4.12 Eulerian Path

```

struct Edge;
typedef list<Edge>::iterator iter;
struct Edge
{
    int next_vertex;
    iter reverse_edge;
    Edge(int next_vertex) :next_vertex(next_vertex) { }
};

const int max_vertices = ;
int num_vertices;
list<Edge> adj[max_vertices]; // adjacency list
vector<int> path;

void find_path(int v)
{
    while(adj[v].size() > 0)
    {
        int vn = adj[v].front().next_vertex;
        adj[vn].erase(adj[v].front().reverse_edge);
        adj[v].pop_front();
        find_path(vn);
    }
    path.push_back(v);
}

void add_edge(int a, int b)
{
    adj[a].push_front(Edge(b));
    iter ita = adj[a].begin();
    adj[b].push_front(Edge(a));
    iter itb = adj[b].begin();
    ita->reverse_edge = itb;
    itb->reverse_edge = ita;
}

```

4.13 Hopcroft-Karp Algorithm

```

#include <vector>

vector<int> g[N];
int r[N], l[N], n, m, e, a, b;

// generate g;

bool dfs(int v)
{
    if(vis[v]) return false;
    vis[v] = true;
    for(int u=0; u<g[v].size(); ++u)
    {
        if(!r[g[v][u]])
        {
            l[v]=g[v][u];
            r[g[v][u]]=v;
            return true;
        }
    }
    for(int u=0; u<g[v].size(); ++u)
    {
        if(dfs(r[g[v][u]]))
        {
            l[v]=g[v][u];
            r[g[v][u]]=v;
            return true;
        }
    }
}

```



```

    }
    return false;
}

void hopcroft_karp()
{
    bool change = true;
    while(change)
    {
        change = false;
        fill(vis, vis+n+1, false);
        for(int i=1; i<=n; ++i)
            if(!l[i])
                change|= dfs(i);
    }
}

```

4.14 2-SAT Kosaraju

```

/*****
 * 2-SAT (TELL WHETHER A SERIES OF STATEMENTS CAN OR CANNOT BE FEASIBLE AT THE SAME TIME)
 *****/

 * Time complexity: O(V+E)
 * Usage: n      -> number of variables, 1-indexed
 *         p = v(i) -> picks the "true" state for variable i
 *         p = nv(i) -> picks the "false" state for variable i, i.e. ~i
 *         add(p, q) -> add clause (p v q) (which also means ~p => q, which also means ~q => p)
 *         run2sat() -> true if possible, false if impossible
 *         val[i]    -> tells if i has to be true or false for that solution
 *****/

int n, vis[2*N], ord[2*N], ordn, cnt, cmp[2*N], val[N];
vector<int> adj[2*N], adjt[2*N];

// for a variable u with idx i
// u is 2*i and !u is 2*i+1
// (a v b) == !a -> b ^ !b -> a

int v(int x) { return 2*x; }
int nv(int x) { return 2*x+1; }

// add clause (a v b)
void add(int a, int b){
    adj[a^1].push_back(b);
    adj[b^1].push_back(a);
    adjt[b].push_back(a^1);
    adjt[a].push_back(b^1);
}

void dfs(int x){
    vis[x] = 1;
    for(auto v : adj[x]) if(!vis[v]) dfs(v);
    ord[ordn++] = x;
}

void dfst(int x){
    cmp[x] = cnt, vis[x] = 0;
    for(auto v : adjt[x]) if(vis[v]) dfst(v);
}

bool run2sat(){
    for(int i = 1; i <= n; i++) {
        if(!vis[v(i)]) dfs(v(i));
        if(!vis[nv(i)]) dfs(nv(i));
    }
    for(int i = ordn-1; i >= 0; i--)
        if(vis[ord[i]]) cnt++, dfst(ord[i]);
    for(int i = 1; i <= n; i++){
        if(cmp[v(i)] == cmp[nv(i)]) return false;
        val[i] = cmp[v(i)] > cmp[nv(i)];
    }
    return true;
}

int main () {
    for (int i = 1; i <= n; i++) {
        if (val[i]); // i-th variable is true
        else       // i-th variable is false
    }
}

```

4.15 Articulation points and bridges

```
// Articulation points and Bridges O(V+E)
```

```

// An articulation (cut vertex), if removed, disconnects the graph
int par[N], art[N], low[N], num[N], ch[N], cnt;

void articulation(int u) {
    low[u] = num[u] = ++cnt;
    for (int v : adj[u]) {
        if (!num[v]) {
            par[v] = u; ch[u]++;
            articulation(v);
            if (low[v] >= num[u]) art[u] = 1;
            if (low[v] > num[u]) { /* u-v bridge */
                low[u] = min(low[u], low[v]);
            }
        } else if (v != par[u]) low[u] = min(low[u], num[v]);
    }
}

for (int i = 0; i < n; ++i) if (!num[i])
    articulation(i), art[i] = ch[i]>1;

```

4.16 Hungarian

```

// Hungarian - O(m*n^2)
// Assignment Problem

int n, m;
int pu[N], pv[N], cost[N][M];
int pairV[N], way[M], minv[M], used[M];

void hungarian() {
    for(int i = 1, j0 = 0; i <= n; i++) {
        pairV[i] = 1;
        memset(minv, 63, sizeof minv);
        memset(used, 0, sizeof used);
        do {
            used[j0] = 1;
            int i0 = pairV[j0], delta = INF, j1;
            for(int j = 1; j <= m; j++) {
                if(used[j]) continue;
                int cur = cost[i0][j] - pu[i0] - pv[j];
                if(cur < minv[j]) minv[j] = cur, way[j] = j0;
                if(minv[j] < delta) delta = minv[j], j1 = j;
            }

            for(int j = 0; j <= m; j++) {
                if(used[j]) pu[pairV[j]] += delta, pv[j] -= delta;
                else minv[j] -= delta;
            }
            j0 = j1;
        } while(pairV[j0] != i);

        do {
            int j1 = way[j0];
            pairV[j0] = pairV[j1];
            j0 = j1;
        } while(j0 != 1);
    }
}

// in main
// for(int j = 1; j <= m; j++)
//     if(pairV[j]) ans += cost[pairV[j]][j];
//

```

4.17 Hungarian Navarro

```

// Hungarian - O(n^2 * m)
template<bool is_max = false, class T = int, bool is_zero_indexed = false>
struct Hungarian {
    bool swap_coord = false;
    int lines, cols;
    T ans;

    vector<int> pairV, way;
    vector<bool> used;
    vector<T> pu, pv, minv;
    vector<vector<T>> cost;

    Hungarian(int _n, int _m) {
        if (_n > _m) {
            swap(_n, _m);
            swap_coord = true;
        }
    }
}

```

```

lines = _n + 1, cols = _m + 1;

clear();
cost.resize(lines);
for (auto& line : cost) line.assign(cols, 0);
}

void clear() {
    pairV.assign(cols, 0);
    way.assign(cols, 0);
    pv.assign(cols, 0);
    pu.assign(lines, 0);
}

void update(int i, int j, T val) {
    if (is_zero_indexed) i++, j++;
    if (is_max) val = -val;
    if (swap_coord) swap(i, j);

    assert(i < lines);
    assert(j < cols);

    cost[i][j] = val;
}

T run() {
    T _INF = numeric_limits<T>::max();
    for (int i = 1, j0 = 0; i < lines; i++) {
        pairV[0] = i;
        minv.assign(cols, _INF);
        used.assign(cols, 0);
        do {
            used[j0] = 1;
            int i0 = pairV[j0], j1;
            T delta = _INF;
            for (int j = 1; j < cols; j++) {
                if (used[j]) continue;
                T cur = cost[i0][j] - pv[j];
                if (cur < minv[j]) minv[j] = cur, way[j] = j0;
                if (minv[j] < delta) delta = minv[j], j1 = j;
            }

            for (int j = 0; j < cols; j++) {
                if (used[j]) pu[pairV[j]] += delta, pv[j] -= delta;
                else minv[j] -= delta;
            }
            j0 = j1;
        } while (pairV[j0]);

        do {
            int j1 = way[j0];
            pairV[j0] = pairV[j1];
            j0 = j1;
        } while (j0);
    }

    ans = 0;
    for (int j = 1; j < cols; j++) if (pairV[j]) ans += cost[pairV[j]][j];

    if (is_max) ans = -ans;
    if (is_zero_indexed) {
        for (int j = 0; j + 1 < cols; j++) pairV[j] = pairV[j + 1], pairV[j]--;
        pairV[cols - 1] = -1;
    }
    if (swap_coord) {
        vector<int> pairV_sub(lines, 0);
        for (int j = 0; j < cols; j++) if (pairV[j] >= 0) pairV_sub[pairV[j]] = j;
        swap(pairV, pairV_sub);
    }

    return ans;
}

template <bool is_max = false, bool is_zero_indexed = false>
struct HungarianMult : public Hungarian<is_max, long double, is_zero_indexed> {
    using super = Hungarian<is_max, long double, is_zero_indexed>;

    HungarianMult(int _n, int _m) : super(_n, _m) {}

    void update(int i, int j, long double x) {
        super::update(i, j, log2(x));
    }
};

