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Stanford University ACM Team Notebook (2011-12)
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Dinic.cc 1/27
// Adjacency list implementation of Dinic's blocking flow algorithm.
// This is very fast in practice, and only loses to push-relabel flow.
//
// Running time:
//
       O(|V|^2 |E|)
```

```
//
// INPUT:
//
       - graph, constructed using AddEdge()
//
       - source
//
       - sink
//
// OUTPUT:
//
       - maximum flow value
//
       - To obtain the actual flow values, look at all edges with
//
          capacity > 0 (zero capacity edges are residual edges).
#include <cmath>
#include <vector>
#include <iostream>
#include <queue>
using namespace std;
const int INF = 20000000000;
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) {}
};
struct Dinic {
  int N;
  vector<vector<Edge>>G;
  vector<Edge *> dad;
  vector<int> Q;
  Dinic(int N): N(N), G(N), dad(N), Q(N) {}
  void AddEdge(int from, int to, int 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, G[from].size() - 1));
  }
  long long BlockingFlow(int s, int t) {
     fill(dad.begin(), dad.end(), (Edge *) NULL);
     dad[s] = &G[0][0] - 1;
```

```
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 (!dad[e.to] && e.cap - e.flow > 0) {
     dad[e.to] = &G[x][i];
     Q[tail++] = e.to;
  }
     }
  if (!dad[t]) return 0;
  long long totflow = 0;
  for (int i = 0; i < G[t].size(); i++) {
     Edge *start = &G[G[t][i].to][G[t][i].index];
     int amt = INF;
     for (Edge *e = start; amt && e != dad[s]; e = dad[e -> from]) {
  if (!e) { amt = 0; break; }
  amt = min(amt, e->cap - e->flow);
     }
     if (amt == 0) continue;
     for (Edge *e = start; amt && e != dad[s]; e = dad[e-sfrom]) {
  e->flow += amt;
  G[e->to][e->index].flow -= amt;
     totflow += amt;
  }
  return totflow;
}
long long GetMaxFlow(int s, int t) {
  long long totflow = 0;
  while (long long flow = BlockingFlow(s, t))
     totflow += flow;
  return totflow;
}
```

};

```
MinCostMaxFlow.cc 2/27
// Implementation of min cost max flow algorithm using 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.
#include <cmath>
#include <vector>
#include <iostream>
using namespace std;
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 \text{ val} = \text{dist}[s] + \text{pi}[s] - \text{pi}[k] + \text{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 \parallel 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);
};
PushRelabel.cc 3/27
// 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.
//
// Running time:
//
       O(|V|^{3})
// INPUT:
//
       - graph, constructed using AddEdge()
//
       - source
//
       - sink
// OUTPUT:
//
       - maximum flow value
//
       - To obtain the actual flow values, look at all edges with
//
          capacity > 0 (zero capacity edges are residual edges).
#include <cmath>
#include <vector>
```

#include <iostream>
#include <queue>

```
using namespace std;
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) {}
};
struct PushRelabel {
  int N;
  vector<vector<Edge> > G;
  vector<LL> excess;
  vector<int> dist, active, count;
  queue<int> Q;
  PushRelabel(int N): N(N), G(N), excess(N), dist(N), active(N), count(2*N) {}
  void AddEdge(int from, int to, int 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, G[from].size() - 1));
  }
  void Enqueue(int v) {
     if (!active[v] && excess[v] > 0) { active[v] = true; Q.push(v); }
  }
  void Push(Edge &e) {
     int amt = int(min(excess[e.from], LL(e.cap - e.flow)));
    if (dist[e.from] \le dist[e.to] \parallel 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(int k) {
     for (int 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(int v) {
  count[dist[v]]--;
  dist[v] = 2*N;
  for (int 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(int v) {
  for (int 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(int s, int t) {
  count[0] = N-1;
  count[N] = 1;
  dist[s] = N;
  active[s] = active[t] = true;
  for (int i = 0; i < G[s].size(); i++) {
     excess[s] += G[s][i].cap;
     Push(G[s][i]);
  }
  while (!Q.empty()) {
     int v = Q.front();
     Q.pop();
     active[v] = false;
     Discharge(v);
  }
  LL totflow = 0;
  for (int i = 0; i < G[s].size(); i++) totflow += G[s][i].flow;
```

```
return totflow;
  }
};
MinCostMatching.cc 4/27
// Min cost bipartite matching via shortest augmenting paths
//
// This is an O(n^3) implementation of a shortest augmenting path
// algorithm for finding min cost perfect matchings in dense
// graphs. In practice, it solves 1000x1000 problems in around 1
// second.
//
//
    cost[i][j] = cost for pairing left node i with right node j
    Lmate[i] = index of right node that left node i pairs with
//
//
    Rmate[j] = index of left node that right node j pairs with
//
// The values in cost[i][j] may be positive or negative. To perform
// maximization, simply negate the cost[][] matrix.
#include <algorithm>
#include <cstdio>
#include <cmath>
#include <vector>
using namespace std;
typedef vector<double> VD;
typedef vector<VD> VVD;
typedef vector<int> VI;
double MinCostMatching(const VVD &cost, VI &Lmate, VI &Rmate) {
  int n = int(cost.size());
  // construct dual feasible solution
  VD u(n);
  VD v(n);
  for (int i = 0; i < n; i++) {
    u[i] = cost[i][0];
```

```
for (int j = 1; j < n; j++) u[i] = min(u[i], cost[i][j]);
}
for (int j = 0; j < n; j++) {
  v[j] = cost[0][j] - u[0];
  for (int i = 1; i < n; i++) v[j] = min(v[j], cost[i][j] - u[i]);
}
// construct primal solution satisfying complementary slackness
Lmate = VI(n, -1);
Rmate = VI(n, -1);
int mated = 0;
for (int i = 0; i < n; i++) {
  for (int j = 0; j < n; j++) {
     if (Rmate[j] != -1) continue;
     if (fabs(cost[i][j] - u[i] - v[j]) < 1e-10) {
  Lmate[i] = j;
  Rmate[j] = i;
  mated++;
  break;
     }
  }
}
VD dist(n);
VI dad(n);
VI seen(n);
// repeat until primal solution is feasible
while (mated < n) {
  // find an unmatched left node
  int s = 0;
  while (Lmate[s] !=-1) s++;
  // initialize Dijkstra
  fill(dad.begin(), dad.end(), -1);
  fill(seen.begin(), seen.end(), 0);
  for (int k = 0; k < n; k++)
     dist[k] = cost[s][k] - u[s] - v[k];
  int j = 0;
  while (true) {
     // find closest
```

```
j = -1;
   for (int k = 0; k < n; k++) {
if (seen[k]) continue;
if (j == -1 || dist[k] < dist[j]) j = k;
   }
   seen[j] = 1;
   // termination condition
   if (Rmate[j] == -1) break;
   // relax neighbors
   const int i = Rmate[j];
   for (int k = 0; k < n; k++) {
if (seen[k]) continue;
const double new_dist = dist[j] + cost[i][k] - u[i] - v[k];
if (dist[k] > new_dist) {
   dist[k] = new_dist;
  dad[k] = j;
}
   }
}
// update dual variables
for (int k = 0; k < n; k++) {
   if (k == j \parallel !seen[k]) continue;
   const int i = Rmate[k];
   v[k] += dist[k] - dist[j];
   u[i] = dist[k] - dist[j];
}
u[s] += dist[j];
// augment along path
while (dad[j] >= 0) {
   const int d = dad[j];
   Rmate[j] = Rmate[d];
   Lmate[Rmate[j]] = j;
  j = d;
}
Rmate[j] = s;
Lmate[s] = j;
mated++;
```

}

