# **Combinatorial optimization**

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## Max flow (C++)

```
// Fattest path network flow algorithm using an adjacency matrix. //  
// Running time: O(|E|^2 \log (|V| * U)), where U is the largest capacity of any edge. If you replace the 'fattest path' search with a minimum number of edges search, the running time becomes O(|E|^2 |V|).
```

```
// INPUT: cap -- a matrix such that cap[i][i] is the capacity of
                 a directed edge from node i to node j
//
//
                 * Note that it is legitimate to create an i->j
                   edge without a corresponding j->i edge.
//
//
//
                 * Note that for an undirected edge, set
//
                   both cap[i][j] and cap[j][i] to the capacity of
//
                   the undirected edge.
//
//
          source -- starting node
//
          sink -- ending node
//
// OUTPUT: value of maximum flow; also, the flow[][] matrix will
           contain both positive and negative integers -- if you
//
//
           want the actual flow assignments, look at the
//
           *positive* flow values only.
//
// To use this, create a MaxFlow object, and call it like this:
//
//
    MaxFlow nf:
// int maxflow = nf.getMaxFlow(cap, source, sink);
typedef vector<int> VI;
typedef vector<VI> VVI;
const int INF = 1000000000;
struct MaxFlow {
  int N:
  VVI cap, flow;
  VI found, dad, dist:
  bool searchFattest(int source, int sink){
    fill (found.begin(), found.end(), false);
    fill (dist.begin(), dist.end(), 0);
    dist[source] = INF;
    while (source != N){
      int best = N;
      found[source] = true;
      if (source == sink) break;
      for (int k = 0; k < N; k++) {
        if (found[k]) continue;
        int possible = min(cap[source][k] - flow[source][k], dist[source]);
        if (dist[k] < possible) {</pre>
          dist[k] = possible;
          dad[k] = source;
        if (dist[k] > dist[best]) best = k;
      source = best;
```

```
return found[sink];
 }
  bool searchShortest(int source, int sink){
   fill (found.begin(), found.end(), false);
   fill (dist.begin(), dist.end(), INF);
   dist[source] = 0;
   while (source != N){
      int best = N;
      found[source] = true;
      if (source == sink) break;
      for (int k = 0; k < N; k++){
        if (found[k]) continue;
        if (cap[source][k] - flow[source][k] > 0){
          if (dist[k] > dist[source] + 1){
            dist[k] = dist[source] + 1;
            dad[k] = source;
          }
        if (dist[k] < dist[best]) best = k;</pre>
      }
      source = best;
    return found[sink];
  int getMaxFlow (const VVI &cap, int source, int sink){
   this->cap = cap;
   N = cap.size();
   found = VI(N);
   flow = VVI(N,VI(N)):
   dist = VI(N+1);
   dad = VI(N);
   int totflow = 0:
   while (searchFattest(source, sink)){
      int amt = INF;
      for (int x = sink; x != source; x = dad[x])
        amt = min (amt, cap[dad[x]][x] - flow[dad[x]][x]);
      for (int x = sink; x != source; x = dad[x]){
        flow[dad[x]][x] += amt;
        flow[x][dad[x]] -= amt;
      totflow += amt;
    return totflow;
};
```

#### Min cost max flow (C++)

```
// Min cost max flow algorithm using an adjacency matrix. If you
// want just regular max flow, setting all edge costs to 1 gives
// running time O(|E|^2 |V|).
// Running time: O(\min(|V|^2 * totflow, |V|^3 * totcost))
// INPUT: cap -- a matrix such that cap[i][j] is the capacity of
//
                 a directed edge from node i to node j
//
//
          cost -- a matrix such that cost[i][j] is the (positive)
//
                  cost of sending one unit of flow along a
//
                  directed edge from node i to node j
//
//
          source -- starting node
//
          sink -- ending node
//
// OUTPUT: max flow and min cost; the matrix flow will contain
           the actual flow values (note that unlike in the MaxFlow
//
//
           code, you don't need to ignore negative flow values -- there
//
           shouldn't be any)
//
// To use this, create a MinCostMaxFlow object, and call it like this:
// MinCostMaxFlow nf;
// int maxflow = nf.getMaxFlow(cap,cost,source,sink);
typedef vector<int> VI;
typedef vector<VI> VVI;
const int INF = 1000000000:
struct MinCostMaxFlow {
  int N;
  VVI cap, flow, cost;
  VI found, dad, dist, pi;
  bool search(int source, int sink) {
    fill(found.begin(), found.end(), false);
    fill(dist.begin(), dist.end(), INF);
    dist[source] = 0;
    while (source != N) {
     int best = N;
      found[source] = true;
      for (int k = 0; k < N; k++) {
       if (found[k]) continue;
        if (flow[k][source]) {
          int val = dist[source] + pi[source] - pi[k] - cost[k][source];
          if (dist[k] > val) {
```

```
dist[k] = val;
          dad[k] = source;
      if (flow[source][k] < cap[source][k]) {</pre>
        int val = dist[source] + pi[source] - pi[k] + cost[source][k];
        if (dist[k] > val) {
          dist[k] = val;
          dad[k] = source;
        }
      if (dist[k] < dist[best]) best = k;</pre>
   source = best;
  for (int k = 0; k < N; k++)
    pi[k] = min(pi[k] + dist[k], INF);
 return found[sink];
pair<int,int> getMaxFlow(const VVI &cap, const VVI &cost, int source, int si
 this->cap = cap;
 this->cost = cost;
 N = cap.size();
 found = VI(N);
 flow = VVI(N, VI(N));
 dist = VI(N+1);
 dad = VI(N);
 pi = VI(N);
 int totflow = 0, totcost = 0;
 while (search(source, sink)) {
    int amt = INF:
    for (int x = sink; x != source; x = dad[x])
      amt = min(amt, flow[x][dad[x]] ? flow[x][dad[x]] :
                cap[dad[x]][x] - flow[dad[x]][x]);
    for (int x = sink; x != source; x = dad[x]) {
      if (flow[x][dad[x]]) {
        flow[x][dad[x]] -= amt;
        totcost -= amt * cost[x][dad[x]];
     } else {
        flow[dad[x]][x] += amt;
        totcost += amt * cost[dad[x]][x];
    totflow += amt;
  return make_pair(totflow, totcost);
```

```
};
(Min cost) maximum matching (C++)
// This code performs maximum bipartite matching and, optionally min-cost match
// It has a heuristic that will give excellent performance on complete graphs
// where rows <= columns.
//
//
     INPUT: w[i][i] = cost from row node i and column node i or NO EDGE
//
     OUTPUT: mr[i] = assignment for row node i or -1 if unassigned
             mc[i] = assignment for column node i or -1 if unassigned
//
//
     BipartiteMatching and MinCostMatching return the number of matches made.
    MatchingCost will give you the cost, if you need it.
// Contributed by Andy Lutomirski.
typedef vector<int> VI;
typedef vector<VI> VVI;
const int NO_EDGE = -(1<<30); // Or any other value.</pre>
bool FindMatch(int i, const VVI &w, VI &mr, VI &mc, VI &seen)
  if (seen[i])
    return false;
  seen[i] = true;
  for (int j = 0; j < w[i].size(); j++) {</pre>
   if (w[i][j] != NO EDGE && mc[j] < 0) {
      mr[i] = i:
      mc[j] = i;
      return true;
   }
  for (int j = 0; j < w[i].size(); j++) {</pre>
   if (w[i][j] != NO EDGE && mr[i] != j) {
      if (mc[j] < 0 \mid | 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)
```