```

4.18 Max Bipartite Cardinality Matching (Kuhn)

```

/*****
 * KUHN'S ALGORITHM (FIND GREATEST NUMBER OF MATCHINGS - BIPARTITE GRAPH)
 * Time complexity: O(VE)
 * Notation: ans:      number of matchings
 *           b[j]:     matching edge b[j] <-> j
 *           adj[i]:    adjacency list for node i
 *           vis:       visited nodes
 *           x:         counter to help reuse vis list
 *****/

// TIP: If too slow, shuffle nodes and try again.
int x, vis[N], b[N], ans;

bool match(int u) {
    if (vis[u] == x) return 0;
    vis[u] = x;
    for (int v : adj[u])
        if (!b[v] || match(b[v])) return b[v] = u;
    return 0;
}

for (int i = 1; i <= n; ++i) ++x, ans += match(i);

// Maximum Independent Set on bipartite graph
MIS + MCBM = V

// Minimum Vertex Cover on bipartite graph
MVC = MCBM

```

4.19 Lowest Common Ancestor

```

const int max_nodes, log_max_nodes;

int num_nodes, log_num_nodes, root;
vector<int> children[max_nodes];
// children[i] contains the children of node i int A[max_nodes][log_max_nodes+1]; // A[i][j] is
// the 2^j-th ancestor of node i, or -1 if that ancestor does not exist int L[max_nodes];
// L[i] is the distance between node i and the root

// floor of the binary logarithm of n
int lb(unsigned int n)
{
    if (n==0) return -1;
    int p = 0;
    if (n >= 1<<16) { n >>= 16; p += 16; }
    if (n >= 1<<8) { n >>= 8; p += 8; }
    if (n >= 1<<4) { n >>= 4; p += 4; }
    if (n >= 1<<2) { n >>= 2; p += 2; }
    if (n >= 1<<1) { p += 1; }
    return p;
}

void DFS(int i, int l)
{
    L[i] = l;
    for (int j = 0; j < children[i].size(); j++)
        DFS(children[i][j], l+1);
}

int LCA(int p, int q) {
    // ensure node p is at least as deep as node q
    if (L[p] < L[q]) swap(p, q);

    // "binary search" for the ancestor of node p situated on the same level as q
    for (int i = log_num_nodes; i >= 0; i--)
        if (L[p] - (1<<i) >= L[q])
            p = A[p][i];
    if (p == q) return p;

    // "binary search" for the LCA
    for (int i = log_num_nodes; i >= 0; i--)
    {
        if (A[p][i] != -1 && A[p][i] != A[q][i])
        {
            p = A[p][i];
            q = A[q][i];
        }
    }
    return A[p][0];
}

int main(int argc, char* argv[])
{
    // read num_nodes, the total number of nodes
    log_num_nodes = lb(num_nodes);
    for (int i = 0; i < num_nodes; i++)
    {
        int p;
        // read p, the parent of node i or -1 if node i is the root
    }
}

```

```

    A[i][0] = p;
    if(p != -1) children[p].push_back(i);
    else root = i;
}
// precompute A using dynamic programming
for(int j = 1; j <= log_num_nodes; j++)
    for(int i = 0; i < num_nodes; i++)
        if(A[i][j-1] != -1) A[i][j] = A[A[i][j-1]][j-1];
        else A[i][j] = -1;
// precompute L
DFS(root, 0);
return 0;
}

```

4.20 Strongly Connected Components

```

#include <memory.h>

struct edge
{
    int e, nxt;
};

int V, E;
edge e[MAXE], er[MAXE];
int sp[MAXV], spr[MAXV];
int group_cnt, group_num[MAXV];
bool v[MAXV];
int stk[MAXV];

void fill_forward(int x)
{
    int i;
    v[x]=true;
    for(i=sp[x];i!=e[i].nxt;
        if(!v[e[i].e]) fill_forward(e[i].e);
    stk[++stk[0]]=x;
}

void fill_backward(int x)
{
    int i;
    v[x]=false;
    group_num[x]=group_cnt;
    for(i=spr[x];i!=er[i].nxt;
        if(v[er[i].e]) fill_backward(er[i].e);
}

void add_edge(int v1, int v2) //add edge v1->v2
{
    e[++E].e=v2;
    e[E].nxt=sp[v1];
    sp[v1]=E;
    er[E].e=v1;
    er[E].nxt=spr[v2];
    spr[v2]=E;
}

void SCC()
{
    int i;
    stk[0]=0;
    memset(v, false, sizeof(v));
    for(i=1;i<=V;i++)
        if(!v[i]) fill_forward(i);

    group_cnt=0;
    for(i=stk[0];i>=1;i--)
        if(v[stk[i]]){group_cnt++; fill_backward(stk[i]);}
}

```

4.21 Strongly Connected Components (Kosaraju)

```

/*****
* KOSARAJU'S ALGORITHM (GET EVERY STRONGLY CONNECTED COMPONENTS (SCC))
* Description: Given a directed graph, the algorithm generates a list of every
* strongly connected components. A SCC is a set of points in which you can reach
* every point regardless of where you start from. For instance, cycles can be
* a SCC themselves or part of a greater SCC.
* This algorithm starts with a DFS and generates an array called "ord" which
* stores vertices according to the finish times (i.e. when it reaches "return").
* Then, it makes a reversed DFS according to "ord" list. The set of points
*****/

```

```

* visited by the reversed DFS defines a new SCC.
* One of the uses of getting all SCC is that you can generate a new DAG (Directed
* Acyclic Graph), easier to work with, in which each SCC being a "supernode" of
* the DAG.
* Time complexity: O(V+E)
* Notation: adj[i]: adjacency list for node i
* adjt[i]: reversed adjacency list for node i
* ord: array of vertices according to their finish time
* ordn: ord counter
* scc[i]: supernode assigned to i
* scc_cnt: amount of supernodes in the graph
*****/
const int N = 2e5 + 5;

vector<int> adj[N], adjt[N];
int n, ordn, scc_cnt, vis[N], ord[N], scc[N];

//Directed Version
void dfs(int u) {
    vis[u] = 1;
    for (auto v : adj[u]) if (!vis[v]) dfs(v);
    ord[ordn++] = u;
}

void dfst(int u) {
    scc[u] = scc_cnt, vis[u] = 0;
    for (auto v : adjt[u]) if (vis[v]) dfst(v);
}

// add edge: u -> v
void add_edge(int u, int v) {
    adj[u].push_back(v);
    adjt[v].push_back(u);
}

//Undirected version:
/*
int par[N];

void dfs(int u) {
    vis[u] = 1;
    for (auto v : adj[u]) if(!vis[v]) par[v] = u, dfs(v);
    ord[ordn++] = u;
}

void dfst(int u) {
    scc[u] = scc_cnt, vis[u] = 0;
    for (auto v : adj[u]) if(vis[v] and u != par[v]) dfst(v);
}

// add edge: u -> v
void add_edge(int u, int v) {
    adj[u].push_back(v);
    adj[v].push_back(u);
}
*/

// run kosaraju
void kosaraju() {
    for (int i = 1; i <= n; ++i) if (!vis[i]) dfs(i);
    for (int i = ordn - 1; i >= 0; --i) if (vis[ord[i]]) scc_cnt++, dfst(ord[i]);
}