```
double value = 0;
  for (int i = 0; i < n; i++)
     value += cost[i][Lmate[i]];
  return value;
}
MaxBipartiteMatching.cc 5/27
// This code performs maximum bipartite matching.
// Running time: O(|E||V|) -- often much faster in practice
//
//
     INPUT: w[i][j] = edge between row node i and column node j
//
     OUTPUT: mr[i] = assignment for row node i, -1 if unassigned
//
               mc[j] = assignment for column node j, -1 if unassigned
//
               function returns number of matches made
#include <vector>
using namespace std;
typedef vector<int> VI;
typedef vector<VI>VVI;
bool FindMatch(int i, const VVI &w, VI &mr, VI &mc, VI &seen) {
  for (int j = 0; j < w[i].size(); j++) {
     if (w[i][j] && !seen[j]) {
       seen[j] = true;
       if (mc[j] < 0 \parallel FindMatch(mc[j], w, mr, mc, seen)) {
          mr[i] = j;
          mc[j] = i;
          return true;
       }
     }
  return false;
}
int BipartiteMatching(const VVI &w, VI &mr, VI &mc) {
  mr = VI(w.size(), -1);
  mc = VI(w[0].size(), -1);
```

```
int ct = 0;
  for (int i = 0; i < w.size(); i++) {
     VI seen(w[0].size());
    if (FindMatch(i, w, mr, mc, seen)) ct++;
  }
  return ct;
}
MinCut.cc 6/27
// Adjacency matrix implementation of Stoer-Wagner min cut algorithm.
//
// Running time:
//
       O(|V|^{3})
//
// INPUT:
//
       - graph, constructed using AddEdge()
//
// OUTPUT:
       - (min cut value, nodes in half of min cut)
#include <cmath>
#include <vector>
#include <iostream>
using namespace std;
typedef vector<int> VI;
typedef vector<VI>VVI;
const int INF = 10000000000;
pair<int, VI> GetMinCut(VVI &weights) {
  int N = weights.size();
  VI used(N), cut, best_cut;
  int best_weight = -1;
  for (int phase = N-1; phase >= 0; phase--) {
     VI w = weights[0];
     VI added = used;
     int prev, last = 0;
```

```
for (int i = 0; i < phase; i++) {
       prev = last;
       last = -1;
       for (int j = 1; j < N; j++)
     if (!added[j] && (last == -1 \parallel w[j] > w[last])) last = j;
       if (i == phase-1) {
     for (int j = 0; j < N; j++) weights[prev][j] += weights[last][j];
     for (int j = 0; j < N; j++) weights[j][prev] = weights[prev][j];
     used[last] = true;
     cut.push_back(last);
     if (best_weight == -1 || w[last] < best_weight) {
       best_cut = cut;
       best_weight = w[last];
       } else {
     for (int j = 0; j < N; j++)
       w[j] += weights[last][j];
     added[last] = true;
       }
     }
  }
  return make_pair(best_weight, best_cut);
}
ConvexHull.cc 7/27
// Compute the 2D convex hull of a set of points using the monotone chain
// algorithm. Eliminate redundant points from the hull if REMOVE_REDUNDANT is
// #defined.
//
// Running time: O(n log n)
//
//
     INPUT:
                 a vector of input points, unordered.
//
     OUTPUT: a vector of points in the convex hull, counterclockwise, starting
//
                with bottommost/leftmost point
#include <cstdio>
#include <cassert>
#include <vector>
#include <algorithm>
```

#include <cmath>

```
using namespace std;
#define REMOVE REDUNDANT
typedef double T;
const T EPS = 1e-7;
struct PT {
  T x, y;
  PT() {}
  PT(T x, T y) : x(x), y(y) \{ \}
  bool operator<(const PT &rhs) const { return make_pair(y,x) < make_pair(rhs.y,rhs.x); }
  bool operator==(const PT &rhs) const { return make_pair(y,x) == make_pair(rhs.y,rhs.x); }
};
T cross(PT p, PT q) { return p.x*q.y-p.y*q.x; }
T area2(PT a, PT b, PT c) { return cross(a,b) + cross(b,c) + cross(c,a); }
#ifdef REMOVE_REDUNDANT
bool between(const PT &a, const PT &b, const PT &c) {
  return (fabs(area2(a,b,c)) < EPS && (a.x-b.x)*(c.x-b.x) <= 0 && (a.y-b.y)*(c.y-b.y) <= 0);
}
#endif
void ConvexHull(vector<PT> &pts) {
  sort(pts.begin(), pts.end());
  pts.erase(unique(pts.begin(), pts.end()), pts.end());
  vector<PT> up, dn;
  for (int i = 0; i < pts.size(); i++) {
     while (up.size() > 1 \&\& area2(up[up.size()-2], up.back(), pts[i]) >= 0) up.pop_back();
     while (dn.size() > 1 && area2(dn[dn.size()-2], dn.back(), pts[i]) <= 0) dn.pop_back();
     up.push_back(pts[i]);
    dn.push_back(pts[i]);
  }
  pts = dn;
  for (int i = (int) \text{ up.size}() - 2; i \ge 1; i--) \text{ pts.push\_back}(\text{up}[i]);
#ifdef REMOVE_REDUNDANT
  if (pts.size() <= 2) return;
  dn.clear();
  dn.push_back(pts[0]);
  dn.push_back(pts[1]);
  for (int i = 2; i < pts.size(); i++) {
     if (between(dn[dn.size()-2], dn[dn.size()-1], pts[i])) dn.pop_back();
```

```
dn.push_back(pts[i]);
  }
  if (dn.size() \ge 3 \&\& between(dn.back(), dn[0], dn[1])) {
     dn[0] = dn.back();
    dn.pop_back();
  }
  pts = dn;
#endif
}
Geometry.cc 8/27
// C++ routines for computational geometry.
#include <iostream>
#include <vector>
#include <cmath>
#include <cassert>
using namespace std;
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; }
ostream & operator << (ostream & os, const PT & p) {
  os << "(" << p.x << "," << p.y << ")";
}
```

```
// 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 \parallel dist2(a, d) < EPS \parallel
       dist2(b, c) < EPS \parallel 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 \le q.y \&\& q.y < p[j].y ||
       p[j].y \le 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 = \operatorname{sqrt}(\operatorname{dist2}(a, b));
  if (d > r + R \parallel 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 1 = (k+1) \% p.size();
       if (i == 1 || j == k) continue;
       if \ (SegmentsIntersect(p[i], p[j], p[k], p[l])) \\
          return false;
     }
  }
```

```
return true;
int main() {
  // expected: (-5,2)
  cerr << RotateCCW90(PT(2,5)) << endl;
  // expected: (5,-2)
  cerr << RotateCW90(PT(2,5)) << endl;
  // expected: (-5,2)
  cerr << RotateCCW(PT(2,5),M_PI/2) << endl;
  // expected: (5,2)
  cerr << ProjectPointLine(PT(-5,-2), PT(10,4), PT(3,7)) << endl;
  // expected: (5,2)(7.5,3)(2.5,1)
  cerr << ProjectPointSegment(PT(-5,-2), PT(10,4), PT(3,7)) << " "
        << ProjectPointSegment(PT(7.5,3), PT(10,4), PT(3,7)) << " "
        << ProjectPointSegment(PT(-5,-2), PT(2.5,1), PT(3,7)) << endl;
  // expected: 6.78903
  cerr << DistancePointPlane(4,-4,3,2,-2,5,-8) << endl;
  // expected: 1 0 1
  cerr << LinesParallel(PT(1,1), PT(3,5), PT(2,1), PT(4,5)) << " "
        << LinesParallel(PT(1,1), PT(3,5), PT(2,0), PT(4,5)) << " "
        << LinesParallel(PT(1,1), PT(3,5), PT(5,9), PT(7,13)) << endl;
  // expected: 0 0 1
  cerr << LinesCollinear(PT(1,1), PT(3,5), PT(2,1), PT(4,5)) << " "
        << LinesCollinear(PT(1,1), PT(3,5), PT(2,0), PT(4,5)) << " "
        << LinesCollinear(PT(1,1), PT(3,5), PT(5,9), PT(7,13)) << endl;
  // expected: 1 1 1 0
  cerr << SegmentsIntersect(PT(0,0), PT(2,4), PT(3,1), PT(-1,3)) << " "
        << SegmentsIntersect(PT(0,0), PT(2,4), PT(4,3), PT(0,5)) << " "
        << SegmentsIntersect(PT(0,0), PT(2,4), PT(2,-1), PT(-2,1)) << " "
        << SegmentsIntersect(PT(0,0), PT(2,4), PT(5,5), PT(1,7)) << endl;
  // expected: (1,2)
  cerr << ComputeLineIntersection(PT(0,0), PT(2,4), PT(3,1), PT(-1,3)) << endl;
```