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mr = VI (w.size(), -1);

```
mc = VI(w[0].size(), -1);
 VI seen(w.size());
 int ct = 0:
  for(int i = 0; i < w.size(); i++)</pre>
      fill(seen.begin(), seen.end(), 0);
      if (FindMatch(i, w, mr, mc, seen)) ct++;
 return ct;
// ----- CUT HERE FOR JUST MAXIMUM MATCHING ------
bool Augment(int start, const VVI &w, VI &mr, VI &mc)
  const int INF = (1<<31) - 1;</pre>
 VI cost(w.size(), INF);
 VI dad(w.size());
 cost[start] = 0;
  int last = 0:
  for(int i = 0; i < w.size() && last != -1 && cost[start] == 0; i++) // Beli</pre>
      last = -1:
      for(int r = 0; r < w.size(); r++)</pre>
          if (cost[r] == INF)
            continue;
          for(int r2 = 0; r2 < w.size(); r2++)
              if (r == r2 || mr[r2] == -1 || w[r][mr[r2]] == NO EDGE) continue
              int val = cost[r] + w[r][mr[r2]] - w[r2][mr[r2]];
              if (val < cost[r2] && val < 0) {</pre>
                cost[r2] = val;
                last = r2;
                dad[r2] = r;
            }
  if (mr[start] == -1)
      int best = -1;
      for(int i = 0; i < w.size(); i++)</pre>
          if (cost[i] < 0 \&\& (best == -1 \mid | cost[i] < cost[best]))
            best = i;
        }
      if (best == -1)
```

```
return false;
      // Augment a non-cycle
      int a = dad[best], b = best;
      VI oldmr = mr;
      mr[best] = -1;
      do {
        mr[a] = oldmr[b];
        mc[mr[a]] = a;
        b = a;
        a = dad[a];
     } while(b != start);
      return true:
  if (last == -1)
    return false:
  if (cost[start] == 0)
    last = start;
  for(int i = 0; i < w.size(); i++)</pre>
    last = dad[last]; // Find a cycle
  // Augment.
  VI oldmr = mr;
  int a = last, b;
  do {
    b = a;
    a = dad[a];
    mr[a] = oldmr[b];
    mc[oldmr[b]] = a;
  } while(a != last);
  for(int i = 0; i < w.size(); i++)</pre>
   if (mr[i] != -1)
      assert(mc[mr[i]] == i);
  return true;
// Returns the size of the matching.
int MinCostMatching(const VVI &w, VI &mr, VI &mc)
  int size = BipartiteMatching(w, mr, mc);
  int blanks = 0;
  for(int start = 0; blanks < w.size(); start = (start+1)%w.size())</pre>
      blanks++:
      if (Augment(start, w, mr, mc))
        blanks = 0;
  return size;
```

```
}
int MatchingCost(const VVI &w, const VI &mr)
  int c = 0;
  for(int i = 0; i < w.size(); i++)</pre>
   if (mr[i] != -1)
      c += w[i][mr[i]];
  return c;
Non-bipartite matching (C++)
/* An implementation of Nick Harvey's algorithm for nonbipartiate max matching
   in undirected, unweighted graphs. Note the algorithm is randomized, so to L
   safe you should run matching() a few times and take the largest one. If yo
   find a bug in this code, e-mail Jelani Nelson (minilek@mit.edu). */
#define FPS 1e-9
// MAXN is the maximum number of vertices
#define MAXN 64
// adj[i][i] = adj[j][i] = 1 iff the edge (i,j) exists (else both are 0)
char adj[MAXN][MAXN];
// number of vertices in graph
int V;
// counts number of bits of x set to 1
int pc(int x) { return !x?0:(x&1)+pc(x>>1); }
typedef struct matrix {
 vector< vector<long double> > a;
 int n. m:
  matrix(int x, int y) {
   n = x, m = y;
   a = vector< vector<long double> >();
   for (int i = 0; i < n; ++i)
      a.push back(vector<long double>(m, 0));
 }
  matrix() {
   n = m = 0;
   a = vector< vector<long double> >(0);
  matrix(const matrix &x) {
   n = x.n, m = x.m, a = x.a;
  void operator=(const matrix& x) {
   n = x.n, m = x.m, a = x.a;
```

```
matrix operator*(const matrix &b) {
    matrix c(n, b.m);
    for (int i = 0; i < c.n; ++i)
      for (int j = 0; j < c.m; ++j)
        for (int k = 0; k < m; ++k)
          c.a[i][j] += a[i][k] * b.a[k][j];
    return c;
  matrix operator+(const matrix &b) {
    matrix c = b;
    for (int i = 0; i < c.n; ++i)
      for (int j = 0; j < c.m; ++j)
        c.a[i][j] += a[i][j];
    return c:
  matrix operator-(const matrix &b) {
    matrix c = b;
    for (int i = 0; i < c.n; ++i)
      for (int j = 0; j < c.m; ++j)
        c.a[i][j] = a[i][j] - c.a[i][j];
    return c;
  matrix operator-() {
    matrix c(n, m);
    for (int i = 0; i < c.n; ++i)
      for (int j = 0; j < c.m; ++j)
        c.a[i][j] = -a[i][j];
    return c;
  long double& operator()(unsigned i, unsigned j) {
    return a[i][j];
  matrix operator()(vector<int> x, vector<int> y) {
    matrix c(x.size(), y.size());
    for (int i = 0; i < c.n; ++i)
      for (int j = 0; j < c.m; ++j)
        c(i, j) = a[x[i]][y[j]];
    return c;
} matrix;
// utility function to print a matrix
void printMatrix(matrix A) {
  for (int i = 0; i < A.n; ++i) {
    for (int j = 0; j < A.m; ++j)
      cout << A(i, j) << " ";
    cout << endl;</pre>
  cout << endl;</pre>
```

```
// returns an LxL identity matrix
matrix identity(int L) {
  matrix I(L, L);
  for (int i = 0; i < L; ++i)
   I(i, i) = 1;
  return I;
}
// returns a maximum size full rank square submatrix of A
// the vector<int> returned is the list of indices of rows used
pair< vector<int>, matrix> max rank submatrix(matrix A) {
  vector<int> indices:
 matrix B = A:
  int at = 0:
  vector<int> perm = vector<int>(A.n);
  for (int i = 0; i < A.n; ++i)
    perm[i] = i;
  for (int i = 0: (at < A.n) && (i < A.m): ++i) {
   int maxrow = at;
    for (int j = at + 1; j < A.n; ++j)
     if (fabs(A(j, i)) > fabs(A(maxrow, i)))
        maxrow = j;
   if (maxrow != at) {
      for (int j = 0; j < A.n; ++j)
        swap(A(at, j), A(maxrow, j));
      swap(perm[at], perm[maxrow]);
    if (fabs(A(at, i)) < EPS)
     continue:
   indices.push back(perm[at]);
   long double c = A(at. i):
   for (int j = i; j < A.m; ++j)
      A(at, j) /= c;
    for (int j = at + 1; j < A.n; ++j) {
     long double c = A(i, i):
     for (int k = i; k < A.m; ++k)
        A(i, k) -= A(at, k) * c;
   }
   ++at;
  sort(indices.begin(), indices.end());
  return make pair(indices, B(indices, indices));
// assumes matrix is non-singular
matrix matrix inverse(matrix A) {
 // input should have A.n = A.m
  matrix B = matrix(A.n, 2 * A.n);
  for (int i = 0; i < A.n; ++i)
   for (int j = 0; j < A.n; ++j)
```

```
B(i, j) = A(i, j);
  for (int i = 0; i < A.n; ++i)
    B(i, i + A.n) = 1;
  for (int i = 0; i < A.n; ++i) {
    int maxrow = i;
    for (int j = i + 1; j < A.n; ++j)
      if (fabs(B(j, i)) > fabs(B(maxrow, i)))
        maxrow = j;
    if (maxrow != i)
      for (int j = 0; j < B.m; ++j)
        swap(B(i, j), B(maxrow, j));
    long double c = B(i, i);
    for (int j = i; j < B.m; ++j)
      B(i, j) /= c;
    for (int j = 0; j < A.n; ++j) if (i != j) {
     long double c = B(j, i);
      for (int k = i; k < B.m; ++k)
        B(j, k) -= B(i, k) * c;
  matrix ret = matrix(A.n, A.n);
  for (int i = 0; i < A.n; ++i)
    for (int j = 0; j < A.n; ++j)
      ret(i, j) = B(i, j + A.n);
  return ret;
}
struct matrix T, N;
void delete edges(vector<int> J) {
  if (J.size() == 2) {
    int i = J[0], i = J[1]:
    if (fabs(T(i, j))>EPS \&\& fabs(N(i,j) + 1./T(i, j))>EPS)
      T(i, j) = T(j, i) = 0;
  } else {
    int C = 4:
    for (int i = 1; i <= C; ++i)</pre>
      for (int j = i + 1; j \le C; ++j) {
        vector<int> Jp;
        for (int k = (i-1)*J.size()/C; k < i*J.size()/C; ++k)
          Jp.push back(J[k]);
        for (int k = (j-1)*J.size()/C; k < j*J.size()/C; ++k)
          Jp.push back(J[k]);
        matrix W = T(Jp, Jp), WHat = N(Jp, Jp);
        delete_edges(Jp);
        matrix w = T(Jp, Jp);
        for (int k = 0; k < Jp.size(); ++k)
          for (int l = 0; l < Jp.size(); ++l)</pre>
            N(Jp[k], Jp[l]) = WHat(k, l);
        matrix X = N(J, J) - N(J, Jp) *
          matrix_inverse(identity(Jp.size()) + (w-W)*WHat)*(w-W)*N(Jp, J);
```