```

4.22 Tarjan

```

// Tarjan for SCC and Edge Biconnected Componentes - O(n + m)
vector<int> adj[N];
stack<int> st;
bool inSt[N];

int id[N], cmp[N];
int cnt, cmpCnt;

void clear() {
    memset(id, 0, sizeof id);
    cnt = cmpCnt = 0;
}

int tarjan(int n) {
    int low;
    id[n] = low = ++cnt;
    st.push(n), inSt[n] = true;

    for(auto x : adj[n]){
        if(id[x] and inSt[x]) low = min(low, id[x]);
        else if(!id[x]) {

```

```

    int lowx = tarjan(x);
    if(inSt[x])
        low = min(low, lowx);
}

if(low == id[n]){
    while(st.size()){
        int x = st.top();
        inSt[x] = false;
        cmp[x] = cmpCnt;

        st.pop();
        if(x == n) break;
    }
    cmpCnt++;
}
return low;
}

```

4.23 Zero One BFS

```

// 0-1 BFS - O(V+E)

const int N = 1e5 + 5;

int dist[N];
vector<pii> adj[N];
deque<pii> dq;

void zero_one_bfs (int x){
    cl(dist, 63);
    dist[x] = 0;
    dq.push_back({x, 0});
    while(!dq.empty()){
        int u = dq.front().st;
        int ud = dq.front().nd;
        dq.pop_front();
        if(dist[u] < ud) continue;
        for(auto x : adj[u]){
            int v = x.st;
            int w = x.nd;
            if(dist[u] + w < dist[v]){
                dist[v] = dist[u] + w;
                if(w) dq.push_back({v, dist[v]});
                else dq.push_front({v, dist[v]});
            }
        }
    }
}

```

4.24 Union-Find Set

```

struct Edge // MST
{
    int a,b,d;
    bool operator < (const Edge &E) const{
        return this->d < E.d;
    }
};

int ranks[M]; int c[N];

int Find(int x)
{
    int y=x;
    while(y!=c[y])
        y=c[y];
    while(x!=c[x])
    {
        int aux=c[x];
        c[x]=y;
        x=aux;
    }
    return y;
}

void Union(int x,int y)
{
    if(ranks[x]>ranks[y])
        c[x]=y;
    else
        c[y]=x;
    if(ranks[x]==ranks[y])

```

```

        ranks[y]++;
    }

```

5 Tree

5.1 Cartesian Tree

```

struct Tr
{
    Tr *l,*r;
    int key,pr,cnt,val,rev;
    long long sum;
    Tr(int new_key,int new_pr,int new_val)
    {
        rev=0;
        key=new_key;
        cnt=1;
        l=r=NULL;
        pr=new_pr;
        val=new_val;
        sum=new_val;
    }
};

#define T Tr*
T R=NULL;

int cnt (T t)
{
    if(!t) return 0;
    return t->cnt;
}

void upd_cnt(T &t)
{
    if(t) t->cnt=cnt(t->l)+cnt(t->r)+1;
}

long long sum(T t)
{
    if(!t) return 0;
    return t->sum;
}

void upd_sum(T &t)
{
    if(t) t->sum=sum(t->l)+sum(t->r)+t->val;
}

void push(T &t)
{
    if(t && t->rev)
    {
        t->rev=0;
        swap(t->l,t->r);
        upd_sum(t);
        if(t->l) t->l->rev^=1;
        if(t->r) t->r->rev^=1;
    }
}

void split(T t,T &l,T &r,int key,int add)
{
    if(!t)
        return void(l=r=NULL);
    push(t);
    upd_cnt(t);
    int current_key=add+cnt(t->l)+1;
    if(key<=current_key)
        split(t->l,l,t->l,key,add),r=t;
    else
        split(t->r,t->r,r,key,current_key),l=t;
    upd_cnt(t);
    upd_sum(t);
}

void merge(T &t,T l,T r)
{
    push(l);
    push(r);
    if(!l || !r)
        t=l?l:r;
    else if(l->pr>r->pr)
        merge(l->r,l->r,r), t=l;
}

```

```

else
    merge(r->l,l,r->l), t=r;
upd_cnt(t);
upd_sum(t);
}

void insert(T &t,T it,int add)
{
    push(t);
    if(!t)
    {
        t=it;
        upd_cnt(t);
        return;
    }
    upd_sum(t);
    if(it->pr > t->pr)
        split(t,it->l,it->r,it->key,add),t=it;
    else if(it->key>add+cnt(t->l)+1)
        insert(t->r,it,add+cnt(t->l)+1);
    else
        insert(t->l,it,add);
    upd_sum(t);
    upd_cnt(t);
}

void print(T t)
{
    if(!t) return;
    print(t->l);
    cout<<t->val<<" ";
    print(t->r);
}

void reverse(int left,int right)
{
    Tr *t1,*t2,*t3;
    t1=t2=t3=NULL;
    split(R,t1,t2,left,0);
    split(t2,t2,t3,right-left+2,0);
    t2->rev^=1;
    merge(R,t1,t2);
    merge(R,R,t3);
}

void get_sum(int left,int right)
{
    Tr *t1,*t2,*t3;
    t1=t2=t3=NULL;
    split(R,t1,t2,left,0);
    split(t2,t2,t3,right-left+2,0);
    cout<<t2->sum<<"\n";
    merge(R,t1,t2);
    merge(R,R,t3);
}

int n,m, q,a,b;
void example()
{
    ios_base::sync_with_stdio(0);
    cin.tie(0);
    freopen("reverse.in","r",stdin);
    freopen("reverse.out","w",stdout);
    srand(time(0));
    cin>>n>>m;
    for(int i=1;i<=n;++i)
    {
        cin>>a;
        T it=new Tr(i,rand()+1,a);
        insert(R,it,0);
    }
    for(int i=0;i<m;++i)
    {
        cin>>q>>a>>b;
        if(q) reverse(a,b);
        else get_sum(a,b);
    }
    return 0;
}

```

5.2 Segment Tree

```

//Segment Tree
#include <iostream>
#define N (1<<18)

```

```

using namespace std;
//int find(vector<int>& C, int x){return (C[x]==x) ? x : C[x]=find(C, C[x]);} C++
//int find(int x){return (C[x]==x)?x:C[x]=find(C[x]);}
typedef pair<int,int> ii;
ii arb[N]={0,0};
int n,m,a,b,v;
char type;

void update(int node,int l,int r,int a,int b,int val,int p)
{
    if(a<=l && r<=b)
    {
        arb[node].first=val;
        arb[node].second=p;
    }
    else
    {
        int mid=(l+r)/2;
        if(a<=mid)
            update(node*2,l,mid,a,b,val,p);
        if(b>mid)
            update(2*node+1,mid+1,r,a,b,val,p);
    }
}

pair<int,int> search(int node,int l,int r,int a)
{
    if(a==l && a==r)
    {
        return arb[node];
    }
    else
    {
        int mid=(l+r)/2;
        ii cur;
        if(a<=mid)
            cur=search(2*node,l,mid,a);
        else
            cur=search(2*node+1,mid+1,r,a);
        if(cur.second<arb[node].second)
            return arb[node];
        else
            return cur;
    }
}

```

6 Math with Numbers

6.1 Extended Euclid's Algorithm

```

#include "GcdLcm.h"

// returns d = gcd(a,b); find x, y such that d = ax + by
int extended_euclid(int a, int b, int &x, int &y){
    int current_x = y = 0;
    int current_y = x = 1;
    while(b)
    {
        int q = a/b;
        int t = b;
        b = a%b;
        a = t;
        t = current_x; current_x = x-q*current_x; x = t;
        t = current_y; current_y = y-q*current_y; y = t;
    }
    return a;
}

```

6.2 Fast Prime Number Sieve

```

// compute prime numbers to N really fast (Sieve)
#define N 10000100
vector<int> primes;
char p[N/8];

void GeneratePrimes()
{
    primes.push_back(2);
    int i,j;