```
// expected: (1,1)
cerr << ComputeCircleCenter(PT(-3,4), PT(6,1), PT(4,5)) << endl;
vector<PT> v;
v.push_back(PT(0,0));
v.push_back(PT(5,0));
v.push_back(PT(5,5));
v.push_back(PT(0,5));
// expected: 1 1 1 0 0
cerr << PointInPolygon(v, PT(2,2)) << " "
      << PointInPolygon(v, PT(2,0)) << " "
      << PointInPolygon(v, PT(0,2)) << " "
      << PointInPolygon(v, PT(5,2)) << " "
      << PointInPolygon(v, PT(2,5)) << endl;
// expected: 0 1 1 1 1
cerr << PointOnPolygon(v, PT(2,2)) << " "
      << PointOnPolygon(v, PT(2,0)) << " "
      << PointOnPolygon(v, PT(0,2)) << " "
      << PointOnPolygon(v, PT(5,2)) << " "
      << PointOnPolygon(v, PT(2,5)) << endl;
// expected: (1,6)
//
               (5,4)(4,5)
//
               blank line
//
               (4,5)(5,4)
//
               blank line
               (4,5)(5,4)
vector<PT> u = CircleLineIntersection(PT(0,6), PT(2,6), PT(1,1), 5);
for (int i = 0; i < u.size(); i++) cerr << u[i] << " "; cerr <math><< endl;
u = CircleLineIntersection(PT(0,9), PT(9,0), PT(1,1), 5);
for (int i = 0; i < u.size(); i++) cerr << u[i] << " "; cerr <math><< endl;
u = CircleCircleIntersection(PT(1,1), PT(10,10), 5, 5);
for (int i = 0; i < u.size(); i++) cerr << u[i] << " "; cerr <math><< endl;
u = CircleCircleIntersection(PT(1,1), PT(8,8), 5, 5);
for (int i = 0; i < u.size(); i++) cerr << u[i] << " "; cerr <math><< endl;
u = CircleCircleIntersection(PT(1,1), PT(4.5,4.5), 10, sqrt(2.0)/2.0);
for (int i = 0; i < u.size(); i++) cerr << u[i] << " "; cerr <math><< endl;
u = CircleCircleIntersection(PT(1,1), PT(4.5,4.5), 5, sqrt(2.0)/2.0);
for (int i = 0; i < u.size(); i++) cerr << u[i] << " "; cerr <math><< endl;
// area should be 5.0
// centroid should be (1.1666666, 1.166666)
```

```
PT pa[] = \{ PT(0,0), PT(5,0), PT(1,1), PT(0,5) \};
  vector<PT> p(pa, pa+4);
  PT c = ComputeCentroid(p);
  cerr << "Area: " << ComputeArea(p) << endl;</pre>
  cerr << "Centroid: " << c << endl;
  return 0;
}
JavaGeometry.java 9/27
// In this example, we read an input file containing three lines, each
// containing an even number of doubles, separated by commas. The first two
// lines represent the coordinates of two polygons, given in counterclockwise
// (or clockwise) order, which we will call "A" and "B".  The last line
// contains a list of points, p[1], p[2], ...
//
// Our goal is to determine:
     (1) whether B - A is a single closed shape (as opposed to multiple shapes)
//
     (2) the area of B - A
//
     (3) whether each p[i] is in the interior of B - A
//
// INPUT:
//
     00100010
     0 0 10 10 10 0
     86
//
     5 1
//
// OUTPUT:
     The area is singular.
//
//
     The area is 25.0
     Point belongs to the area.
//
     Point does not belong to the area.
import java.util.*;
import java.awt.geom.*;
import java.io.*;
public class JavaGeometry {
     // make an array of doubles from a string
     static double[] readPoints(String s) {
```

```
String[] arr = s.trim().split("\s++");
     double[] ret = new double[arr.length];
     for (int i = 0; i < arr.length; i++) ret[i] = Double.parseDouble(arr[i]);
     return ret;
}
// make an Area object from the coordinates of a polygon
static Area makeArea(double[] pts) {
     Path2D.Double p = new Path2D.Double();
     p.moveTo(pts[0], pts[1]);
     for (int i = 2; i < pts.length; i += 2) p.lineTo(pts[i], pts[i+1]);
     p.closePath();
     return new Area(p);
}
// compute area of polygon
static double computePolygonArea(ArrayList<Point2D.Double> points) {
     Point2D.Double[] pts = points.toArray(new Point2D.Double[points.size()]);
     double area = 0;
    for (int i = 0; i < pts.length; i++){
          int j = (i+1) % pts.length;
          area += pts[i].x * pts[j].y - pts[j].x * pts[i].y;
     return Math.abs(area)/2;
}
// compute the area of an Area object containing several disjoint polygons
static double computeArea(Area area) {
     double totArea = 0;
     PathIterator iter = area.getPathIterator(null);
     ArrayList<Point2D.Double> points = new ArrayList<Point2D.Double>();
     while (!iter.isDone()) {
          double[] buffer = new double[6];
          switch (iter.currentSegment(buffer)) {
          case PathIterator.SEG_MOVETO:
          case PathIterator.SEG_LINETO:
               points.add(new Point2D.Double(buffer[0], buffer[1]));
              break;
          case PathIterator.SEG_CLOSE:
              totArea += computePolygonArea(points);
              points.clear();
              break;
          }
```

```
iter.next();
     }
     return totArea;
}
// notice that the main() throws an Exception -- necessary to
// avoid wrapping the Scanner object for file reading in a
// try { ... } catch block.
public static void main(String args[]) throws Exception {
     Scanner scanner = new Scanner(new File("input.txt"));
     // also,
          Scanner scanner = new Scanner (System.in);
     double[] pointsA = readPoints(scanner.nextLine());
     double[] pointsB = readPoints(scanner.nextLine());
     Area areaA = makeArea(pointsA);
     Area areaB = makeArea(pointsB);
     areaB.subtract(areaA);
     // also,
          areaB.exclusiveOr (areaA);
     //
          areaB.add (areaA);
          areaB.intersect (areaA);
     // (1) determine whether B - A is a single closed shape (as
            opposed to multiple shapes)
     boolean isSingle = areaB.isSingular();
     // also,
          areaB.isEmpty();
     if (isSingle)
          System.out.println("The area is singular.");
     else
          System.out.println("The area is not singular.");
     // (2) compute the area of B - A
     System.out.println("The area is " + computeArea(areaB) + ".");
     // (3) determine whether each p[i] is in the interior of B - A
     while (scanner.hasNextDouble()) {
          double x = scanner.nextDouble();
          assert(scanner.hasNextDouble());
          double y = scanner.nextDouble();
```