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```
for (int k = 0; k < J.size(); ++k)
          for (int l = 0; l < J.size(); ++l)</pre>
            N(J[k], J[l]) = X(k, l);
// returns a vector<int> v
//v[i] = -1 if i isn't matched, else v[i] is the vertex i is matched to
#define MAXRAND 10000
vector<int> matching() {
 T = matrix(V, V);
 for (int i = 0; i < V; ++i)
   for (int j = i + 1; j < V; ++j)
      if (adj[i][j])
        T(i, j) = (rand() % MAXRAND) + 1;
  for (int i = 0; i < V; ++i)
    for (int j = i; j < V; ++j)
      T(j, i) = -T(i, j);
  pair< vector<int>, matrix > x = max rank submatrix(T);
  vector<int> indices = x.first;
  if (indices.size() == 0)
   return vector<int>(V, -1);
 // make the number of vertices in T a power of 2 while keeping
 // full rank (put the new vertices in a clique)
 // Nick's algorithm assumes #vertices is a power of 2
 T = x.second;
 int newLength = T.n;
  while (pc(newLength) > 1)
   ++newLenath:
  matrix newT = matrix(newLength, newLength);
  for (int i = 0; i < T.n: ++i)
   for (int j = 0; j < T.n; ++j)
      newT(i, j) = T(i, j);
  for (int i = T.n; i < newT.n; ++i)</pre>
    for (int j = i + 1; j < newT.n; ++j) {</pre>
      newT(i, j) = (rand() % MAXRAND) + 1;
      newT(j, i) = -newT(i, j);
 T = newT;
 N = matrix_inverse(T);
  vector<int> vertices;
  for (int i = 0; i < T.n; ++i)
   vertices.push back(i);
  delete edges(vertices); // deletes edges until left with a perf. matching
  vector<int> v = vector<int>(V, -1);
  for (int i = 0; i < indices.size(); ++i)</pre>
   for (int j = 0; j < indices.size(); ++j)</pre>
```

```
if (fabs(T(i, j)) > EPS)
        v[indices[i]] = indices[j];
  // make sure this is a valid matching
  for (int i = 0; i < v.size(); ++i)</pre>
   if (v[i] != -1) {
     if (v[v[i]] != i)
        cout << "failed at " << i << endl;</pre>
      assert(v[v[i]] == i);
  return v;
int main() {
  timeval tp:
  gettimeofday(&tp, NULL);
  srand(tp.tv_usec);
  /* EXAMPLE USAGE */
  // set the adj array and #vertices here
  V = 64;
  memset(adj, 0, sizeof(adj));
  for (int i = 0; i < V; i ++)
    for (int j = i + 1; j < V; ++j)
      if (rand()%50 == 0) // making it somewhat sparse so there's no perf. mat
        adi[i][i] = adi[i][i] = 1;
  // find the max matching
  vector<int> matches = matching();
  // make sure max matching only used real edges!
  for (int i = 0; i < V; ++i)
   if (matches[i] != -1)
      assert(adj[i][matches[i]]);
  for (int i = 0; i < matches.size(); ++i)</pre>
    cout << i << ": " << matches[i] << endl;</pre>
  /* END OF EXAMPLE */
  return 0;
Convex hull (C++)
```

```
// 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
using namespace std;
#define REMOVE REDUNDANT
typedef double T;
typedef pair<T,T> PT;
typedef vector<PT> VPT;
const double EPS = 1e-7;
T det (const PT &a, const PT &b){
  return a.first * b.second - a.second * b.first;
T area2 (const PT &a, const PT &b, const PT &c){
  return det(a,b) + det(b,c) + det(c,a);
#ifdef REMOVE REDUNDANT
// return true if point b is between points a and c
bool between (const PT &a, const PT &b, const PT &c){
  return (fabs(area2(a,b,c)) < EPS &&</pre>
          (a.first - b.first) * (c.first - b.first) <= 0 &&
          (a.second - b.second) * (c.second - b.second) <= 0);</pre>
#endif
void convex hull (VPT &pts){
  sort (pts.begin(), pts.end());
  pts.erase (unique (pts.begin(), pts.end()), pts.end());
 VPT up, dn;
  for (int i = 0; i < pts.size(); i++){</pre>
    while (up.size() > 1 \& area2(up[up.size()-2], up.back(), pts[i]) >= 0)
    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 >= 1; i--) \text{ pts.push back}(up[i]);
```