```

```

for (i=1; ((i+1)<<1)+(i<<1)<N; ++i)
{
    if ((p[i]>>3] & (1<<(i&7)))==0)
    {
        primes.push_back((i<<1)+1);
        for (j=((i+1)<<1)+(i<<1), k=(i<<1)+1; (j<<1)+1<N; j+=k)
        {
            p[j]>>3] |= (1<<(j&7));
        }
    }
}

for (; (i<<1)+1<N; ++i)
{
    if ((p[i]>>3] & (1<<(i&7)))==0)
    {
        primes.push_back((i<<1)+1);
    }
}
}

```

6.3 GCD and LCM

```

// return a*b
int mod(int a, int b)
{
    return ((a%b)+b)%b;
}

// computes gcd(a,b)
int gcd(int a, int b)
{
    int tmp;
    while(b)
    {
        a %= b;
        tmp = a;
        a = b;
        b = tmp;
    }
    return a;
}

// computes lcm(a,b)
int lcm(int a, int b)
{
    return a/gcd(a,b)*b;
}

```

6.4 Euler Phi/Totient

```

// Euler phi (totient)
int ind = 0, pf = primes[0], ans = n;
while (1ll*pf*pf <= n) {
    if (n%pf==0) ans -= ans/pf;
    while (n%pf==0) n /= pf;
    pf = primes[++ind];
}
if (n != 1) ans -= ans/n;

// IME2014
int phi[N];
void totient() {
    for (int i = 1; i < N; ++i) phi[i]=i;
    for (int i = 2; i < N; i+=2) phi[i]>>=1;
    for (int j = 3; j < N; j+=2) if (phi[j]==j) {
        phi[j]--;
        for (int i = 2*j; i < N; i+=j) phi[i]=phi[i]/j*(j-1);
    }
}

```

7 Math with Matrices?

7.1 Modular Linear Equation Solver

```

// Stanford Notebook
#include <vector>
#include "ExtendedEuclid.h"

// find all solutions to ax = b (mod n)
vector<int> modular_linear_equation_solver(int a, int b, int n)
{
    int x,y;
    vector<int> sol;
    int d = extended_euclid(a, n, x, y);
    if (!(b%d)) {
        x = mod(x%(b/d), n);
        for (int i=0; i < d; ++i)
            sol.push_back(mod(x + i*(n/d), n));
    }
    return sol;
}

// computes b such that ab = 1(mod n), returns -1 on failure
int mod_inverse(int a, int n)
{
    int x, y;
    int d = extended_euclid(a, n, x, y);
    if (d>1) return -1;
    return mod(x,n);
}

// computes x and y such that ax + by =c; on failure, x = y =-1
void linear_diophantine(int a,int b, int c,int &x, int &y)
{
    int d = gcd(a,b);
    if (c%d) x = y = -1;
    else
    {
        x = c/d * mod_inverse(a/d,b/d);
        y = (c-a*x)/b;
    }
}

```

7.2 Fast Fourier Transform

```

struct cpx
{
    cpx() {}
    cpx(double aa): a(aa) {}
    cpx(double aa, double bb): a(aa), b(bb) {}
    double a;
    double b;
    double modsq(void) const { return a*a+b*b; }
    cpx bar(void) const { return cpx(a,-b); }
};

cpx b[N+100], c[N+100], B[N+100], C[N+100];
int a[N+100], int x[N+100];
double coss[N+100], sins[N+100];
int n,m,p;

cpx operator +(cpx a, cpx b) { return cpx(a.a+b.a, a.b+b.b); }
cpx operator *(cpx a, cpx b) { return cpx(a.a*b.a-a.b*b.b, a.a*b.b+a.b*b.a); }
cpx operator /(cpx a, cpx b) { cpx r = a*b.bar(); return cpx(r.a/b.modsq(), r.b/b.modsq()); }
cpx EXP(int i, int dir) { return cpx(coss[i], sins[i]*dir); }

const double two_pi = 4 * acos(0);

void FFT(cpx *in, cpx *out, int step, int size, int dir)
{
    if (size<1) return;
    if (size==1)
    {
        out[0]=in[0];
        return;
    }
    FFT(in, out, step*2, size/2, dir);
    FFT(in+step, out+size/2, step*2, size/2, dir);
    for (int i=0; i<size/2; ++i)
    {
        cpx even=out[i];
        cpx odd=out[i+size/2];
        out[i] = even+EXP(i*step, dir)*odd;
        out[i+size/2]=even+EXP((i+size/2)*step, dir)*odd;
    }
}

void exemple
{
    for (int i=0; i<=N; ++i)

```

```

{
    coss[i]=cos(two_pi*i/N);
    sins[i]=sin(two_pi*i/N);
}
while(scanf("%d",&n)==1)
{
    fill(x,x+N+100,0);
    fill(a,a+N+100,0);
    for(int i=0;i<n;++i)
    {
        scanf("%d",&p);
        x[p]=1;
    }
    for(int i=0;i<N+100;++i)
    {
        b[i]=cpx(x[i],0);
    }
    scanf("%d",&m);
    for(int i=0;i<m;++i)
    {
        scanf("%d",&a[i]);
    }
    FFT(b,B,1,N,1);
    for(int i=0;i<N;++i)
        C[i]=B[i]*B[i];
    FFT(C,c,1,N,-1);
    for(int i=0;i<N;++i)
        c[i]=c[i]/N;
    int cnt=0;
    for(int i=0;i<16;++i)
        cout<<c[i].a<<" ";
    for(int i=0;i<m;++i)
        if(c[a[i]].a>0.5 || x[a[i]])
            cnt++;
    printf("%d\n",cnt);
}
}

```

7.3 Gauss-Jordan Elimination

```

/*
Stanford notebook
-----
GaussJordan Algorithm (elimination with full pivoting)
Uses:
(1) solving systems of linear equations (AX=B)
(2) inverting matrices (AX=I)
(3) computing determinants of square matrices
Running time: O(n^3)
INPUT:   a[][] = an nxn matrix
         b[][] = an nxm matrix
OUTPUT:  X      = an nxm matrix (stored in b[][])
         A^{-1} = an nxn matrix (stored in a[][])
         returns determinant of a[][]
-----
*/

const double EPS = 1e-10;
typedef vector<int> VI;
typedef double T;
typedef vector<T> VT;
typedef vector<VT> VVT;

T GaussJordan(VVT &a, VVT &b)
{
    const int n = a.size();
    const int m = b[0].size();
    VI irow(n), icol(n), ipiv(n);
    T det = 1;

    for (int i = 0; i < n; i++)
    {
        int pj = -1, pk = -1;
        for (int j = 0; j < n; j++)
            if (!ipiv[j])
                for (int k = 0; k < n; k++)
                    if (!ipiv[k])
                        if (pj == -1 || fabs(a[j][k]) > fabs(a[pj][pk])) { pj = j; pk = k; }

        if (fabs(a[pj][pk]) < EPS)
        {
            cerr << "Matrix is singular." << endl;
            exit(0);
        }
        ipiv[pj]++;
        swap(a[pj], a[pk]);
        swap(b[pj], b[pk]);
        if (pj != pk) det *= -1;
    }
}