```
if (areaB.contains(x,y)) {
                    System.out.println ("Point belongs to the area.");
               } else {
                   System.out.println ("Point does not belong to the area.");
               }
          }
          // Finally, some useful things we didn't use in this example:
          //
               Ellipse2D.Double ellipse = new Ellipse2D.Double (double x, double y,
          //
          //
                                                                             double w, double h);
          //
          //
                 creates an ellipse inscribed in box with bottom-left corner (x,y)
                 and upper-right corner (x+y,w+h)
          //
          //
               Rectangle2D.Double rect = new Rectangle2D.Double (double x, double y,
          //
                                                                              double w, double h);
          //
          //
                 creates a box with bottom-left corner (x,y) and upper-right
          //
                 corner (x+y,w+h)
          //
         // Each of these can be embedded in an Area object (e.g., new Area (rect)).
     }
}
Geom3D.java 10/27
public class Geom3D {
  // distance from point (x, y, z) to plane aX + bY + cZ + d = 0
  public static double ptPlaneDist(double x, double y, double z,
       double a, double b, double c, double d) {
     return Math.abs(a*x + b*y + c*z + d) / Math.sqrt(a*a + b*b + c*c);
  }
  // distance between parallel planes aX + bY + cZ + d1 = 0 and
  // aX + bY + cZ + d2 = 0
  public static double planePlaneDist(double a, double b, double c,
       double d1, double d2) {
     return Math.abs(d1 - d2) / Math.sqrt(a*a + b*b + c*c);
  }
```

```
// distance from point (px, py, pz) to line (x1, y1, z1)-(x2, y2, z2)
// (or ray, or segment; in the case of the ray, the endpoint is the
// first point)
public static final int LINE = 0;
public static final int SEGMENT = 1;
public static final int RAY = 2;
public static double ptLineDistSq(double x1, double y1, double z1,
     double x2, double y2, double z2, double px, double py, double pz,
     int type) {
  double pd2 = (x1-x2)*(x1-x2) + (y1-y2)*(y1-y2) + (z1-z2)*(z1-z2);
  double x, y, z;
  if (pd2 == 0) {
     x = x1;
     y = y1;
     z = z1;
  } else {
     double u = ((px-x1)*(x2-x1) + (py-y1)*(y2-y1) + (pz-z1)*(z2-z1)) / pd2;
     x = x1 + u * (x2 - x1);
     y = y1 + u * (y2 - y1);
     z = z1 + u * (z2 - z1);
     if (type != LINE && u < 0) {
       x = x1;
       y = y1;
       z = z1;
     if (type == SEGMENT && u > 1.0) {
       x = x2;
       y = y2;
       z = z2;
     }
  }
  return (x-px)*(x-px) + (y-py)*(y-py) + (z-pz)*(z-pz);
public static double ptLineDist(double x1, double y1, double z1,
     double x2, double y2, double z2, double px, double py, double pz,
     int type) {
  return Math.sqrt(ptLineDistSq(x1, y1, z1, x2, y2, z2, px, py, pz, type));
}
```

}

```
Delaunay.cc 11/27
// Slow but simple Delaunay triangulation. Does not handle
// degenerate cases (from O'Rourke, Computational Geometry in C)
// Running time: O(n^4)
//
// INPUT:
               x[] = x-coordinates
//
               y[] = y-coordinates
// OUTPUT:
                triples = a vector containing m triples of indices
//
                            corresponding to triangle vertices
#include<vector>
using namespace std;
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 (i == 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
     //
                   032
     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;
}
Euclid.cc 12/27
// This is a collection of useful code for solving problems that
// involve modular linear equations. Note that all of the
// algorithms described here work on nonnegative integers.
#include <iostream>
#include <vector>
#include <algorithm>
using namespace std;
typedef vector<int> VI;
typedef pair<int,int> PII;
// return a % b (positive value)
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;
}
// returns d = gcd(a,b); finds x,y such that d = ax + by
int extended_euclid(int a, int b, int &x, int &y) {
  int xx = y = 0;
  int yy = x = 1;
  while (b) {
     int q = a/b;
     int t = b; b = a\%b; a = t;
     t = xx; xx = x-q*xx; x = t;
     t = yy; yy = y-q*yy; y = t;
  return a;
}
// finds all solutions to ax = b \pmod{n}
VI modular_linear_equation_solver(int a, int b, int n) {
  int x, y;
  VI solutions;
  int d = extended_euclid(a, n, x, y);
  if (!(b%d)) {
     x = mod(x*(b/d), n);
     for (int i = 0; i < d; i++)
        solutions.push\_back(mod(x + i*(n/d), n));
  }
  return solutions;
}
// computes b such that ab = 1 \pmod{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);
}
// Chinese remainder theorem (special case): find z such that
// z % x = a, z % y = b. Here, z is unique modulo M = lcm(x,y).
// Return (z,M). On failure, M = -1.
PII chinese_remainder_theorem(int x, int a, int y, int b) {
  int s. t:
  int d = \text{extended\_euclid}(x, y, s, t);
  if (a\%d != b\%d) return make_pair(0, -1);
  return make_pair(mod(s*b*x+t*a*y,x*y)/d, x*y/d);
}
// Chinese remainder theorem: find z such that
// z % x[i] = a[i] for all i. Note that the solution is
// unique modulo M = lcm_i (x[i]). Return (z,M). On
// failure, M = -1. Note that we do not require the a[i]'s
// to be relatively prime.
PII chinese_remainder_theorem(const VI &x, const VI &a) {
  PII ret = make_pair(a[0], x[0]);
  for (int i = 1; i < x.size(); i++) {
     ret = chinese_remainder_theorem(ret.first, ret.second, x[i], a[i]);
     if (ret.second == -1) break;
  }
  return ret;
}
// 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;
  }
}
int main() {
  // expected: 2
  cout << gcd(14, 30) << endl;
```

```
// expected: 2 -2 1
  int x, y;
  int d = \text{extended\_euclid}(14, 30, x, y);
  cout << d << " \ " << x << " \ " << y << endl;
  // expected: 95 45
  VI sols = modular_linear_equation_solver(14, 30, 100);
  for (int i = 0; i < (int) sols.size(); i++) cout << sols[i] << " ";
  cout << endl;
  // expected: 8
  cout << mod_inverse(8, 9) << endl;</pre>
  // expected: 23 56
  //
                  11 12
  int xs[] = \{3, 5, 7, 4, 6\};
  int as[] = \{2, 3, 2, 3, 5\};
  PII ret = chinese_remainder_theorem(VI (xs, xs+3), VI(as, as+3));
  cout << ret.first << " " << ret.second << endl;</pre>
  ret = chinese\_remainder\_theorem (VI(xs+3, xs+5), VI(as+3, as+5));
  cout << ret.first << " " << ret.second << endl;</pre>
  // expected: 5 -15
  linear_diophantine(7, 2, 5, x, y);
  cout << x << " " << y << endl;
}
GaussJordan.cc 13/27
// Gauss-Jordan 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[][]
#include <iostream>
#include <vector>
#include <cmath>
using namespace std;
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 p_i = -1, p_i = -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[pk]++;
     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;</pre>
  // 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;</pre>
  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;</pre>
```

```
for (int i = 0; i < n; i++) {
     for (int j = 0; j < m; j++)
        cout << b[i][j] << ' ';
     cout << endl;
  }
}
```