```
#ifdef REMOVE REDUNDANT
  if (pts.size() <= 2) return;</pre>
  dn.clear();
  dn.push back (pts[0]);
  dn.push back (pts[1]);
  for (int i = 2; i < pts.size(); i++){</pre>
   if (between (dn[dn.size()-2], dn[dn.size()-1], pts[i])) dn.pop back();
    dn.push back (pts[i]);
  if (dn.size() >= 3 \&\& between (dn.back(), dn[0], dn[1])){
    dn[0] = dn.back():
    dn.pop back();
  pts = dn;
#endif
Area and centroid (C++)
// This code computes the area or centroid of a polygon,
// assuming that the coordinates are listed in a clockwise
// or counterclockwise fashion.
// Running time: O(n)
// INPUT: list of x[] and y[] coordinates
// OUTPUTS: (signed) area or centroid
// Note that the centroid is often known as the
// "center of gravity" or "center of mass".
typedef vector<double> VD:
typedef pair<double, double> PD;
double ComputeSignedArea (const VD &x, const VD &y){
  double area = 0:
  for (int i = 0; i < x.size(); i++){</pre>
   int j = (i+1) % x.size();
    area += x[i]*y[j] - x[j]*y[i];
  return area / 2.0;
double ComputeArea (const VD &x, const VD &y){
  return fabs (ComputeSignedArea (x, y));
```

```
PD ComputeCentroid (const VD &x, const VD &y){
    double cx = 0, cy = 0;
    double scale = 6.0 * ComputeSignedArea (x, y);
    for (int i = 0; i < x.size(); i++){
        int j = (i+1) % x.size();
        cx += (x[i]+x[j])*(x[i]*y[j]-x[j]*y[i]);
        cy += (y[i]+y[j])*(x[i]*y[j]-x[j]*y[i]);
    }
    return make_pair (cx/scale, cy/scale);
}</pre>
```

## Misc geometry (C++)

```
// C++ routines for computational geometry.
double INF = 1e100:
double EPS = 1e-7:
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){ return PT(x-p.x,y-p.y); }
  PT operator+ (const PT &p){ return PT(x+p.x,y+p.y); }
  PT operator* (double c){ return PT(x*c,y*c); }
  PT operator/ (double c) { return PT(x/c,y/c); }
};
double dot (PT p, PT a){ return p,x*a,x+p,v*a,v: }
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;</pre>
  r = dot(c-a.b-a)/r:
  if (r < 0) return a;</pre>
  if (r > 1) return b;
  return a + (b-a)*r;
// 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 two lines are parallel or collinear
bool LinesParallel (PT a, PT b, PT c, PT d){
  return fabs(cross(b-a,c-d)) < EPS;</pre>
bool LinesCollinear (PT a, PT b, PT c, PT d){
  return LinesParallel(a,b,c,d) && fabs(cross(a-c,d-c)) < EPS;</pre>
}
// 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 (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
PT ComputeLineIntersection (PT a, PT b, PT c, PT d){
  b=b-a; d=c-d; c=c-a;
  if (dot(b,b) < EPS) return a;</pre>
  if (dot(d,d) < EPS) return c;</pre>
  return a + b*cross(c,d)/cross(b,d);
```

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```
// 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++){</pre>
   int j = (i+1)%p.size();
   if ((p[i].y <= 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++)</pre>
   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:
 PT d = b-a:
 double D = cross(a-c,b-c);
  double e = r*r*dot(d,d)-D*D;
  if (e < 0) return ret;</pre>
  e = sqrt(e);
```

```
ret.push back (c+PT(D*d.y+(d.y)=0?1:-1)*d.x*e,-D*d.x+fabs(d.y)*e)/dot(d,d));
  if (e > 0)
    ret.push back (c+PT(D*d.y-(d.y)=0?1:-1)*d.x*e,-D*d.x-fabs(d.y)*e)/dot(d,d)
  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 \mid\mid d+min(r,R) < max(r,R)) return ret;
  double x = (d*d-R*R+r*r)/(2*d);
  double v = sart(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;
}
```

file:///mit/ecprice/acm/acm08/notebook.html#file7

#### Voronoi diagrams (C++)

```
#include "Geometry.cc"
#define MAXN 1024
#define INF 1000000
//Voronoi diagrams: O(N^2*LogN)
//Convex hull: O(N*LoaN)
typedef struct {
  int id;
  double x;
  double v:
  double ang:
} chp;
int n;
double x[MAXN], y[MAXN]; // Input points
chp inv[2*MAXN]; // Points after inversion (to be given to Convex Hull)
int vors:
int vor[MAXN]; // Set of points in convex hull;
               //starts at lefmost; last same as first!!
PT ans[MAXN][2];
int chpcmp(const void *aa, const void *bb) {
  double a = ((chp *)aa)->ang;
  double b = ((chp *)bb)->ang;
  if (a<b) return -1;</pre>
```

```
else if (a>b) return 1;
  else return 0; // Might be better to include a
                 // tie-breaker on distance, instead of the cheap hack below
int orient(chp *a, chp *b, chp *c) {
 double s = a->x*(b->y-c->y) + b->x*(c->y-a->y) + c->x*(a->y-b->y);
 if (s>0) return 1;
  else if (s<0) return -1:
  else if (a->ang=b->ang && a->ang=ec->ang) return -1; // Cheap hack
           //for points with same angles
  else return 0;
}
//the pt argument must have the points with precomputed angles (atan2()'s)
//with respect to a point on the inside (e.g. the center of mass)
int convexHull(int n, chp *pt, int *ans) {
 int i, j, st, anses=0;
 qsort(pt, n, sizeof(chp), chpcmp);
  for (i=0; i<n; i++) pt[n+i] = pt[i];</pre>
  st = 0:
  for (i=1; i<n; i++) { // Pick leftmost (bottommost)</pre>
                        //point to make sure it's on the convex hull
    if (pt[i].x<pt[st].x || (pt[i].x==pt[st].x && pt[i].y<pt[st].y)) st = i;</pre>
  ans[anses++] = st;
  for (i=st+1; i<=st+n; i++) {</pre>
    for (j=anses-1; j; j--) {
     if (orient(pt+ans[j-1], pt+ans[j], pt+i)>=0) break;
      // Should change the above to strictly greater,
      // if you don't want points that lie on the side (not on a vertex) of the
      // If you really want them, you might also put an epsilon in orient
    ans[j+1] = i;
    anses = i+2:
  for (i=0; i<anses; i++) ans[i] = pt[ans[i]].id;</pre>
  return anses;
int main(void) {
 int i, j, jj;
 double tmp;
  scanf("%d", &n);
  for (i=0; i<n; i++) scanf("%lf %lf", &x[i], &y[i]);</pre>
  for (i=0; i<n; i++) {</pre>
   x[n] = 2*(-INF)-x[i]; y[n] = y[i];
   x[n+1] = x[i]; y[n+1] = 2*INF-y[i];
    x[n+2] = 2*INF-x[i]; y[n+2] = y[i];
```

```
x[n+3] = x[i]; y[n+3] = 2*(-INF)-y[i];
  for (j=0; j<n+4; j++) if (j!=i) {
   jj = j - (j>i);
   inv[jj].id = j;
   tmp = (x[j]-x[i])*(x[j]-x[i]) + (y[j]-y[i])*(y[j]-y[i]);
   inv[jj].x = (x[j]-x[i])/tmp;
   inv[jj].y = (y[j]-y[i])/tmp;
   inv[jj].ang = atan2(inv[jj].y, inv[jj].x);
  vors = convexHull(n+3, inv, vor);
  // Build bisectors
  for (j=0; j<vors; j++) {</pre>
   ans[j][0].x = (x[i]+x[vor[j]])/2;
    ans[j][0].y = (y[i]+y[vor[j]])/2;
    ans[j][1].x = ans[j][0].x - (y[vor[j]]-y[i]);
   ans[j][1].y = ans[j][0].y + (x[vor[j]]-x[i]);
 printf("Around (%lf, %lf)\n", x[i], y[i]);
 // List all intersections of the bisectors
  for (j=1; j<vors; j++) {</pre>
   PT vv:
   vv = ComputeLineIntersection(ans[j-1][0], ans[j-1][1],
                                 ans[i][0], ans[i][1]);
   printf("%lf, %lf\n", vv.x, vv.y);
 printf("\n");
return 0;
```