```

```

irow[i] = pj;
icol[i] = pk;

T c = 1.0 / a[pk][pk];
det *= a[pk][pk];
a[pk][pk] = 1.0;

for (int p = 0; p < n; p++)
    a[pk][p] *= c;
for (int p = 0; p < m; p++)
    b[pk][p] *= c;
for (int p = 0; p < n; p++)
{
    if (p != pk)
    {
        c = a[p][pk];
        a[p][pk] = 0;
        for (int q = 0; q < n; q++)
            a[p][q] -= a[pk][q] * c;
        for (int q = 0; q < m; q++)
            b[p][q] -= b[pk][q] * c;
    }
}

for (int p = n-1; p >= 0; p--)
{
    if (irow[p] != icol[p])
    {
        for (int k = 0; k < n; k++)
            swap(a[k][irow[p]], a[k][icol[p]]);
    }
    return det;
}

int main()
{
    const int n = 4;
    const int m = 2;
    double A[n][n] = { {1,2,3,4},{1,0,1,0},{5,3,2,4},{6,1,4,6} };
    double B[n][m] = { {1,2},{4,3},{5,6},{8,7} };
    VVT a(n), b(n);
    for (int i = 0; i < n; i++)
    {
        a[i] = VT(A[i], A[i] + n);
        b[i] = VT(B[i], B[i] + m);
    }
    double det = GaussJordan(a, b);
    // expected: 60
    cout << "Determinant: " << det << endl;
    // expected: -0.233333 0.166667 0.133333 0.0666667
    // 0.166667 0.166667 0.333333 -0.333333
    // 0.233333 0.833333 -0.133333 -0.0666667
    // 0.05 -0.75 -0.1 0.2
    cout << "Inverse: " << endl;
    for (int i = 0; i < n; i++)
    {
        for (int j = 0; j < n; j++)
            cout << a[i][j] << ' ';
        cout << endl;
    }
    // expected: 1.63333 1.3
    // -0.166667 0.5
    // 2.36667 1.7
    // -1.85 -1.35
    cout << "Solution: " << endl;
    for (int i = 0; i < n; i++)
    {
        for (int j = 0; j < m; j++)
            cout << b[i][j] << ' ';
        cout << endl;
    }
}

```

8 Strings

8.1 Aho-Corasick Algorithm

```

/*
Implementation - Benoit Chabod
-----
Aho Corasick algorithm
-----

```

8.2 Knuth-Morris-Pratt Algorithm

```

*/
struct node
{
    int f;
    map<char, int> g;
    vector<short> out;
    node(int fail = -1): f(fail) {}
};

vector<node> nodes;

void add_str(const string & s, int num)
{
    int cur = 0;
    int n = s.size();
    for(int i = 0; i < n; i++)
    {
        auto it = nodes[cur].g.find(s[i]);
        if(it == nodes[cur].g.end())
        {
            nodes[cur].g[s[i]] = nodes.size();
            cur = nodes.size();
            nodes.push_back(node());
        }
        else
        {
            cur = it->second;
        }
    }
    nodes[cur].out.push_back(num);
}

void init_fail()
{
    int cur = 0;
    queue<int> q;
    q.push(cur);

    while( !q.empty() )
    {
        cur = q.front();
        map<char, int>::iterator it;
        for(it = nodes[cur].g.begin(); it != nodes[cur].g.end(); it++)
        {
            int child = it->second;
            int pfail = nodes[cur].f;
            char ch = it->first;
            map<char, int>::iterator f;
            while( pfail != -1 && ((f = nodes[pfail].g.find(ch)) == nodes[pfail].g.end()) )
            {
                pfail = nodes[pfail].f;
            }
            nodes[child].f = (pfail == -1)? 0 : f->second;
            pfail = nodes[child].f;
            nodes[child].out.insert(nodes[child].out.end(), nodes[pfail].out.begin(), nodes[pfail].out.end());
            q.push(child);
        }
        q.pop();
    }

    // Usage
    void usage()
    {
        nodes.push_back(node());
        for [each word] add_str(word,i)
        init_fail();
        for [each letter]
        {
            map<char, int>::iterator f;
            while( cur != -1 && ((f = nodes[cur].g.find(letter)) == nodes[cur].g.end()) )
            {
                cur = nodes[cur].f;
                if( cur == -1 )
                {
                    cur = 0;
                    continue;
                }
            }
            cur = f->second;
            for(auto v : nodes[cur].out)
            {
                // Word v was found
            }
        }
    }
}

```

```

/*
-----
KMP/Pi function
Note : cin>>(s+1) (the operations in the pi-function start at 1)
-----
*/
void preKmp()
{
    int k;
    k=kmpNext[1]=0;
    for(int i=2; i<=n; ++i)
    {
        while(k && p[k+1]!=p[i]) k=kmpNext[k];
        if(p[k+1]==p[i])
            k++;
        kmpNext[i]=k;
    }
}

void KMP()
{
    preKmp();
    int k=0;

    for(int i=1; i<=m; ++i)
    {
        while(k && p[k+1]!=s[i])
            k=kmpNext[k];
        if(p[k+1]==s[i])
            k++;
        if(k==n)
        {
            // here we have a match
            k=kmpNext[k];
        }
    }
}

```

8.3 Suffix Array

```

//Suffix Array
struct SuffixArray
{
    const int L;
    string s;
    vector<vector<int>> > P;
    vector<pair<pair<int,int>,int>> > M;
    SuffixArray(const string &s) : L(s.length()), s(s), P(1, vector<int>(L, 0)), M(L)
    {
        for (int i = 0; i < L; i++)
            P[0][i] = int(s[i]);
        for (int skip = 1, level = 1; skip < L; skip *= 2, level++)
        {
            P.push_back(vector<int>(L, 0));
            for (int i = 0; i < L; i++)
            {
                M[i] = make_pair(make_pair(P[level-1][i], i + skip < L ? P[level-1][i]
                    + skip : -1000), i);
            }
            sort(M.begin(), M.end());
            for (int i = 0; i < L; i++)
                P[level][M[i].second] = (i > 0 && M[i].first == M[i-1].first) ? P[
                    level][M[i-1].second] : i;
        }
    }

    vector<int> GetSuffixArray()
    {
        return P.back();
    }

    // returns the length of the longest common prefix of s[i...L-1] and s[j...L-1]
    int LongestCommonPrefix(int i, int j)
    {
        int len = 0;
        if (i == j) return L - i;
        for (int k = P.size() - 1; k >= 0 && i < L && j < L; k--)
        {
            if (P[k][i] == P[k][j])
            {
                i += 1 << k;
                j += 1 << k;
                len += 1 << k;
            }
        }
    }
}

```



```

        return len;
    }
};

int main()
{
    // bobocel is the 0'th suffix
    // obocel is the 5'th suffix
    // bocel is the 1'st suffix
    // ocel is the 6'th suffix
    // cel is the 2'nd suffix
    // el is the 3'rd suffix
    // l is the 4'th suffix
    SuffixArray suffix("bobocel");
    vector<int> v = suffix.GetSuffixArray();
    // Expected output: 0 5 1 6 2 3 4
    //                2
    for (int i = 0; i < v.size(); i++)
        cout << v[i] << " ";
    cout << endl;
    cout << suffix.LongestCommonPrefix(0, 2) << endl;
}

```

8.4 Manacher (Longest Palindromic Substring)

```

// Manacher (Longest Palindromic String) - O(n)
int lps[2*N+5];
char s[N];

int manacher() {
    int n = strlen(s);

    string p(2*n+3, '#');
    p[0] = '^';
    for (int i = 0; i < n; i++) p[2*(i+1)] = s[i];
    p[2*n+2] = '$';

    int k = 0, r = 0, m = 0;
    int l = p.length();
    for (int i = 1; i < l; i++) {
        int o = 2*k - i;
        lps[i] = (r > i) ? min(r-i, lps[o]) : 0;
        while (p[i+1+lps[i]] == p[i-1-lps[i]]) lps[i]++;
        if (i+lps[i] > r) k = i, r = i+lps[i];
        m = max(m, lps[i]);
    }
    return m;
}

```

8.5 Z Function

```

// Z-Function - O(n)
// Array where ith element is greatest number of characters starting at i matching the prefix

vector<int> zfunction(const string& s){
    vector<int> z (s.size());
    for (int i = 1, l = 0, r = 0, n = s.size(); i < n; i++){
        if (i <= r) z[i] = min(z[i-l], r - i + 1);
        while (i + z[i] < n and s[z[i]] == s[z[i] + i]) z[i]++;
        if (i + z[i] - 1 > r) l = i, r = i + z[i] - 1;
    }
    return z;
}