```
ReducedRowEchelonForm.cc 14/27
// Reduced row echelon form via Gauss-Jordan elimination
// with partial pivoting. This can be used for computing
// the rank of a matrix.
//
// Running time: O(n^3)
//
// INPUT:
              a[][] = an nxn matrix
//
// OUTPUT: rref[][] = an nxm matrix (stored in a[][])
              returns rank of a[][]
#include <iostream>
#include <vector>
#include <cmath>
using namespace std;
const double EPSILON = 1e-10;
typedef double T;
typedef vector<T> VT;
typedef vector<VT> VVT;
int rref(VVT &a) {
  int n = a.size();
  int m = a[0].size();
  int r = 0;
  for (int c = 0; c < m; c++) {
    int j = r;
    for (int i = r+1; i < n; i++)
       if (fabs(a[i][c]) > fabs(a[j][c])) j = i;
     if (fabs(a[j][c]) < EPSILON) continue;
     swap(a[j], a[r]);
```

```
T s = 1.0 / a[r][c];
     for (int j = 0; j < m; j++) a[r][j] *= s;
     for \ (int \ i = 0; \ i < n; \ i++) \ if \ (i \ != r) \ \{
        T t = a[i][c];
        for (int j = 0; j < m; j++) a[i][j] -= t * a[r][j];
     }
     r++;
  return r;
}
int main(){
  const int n = 5;
  const int m = 4;
  double\ A[n][m] = \{\ \{16,2,3,13\}, \{5,11,10,8\}, \{9,7,6,12\}, \{4,14,15,1\}, \{13,21,21,13\}\ \};
  VVT a(n);
  for (int i = 0; i < n; i++)
     a[i] = VT(A[i], A[i] + n);
  int rank = rref(a);
  // expected: 4
  cout << "Rank: " << rank << endl;
  // expected: 1 0 0 1
  //
                   0\,1\,0\,3
                   001-3
  //
  //
                   0 0 0 2.78206e-15
                   0 0 0 3.22398e-15
  cout << "rref: " << endl;
  for (int i = 0; i < 5; i++){
     for (int j = 0; j < 4; j++)
        cout << a[i][j] << ' ';
     cout << endl;
  }
}
```

```
FFT_new.cpp 15/27 #include <cassert>
```

```
#include <cstdio>
#include <cmath>
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 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(double theta)
{
  return cpx(cos(theta),sin(theta));
const double two_pi = 4 * acos(0);
// in:
           input array
```

```
// out:
           output array
// step:
          {SET TO 1} (used internally)
// size:
          length of the input/output {MUST BE A POWER OF 2}
// dir:
          either plus or minus one (direction of the FFT)
// RESULT: out[k] = \sum_{j=0}^{size - 1} in[j] * exp(dir * 2pi * i * j * k / size)
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(dir * two_pi * i / size) * odd;
     out[i + size / 2] = even + EXP(dir * two_pi * (i + size / 2) / size) * odd;
  }
}
// Usage:
// f[0...N-1] and g[0...N-1] are numbers
// Want to compute the convolution h, defined by
// h[n] = sum of f[k]g[n-k] (k = 0, ..., N-1).
// Here, the index is cyclic; f[-1] = f[N-1], f[-2] = f[N-2], etc.
// Let F[0...N-1] be FFT(f), and similarly, define G and H.
// The convolution theorem says H[n] = F[n]G[n] (element-wise product).
// To compute h[] in O(N log N) time, do the following:
//
     1. Compute F and G (pass dir = 1 as the argument).
     2. Get H by element-wise multiplying F and G.
//
     3. Get h by taking the inverse FFT (use dir = -1 as the argument)
//
         and *dividing by N*. DO NOT FORGET THIS SCALING FACTOR.
int main(void)
  printf("If rows come in identical pairs, then everything works.\n");
  cpx a[8] = \{0, 1, cpx(1,3), cpx(0,5), 1, 0, 2, 0\};
  cpx b[8] = \{1, cpx(0,-2), cpx(0,1), 3, -1, -3, 1, -2\};
  cpx A[8];
```

```
cpx B[8];
FFT(a, A, 1, 8, 1);
FFT(b, B, 1, 8, 1);
for(int i = 0; i < 8; i++)
  printf("%7.21f%7.21f", A[i].a, A[i].b);
printf("\n");
for(int i = 0; i < 8; i++)
  cpx Ai(0,0);
  for(int j = 0; j < 8; j++)
     Ai = Ai + a[j] * EXP(j * i * two_pi / 8);
  printf("%7.21f%7.21f", Ai.a, Ai.b);
printf("\n");
cpx AB[8];
for(int i = 0; i < 8; i++)
  AB[i] = A[i] * B[i];
cpx aconvb[8];
FFT(AB, aconvb, 1, 8, -1);
for(int i = 0; i < 8; i++)
  aconvb[i] = aconvb[i] / 8;
for(int i = 0; i < 8; i++)
  printf("%7.2lf%7.2lf", aconvb[i].a, aconvb[i].b);
printf("\n");
for(int i = 0; i < 8; i++)
  cpx aconvbi(0,0);
  for(int j = 0; j < 8; j++)
     aconvbi = aconvbi + a[j] * b[(8 + i - j) % 8];
  printf("%7.2lf%7.2lf", aconvbi.a, aconvbi.b);
printf("\n");
return 0;
```

```
Simplex.cc 16/27
// Two-phase simplex algorithm for solving linear programs of the form
//
                      c^T x
//
       maximize
       subject to
                    Ax \le b
//
                      x >= 0
// INPUT: A -- an m x n matrix
          b -- an m-dimensional vector
//
          c -- an n-dimensional vector
//
          x -- a vector where the optimal solution will be stored
//
// OUTPUT: value of the optimal solution (infinity if unbounded
//
            above, nan if infeasible)
//
// To use this code, create an LPSolver object with A, b, and c as
// arguments. Then, call Solve(x).
#include <iostream>
#include <iomanip>
#include <vector>
#include <cmath>
#include imits>
using namespace std;
typedef long double DOUBLE;
typedef vector<DOUBLE> VD;
typedef vector<VD> VVD;
typedef vector<int> VI;
const DOUBLE EPS = 1e-9;
struct LPSolver {
  int m, n;
  VIB, N;
  VVD D;
```

```
LPSolver(const VVD &A, const VD &b, const VD &c):
  m(b.size()), n(c.size()), N(n+1), B(m), D(m+2, VD(n+2)) 
  for (int i = 0; i < m; i++) for (int j = 0; j < n; j++) D[i][j] = A[i][j];
  for (int i = 0; i < m; i++) { B[i] = n+i; D[i][n] = -1; D[i][n+1] = b[i]; }
  for (int j = 0; j < n; j++) { N[j] = j; D[m][j] = -c[j]; }
  N[n] = -1; D[m+1][n] = 1;
}
void Pivot(int r, int s) {
  for (int i = 0; i < m+2; i++) if (i != r)
     for (int j = 0; j < n+2; j++) if (j != s)
  D[i][j] = D[r][j] * D[i][s] / D[r][s];
  for (int j = 0; j < n+2; j++) if (j != s) D[r][j] /= D[r][s];
  for (int i = 0; i < m+2; i++) if (i != r) D[i][s] /= -D[r][s];
  D[r][s] = 1.0 / D[r][s];
  swap(B[r], N[s]);
}
bool Simplex(int phase) {
  int x = phase == 1 ? m+1 : m;
  while (true) {
     int s = -1;
     for (int j = 0; j \le n; j++) {
  if (phase == 2 \&\& N[j] == -1) continue;
  if (s == -1 \parallel D[x][j] < D[x][s] \parallel D[x][j] == D[x][s] \&\& N[j] < N[s]) s = j;
     if (D[x][s] \ge -EPS) return true;
     int r = -1;
     for (int i = 0; i < m; i++) {
  if (D[i][s] \le 0) continue;
  if (r == -1 \parallel D[i][n+1] / D[i][s] < D[r][n+1] / D[r][s] \parallel
        D[i][n+1] / D[i][s] == D[r][n+1] / D[r][s] && B[i] < B[r]) r = i;
     if (r == -1) return false;
     Pivot(r, s);
  }
}
DOUBLE Solve(VD &x) {
  int r = 0;
  for (int i = 1; i < m; i++) if (D[i][n+1] < D[r][n+1]) r = i;
  if (D[r][n+1] \le -EPS) {
     Pivot(r, n);
     if (!Simplex(1) || D[m+1][n+1] < -EPS) return -numeric_limits<DOUBLE>::infinity();
```