file:///mit/ecprice/acm/acm08/notebook.html#file7

## Euclid's algorithm, etc. (C++)

```
// 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.
using namespace std;
typedef vector<int> VI;
typedef pair<int,int> PII;
// return a % b (positive value)
int mod (int a, int b) {
  int ret = a % b;
  if (ret < 0) ret += b;</pre>
  return ret;
// computes gcd(a,b)
```

```
int gcd (int a, int b){
 if (b == 0) return a;
  return gcd (b, a % b);
// 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 == 0){
   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 = 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(x[0], a[0]);
  for (int i = 1; i < x.size(); i++){</pre>
    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;
}
```

file:///mit/ecprice/acm/acm08/notebook.html#file7

#### Linear systems, matrix inverse (Stanford) (C++)

```
// Gauss-Jordan elimination with partial 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[][])
             returns determinant of a[][]
//
const double EPSILON = 1e-7;
typedef vector<double> VD;
typedef vector<VD> VVD;
// Gauss-Jordan elimination with partial pivoting
double GaussJordan (VVD &a, VVD &b){
 double det = 1;
  int i,j,k;
  int n = a.size();
  int m = b[0].size();
  for (k=0; k<n; k++) {
   j=k;
    for (i=k+1;i<n;i++) if (fabs(a[i][k])>fabs(a[j][k])) j = i;
    if (fabs(a[j][k]) < EPSILON) { cerr << "Matrix is singular." << endl; exit(1)</pre>
    for (i=0;i<n;i++) swap (a[j][i],a[k][i]);</pre>
    for (i=0;i<m;i++) swap (b[i][i],b[k][i]);</pre>
    if (j!=k) det *= -1;
    double s = a[k][k];
    for (j=0;j<n;j++) a[k][j] /= s;
    for (j=0;j<m;j++) b[k][j] /= s;
    det *= s;
    for (i=0;i<n;i++) if (i != k){</pre>
      double t = a[i][k];
      for (j=0;j<n;j++) a[i][j] -= t*a[k][j];</pre>
      for (j=0;j<m;j++) b[i][j] -= t*b[k][j];
   }
 }
  return det;
RREF, matrix rank (C++)
// Reduced row echelon form via Gauss-Jordan elimination
// with partial pivoting. This can be used for computing
// the rank of a matrix.
// Running time: 0(n^3)
//
// INPUT:
             a[][] = an nxn matrix
```

// OUTPUT: rref[][] = an nxm matrix (stored in a[][])

