PeerSim HOWTO 2: build a topology generator

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1 Introduction

This tutorial teaches you how to build from scratch a new peersim (peersim project page: `http://sourceforge.net/projects/peersim`) topology generator. In order to understand this tutorial, the reader is encouraged to start reading the first peersim tutorial (`http://peersim.sourceforge.net/peersim_HOWTO.html`) to have an idea of the basic concepts that will not be discussed any further in this document.

The aim of this tutorial is to be as practical as possible; the goal is to give the reader ideas about technical or intermediate level features of peersim and to encourage him/her to experiment further. The full source code discussed in this document is available via CVS at peersim project page in the *peersim.example.hot* class package.

2 What is a topology?

The network abstraction in peersim is a (sometimes huge) array of *Node* structures (interfaces); because of the size of the network and to overcome scalability problems, usually in large P2P networks each node knows about the existence of a very small subset of other nodes (ex: order of $\log(N)$ where N is the whole network size). Thus each node has a short list of other node references, usually called "neighbors", build accordingly to some kind of strategy or rule.

Thus, we can say that a topology is how nodes are arranged (linked) together and clearly this depends upon the particular chosen rule. Examples of topology are the following (not exhaustive at all):

- random graphs

- Watts-Strogatz model graph
- star model
- ring model
- lattice model
- ...

2.1 Which rule to choose?

In this document, we have chosen to code a particular topology generator to build Internet-like tree topologies. The building process is based on the *preferential attachment* approach. The rule applied is quite simple and takes into account geometric and network constraints to better mimic real world network. The preferential attachment choice can be affected by a parameter (α) that amplifies or reduces the geometric location influence in favor of the path distance.

The rule strategy is the following: we consider a square unit region D, then we start with node $x(0)$ chosen at random and we set $W(x(0)) = 0$ (it is the root node). For each i with $i = 1...n-1$ we choose a new node $x(i)$ in the region D and we connect it to an **early inserted** node $x(j)$ that minimize the following formula:

$$W(x(j)) + \alpha \cdot dist(x(i), x(j)) \text{ with } 0 \leq \text{j} < \text{i}$$

where:

- $W(x(j))$ is the distance in terms of hops (the path distance from node $x(j)$ to the root node);
- $dist(...)$ is the usual Euclidean distance;
- α is a weight parameter that can minimize or maximize the geometric distance influence;

After having chosen a node $x(j)$, we set $W(x(i)) = W(x(j)) + 1$. At the end we obtain a tree rooted in $x(0)$.

We have extended this model to improve robustness allowing every node to have exactly d outbound neighbors instead of only one. This means that, at the time of joining the network, each node should have at least d candidates to be selected as neighbors. To achieve this property, as a first step we select

at random exactly d root nodes and we connect them together in a ring fashion (a doubly linked list). In this way each ordinary node has at least d nodes (the d roots) to choose from in order to select its neighbors. In other words, each node has to select the best d nodes that minimize the function above.

To get further details about this model, we suggest the following readings:

1. "Heuristically Optimized Trade-offs: A New Paradigm for Power Laws in the Internet"
   (`http://cgi.di.uoa.gr/~elias/publications/paper-fkp02.pdf`)
2. "Degree distributions of the FKP network model"
   (`http://research.microsoft.com/~jchayes/Papers/FKPdgrees.pdf`)
3. "On Power-Law Relationships of the Internet Topology"
   (`http://www.cs.ucr.edu/~michalis/CAMERA.ps`)

The model should generate a topology that exhibits a power-law bound on the in-degree sequence of nodes; but, as stated in the second previously listed paper, this power-law prediction is not true.

3 What we need to code

In order to run this model in peersim we need to write Java classes extending some peersim typical interfaces. Very in brief, we need:

- the protocol class: the protocol itself does nothing because we want something that auto-magically builds this topology model from a raw list of unconnected nodes; we are not interested in running any piece of code over time (cycles in peersim terminology). This class is a sort of structure to collect some needed values such as the node space coordinates, the hop distance, the in degree counter and so on. The reader can think to this class as "glue code".
- the initializer class: it extends the *Dynamics* interface and deal with all the initialization process. As we'll see further, the initializer code itself is very compact because all the building process complexity is hidden in a custom made factory pattern class. Please note that we don't need explicitly an *Observer* object because we are not interested in observing any behavior over time; due to that, the code to track down informations about the actual generated tree (coordinates and in-degree distribution) are embedded into the initializer.