```
for (int i = 0; i < m; i++) if (B[i] == -1) {
     int s = -1;
     for (int j = 0; j \le n; j++)
       if \ (s == -1 \parallel D[i][j] < D[i][s] \parallel D[i][j] == D[i][s] \ \&\& \ N[j] < N[s]) \ s = j;
     Pivot(i, s);
       }
     if (!Simplex(2)) return numeric_limits<DOUBLE>::infinity();
     x = VD(n);
     for (int i = 0; i < m; i++) if (B[i] < n) \times [B[i]] = D[i][n+1];
     return D[m][n+1];
  }
};
int main() {
  const int m = 4;
  const int n = 3;
  DOUBLE A[m][n] = {
     \{6, -1, 0\},\
     \{-1, -5, 0\},\
     { 1, 5, 1 },
     \{-1, -5, -1\}
  };
  DOUBLE _b[m] = \{ 10, -4, 5, -5 \};
  DOUBLE _{c}[n] = \{ 1, -1, 0 \};
  VVDA(m);
  VD b(_b, _b + m);
  VD c(_c, _c + n);
  for (int i = 0; i < m; i++) A[i] = VD(\_A[i], \_A[i] + n);
  LPSolver solver(A, b, c);
  VD x;
  DOUBLE value = solver.Solve(x);
  cerr << "VALUE: "<< value << endl;
  cerr << "SOLUTION:";
  for (size_t i = 0; i < x.size(); i++) cerr << " " << x[i];
  cerr << endl;
  return 0;
}
```

.....

```
FastDijkstra.cc 17/27
// Implementation of Dijkstra's algorithm using adjacency lists
// and priority queue for efficiency.
//
// Running time: O(|E| log |V|)
#include <queue>
#include <stdio.h>
using namespace std;
const int INF = 20000000000;
typedef pair<int,int> PII;
int main(){
  int N, s, t;
  scanf ("%d%d%d", &N, &s, &t);
  vector < vector < PII > edges(N);
  for (int i = 0; i < N; i++)
     int M;
     scanf ("%d", &M);
     for (int j = 0; j < M; j++){
       int vertex, dist;
       scanf ("%d%d", &vertex, &dist);
       edges[i].push_back (make_pair (dist, vertex)); // note order of arguments here
  }
  // use priority queue in which top element has the "smallest" priority
  priority_queue<PII, vector<PII>, greater<PII> > Q;
  vector<int> dist(N, INF), dad(N, -1);
  Q.push (make_pair (0, s));
  dist[s] = 0;
  while (!Q.empty()){
     PII p = Q.top();
     if (p.second == t) break;
     Q.pop();
     int here = p.second;
     for (vector<PII>::iterator it=edges[here].begin(); it!=edges[here].end(); it++){
```

```
if (dist[here] + it->first < dist[it->second]){
          dist[it->second] = dist[here] + it->first;
          dad[it->second] = here;
          Q.push (make_pair (dist[it->second], it->second));
       }
     }
  printf ("%d\n", dist[t]);
  if (dist[t] < INF)
     for(int i=t;i!=-1;i=dad[i])
       printf ("%d%c", i, (i==s?'\n':' '));
  return 0;
}
SCC.cc 18/27
#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;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;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]);}
Suffix Array.cc 19/27
// Suffix array construction in O(L log^2 L) time. Routine for
// computing the length of the longest common prefix of any two
// suffixes in O(log L) time.
//
// INPUT:
              string s
// OUTPUT: array suffix[] such that suffix[i] = index (from 0 to L-1)
//
              of substring s[i...L-1] in the list of sorted suffixes.
//
              That is, if we take the inverse of the permutation suffix[],
              we get the actual suffix array.
//
#include <vector>
#include <iostream>
#include <string>
using namespace std;
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 \ge 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
  //
            1 is the 4'th suffix
  SuffixArray suffix("bobocel");
  vector<int> v = suffix.GetSuffixArray();
  // Expected output: 0 5 1 6 2 3 4
  for (int i = 0; i < v.size(); i++) cout << v[i] << " ";
  cout << endl;
  cout << suffix.LongestCommonPrefix(0, 2) << endl;</pre>
```

}

.....

```
BIT.cc 20/27
#include <iostream>
using namespace std;
#define LOGSZ 17
int tree[(1<<LOGSZ)+1];</pre>
int N = (1 << LOGSZ);
// add v to value at x
void set(int x, int v) {
  while(x \le N) {
     tree[x] += v;
     x += (x \& -x);
  }
}
// get cumulative sum up to and including \boldsymbol{x}
int get(int x) {
  int res = 0;
  while(x) {
     res += tree[x];
     x = (x \& -x);
  }
  return res;
}
// get largest value with cumulative sum less than or equal to x;
// for smallest, pass x-1 and add 1 to result
int getind(int x) {
  int idx = 0, mask = N;
  while (mask && idx < N) {
     int t = idx + mask;
     if(x \ge tree[t]) {
       idx = t;
        x = tree[t];
     }
     mask >>= 1;
   }
  return idx;
```

```
}
UnionFind.cc 21/27
//union-find set: the vector/array contains the parent of each node
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]);} //C
KDTree.cc 22/27
// -----
// A straightforward, but probably sub-optimal KD-tree implmentation that's
// probably good enough for most things (current it's a 2D-tree)
//
// - constructs from n points in O(n lg^2 n) time
// - handles nearest-neighbor query in O(lg n) if points are well distributed
// - worst case for nearest-neighbor may be linear in pathological case
//
// Sonny Chan, Stanford University, April 2009
// -----
#include <iostream>
#include <vector>
#include inits>
#include <cstdlib>
using namespace std;
// number type for coordinates, and its maximum value
typedef long long ntype;
const ntype sentry = numeric_limits<ntype>::max();
// point structure for 2D-tree, can be extended to 3D
struct point {
    ntype x, y;
    point(ntype xx = 0, ntype yy = 0): x(xx), y(yy) {}
```

};

bool operator==(const point &a, const point &b)

```
return a.x == b.x && a.y == b.y;
}
// sorts points on x-coordinate
bool on_x(const point &a, const point &b)
{
     return a.x < b.x;
}
// sorts points on y-coordinate
bool on_y(const point &a, const point &b)
{
    return a.y < b.y;
}
// squared distance between points
ntype pdist2(const point &a, const point &b)
{
    ntype dx = a.x-b.x, dy = a.y-b.y;
    return dx*dx + dy*dy;
}
// bounding box for a set of points
struct bbox
{
    ntype x0, x1, y0, y1;
    bbox(): x0(sentry), x1(-sentry), y0(sentry), y1(-sentry) {}
    // computes bounding box from a bunch of points
     void compute(const vector<point> &v) {
          for (int i = 0; i < v.size(); ++i) {
               x0 = min(x0, v[i].x); x1 = max(x1, v[i].x);
               y0 = min(y0, v[i].y); y1 = max(y1, v[i].y);
          }
     }
    // squared distance between a point and this bbox, 0 if inside
     ntype distance(const point &p) {
          if (p.x < x0) {
               if (p.y < y0)
                                    return pdist2(point(x0, y0), p);
               else if (p.y > y1) return pdist2(point(x0, y1), p);
               else
                                       return pdist2(point(x0, p.y), p);
          }
```