returns rank of a[][]

```
const double EPSILON = 1e-7:
typedef vector<double> VD;
typedef vector<VD> VVD;
// returns rank
int rref (VVD &a){
  int i, j, r, c;
  int n = a.size();
  int m = a[0].size();
  for (r=c=0;c<m;c++){
    j=r;
    for (i=r+1;i<n;i++) if (fabs(a[i][c])>fabs(a[j][c])) j = i;
    if (fabs(a[j][c])<EPSILON) continue;</pre>
    for (i=0;i<m;i++) swap (a[j][i],a[r][i]);</pre>
    double s = a[r][c];
    for (j=0;j<m;j++) a[r][j] /= s;</pre>
    for (i=0;i<n;i++) if (i != r){</pre>
      double t = a[i][c];
      for (j=0;j<m;j++) a[i][j] -= t*a[r][j];</pre>
    r++;
  return r;
Simplex (C++)
// This is a simple simplex solver. It solves:
// Maximize obi[0] + obi[1]*x*1 + ... + obi[n]*x n
// Subject to
// x 1 >= 0, ..., x n >= 0
// for each i, c[i][0] + c[i][1]*x_1 + ... + c[i][n]*x_n >= 0
// DO NOT TRY TO REUSE LP OBJECTS!!!!! (INFEASIBLE corrupts them.)
// You should consider calling srand() first.
typedef vector<double> VD;
typedef vector<VD> VVD:
typedef vector<int> VI;
class LP
public:
  int nvars, ncons; // # decision vars and # constraints
  VD obj;
                     // [cols]
  VVD c;
                     // ncons * cols (left column is constant)
```

```
// Results in intelligible form
  double objval;
  VD assignments;
  enum Result {FAILED, INFEASIBLE, UNBOUNDED, FEASIBLE, OPTIMAL};
private:
 int cols:
                     // width of the constraint matrix
 VI nonbasic orig; // [nvars]
 VI basic_orig; // [ncons]
public:
 LP(int nvars, int ncons) : nvars(nvars), ncons(ncons),
                             cols(1 + nvars).
                             obj(1 + nvars, 0.0)
  {
    c = VVD(ncons, VD(cols, 0.0));
    for(int i = 0; i < nvars; i++)</pre>
      nonbasic orig.push back(i);
    for(int i = 0; i < ncons; i++)</pre>
      basic_orig.push_back(i + nvars);
  void pivot(int col, int row)
    // Enforce that the old col remains nonnegative.
      double val = 1.0 / c[row][col];
      for (int i = 0; i < cols; i++)</pre>
        c[row][i] *= -val;
      c[row][col] = val;
   // Subtract the extra stuff the pivot row brings along.
    for (int i = 0: i < ncons: i++) {</pre>
      if (i == row) continue:
      double coeff = c[i][col];
      c[i][col] = 0.0;
      for (int j = 0; j < cols; j++)
        c[i][j] += coeff * c[row][j];
    double coeff = obj[col];
    obj[col] = 0.0;
    for (int j = 0; j < cols; j++)
      obj[j] += coeff * c[row][j];
   // Update maps to original indices.
    swap(nonbasic_orig[col - 1], basic_orig[row]);
```

```
// Returns true if successful, false if unbounded
bool simplex()
{
 // Bland's rule: pick an arbitrary column and
 // do the pivot that will change it the least.
 while (true) {
   // Pick a random nonbasic column to pivot.
   int offset = rand() % (cols - 1), col = -1;
    for (int i = 0; i < cols - 1; i++) {</pre>
     int c = (offset + i) % (cols - 1) + 1;
     if (obj[c] > 1e-8) {
       col = c;
        break:
     }
    if (col == -1)
     break; // This basis is optimal.
   // Find the row that will hit zero first.
    double min change = 1e100;
    int best row = -1;
    for (int row = 0; row < ncons; row++) {</pre>
     if (c[row][col] >= -1e-8) continue;
      double change = -c[row][0] / c[row][col];
     if (change < min change) {</pre>
        min change = change;
        best row = row;
    if (best row == -1) // Unbounded!
     return false:
    pivot(col, best_row);
 }
 // Produce output
  objval = obj[0];
 assignments.resize(ncons + nvars);
  for (int i = 0; i < ncons; i++)
    assignments[basic_orig[i]] = c[i][0];
  for (int i = 0; i < nvars; i++)</pre>
    assignments[nonbasic orig[i]] = 0.0;
  return true;
Result phase1()
 // Find equation with minimum b
 int worst row = 0;
 for (int i = 1; i < ncons; i++)</pre>
```

```
if (c[i][0] < c[worst row][0])
    worst row = i;
if (c[worst row][0] >= -1e-8)
  return FEASIBLE;
// Add a new variable epsilon, which we minimize.
for (int i = 0; i < ncons; i++)</pre>
  c[i].push back(1.0);
VD orig obj = obj;
obj = VD(cols, 0.0);
obj.push back(-1.0);
int eps var = nvars + ncons;
nonbasic orig.push back(eps var);
nvars++:
cols++;
// We started out infeasible, so pivot epsilon into the basis.
pivot(cols-1, worst row);
if (!simplex())
  return FAILED; // Unbounded phase 1 here is bad.
if (objval < -1e-9)
  return INFEASIBLE; // Epsilon must be nonpositive.
// Force epsilon out of the basis
// (It's zero anyway within our precision).
for (int i = 0; i < ncons; i++) {</pre>
  if (basic_orig[i] == eps_var) {
    pivot(1, i);
    break;
// Find epsilon's column.
int eps col = -1;
for (int i = 0: i < nvars: i++)
  if (nonbasic orig[i] == eps var)
    eps col = i+1;
// Epsilon is nonbasic and thus zero, so we can remove it.
for (int i = 0; i < ncons; i++) {</pre>
  c[i][eps col] = c[i][cols-1];
  c[i].pop back();
nonbasic_orig[eps_col - 1] = nonbasic_orig.back();
nonbasic orig.pop back();
cols--:
nvars--;
// Restore the original objective.
obj = VD(cols, 0.0);
```

```
obi[0] = orig obi[0];
  for (int i = 0; i < nvars; i++) {</pre>
   if (nonbasic orig[i] < nvars)</pre>
      obj[i+1] = orig obj[nonbasic orig[i] + 1];
  for (int i = 0; i < ncons; i++) {</pre>
    if (basic orig[i] < nvars)</pre>
      for (int j = 0; j < cols; j++)
        obj[j] += orig_obj[basic_orig[i] + 1] * c[i][j];
 return FEASIBLE;
Result solve()
  Result p1_res = phase1();
 if (p1_res != FEASIBLE)
    return p1_res;
  assignments.clear(); // Poison it.
  if (!simplex())
    return UNBOUNDED:
  return OPTIMAL;
void printState()
  printf("Maximize %lf ", obj[0]);
  for(int i = 1; i < cols; i++)</pre>
   printf(" + %lf*x %d", obj[i], nonbasic orig[i-1]);
  printf("\nSubject to\n
  for(int i = 1; i < cols; i++) {</pre>
   printf("x%-7d", nonbasic orig[i-1]);
  for(int i = 0; i < ncons; i++)
      printf("\nx%-5d", basic orig[i]);
      for(int j = 1; j < cols; j++)</pre>
        printf("%8.4lf", c[i][j]);
      printf(" + %lf >= 0", c[i][0]);
  printf("\n\n");
void printResult()
  printf("Objective = %.6lf\n", objval);
  for (int i = 0; i < nvars; i++)</pre>
    printf(" x%d = %.6lf\n", i, assignments[i]);
  for (int i = 0; i < ncons; i++)
    printf(" r%d = %.6lf\n", i, assignments[nvars + i]);
```

```
};
```

```
FFT(C++)
typedef vector<int> VI;
double PI = acos(0) * 2;
class complex
public:
        double a, b;
        complex() \{a = 0.0; b = 0.0; \}
        complex(double na, double nb) {a = na; b = nb;}
        const complex operator+(const complex &c) const
                {return complex(a + c.a, b + c.b);}
        const complex operator-(const complex &c) const
                {return complex(a - c.a, b - c.b);}
        const complex operator*(const complex &c) const
                {return complex(a*c.a - b*c.b, a*c.b + b*c.a);}
        double magnitude() {return sqrt(a*a+b*b);}
        void print() {printf("(%.3f %.3f)\n", a, b);}
};
class FFT
{
public:
        vector<complex> data;
        vector<complex> roots;
        VI rev;
        int s, n;
        void setSize(int ns)
                s = ns;
                n = (1 << s);
                int i, j;
                rev = VI(n);
                data = vector<complex> (n);
                roots = vector<complex> (n+1);
                for (i = 0; i < n; i++)
                        for (j = 0; j < s; j++)
                                if ((i & (1 << j)) != 0)
                                        rev[i] += (1 << (s-j-1));
                roots[0] = complex(1, 0);
                complex mult = complex(cos(2*PI/n), sin(2*PI/n));
                for (i = 1; i <= n; i++)
                        roots[i] = roots[i-1] * mult;
        }
        void bitReverse(vector<complex> &array)
```

```
{
        vector<complex> temp(n);
        int i;
        for (i = 0; i < n; i++)
                temp[i] = array[rev[i]];
        for (i = 0; i < n; i++)
                array[i] = temp[i];
}
void transform(bool inverse = false)
        bitReverse(data);
        int i, j, k;
        for (i = 1; i <= s; i++) {
                int m = (1 << i), md2 = m / 2;
                int start = 0, increment = (1 << (s-i));</pre>
                if (inverse) {
                        start = n;
                        increment *= -1:
                complex t, u;
                for (k = 0; k < n; k += m) {
                        int index = start;
                        for (j = k; j < md2+k; j++) {
                                t = roots[index] * data[j+md2];
                                index += increment;
                                data[i+md2] = data[i] - t;
                                data[j] = data[j] + t;
                        }
                }
        if (inverse)
                for (i = 0; i < n; i++) {
                        data[i].a /= n;
                        data[i].b /= n;
                }
}
static VI convolution(VI &a, VI &b)
        int alen = a.size(), blen = b.size();
        int resn = alen + blen - 1;
                                       // size of the resulting array
        int s = 0. i:
        while ((1 << s) < resn) s++; // n = 2^s
        int n = 1 << s; // round up the the nearest power of two</pre>
        FFT pga, pgb;
        pga.setSize(s); // fill and transform first array
        for (i = 0; i < alen; i++) pga.data[i] = complex(a[i], 0);
        for (i = alen; i < n; i++)</pre>
                                        pga.data[i] = complex(0, 0);
        pga.transform();
```