4 Code writing

4.1 Protocol class

As we stated so far, the protocol code is minimal:

```
import peersim.core.IdleProtocol;

public class InetNodeProtocol extends IdleProtocol {
    // coordinates in space:
    public double x;
    public double y;

    public int in_degree;
    public int hops;
    public boolean isroot;

    /** Creates a new instance of hotNodeProtocol */
    public InetNodeProtocol(String prefix, Object obj) {
   super(prefix);
        in_degree = 0;
        hops = 0;
        isroot = false;
    }

    public Object clone() throws CloneNotSupportedException {
        InetNodeProtocol af = (InetNodeProtocol) super.clone();
        ...
return af;
    }

    public void nextCycle(peersim.core.Node, int protocolID) {
    }
```

The *nextCycle()* method is empty, so it's presence is completely optional. The class is basically a structure encapsulated in an object.

4.2 Initializer class

This initializer can be considered as a public interface to the model. From the outside (means from the peersim configuration file) only few main parameters are needed, such as the *Linkable* enabled protocol, the outbound degree and the α parameter. The constructor method takes care of collecting these

parameters or to set up the corresponding default values The actual structure generation is performed elsewhere.

```
public class InetInitializer implements peersim.dynamics.Dynamics {
     /**
      * String name of the parameter that defines the protocol to initialize.
     * Parameter read will has the full name
     * <tt>prefix+"."+PAR_PROT</tt>
     */
     public static final String PAR_PROT = "protocol";

     /**
      * String name of the parameter about the out degree value.
     */
     public static final String PAR_OUTDEGREE = "d";

     /**
      * String name of the parameter used as a weight.
     */
     public static final String PAR_ALFA = "alfa";

     /**
      * String name of the parameter used as a maximum x or y coordinate. All the
      * nodes are on a square region.
     */
     public static final String PAR_MAX_COORD = "max_coord";

     ....

     /** Creates a new instance of InetInitializer */
     public InetInitializer(String prefix) {
         // super(prefix);
         pid = Configuration.getPid(prefix+"."+PAR_PROT);
         d = Configuration.getInt(prefix+"."+PAR_OUTDEGREE);
         alfa = Configuration.getDouble(prefix+"."+PAR_ALFA);
         graph_filename = "cmplxnet_d"+ d + "_alfa"+alfa+".dat";
         dg_filename = "degree_d"+ d + "_alfa"+alfa+".dat";

         maxcoord = Configuration.getDouble(prefix + "." + PAR_MAX_COORD, 1.0);
         if ( !graph_filename.equals("") ) {
             try { graph_fileout = new PrintWriter(new FileWriter(graph_filename));
                 System.out.println(prefix +" filename: "+graph_filename +" selected");}
             catch (Exception e) {;}
         }
         if ( !dg_filename.equals("") ) {
             try { dg_fileout = new PrintWriter(new FileWriter(dg_filename));
                 System.out.println(prefix +" filename: "+dg_filename +" selected");}
             catch (Exception e) {;}
         }
```

```
    }
}
```

The *modify()* method is defined by the *Dynamics* interface; it invokes a specialized factory object that works on top of a *OverlayGraph* type object (actual implementation of the *Graph* interface). This object provides the high level abstraction of a graph on the simulator overlay network; in this way, it allows the application to use many well known graph algorithms and operations (many operations can be found in *peersim.graphGraphAlgorithms* class). The operations performed on it are reflected on the actual topology (ex: adding an edge).

```
    public void modify() {
        OverlayGraph ogr = new OverlayGraph(pid);
        InetFactory.InetTree(ogr, CommonRandom.r, pid, maxcoord, d, alfa );
        graphToFile(ogr);
        dgDistribToFile(ogr);
        }
    }
```

The other two method invocation in the *modify()* method are used to write data on stable storage. In the first one, for each node n the x and y coordinates are collected and then for each neighbor i of node n the coordinates are written in the following format:

```
n.neighbor(i).x n.neighbor(i).y \newline
n.x n.y \newline
\newline}
```