```
else if (p.x > x1) {
               if (p.y < y0)
                                     return pdist2(point(x1, y0), p);
               else if (p.y > y1) return pdist2(point(x1, y1), p);
                                        return pdist2(point(x1, p.y), p);
               else
          }
          else {
               if (p.y < y0)
                                     return pdist2(point(p.x, y0), p);
               else if (p.y > y1) return pdist2(point(p.x, y1), p);
               else
                                        return 0:
          }
     }
};
// stores a single node of the kd-tree, either internal or leaf
struct kdnode
     bool leaf;
                      // true if this is a leaf node (has one point)
     point pt;
                      // the single point of this is a leaf
     bbox bound;
                        // bounding box for set of points in children
     kdnode *first, *second; // two children of this kd-node
     kdnode(): leaf(false), first(0), second(0) {}
     ~kdnode() { if (first) delete first; if (second) delete second; }
     // intersect a point with this node (returns squared distance)
     ntype intersect(const point &p) {
          return bound.distance(p);
     }
    // recursively builds a kd-tree from a given cloud of points
     void construct(vector<point> &vp)
     {
          // compute bounding box for points at this node
          bound.compute(vp);
          // if we're down to one point, then we're a leaf node
          if (vp.size() == 1) {
               leaf = true;
               pt = vp[0];
          }
          else {
               // split on x if the bbox is wider than high (not best heuristic...)
               if (bound.x1-bound.x0 >= bound.y1-bound.y0)
```

```
sort(vp.begin(), vp.end(), on_x);
               // otherwise split on y-coordinate
               else
                    sort(vp.begin(), vp.end(), on_y);
               // divide by taking half the array for each child
               // (not best performance if many duplicates in the middle)
               int half = vp.size()/2;
               vector<point> vl(vp.begin(), vp.begin()+half);
               vector<point> vr(vp.begin()+half, vp.end());
               first = new kdnode();
                                       first->construct(vl);
               second = new kdnode(); second->construct(vr);
          }
     }
};
// simple kd-tree class to hold the tree and handle queries
struct kdtree
     kdnode *root;
    // constructs a kd-tree from a points (copied here, as it sorts them)
     kdtree(const vector<point> &vp) {
          vector<point> v(vp.begin(), vp.end());
          root = new kdnode();
          root->construct(v);
     }
     ~kdtree() { delete root; }
     // recursive search method returns squared distance to nearest point
     ntype search(kdnode *node, const point &p)
     {
          if (node->leaf) {
               // commented special case tells a point not to find itself
//
                if (p == node->pt) return sentry;
                else
                    return pdist2(p, node->pt);
          }
          ntype bfirst = node->first->intersect(p);
          ntype bsecond = node->second->intersect(p);
          // choose the side with the closest bounding box to search first
          // (note that the other side is also searched if needed)
```

```
if (bfirst < bsecond) {
               ntype best = search(node->first, p);
               if (bsecond < best)
                    best = min(best, search(node->second, p));
               return best;
          }
          else {
               ntype best = search(node->second, p);
               if (bfirst < best)
                    best = min(best, search(node->first, p));
               return best;
          }
     }
     // squared distance to the nearest
     ntype nearest(const point &p) {
          return search(root, p);
     }
};
// some basic test code here
int main()
{
    // generate some random points for a kd-tree
     vector<point> vp;
     for (int i = 0; i < 100000; ++i) {
          vp.push_back(point(rand()%100000, rand()%100000));
     kdtree tree(vp);
    // query some points
     for (int i = 0; i < 10; ++i) {
          point q(rand()% 100000, rand()% 100000);
          cout << "Closest squared distance to (" << q.x << ", " << q.y << ")"
                << " is " << tree.nearest(q) << endl;
     }
     return 0;
}
```

.....

```
LongestIncreasingSubsequence.cc 23/27
// Given a list of numbers of length n, this routine extracts a
// longest increasing subsequence.
// Running time: O(n log n)
//
//
     INPUT: a vector of integers
    OUTPUT: a vector containing the longest increasing subsequence
#include <iostream>
#include <vector>
#include <algorithm>
using namespace std;
typedef vector<int> VI;
typedef pair<int,int> PII;
typedef vector<PII> VPII;
#define STRICTLY_INCREASNG
VI LongestIncreasingSubsequence(VI v) {
  VPII best;
  VI dad(v.size(), -1);
  for (int i = 0; i < v.size(); i++) {
#ifdef STRICTLY INCREASIG
    PII item = make_pair(v[i], 0);
     VPII::iterator it = lower_bound(best.begin(), best.end(), item);
     item.second = i;
#else
     PII item = make_pair(v[i], i);
     VPII::iterator it = upper_bound(best.begin(), best.end(), item);
#endif
    if (it == best.end()) {
       dad[i] = (best.size() == 0 ? -1 : best.back().second);
       best.push_back(item);
     } else {
       dad[i] = dad[it->second];
       *it = item;
     }
```

```
}
  VI ret;
  for (int i = best.back().second; i >= 0; i = dad[i])
     ret.push_back(v[i]);
  reverse(ret.begin(), ret.end());
  return ret;
}
Dates.cc 24/27
// Routines for performing computations on dates. In these routines,
// months are expressed as integers from 1 to 12, days are expressed
// as integers from 1 to 31, and years are expressed as 4-digit
// integers.
#include <iostream>
#include <string>
using namespace std;
string dayOfWeek[] = {"Mon", "Tue", "Wed", "Thu", "Fri", "Sat", "Sun"};
// converts Gregorian date to integer (Julian day number)
int dateToInt (int m, int d, int y){
  return
     1461*(y+4800+(m-14)/12)/4+
     367*(m-2-(m-14)/12*12)/12-
    3*((y+4900+(m-14)/12)/100)/4+
    d - 32075;
}
// converts integer (Julian day number) to Gregorian date: month/day/year
void intToDate (int jd, int &m, int &d, int &y){
  int x, n, i, j;
  x = id + 68569;
  n = 4 * x / 146097;
  x = (146097 * n + 3) / 4;
  i = (4000 * (x + 1)) / 1461001;
  x = 1461 * i / 4 - 31;
  j = 80 * x / 2447;
```

```
d = x - 2447 * j / 80;
  x = j / 11;
  m = j + 2 - 12 * x;
  y = 100 * (n - 49) + i + x;
}
// converts integer (Julian day number) to day of week
string intToDay (int jd){
  return dayOfWeek[jd % 7];
}
int main (int argc, char **argv){
  int jd = dateToInt (3, 24, 2004);
  int m, d, y;
  intToDate (jd, m, d, y);
  string day = intToDay (jd);
  // expected output:
         2453089
  //
  //
         3/24/2004
         Wed
  cout << jd << endl
     << m << "/" << d << "/" << y << endl
     << day << endl;
}
LogLan.java 25/27
// Code which demonstrates the use of Java's regular expression libraries.
// This is a solution for
//
//
     Loglan: a logical language
//
     http://acm.uva.es/p/v1/134.html
// In this problem, we are given a regular language, whose rules can be
// inferred directly from the code. For each sentence in the input, we must
// determine whether the sentence matches the regular expression or not. The
// code consists of (1) building the regular expression (which is fairly
// complex) and (2) using the regex to match sentences.
import java.util.*;
import java.util.regex.*;
```