file:///mit/ecprice/acm/acm08/notebook.html#file7

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```
pgb.setSize(s); // fill and transform second array
                for (i = 0; i < blen; i++)</pre>
                                                 pgb.data[i] = complex(b[i], 0)
                for (i = blen; i < n; i++)</pre>
                                                 pgb.data[i] = complex(0, 0);
                pgb.transform();
                for (i = 0; i < n; i++) pga.data[i] = pga.data[i] * pgb.data[i]
                pga.transform(true); // inverse transform
                VI result = VI (resn); // round to nearest integer
                for (i = 0; i < resn; i++)</pre>
                                                 result[i] = (int) (pga.data[i]
                int actualSize = resn - 1;  // find proper size of array
                while (result[actualSize] == 0)
                        actualSize--:
                if (actualSize < 0) actualSize = 0:</pre>
                result.resize(actualSize+1);
                return result;
        }
};
int main()
{
        VI a = VI (10);
        for (int i = 0; i < 10; i++)
                a[i] = (i+1)*(i+1);
        VI b = FFT::convolution(a, a);
        /* 1 8 34 104 259 560 1092 1968 3333
        5368 8052 11120 14259 17104 19234 20168 19361 16200 10000*/
        for (int i = 0; i < b.size(); i++)</pre>
                printf("%d ", b[i]);
        return 0:
```

#### Dense Dijkstra's (C++)

```
void Dijkstra (const VVT &w, VT &dist, VI &prev, int start){
  int n = w.size();
  VI found (n);
  prev = VI(n, -1);
  dist = VT(n, 1000000000);
  dist[start] = 0;

while (start != -1){
  found[start] = true;
  int best = -1;
  for (int k = 0; k < n; k++) if (!found[k]){
    if (dist[k] > dist[start] + w[start][k]){
      dist[k] = dist[start] + w[start][k];
      prev[k] = start;
  }
```

```
if (best == -1 || dist[k] < dist[best]) best = k;
}
start = best;
}</pre>
```

```
Topological sort (C++)
// This function uses performs a non-recursive topological sort.
// Running time: O(|V|^2). If you use adjacency lists (vector<map<int> >),
                 the running time is reduced to O(|E|).
//
//
    INPUT: w[i][i] = 1 if i should come before i. 0 otherwise
//
    OUTPUT: a permutation of 0, \ldots, n-1 (stored in a vector)
              which represents an ordering of the nodes which
//
//
              is consistent with w
//
// If no ordering is possible, false is returned.
typedef double TYPE;
typedef vector<TYPE> VT;
typedef vector<VT> VVT;
typedef vector<int> VI:
typedef vector<VI> VVI;
bool TopologicalSort (const VVI &w, VI &order){
  int n = w.size();
  VI parents (n):
  aueue<int> a:
  order.clear();
  for (int i = 0; i < n; i++){
    for (int j = 0; j < n; j++)
     if (w[j][i]) parents[i]++;
    if (parents[i] == 0) q.push (i);
  while (q.size() > 0){
    int i = q.front();
    q.pop();
    order.push_back (i);
    for (int j = 0; j < n; j++) if (w[i][j]){
     parents[i]--;
      if (parents[j] == 0) q.push (j);
```

return (order.size() == n);

}

```
Kruskal's (C++)
Uses Kruskal's Algorithm to calculate the weight of the minimum spanning
forest (union of minimum spanning trees of each connected component) of
a possibly disjoint graph, given in the form of a matrix of edge weights
(-1 if no edge exists). Returns the weight of the minimum spanning
forest (also calculates the actual edges - stored in T). Note: uses a
disjoint-set data structure with amortized (effectively) constant time per
union/find. Runs in O(E*log(E)) time.
typedef int TYPE;
struct edge
    int u. v:
    TYPE d;
};
struct edgeCmp
{
    int operator()(const edge& a, const edge& b) { return a.d > b.d; }
};
int find(vector <int>& C, int x) { return (C[x] == x) ? x : C[x] = find(C, C[x])
TYPE Kruskal (vector <vector <TYPE> >& w)
     int n = w.size();
     TYPE weight = 0;
     vector <int> C(n), R(n):
     for(int i=0; i<n; i++) { C[i] = i; R[i] = 0; }</pre>
     vector <edge> T;
     priority_queue <edge, vector <edge>, edgeCmp> E;
     for(int i=0; i<n; i++)</pre>
         for(int j=i+1; j<n; j++)</pre>
            if(w[i][j] >= 0)
                edge e;
                e.u = i; e.v = j; e.d = w[i][j];
                E.push(e);
     while (T.size() < n-1 \&\& !E.empty())
```

```
{
    edge cur = E.top(); E.pop();
    int uc = find(C, cur.u), vc = find(C, cur.v);
    if(uc != vc)
    {
        T.push_back(cur); weight += cur.d;
        if(R[uc] > R[vc]) C[vc] = uc;
        else if(R[vc] > R[uc]) C[uc] = vc;
        else { C[vc] = uc; R[uc]++; }
    }
}
return weight;
}
```

# **Longest Increasing Subsequence (C++)**

```
// 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
typedef vector<int> VI;
typedef pair<int,int> PII;
typedef vector<PII> VPII:
#define STRICTLY INCREASING
VI LongestIncreasingSubsequence(VI v) {
  VPII best:
  VI dad(v.size(), -1):
  for (int i = 0; i < v.size(); i++) {</pre>
#ifdef STRICTLY INCREASIG
    PII item = make pair(v[i], 0);
    VPII::iterator iter = lower bound(best.begin(), best.end(), item);
    item.second = i:
#else
    PII item = make_pair(v[i], i);
    VPII::iterator iter = upper_bound(best.begin(), best.end(), item);
    if (iter == best.end()) {
      dad[i] = (best.size() == 0 ? -1 : best.back().second);
      best.push back(item);
   } else {
```

```
dad[i] = dad[iter->second];
   *iter = 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 (C++)

```
// Routines for performing computations on dates. In these routines,
// months are exprsesed as integers from 1 to 12, days are expressed
// as integers from 1 to 31, and years are expressed as 4-digit
// integers.
string dayOfWeek[] = {"Mo", "Tu", "We", "Th", "Fr", "Sa", "Su"};
// 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;
  i = 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];
Knuth-Morris-Pratt (C++)
Searches for the string w in the string s (of length k). Returns the
O-based index of the first match (k if no match is found). Algorithm
runs in O(k) time.
void buildTable(string& w, vector <int>& t)
  t = vector <int>(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;
  vector <int> t;
  buildTable(w, t);
  while(m+i < s.length()) {</pre>
   if(w[i] == s[m+i]) {
      i++;
      if(i == w.length()) return m;
   } else {
      m += i-t[i];
      if(i > 0) i = t[i];
```

## Hashed strstr (C++)

return s.length();

```
const char *fast_strstr(const char *haystack, const char *needle)
{
  unsigned target = 0, power = 1, hash = 0;
  size t nlen = strlen(needle), hlen = strlen(haystack);
```

```
if (hlen < nlen || !*needle)
  return 0;
for(int i = 0; i < nlen; i++) {
  target = target * 257 + needle[i];
  hash = hash * 257 + haystack[i];
  power = power * 257;
}
for(int i = nlen; i <= hlen; i++) {
  if (hash == target && !memcmp(haystack + i - nlen, needle, nlen))
    return haystack + i - nlen;
  hash = hash * 257 + haystack[i] - power * haystack[i-nlen];
}
return 0;
}</pre>
```