```

9 Python Stuff...

10 Geometry

10.1 Convex Hull

```

def convex_hull(points):
    """Computes the convex hull of a set of 2D points.

```

Input: an iterable sequence of (x, y) pairs representing the points.
Output: a list of vertices of the convex hull in counter-clockwise order,
starting from the vertex with the lexicographically smallest coordinates.
Implements Andrew's monotone chain algorithm. $O(n \log n)$ complexity.
"""

```

# Sort the points lexicographically (tuples are compared lexicographically).
# Remove duplicates to detect the case we have just one unique point.
points = sorted(set(points))

# Boring case: no points or a single point, possibly repeated multiple times.
if len(points) <= 1:
    return points

# 2D cross product of OA and OB vectors, i.e. z-component of their 3D cross product.
# Returns a positive value, if OAB makes a counter-clockwise turn,
# negative for clockwise turn, and zero if the points are collinear.
def cross(o, a, b):
    return (a[0] - o[0]) * (b[1] - o[1]) - (a[1] - o[1]) * (b[0] - o[0])

# Build lower hull
lower = []
for p in points:
    while len(lower) >= 2 and cross(lower[-2], lower[-1], p) <= 0:
        lower.pop()
    lower.append(p)

# Build upper hull
upper = []
for p in reversed(points):
    while len(upper) >= 2 and cross(upper[-2], upper[-1], p) <= 0:
        upper.pop()
    upper.append(p)

# Concatenation of the lower and upper hulls gives the convex hull.
# Last point of each list is omitted because it is repeated at the beginning of the other list.
return lower[:-1] + upper[:-1]

# Example: convex hull of a 10-by-10 grid.
assert convex_hull([(i//10, i%10) for i in range(100)]) == [(0, 0), (9, 0), (9, 9), (0, 9)]

```

11 Dynamic Programming

11.1 Max Sum Subarray (Kadane's Algorithm)

```

def maxSubArraySum(a, size):
    max_so_far = 0
    max_ending_here = 0
    for i in range(0, size):
        max_ending_here = max_ending_here + a[i]
        if max_ending_here < 0:
            max_ending_here = 0
        elif (max_so_far < max_ending_here):
            max_so_far = max_ending_here
    return max_so_far

```

11.2 Longest Common Subsequence

```

def lcs(X, Y):
    # find the length of the strings
    m = len(X)
    n = len(Y)

    # declaring the array for storing the dp values
    L = [[None]*(n+1) for i in xrange(m+1)]

    """Following steps build L[m+1][n+1] in bottom up fashion
    Note: L[i][j] contains length of LCS of X[0..i-1]
    and Y[0..j-1]"""
    for i in range(m+1):
        for j in range(n+1):
            if i == 0 or j == 0 :
                L[i][j] = 0
            elif X[i-1] == Y[j-1]:
                L[i][j] = L[i-1][j-1] + 1
            else:
                L[i][j] = max(L[i-1][j], L[i][j-1])

    # L[m][n] contains the length of LCS of X[0..n-1] & Y[0..m-1]
    return L[m][n]

```

11.3 Levenshtein Distance

```
def levenshtein(s1, s2):
    if len(s1) < len(s2):
        return levenshtein(s2, s1)

    # len(s1) >= len(s2)
    if len(s2) == 0:
        return len(s1)

    previous_row = range(len(s2) + 1)
    for i, c1 in enumerate(s1):
        current_row = [i + 1]
        for j, c2 in enumerate(s2):
            insertions = previous_row[j + 1] + 1 # j+1 instead of j since previous_row and current_row
            # are one character longer
            deletions = current_row[j] + 1 # than s2
            substitutions = previous_row[j] + (c1 != c2)
            current_row.append(min(insertions, deletions, substitutions))
        previous_row = current_row
    return previous_row[-1]
```

11.4 Longest Increasing Subsequence

```
def lis(arr):
    n = len(arr)

    # Declare the list (array) for LIS and initialize LIS
    # values for all indexes
    lis = [1]*n

    # Compute optimized LIS values in bottom up manner
    for i in range(1, n):
        for j in range(0, i):
            if arr[i] > arr[j] and lis[i] < lis[j] + 1 :
                lis[i] = lis[j] + 1

    # Initialize maximum to 0 to get the maximum of all
    # LIS
    maximum = 0

    # Pick maximum of all LIS values
    for i in range(n):
        maximum = max(maximum, lis[i])
    return maximum
```

12 Graphs/Trees

12.1 Graph structure example for our DFS and BFS algorithms

```
graph = {'A': set(['B', 'C']),
         'B': set(['A', 'D', 'E']),
         'C': set(['A', 'F']),
         'D': set(['B']),
         'E': set(['B', 'F']),
         'F': set(['C', 'E'])}
```

12.2 Breadth-First Search

```
def bfs(graph, start):
    visited, queue = set(), [start]
    while queue:
        vertex = queue.pop(0)
        if vertex not in visited:
            visited.add(vertex)
            queue.extend(graph[vertex] - visited)
    return visited

bfs(graph, 'A') # {'B', 'C', 'A', 'F', 'D', 'E'}
```

12.3 Breadth-First Search Paths

```
def bfs_paths(graph, start, goal):
    queue = [(start, [start])]
    while queue:
        (vertex, path) = queue.pop(0)
        for next in graph[vertex] - set(path):
            if next == goal:
                yield path + [next]
            else:
                queue.append((next, path + [next]))

list(bfs_paths(graph, 'A', 'F')) # [['A', 'C', 'F'], ['A', 'B', 'E', 'F']]
```

12.4 Breadth-First Search Shortest Path

```
def shortest_path(graph, start, goal):
    try:
        return next(bfs_paths(graph, start, goal))
    except StopIteration:
        return None

shortest_path(graph, 'A', 'F') # ['A', 'C', 'F']
```

12.5 Depth-First Search

```
def dfs(graph, start):
    visited, stack = set(), [start]
    while stack:
        vertex = stack.pop()
        if vertex not in visited:
            visited.add(vertex)
            stack.extend(graph[vertex] - visited)
    return visited

dfs(graph, 'A') # {'E', 'D', 'F', 'A', 'C', 'B'}
```

12.6 Depth-First Search Paths

```
#Returns all paths from start to goal
def dfs_paths(graph, start, goal):
    stack = [(start, [start])]
    while stack:
        (vertex, path) = stack.pop()
        for next in graph[vertex] - set(path):
            if next == goal:
                yield path + [next]
            else:
                stack.append((next, path + [next]))

list(dfs_paths(graph, 'A', 'F')) # [['A', 'C', 'F'], ['A', 'B', 'E', 'F']]
```

12.7 Dijkstra's Algorithm

```
from collections import defaultdict
from heapq import *

def dijkstra(edges, f, t):
    g = defaultdict(list)
    for l,r,c in edges:
        g[l].append((c,r))

    q, seen = [(0,f,())], set()
    while q:
        (cost,v1,path) = heappop(q)
        if v1 not in seen:
            seen.add(v1)
            path = (v1, path)
            if v1 == t: return (cost, path)
            for c, v2 in g.get(v1, ()):
                heappush(q, (cost+c, v2, path+(v2,)))
```

```

        if v2 not in seen:
            heappush(q, (cost+c, v2, path))
    return float("inf")

#Code example
edges = [("A", "B", 7), ("A", "D", 5), ("B", "C", 8),
         ("B", "D", 9), ("B", "E", 7), ("C", "E", 5)]
print "A -> E:"
print dijkstra(edges, "A", "E")    #(14, ('E', ('B', ('A', ())))

```

12.8 Kruskal's Algorithm (including Merge-Find set)

```

parent = dict()
rank = dict()

def make_set(vertice):
    parent[vertice] = vertice
    rank[vertice] = 0

def find(vertice):
    if parent[vertice] != vertice:
        parent[vertice] = find(parent[vertice])
    return parent[vertice]

def union(vertice1, vertice2):
    root1 = find(vertice1)
    root2 = find(vertice2)
    if root1 != root2:
        if rank[root1] > rank[root2]:
            parent[root2] = root1
        else:
            parent[root1] = root2
        if rank[root1] == rank[root2]: rank[root2] += 1

def kruskal(graph):
    for vertice in graph['vertices']:
        make_set(vertice)
    minimum_spanning_tree = set()
    edges = list(graph['edges'])
    edges.sort()
    #print edges
    for edge in edges:
        weight, vertice1, vertice2 = edge
        if find(vertice1) != find(vertice2):
            union(vertice1, vertice2)
            minimum_spanning_tree.add(edge)

    return sorted(minimum_spanning_tree)