The particular line triplet formatting order suits the gluplot needs. Please note that the for loop starts from index d, not from 0; this means that the root node(s) is not directly considered because it has not any outbound connections, but only inbound connections. Nevertheless the root node(s) are plotted inspecting the nodes that are linked directly to it.

```
    private void graphToFile(peersim.graph.Graph g) {
        if (graph_fileout != null) {
            try {
                for (int i = d ; i < g.size() ; i++ ) {
                    Node current = (Node)g.getNode(i);
                    double x_to = ((InetNodeProtocol)current.getProtocol(pid)).x;
                    double y_to = ((InetNodeProtocol)current.getProtocol(pid)).y;
                    Iterator it = (Iterator)g.getNeighbours(i).iterator();
                    while (it.hasNext()) {
```

```
                        int index = ((Integer)it.next()).intValue();
                        Node n = (Node)g.getNode(index);
                        double x_from = ((InetNodeProtocol)n.getProtocol(pid)).x;
                        double y_from = ((InetNodeProtocol)n.getProtocol(pid)).y;

                        graph_fileout.println(x_from+" "+y_from);
                        graph_fileout.println(x_to+" "+y_to);
                        graph_fileout.println("");
                    }
                }
                graph_fileout.close();
            }
            catch (Exception e) {;}
        }
    }
```

The second data collecting method builds an array of in-degree frequencies and an array of in-degree probability and dumps the second collection to file.

```
private void dgDistribToFile(peersim.graph.Graph g) {
        if (dg_fileout != null) {
            int size = g.size();
            try {
                int[] dgfrq = new int[size];
                double[] dgprob = new double[size];
                for (int i = 0 ; i < size ; i++) { // do not plot leaves
                    Node n = (Node)g.getNode(i);
                    InetNodeProtocol protocol = (InetNodeProtocol)n.getProtocol(pid);
                    int degree = protocol.in_degree;
                    dgfrq[degree]++;
                }
                double sum = 0;
                for (int i = size-1 ; i > 0 ; i--) {
                    dgprob[i] = (dgfrq[i] + sum)/size;
                    sum += dgfrq[i];
                }
                // do not count index 0: 'cos the leafs degree is clearly 0!
                for (int i = 0 ; i < dgprob.length ; i++ ) {
                    double k = (double)i/size;
                    dg_fileout.println(k+" "+dgprob[i]);
                }
                dg_fileout.close();
            }
            catch (Exception e) {e.printStackTrace();
                System.out.println(e);}
        }
    }
}
```

5 Factory class

This class is the core one. The actual topology initialization is performed here. Because of the factory pattern, all the methods in this class are static. There is no need of getting parameters from the configuration file, because they have been already collected by the initializer class; thus the factory is completely hidden.

This implementation is an extension of the peersim standard topology factory (*peersim.graph.Graph Factory*).

The only public method is the one that actually builds the topology: *InetTree()*; as parameters it gets all the parameters that the initializer class has collected. The steps performed are the following:

1. set the correct values for the d (at least one) roots, including coordinates; if there is only one root, its coordinates are centered on the square (edge size 1.0 by default), otherwise the coordinates are random (as any ordinary node).

2. initialize the coordinates and the in-degree counter for each ordinary node.

3. if there are more than one root node, than these root nodes are joined together in a ring (connections are non oriented). As an exercise, the reader can change this choice implementing something else (ex: putting the root nodes in a fully connected topology or a random graph or whatever).

4. for each node n other than the root, take exactly d nodes that minimizes the formula (see Section 2.1) and connect node n to those d nodes.

The other present methods are all private and can be considered as utility methods. In fact their function is quite straightforward and can be guessed by the method signature; nevertheless a few comments are presented in the following table:

getParents()	get the the current node best d candidates to connect to
hops()	return the graph distance in terms of hops from the root of the node given as a parameter
minHops()	return the minimum hop valued node between the specified nodes
distance()	get the standard Euclidean distance between two nodes

```
import hot.HotNodeProtocol;
import peersim.graph.*;
import peersim.core.Node;
import peersim.core.Linkable;
import peersim.core.Network;
import peersim.config.Configuration;
import java.util.Random;
import java.util.ArrayList;
import java.util.Arrays;

public class InetFactory extends peersim.graph.GraphFactory {

    private static final String DEBUG_STRING = "inet.InetFactory: ";

    public InetFactory() {
        super();
    }

    public static Graph InetTree(Graph g, Random rnd, int pid, double maxcoord,
    int outdegree, double alfa) {
        int size = g.size(); // size of the network
        System.out.println(DEBUG_STRING+"size: "+size+" outdegree: "+outdegree);

        // build outdegree roots
        System.out.println(DEBUG_STRING+"Generating "+outdegree+" root(s),
means out degree "+outdegree+"...");
        for(int i = 0 ; i < outdegree ; ++i) {
            Node n = (Node)g.getNode(i);
            HotNodeProtocol prot = (HotNodeProtocol)n.getProtocol(pid);
            prot.isroot = true;
            prot.hops = 0;
            prot.in_degree = 0;
            if (outdegree == 1 ) {
                prot.x = maxcoord/2;
                prot.y = maxcoord/2;
            }
            else { // more than one root
                if (rnd.nextBoolean() ) {
                    prot.x = maxcoord/2 + (rnd.nextDouble() * 0.1);
                }
                else {
                    prot.x = maxcoord/2 - (rnd.nextDouble() * 0.1);
                }
                if (rnd.nextBoolean() ) {
                    prot.y = maxcoord/2 + (rnd.nextDouble() * 0.1);
                }
                else {
                    prot.y = maxcoord/2 - (rnd.nextDouble() * 0.1);
```

```
        }
        System.out.println("root coord: "+prot.x+" "+prot.y);
    }
}

// Set coordinates x,y and set indegree 0
System.out.println(DEBUG_STRING+"Generating random cordinates for nodes...");
for (int i = outdegree ; i < size ; i++) {
    Node n = (Node)g.getNode(i);
    HotNodeProtocol prot = (HotNodeProtocol)n.getProtocol(pid);
    if (maxcoord == 1.0) {
        prot.x = rnd.nextDouble();
        prot.y = rnd.nextDouble();
    }
    else {
        prot.x = rnd.nextInt((int)maxcoord);
        prot.y = rnd.nextInt((int)maxcoord);
    }
    prot.in_degree = 0;
}

// Connect the roots in a ring if needed (thus, if there are more than 1
// root nodes.
if (outdegree > 1) {
    System.out.println(DEBUG_STRING+"Putting roots in a ring...");
    for (int i = 0 ; i < outdegree ; i++) {
        Node n = (Node)g.getNode(i);
        ((HotNodeProtocol)n.getProtocol(pid)).in_degree++;
        n = (Node)g.getNode(i+1);
        ((HotNodeProtocol)n.getProtocol(pid)).in_degree++;

        g.setEdge(i, i+1);
        g.setEdge(i+1, i);
    }
    Node n = (Node)g.getNode(0);
    ((HotNodeProtocol)n.getProtocol(pid)).in_degree++;
    n = (Node)g.getNode(outdegree);
    ((HotNodeProtocol)n.getProtocol(pid)).in_degree++;
    g.setEdge(0, outdegree);
    g.setEdge(outdegree, 0);
}

// for all the nodes other than root(s), connect them!
for (int i = outdegree ; i < size ; ++i ) {
    Node n = (Node)g.getNode(i);
    InetNodeProtocol prot = (InetNodeProtocol)n.getProtocol(pid);

    prot.isroot = false;
```

```
        // look for a siutable parent node between those allready part of the
        // overlay topology: alias FIND THE MINIMUM!
        Node candidate = null;
        int candidate_index = 0;
        double min = Double.POSITIVE_INFINITY;
        if (outdegree > 1) {
                int candidates[] = getParents(g, pid, i, outdegree, alfa);
                for (int s = 0 ; s < candidates.length ; s++) {
                    g.setEdge(i, candidates[s]);
                    System.out.print(i+", ");
                }
                prot.hops = minHop(g, candidates, pid) + 1;
        }
        else { // degree 1:
            for (int j = 0 ; j < i ; j++) {
                Node parent = (Node)g.getNode(j);
                InetNodeProtocol prot_parent = (InetNodeProtocol)parent.getProtocol(pid);

                double value = hops(parent, pid) +
                    (alfa * distance(n, parent, pid));
                if (value < min) {
                    candidate = parent; // best parent node to connect to
                    min = value;
                    candidate_index = j;
                }
            }
            prot.hops = ((InetNodeProtocol)candidate.getProtocol(pid)).hops + 1;
            g.setEdge(i, candidate_index);
            ((HotNodeProtocol)candidate.getProtocol(pid)).in_degree++;
        }

    }
    System.out.println(DEBUG_STRING+"Graph generation finished!");
    return g;
}

private static int[] getParents(Graph g, int pid, int cur_node_index,
int how_many, double alfa) {
    int result[] = new int[how_many];
    ArrayList net_copy = new ArrayList(cur_node_index);
    // fill up the sub net copy:
    for (int j = 0 ; j < cur_node_index ; j++) {
        net_copy.add(j, (Node)g.getNode(j));
    }

    // it needs exactly how_many minimums!
    for (int k = 0 ; k < how_many ; k++) {
        int candidate_index = 0;
        double min = Double.POSITIVE_INFINITY;
```

```
            // for all the elements in the copy...
            for (int j = 0 ; j < net_copy.size() ; j++) {
                Node parent = (Node)net_copy.get(j);
                HotNodeProtocol prot_parent = (HotNodeProtocol)parent.getProtocol(pid);
                double value = hops(parent, pid) + (alfa *
                    distance((Node)g.getNode(cur_node_index), parent, pid));

                if (value < min) {
                    min = value;
                    candidate_index = j;
                }
            }
            result[k] = candidate_index;        // collect the parent node
            net_copy.remove(candidate_index); // delete the min from the net copy
        }
        return result;
    }


    private static int hops(Node node, int pid) {
        return ((HotNodeProtocol)node.getProtocol(pid)).hops;
    }

    private static int minHop(Graph g, int[] indexes, int pid) {
        int min = Integer.MAX_VALUE;
        for (int s = 0 ; s < indexes.length ; s++) {
            Node parent = (Node)g.getNode(indexes[s]);
            int value = ((HotNodeProtocol)parent.getProtocol(pid)).hops;
            if (value < min) {
                min = value;
            }
        }
        return min;
    }

    private static double distance(Node new_node, Node old_node, int pid) {
        // Eucleidian distance code...
    }
}
```