## Java formatting (Java)

```
// examples for printing floating point numbers
import java.util.*;
import java.io.*;
import java.text.DecimalFormat;
public class DecFormat {
   public static void main(String[] args) {
        DecimalFormat fmt;
        // round to at most 2 digits, leave of digits if not needed
        fmt = new DecimalFormat("#.##");
        System.out.println(fmt.format(12345.6789)); // produces 12345.68
        System.out.println(fmt.format(12345.0)): // produces 12345
        System.out.println(fmt.format(0.0)); // produces 0
        System.out.println(fmt.format(0.01)); // produces .1
        // round to precisely 2 digits
        fmt = new DecimalFormat("#.00"):
        System.out.println(fmt.format(12345.6789)); // produces 12345.68
        System.out.println(fmt.format(12345.0)); // produces 12345.00
        System.out.println(fmt.format(0.0)); // produces .00
        // round to precisely 2 digits, force leading zero
        fmt = new DecimalFormat("0.00"):
        System.out.println(fmt.format(12345.6789)); // produces 12345.68
        System.out.println(fmt.format(12345.0)); // produces 12345.00
        System.out.println(fmt.format(0.0)); // produces 0.00
        // round to precisely 2 digits, force leading zeros
        fmt = new DecimalFormat("000000000.00");
        System.out.println(fmt.format(12345.6789)); // produces 000012345.68
        System.out.println(fmt.format(12345.0)); // produces 000012345.00
```

```
System.out.println(fmt.format(0.0)); // produces 000000000.00
    // force leading '+'
    fmt = new DecimalFormat("+0; -0");
    System.out.println(fmt.format(12345.6789)); // produces +12346
    System.out.println(fmt.format(-12345.6789)); // produces -12346
    System.out.println(fmt.format(0)); // produces +0
    // force leading positive/negative, pad to 2
    fmt = new DecimalFormat("positive 00; negative 0");
    System.out.println(fmt.format(1)); // produces "positive 01"
    System.out.println(fmt.format(-1)); // produces "negative 01"
    // goute special chars (#)
    fmt = new DecimalFormat("text with '#' followed by #");
    System.out.println(fmt.format(12.34)); // produces "text with # follow
    // always show "."
    fmt = new DecimalFormat("#.#"):
    fmt.setDecimalSeparatorAlwaysShown(true);
    System.out.println(fmt.format(12.34)); // produces "12.3"
    System.out.println(fmt.format(12)); // produces "12."
    System.out.println(fmt.format(0.34)); // produces "0.3"
    // different grouping distances:
    fmt = new DecimalFormat("#,####.###");
    System.out.println(fmt.format(123456789.123)); // produces "1,2345,678
    // scientific:
    fmt = new DecimalFormat("0.000E00");
    System.out.println(fmt.format(123456789.123)); // produces "1.235E08"
    System.out.println(fmt.format(-0.000234)): // produces "-2.34E-04"
    // using variable number of digits:
    fmt = new DecimalFormat("0");
    System.out.println(fmt.format(123.123)); // produces "123"
    fmt.setMinimumFractionDigits(8);
    System.out.println(fmt.format(123.123)); // produces "123.12300000"
    fmt.setMaximumFractionDigits(0);
    System.out.println(fmt.format(123.123)); // produces "123"
    // note: to pad with spaces, you need to do it yourself:
    // String out = fmt.format(...)
    // while (out.length() < targlength) out = " "+out;</pre>
}
```

## Complicated regex example (Java)

// Code which demonstrates the use of Java's regular expression libraries.

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}

```
// This is a solution for
//
//
   Loglan: a logical language
// http://acm.uva.es/p/v1/134.html
import java.util.*;
import java.util.regex.*;
public class LogLan {
    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, when
            // line is terminated by a period. The code below reads lines unt
            // encountering a line whose final character is a '.'. Note the \( \text{\chi} \)
            //
            //
                 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
            //
            // Other useful String manipulation methods include
            //
            // s.compareTo(t) < 0 if s < t, lexicographically
            // s.indexOf("apple") returns index of first occurrence of "apple"
                 s.lastIndexOf("apple") returns index of last occurrence of '
                 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/upperc
            //
            //
                 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)
            if (pattern.matcher (removed period).find()){
```

```
System.out.println ("Good");
            } else {
                System.out.println ("Bad!");
        }
   }
}
```

## Java geometry (Java)

```
// 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:
// 0 0 10 0 0 10
// 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.*;
import iava.awt.geom.*:
import java.math.*;
import java.io.*;
public class JavaGeometrv {
   // make a list of doubles from a string
    static ArrayList<Double> readPoints (String s){
        StringTokenizer st = new StringTokenizer (s);
        ArrayList<Double> ret = new ArrayList<Double>();
        while (st.hasMoreTokens())
            ret.add (Double.parseDouble (st.nextToken()));
        return ret;
```

```
// make an Area object from the coordinates of a polygon
static Area makeArea (ArrayList<Double> points){
    // note that the GeneralPath object does not allow construct
    // of polygons based on doubles -- we must use floats.
    GeneralPath gp = new GeneralPath();
    qp.moveTo ((float) points.get(0).doubleValue(),
               (float) points.get(1).doubleValue());
    for (int i = 2; i < points.size(); i += 2)</pre>
        gp.lineTo ((float) points.get(i).doubleValue(),
                   (float) points.get(i+1).doubleValue());
    gp.closePath();
    return new Area (gp);
}
// compute area of polygon
static double computePolygonArea (ArrayList<Point2D.Double> points){
    // convert to array, for convenience
    Point2D.Double[] pts = points.toArray (new Point2D.Double[0]);
    double area = 0;
    for (int i = 0; i < pts.length; i++){</pre>
        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);
    ArrayList<Double> pointsA = readPoints (scanner.nextLine());
    ArrayList<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\ )
        //
                                                               double w, double
        //
               creates an ellipse inscribed in box with bottom-left corner (x,
        11
        11
               and upper-right corner (x+y,w+h)
        11
             Rectangle2D.Double rect = new Rectangle2D.Double (double x, doub)
        //
        //
                                                                double w, doubl
        //
               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 (re
}
```

## 3D geom (Java)

```
public class Geom3D {
 // distance from point (x, y, z) to plane aX + bY + cZ + d = 0
 public static double ptPlaneDist(double x, double v, 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) {</pre>
        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-py)*(z-py);
  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));
}
```

```
### .emacs
(setq column-number-mode t)
(setq inhibit-startup-message t)
(transient-mark-mode t)
(global-font-lock-mode t)
(show-paren-mode t)
(global-set-key "\C-cg" 'goto-line)
(defun previous-6-lines nil
  (interactive)
  (previous-line 6))
(defun next-6-lines nil
  (interactive)
  (next-line 6))
(global-set-key [(control up)] 'previous-6-lines)
(global-set-key [(control down)] 'next-6-lines)
### .bashrc
export CXXFLAGS="-Wall -g"
PS1='\u@\h:\W\$'
export PYTHONSTARTUP="$HOME/.pythonrc"
alias "l=ls -lh --color=auto"
### .pythonrc
import readline
import rlcompleter
readline.parse_and_bind("tab: complete")
### .vimrc
runtime! debian.vim
```