```

12.9 Bellman-Ford Algorithm

```

# Step 1: For each node prepare the destination and predecessor
def initialize(graph, source):
    d = {} # Stands for destination
    p = {} # Stands for predecessor
    for node in graph:
        d[node] = float('Inf') # We start admitting that the rest of nodes are very very far
        p[node] = None
    d[source] = 0 # For the source we know how to reach
    return d, p

def relax(node, neighbour, graph, d, p):
    # If the distance between the node and the neighbour is lower than the one I have now
    if d[neighbour] > d[node] + graph[node][neighbour]:
        # Record this lower distance
        d[neighbour] = d[node] + graph[node][neighbour]
        p[neighbour] = node

def bellman_ford(graph, source):
    d, p = initialize(graph, source)
    for i in range(len(graph)-1): #Run this until is converges
        for u in graph:
            for v in graph[u]:
                relax(u, v, graph, d, p) #Lets relax it

# Step 3: check for negative-weight cycles
for u in graph:
    for v in graph[u]:
        assert d[v] <= d[u] + graph[u][v]

return d, p

def test():

```

```

graph = {
    'a': {'b': -1, 'c': 4},
    'b': {'c': 3, 'd': 2, 'e': 2},
    'c': {},
    'd': {'b': 1, 'c': 5},
    'e': {'d': -3}
}
d, p = bellman_ford(graph, 'a')
# d = {'a':0, 'b':-1, 'c':2, 'd':-2, 'e':1},
# p = {'a':None, 'b':'a', 'c':'b', 'd':'e', 'e':'b'}

```

12.10 Floyd-Warshall Algorithm

```

# Number of vertices in the graph
V = 4

# Define infinity as the large enough value. This value will be
# used for vertices not connected to each other
INF = 99999

# Solves all pair shortest path via Floyd Warshall Algorithm
def floydWarshall(graph):
    """ dist[][] will be the output matrix that will finally
        have the shortest distances between every pair of vertices """
    """ initializing the solution matrix same as input graph matrix
    OR we can say that the initial values of shortest distances
    are based on shortest paths considering no
    intermediate vertices """
    dist = map(lambda i : map(lambda j : j , i) , graph)

    """ Add all vertices one by one to the set of intermediate
    vertices.
    ---> Before start of a iteration, we have shortest distances
    between all pairs of vertices such that the shortest
    distances consider only the vertices in set
    {0, 1, 2, .. k-1} as intermediate vertices.
    ----> After the end of a iteration, vertex no. k is
    added to the set of intermediate vertices and the
    set becomes {0, 1, 2, .. k}
    """
    for k in range(V):

        # pick all vertices as source one by one
        for i in range(V):

            # Pick all vertices as destination for the
            # above picked source
            for j in range(V):

                # If vertex k is on the shortest path from
                # i to j, then update the value of dist[i][j]
                dist[i][j] = min(dist[i][j] ,
                                dist[i][k]+ dist[k][j]
                                )

    printSolution(dist)

"""
    10
(0)----->(3)
|           /\
5 |         |
|         | 1
|/\       |
(1)----->(2)
      3      """
graph = [[0,5,INF,10],
         [INF,0,3,INF],
         [INF, INF, 0, 1],
         [INF, INF, INF, 0]]

floydWarshall(graph) # [[0,5,8,9],[INF,0,3,4],[INF,INF,0,1],[INF,INF,INF,0]]

```

12.11 Max Flow (Ford-Fulkerson Algorithm)

```

from collections import defaultdict

#This class represents a directed graph using adjacency matrix representation
class Graph:

    def __init__(self,graph):
        self.graph = graph # residual graph

```

12.12 Segment Tree

```

self.ROW = len(graph)
#self.COL = len(gr[0])

'''Returns true if there is a path from source 's' to sink 't' in
residual graph. Also fills parent[] to store the path'''
def BFS(self, s, t, parent):

    # Mark all the vertices as not visited
    visited = [False] * (self.ROW)

    # Create a queue for BFS
    queue = []

    # Mark the source node as visited and enqueue it
    queue.append(s)
    visited[s] = True

    # Standard BFS Loop
    while queue:

        # Dequeue a vertex from queue and print it
        u = queue.pop(0)

        # Get all adjacent vertices of the dequeued vertex u
        # If a adjacent has not been visited, then mark it
        # visited and enqueue it
        for ind, val in enumerate(self.graph[u]):
            if visited[ind] == False and val > 0:
                queue.append(ind)
                visited[ind] = True
                parent[ind] = u

    # If we reached sink in BFS starting from source, then return
    # true, else false
    return True if visited[t] else False

# Returns the maximum flow from s to t in the given graph
def FordFulkerson(self, source, sink):

    # This array is filled by BFS and to store path
    parent = [-1] * (self.ROW)

    max_flow = 0 # There is no flow initially

    # Augment the flow while there is path from source to sink
    while self.BFS(source, sink, parent):

        # Find minimum residual capacity of the edges along the
        # path filled by BFS. Or we can say find the maximum flow
        # through the path found.
        path_flow = float("Inf")
        s = sink
        while(s != source):
            path_flow = min(path_flow, self.graph[parent[s]][s])
            s = parent[s]

        # Add path flow to overall flow
        max_flow += path_flow

        # update residual capacities of the edges and reverse edges
        # along the path
        v = sink
        while(v != source):
            u = parent[v]
            self.graph[u][v] -= path_flow
            self.graph[v][u] += path_flow
            v = parent[v]

    return max_flow

# Create a graph given in the above diagram
graph = [[0, 16, 13, 0, 0, 0],
         [0, 0, 10, 12, 0, 0],
         [0, 4, 0, 0, 14, 0],
         [0, 0, 9, 0, 0, 20],
         [0, 0, 0, 7, 0, 4],
         [0, 0, 0, 0, 0, 0]]

g = Graph(graph)

source = 0; sink = 5

print ("The maximum possible flow is %d" % g.FordFulkerson(source, sink))

```

```

#encoding:utf-8
class SegmentTree(object):
    def __init__(self, start, end):
        self.start = start
        self.end = end
        self.max_value = {}
        self.sum_value = {}
        self.len_value = {}
        self._init(start, end)

    def add(self, start, end, weight=1):
        start = max(start, self.start)
        end = min(end, self.end)
        self._add(start, end, weight, self.start, self.end)
        return True

    def query_max(self, start, end):
        return self._query_max(start, end, self.start, self.end)

    def query_sum(self, start, end):
        return self._query_sum(start, end, self.start, self.end)

    def query_len(self, start, end):
        return self._query_len(start, end, self.start, self.end)

    #####
    def _init(self, start, end):
        self.max_value[(start, end)] = 0
        self.sum_value[(start, end)] = 0
        self.len_value[(start, end)] = 0
        if start < end:
            mid = start + int((end - start) / 2)
            self._init(start, mid)
            self._init(mid+1, end)

    def _add(self, start, end, weight, in_start, in_end):
        key = (in_start, in_end)
        if in_start == in_end:
            self.max_value[key] += weight
            self.sum_value[key] += weight
            self.len_value[key] = 1 if self.sum_value[key] > 0 else 0
            return

        mid = in_start + int((in_end - in_start) / 2)
        if mid >= end:
            self._add(start, end, weight, in_start, mid)
        elif mid+1 <= start:
            self._add(start, end, weight, mid+1, in_end)
        else:
            self._add(start, mid, weight, in_start, mid)
            self._add(mid+1, end, weight, mid+1, in_end)
            self.max_value[key] = max(self.max_value[(in_start, mid)], self.max_value[(mid+1, in_end)])
            self.sum_value[key] = self.sum_value[(in_start, mid)] + self.sum_value[(mid+1, in_end)]
            self.len_value[key] = self.len_value[(in_start, mid)] + self.len_value[(mid+1, in_end)]

    def _query_max(self, start, end, in_start, in_end):
        if start == in_start and end == in_end:
            ans = self.max_value[(start, end)]
        else:
            mid = in_start + int((in_end - in_start) / 2)
            if mid >= end:
                ans = self._query_max(start, end, in_start, mid)
            elif mid+1 <= start:
                ans = self._query_max(start, end, mid+1, in_end)
            else:
                ans = max(self._query_max(start, mid, in_start, mid),
                           self._query_max(mid+1, end, mid+1, in_end))
        #print start, end, in_start, in_end, ans
        return ans

    def _query_sum(self, start, end, in_start, in_end):
        if start == in_start and end == in_end:
            ans = self.sum_value[(start, end)]
        else:
            mid = in_start + int((in_end - in_start) / 2)
            if mid >= end:
                ans = self._query_sum(start, end, in_start, mid)
            elif mid+1 <= start:
                ans = self._query_sum(start, end, mid+1, in_end)
            else:
                ans = self._query_sum(start, mid, in_start, mid) + self._query_sum(mid+1, end, mid+1, in_end)
        return ans

    def _query_len(self, start, end, in_start, in_end):
        if start == in_start and end == in_end:
            ans = self.len_value[(start, end)]
        else:
            mid = in_start + int((in_end - in_start) / 2)