6 Experiments

In order to make the model run, a proper peersim configuration file is needed.
The one presented in the following lines may suits the reader needs:

```
# Complex Network file:
random.seed 1234567890
```

```
simulation.cycles 1

overlay.size 10000
overlay.maxSize 10000

protocol.0 hot.InetNodeProtocol
#protocol.0.maxcoord 1000

init.0 hot.InetInitializer
init.0.alfa 20
init.0.protocol 0
init.0.d 1
```

It produces a 10000 node overlay network with the parameters listed in the `init.0` section.

The presented figures show the produced topology and highlight the parameter α importance. In fact, it affects the clustering behaviour of the system and it is tightly correlated to the size of the network. If α il lower than $\sqrt{netsize}$, the topology becomes more and more clustered (as show in the first two figures); with extremely low α, the topology becomes a star. On the other end, if α is grater than $\sqrt{netsize}$, the topology tends to be random and not clustered at all (the second row of images). For deeper details, please consult the previously listed papers.

All the images has been produced using only one root node and only one outbound connection per node. Using two or more outbound connection per node leads to a massively crowded plot, so it is not a very nice picture to plot!

The degree distribution related to the generated star topology (Figure 1) is not shown (it is simply a straight line). Clearly the plots show that there is not any evidence about in-degree power-law distribution; only in the case of $\alpha = 4$, the corresponding plot exhibits a power-law like behavior at least for a subset of the nodes, but this is very different from what first listed paper was talking about.