```
public class LogLan {
     public static String BuildRegex (){
     String space = " +";
     String A = "([aeiou])";
    String C = "([a-z\&\&[^aeiou]])";
    String MOD = "(g" + A + ")";
     String BA = "(b" + A + ")";
     String DA = "(d" + A + ")";
     String LA = "(1" + A + ")";
    String NAM = "([a-z]^*" + C + ")";
     String PREDA = "(" + C + C + A + C + A + "|" + C + A + C + C + A + ")";
     String predstring = "(" + PREDA + "(" + space + PREDA + ")*)";
     String predname = "(" + LA + space + predstring + "|" + NAM + ")";
     String preds = "(" + predstring + "(" + space + A + space + predstring + ")*)";
     String predclaim = "(" + predname + space + BA + space + preds + "|" + DA + space +
               preds + ")";
     String verbpred = "(" + MOD + space + predstring + ")";
     String statement = "(" + predname + space + verbpred + space + predname + "|" +
               predname + space + verbpred + ")";
     String sentence = "(" + statement + "|" + predclaim + ")";
     return "^" + sentence + "$";
     }
     public static void main (String args[]){
     String regex = BuildRegex();
     Pattern pattern = Pattern.compile (regex);
     Scanner s = new Scanner(System.in);
     while (true) {
               // In this problem, each sentence consists of multiple lines, where the last
          // line is terminated by a period. The code below reads lines until
          // encountering a line whose final character is a '.'. Note the use of
              //
              //
                     s.length() to get length of string
              //
                     s.charAt() to extract characters from a Java string
              //
                     s.trim() to remove whitespace from the beginning and end of Java string
               //
```

```
//
               //
                     s.compareTo(t) < 0 if s < t, lexicographically
               //
                     s.indexOf("apple") returns index of first occurrence of "apple" in s
               //
                     s.lastIndexOf("apple") returns index of last occurrence of "apple" in s
                     s.replace(c,d) replaces occurrences of character c with d
               //
                     s.startsWith("apple) returns (s.indexOf("apple") == 0)
               //
                     s.toLowerCase() / s.toUpperCase() returns a new lower/uppercased string
               //
               //
                     Integer.parseInt(s) converts s to an integer (32-bit)
               //
                     Long.parseLong(s) converts s to a long (64-bit)
               //
                     Double.parseDouble(s) converts s to a double
          String sentence = "";
          while (true){
          sentence = (sentence + " " + s.nextLine()).trim();
          if (sentence.equals("#")) return;
          if (sentence.charAt(sentence.length()-1) == '.') break;
          }
               // now, we remove the period, and match the regular expression
               String removed_period = sentence.substring(0, sentence.length()-1).trim();
          if (pattern.matcher (removed_period).find()){
          System.out.println ("Good");
          } else {
          System.out.println ("Bad!");
}
Primes.cc 26/27
// O(sqrt(x)) Exhaustive Primality Test
#include <cmath>
#define EPS 1e-7
typedef long long LL;
bool IsPrimeSlow (LL x)
  if(x \le 1) return false;
  if(x \le 3) return true;
```

// Other useful String manipulation methods include

```
if (!(x\%2) || !(x\%3)) return false;
  LL s=(LL)(sqrt((double)(x))+EPS);
  for(LL i=5;i \le s;i+=6)
  {
    if (!(x\%i) || !(x\%(i+2))) return false;
  }
  return true;
}
// Primes less than 1000:
        2
               3
                             7
                                                       19
                                                                     29
                                                                            31
                                                                                   37
//
                      5
                                  11
                                         13
                                                17
                                                              23
//
      41
              43
                     47
                            53
                                  59
                                         61
                                                67
                                                       71
                                                              73
                                                                     79
                                                                            83
                                                                                   89
//
      97
             101
                    103
                           107
                                 109
                                        113
                                               127
                                                      131
                                                             137
                                                                    139
                                                                           149
                                                                                  151
//
      157
             163
                   167
                          173
                                 179
                                        181
                                               191
                                                      193
                                                             197
                                                                    199
                                                                           211
                                                                                  223
//
     227
            229
                   233
                          239
                                 241
                                        251
                                               257
                                                      263
                                                             269
                                                                    271
                                                                           277
                                                                                  281
//
     283
            293
                   307
                          311
                                        317
                                               331
                                                      337
                                                             347
                                                                    349
                                                                           353
                                                                                  359
                                 313
//
                   379
                                        397
                                                                                  433
     367
            373
                          383
                                 389
                                               401
                                                      409
                                                             419
                                                                    421
                                                                           431
//
     439
            443
                   449
                          457
                                        463
                                               467
                                                      479
                                                             487
                                                                    491
                                                                           499
                                                                                  503
                                 461
//
     509
            521
                   523
                          541
                                 547
                                        557
                                               563
                                                      569
                                                             571
                                                                    577
                                                                           587
                                                                                  593
     599
//
            601
                   607
                          613
                                 617
                                        619
                                               631
                                                             643
                                                                    647
                                                                           653
                                                                                  659
                                                      641
//
                                        701
                                               709
                                                             727
                                                                    733
                                                                           739
                                                                                  743
     661
            673
                   677
                          683
                                 691
                                                      719
//
     751
                                        787
                                               797
                                                      809
            757
                   761
                          769
                                 773
                                                             811
                                                                    821
                                                                           823
                                                                                  827
//
                                                                           907
                                                                                  911
     829
            839
                   853
                          857
                                 859
                                        863
                                               877
                                                      881
                                                             883
                                                                    887
                                                                                  997
//
     919
            929
                   937
                          941
                                 947
                                        953
                                               967
                                                      971
                                                             977
                                                                    983
                                                                           991
// Other primes:
//
     The largest prime smaller than 10 is 7.
//
     The largest prime smaller than 100 is 97.
//
     The largest prime smaller than 1000 is 997.
//
     The largest prime smaller than 10000 is 9973.
//
     The largest prime smaller than 100000 is 99991.
//
     The largest prime smaller than 1000000 is 999983.
     The largest prime smaller than 10000000 is 9999991.
//
//
     The largest prime smaller than 100000000 is 99999989.
     The largest prime smaller than 1000000000 is 999999937.
//
//
     The largest prime smaller than 10000000000 is 9999999967.
//
     The largest prime smaller than 10000000000 is 9999999977.
     The largest prime smaller than 100000000000 is 99999999999.
//
     The largest prime smaller than 1000000000000 is 999999999971.
//
     The largest prime smaller than 10000000000000 is 9999999999973.
//
     The largest prime smaller than 10000000000000 is 9999999999999989.
//
//
     The largest prime smaller than 100000000000000 is 99999999999937.
//
     The largest prime smaller than 1000000000000000 is 999999999999997.
```

//

```
KMP.cpp 27/27
/*
Searches for the string w in the string s (of length k). Returns the
0-based index of the first match (k if no match is found). Algorithm
runs in O(k) time.
*/
#include <iostream>
#include <string>
#include <vector>
using namespace std;
typedef vector<int> VI;
void buildTable(string& w, VI& t)
  t = VI(w.length());
  int i = 2, j = 0;
  t[0] = -1; t[1] = 0;
  while(i < w.length())</pre>
     if(w[i-1] == w[j]) \{ t[i] = j+1; i++; j++; \}
     else if(j > 0) j = t[j];
     else { t[i] = 0; i++; }
  }
}
int KMP(string& s, string& w)
  int m = 0, i = 0;
  VI t;
  buildTable(w, t);
  while(m+i < s.length())
     if(w[i] == s[m+i])
     {
       i++;
       if(i == w.length()) return m;
```

```
}
     else
     {
       m += i-t[i];
       if(i > 0) i = t[i];
     }
  return s.length();
}
int main()
{
  string a = (string) "The example above illustrates the general technique for assembling "+
     "the table with a minimum of fuss. The principle is that of the overall search: "+
     "most of the work was already done in getting to the current position, so very "+
     "little needs to be done in leaving it. The only minor complication is that the "+
     "logic which is correct late in the string erroneously gives non-proper "+
     "substrings at the beginning. This necessitates some initialization code.";
  string b = "table";
  int p = KMP(a, b);
  cout << p << ":" << a.substr(p, b.length()) << "" << b << endl;\\
}
```

Generated by GNU enscript 1.6.1.