```

```

    if mid >= end:
        ans = self._query_len(start, end, in_start, mid)
    elif mid+1 <= start:
        ans = self._query_len(start, end, mid+1, in_end)
    else:
        ans = self._query_len(start, mid, in_start, mid) + self._query_len(mid+1, end, mid+1, in_end)

# print start, end, in_start, in_end, ans
return ans

```

13 Math

13.1 Prime Number Sieve (generator)

```

from itertools import count

def postponed_sieve():
    yield 2; yield 3; yield 5; yield 7;
    sieve = {}
    ps = postponed_sieve()
    p = next(ps) and next(ps)
    q = p*p
    for c in count(9,2):
        if c in sieve:
            s = sieve.pop(c)
            # c's a multiple of some base prime
            # i.e. a composite ; or
            elif c < q:
                yield c
                continue
            else:
                # (c==q):
                # or the next base prime's square:
                s = count(q+2*p, 2*p)
                p = next(ps)
                q = p*p
                # (25)
            # the next multiple
            if m not in sieve:
                break
            sieve[m] = s
    # original test entry: ideone.com/WFv4f

```

13.2 GCD and Euler's Totient Function

```

# Function to return gcd of a and b
def gcd(a, b):
    if a == 0:
        return b
    return gcd(b%a, a)

# A simple method to evaluate Euler Totient Function
def phi(n):
    result = 1
    for i in range(2, n):
        if gcd(i, n) == 1:
            result = result + 1
    return result

```

13.3 Miller-Rabin Primality Test

```

def miller_rabin(n, k):
    # The optimal number of rounds (k) for this test is 40
    # for justification

    if n == 2:
        return True
    if n % 2 == 0:
        return False
    r, s = 0, n - 1
    while s % 2 == 0:
        r += 1
        s //= 2
    for _ in xrange(k):
        a = random.randrange(2, n - 1)
        x = pow(a, s, n)
        if x == 1 or x == n - 1:
            continue
        for _ in xrange(r - 1):
            x = pow(x, 2, n)

```

```

    if x == n - 1:
        break
    else:
        return False
    return True

```

13.4 Gauss-Jordan Elimination (Matrix inversion and linear system solving)

```

def gauss_jordan(m, eps = 1.0/(10**10)):
    """Puts given matrix (2D array) into the Reduced Row Echelon Form.
    Returns True if successful, False if 'm' is singular.
    NOTE: make sure all the matrix items support fractions! Int matrix will NOT work!
    Written by Jarno Elonen in April 2005, released into Public Domain"""
    (h, w) = (len(m), len(m[0]))
    for y in range(0, h):
        maxrow = y
        for y2 in range(y+1, h):
            if abs(m[y2][y]) > abs(m[maxrow][y]):
                maxrow = y2
        (m[y], m[maxrow]) = (m[maxrow], m[y])
        if abs(m[y][y]) <= eps:
            # Singular?
            return False
        for y2 in range(y+1, h):
            # Eliminate column y
            c = m[y2][y] / m[y][y]
            for x in range(y, w):
                m[y2][x] -= m[y][x] * c
        for y in range(h-1, 0-1, -1): # Backsubstitute
            c = m[y][y]
            for y2 in range(0, y):
                for x in range(w-1, y-1, -1):
                    m[y2][x] -= m[y][x] * m[y2][y] / c
            m[y][y] /= c
        for x in range(h, w):
            # Normalize row y
            m[y][x] /= c
    return True

def solve(M, b):
    """
    solves M*x = b
    return vector x so that M*x = b
    :param M: a matrix in the form of a list of list
    :param b: a vector in the form of a simple list of scalars
    """
    m2 = [row[:] + [right] for row, right in zip(M, b)]
    return [row[-1] for row in m2] if gauss_jordan(m2) else None

def inv(M):
    """
    return the inv of the matrix M
    """
    # clone the matrix and append the identity matrix
    # [int(i==j) for j in range_M] is nothing but the i(th row of the identity matrix
    m2 = [row[:] + [int(i==j) for j in range(len(M))] for i, row in enumerate(M)]
    # extract the appended matrix (kind of m2[m:,...])
    return [row[len(M):] for row in m2] if gauss_jordan(m2) else None

def zeros(s, zero=0):
    """
    return a matrix of size 'size'
    :param size: a tuple containing dimensions of the matrix
    :param zero: the value to use to fill the matrix (by default it's zero)
    """
    return [zeros(s[1:]) for i in range(s[0])] if not len(s) else zero

```

14 Strings

14.1 Knuth-Morris-Pratt Algorithm (fast pattern matching)

```

def KnuthMorrisPratt(text, pattern):
    """Yields all starting positions of copies of the pattern in the text.
    Calling conventions are similar to string.find, but its arguments can be
    lists or iterators, not just strings, it returns all matches, not just
    the first one, and it does not need the whole text in memory at once.
    Whenever it yields, it will have read the text exactly up to and including

```

```

the match that caused the yield.'''

# allow indexing into pattern and protect against change during yield
pattern = list(pattern)

# build table of shift amounts
shifts = [1] * (len(pattern) + 1)
shift = 1
for pos in range(len(pattern)):
    while shift <= pos and pattern[pos] != pattern[pos-shift]:
        shift += shifts[pos-shift]
    shifts[pos+1] = shift

# do the actual search
startPos = 0
matchLen = 0
for c in text:
    while matchLen == len(pattern) or \
          matchLen >= 0 and pattern[matchLen] != c:
        startPos += shifts[matchLen]
        matchLen -= shifts[matchLen]
    matchLen += 1
    if matchLen == len(pattern):
        yield startPos

```

14.2 Rabin-Karp Algorithm (multiple pattern matching)

```

# d is the number of characters in input alphabet
d = 256

# pat -> pattern
# txt -> text
# q -> A prime number

def search(pat, txt, q):
    M = len(pat)
    N = len(txt)
    i = 0
    j = 0
    p = 0 # hash value for pattern
    t = 0 # hash value for txt

```

```

h = 1

# The value of h would be "pow(d, M-1)%q"
for i in xrange(M-1):
    h = (h*d)%q

# Calculate the hash value of pattern and first window
# of text
for i in xrange(M):
    p = (d*p + ord(pat[i]))%q
    t = (d*t + ord(txt[i]))%q

# Slide the pattern over text one by one
for i in xrange(N-M+1):
    # Check the hash values of current window of text and
    # pattern if the hash values match then only check
    # for characters on by one
    if p==t:
        # Check for characters one by one
        for j in xrange(M):
            if txt[i+j] != pat[j]:
                break

        j+=1
        # if p == t and pat[0...M-1] = txt[i, i+1, ...i+M-1]
        if j==M:
            print "Pattern found at index " + str(i)

# Calculate hash value for next window of text: Remove
# leading digit, add trailing digit
if i < N-M:
    t = (d*(t-ord(txt[i])*h) + ord(txt[i+M]))%q

# We might get negative values of t, converting it to
# positive
if t < 0:
    t = t+q

# Driver program to test the above function
txt = "GEEKS FOR GEEKS"
pat = "GEEK"
q = 101 # A prime number
search(pat,txt,q)

```