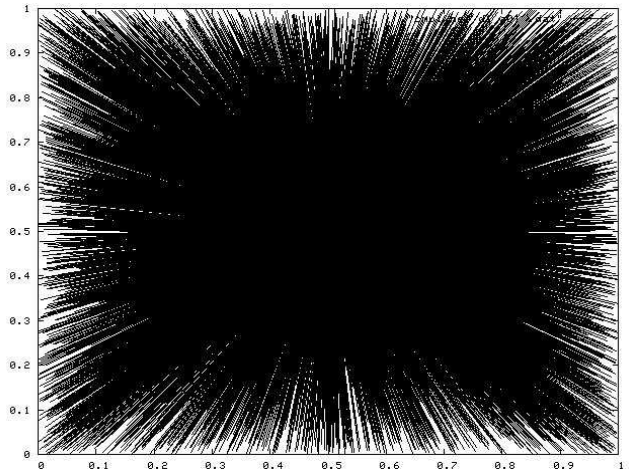

Figure 1: Topology with α 0.1

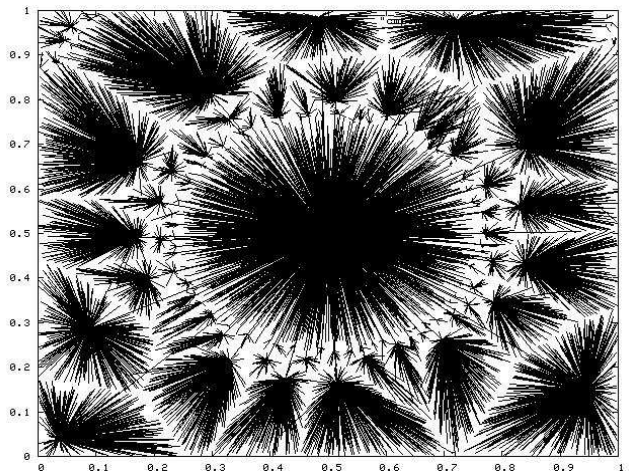

Figure 2: Topology with α 4

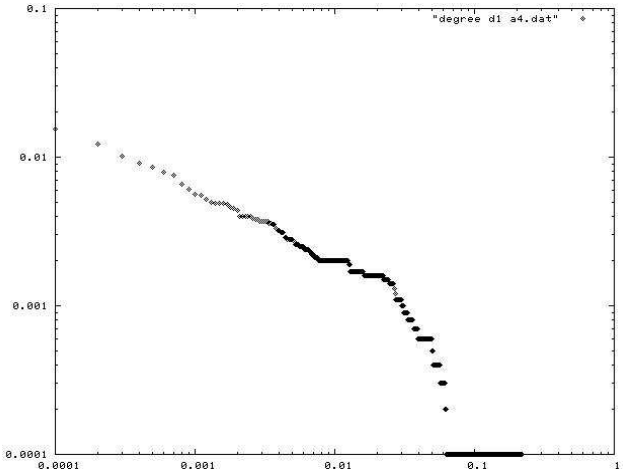

Figure 3: In-degree distribution with α 4

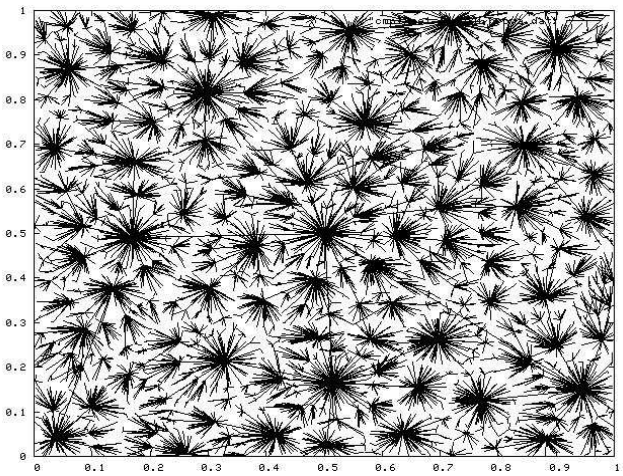

Figure 4: Topology with α 20

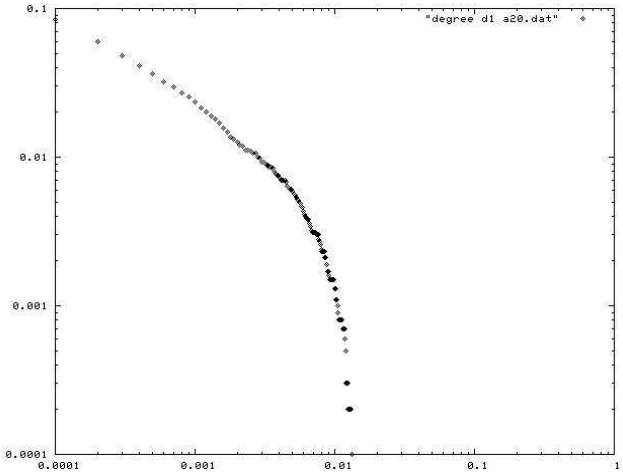

Figure 5: In-degree distribution with α 20

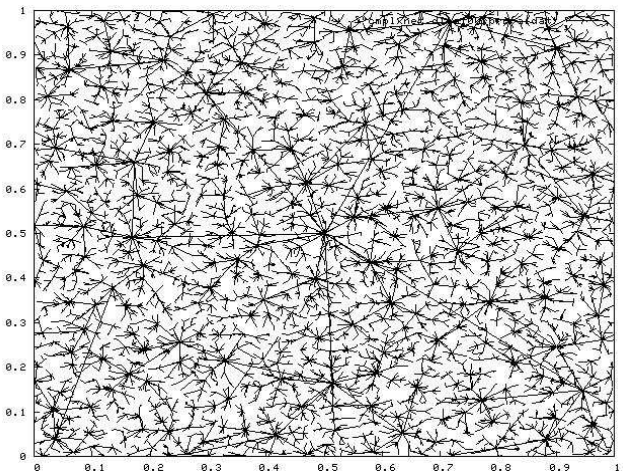

Figure 6: Topology with α 100

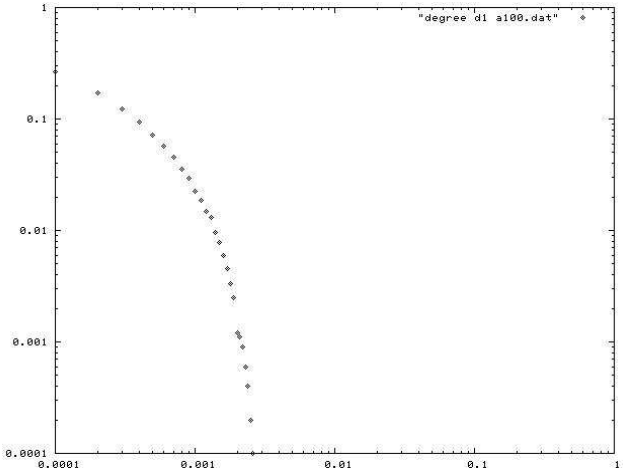

Figure 7: In-degree distribution with α 100

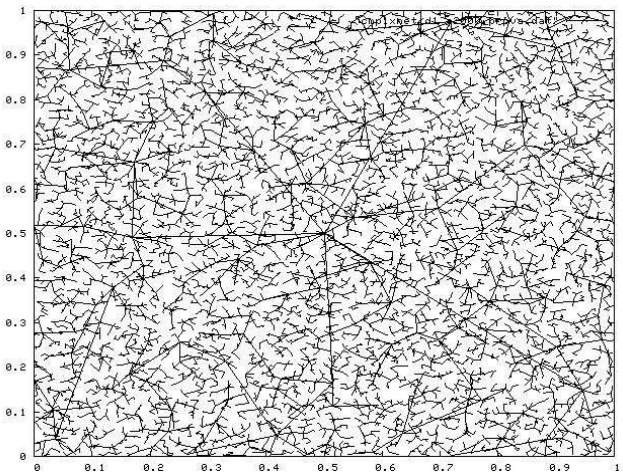

Figure 8: Topology with α 2000

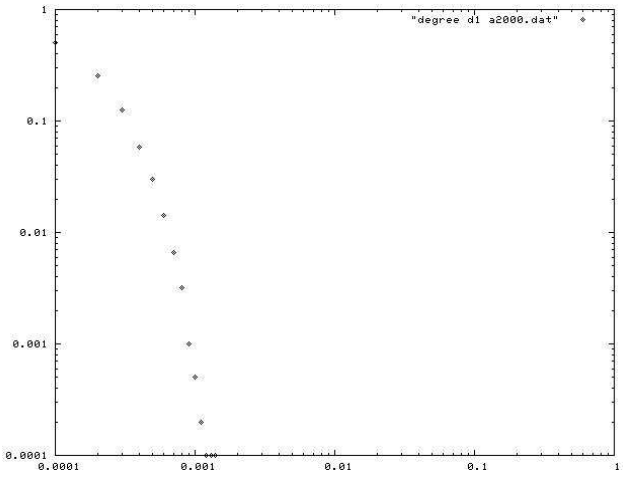

Figure 9: In-degree distribution with α 2000