

# The Percolation Behavior of Current Through Square and Cubic Resistor Networks

Peter Gilbert

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## Abstract

This paper begins with a discussion of elementary percolation theory, especially the "critical phenomenon." This theory is used to study the behavior of current flowing across two and three dimensional disordered electrical networks composed of random resistors. Resistors are assigned resistance values according to a given probability, and the map taking probabilities to expected values of the effective conductance of the network is studied. Given a probability  $p$ , the forward problem is to estimate the value of this map. Given an object composed of a random mixture of two materials each having constant conductance, the inverse problem is to estimate the proportion  $p$  of materials from boundary measurements of the effective conductance of the object. Algorithms are presented which solve the forward and inverse problems and estimate the critical probability in two and three dimensions.

# 1 Introduction to Percolation Theory

Suppose that a large cubic chunk of turf is immersed in water. What is the probability that a grain of dirt near the center of the chunk of turf is wetted? Percolation theory is concerned with questions like this. In the standard "percolation model," let  $Z^d$  be the  $d$  dimensional lattice. Let  $p$  be a number satisfying  $0 \leq p \leq 1$ . Declare each edge of the lattice to be open with probability  $p$  and closed with probability  $1 - p$ . Water can flow through an open edge and not through a closed edge; hence  $p$  is the proportion of passages which are broad enough to allow water to flow along them. Objects such as the chunk of turf have large networks of fissures which can be approximated by the infinite lattice  $Z^3$ . Approximate the chunk of turf by  $Z^3$ , and let  $x$  be a vertex of  $Z^3$  near the center of the chunk of turf. Then the vertex  $x$  is wetted if and only if there is a path from  $x$  to a vertex on the boundary of the chunk of turf, using open edges only (see Figure 1). Percolation theory is concerned with the existence of such open paths.

## 1.1 The Critical Phenomenon

One of the most interesting aspects of percolation theory is the occurrence of a "critical phenomenon." On the infinite lattice  $Z^d$ , define an open cluster to be a set of open edges in which for any two edges in the cluster, there exists a path of open edges connecting them. Define the percolation probability  $\theta^d(p)$  to be the probability that the origin of the lattice  $Z^d$  belongs to an infinite open cluster. A fundamental theorem of percolation theory is that for dimension  $d \geq 1$ , there exists a critical probability  $p_c^d$  such that

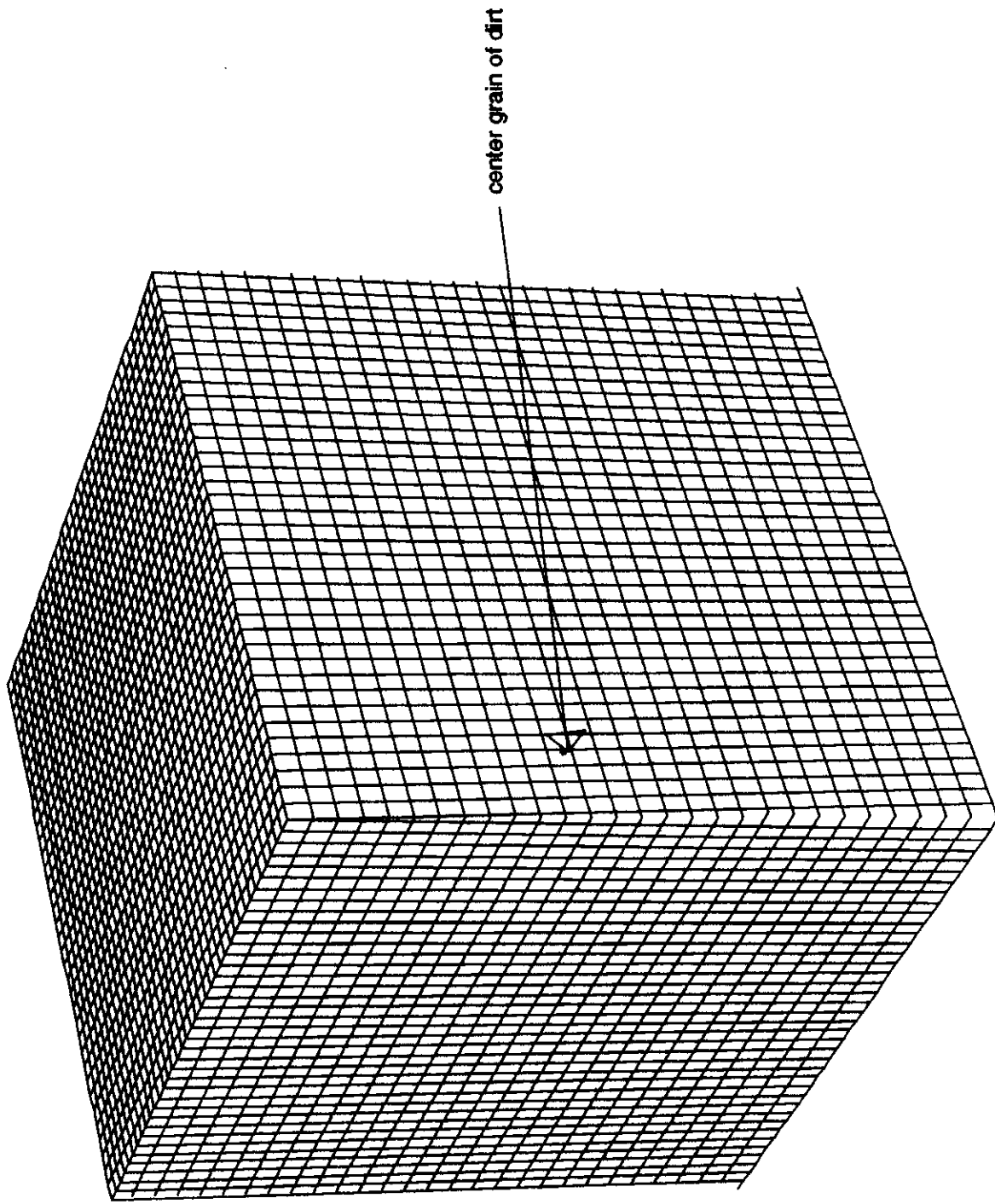
$$\begin{aligned}\theta^d(p) &= 0 \text{ if } p < p_c \\ &> 0 \text{ if } p > p_c\end{aligned}$$

Formally the critical probability  $p_c^d$  is defined by

$$p_c^d = \sup\{p : \theta^d(p) = 0\}$$

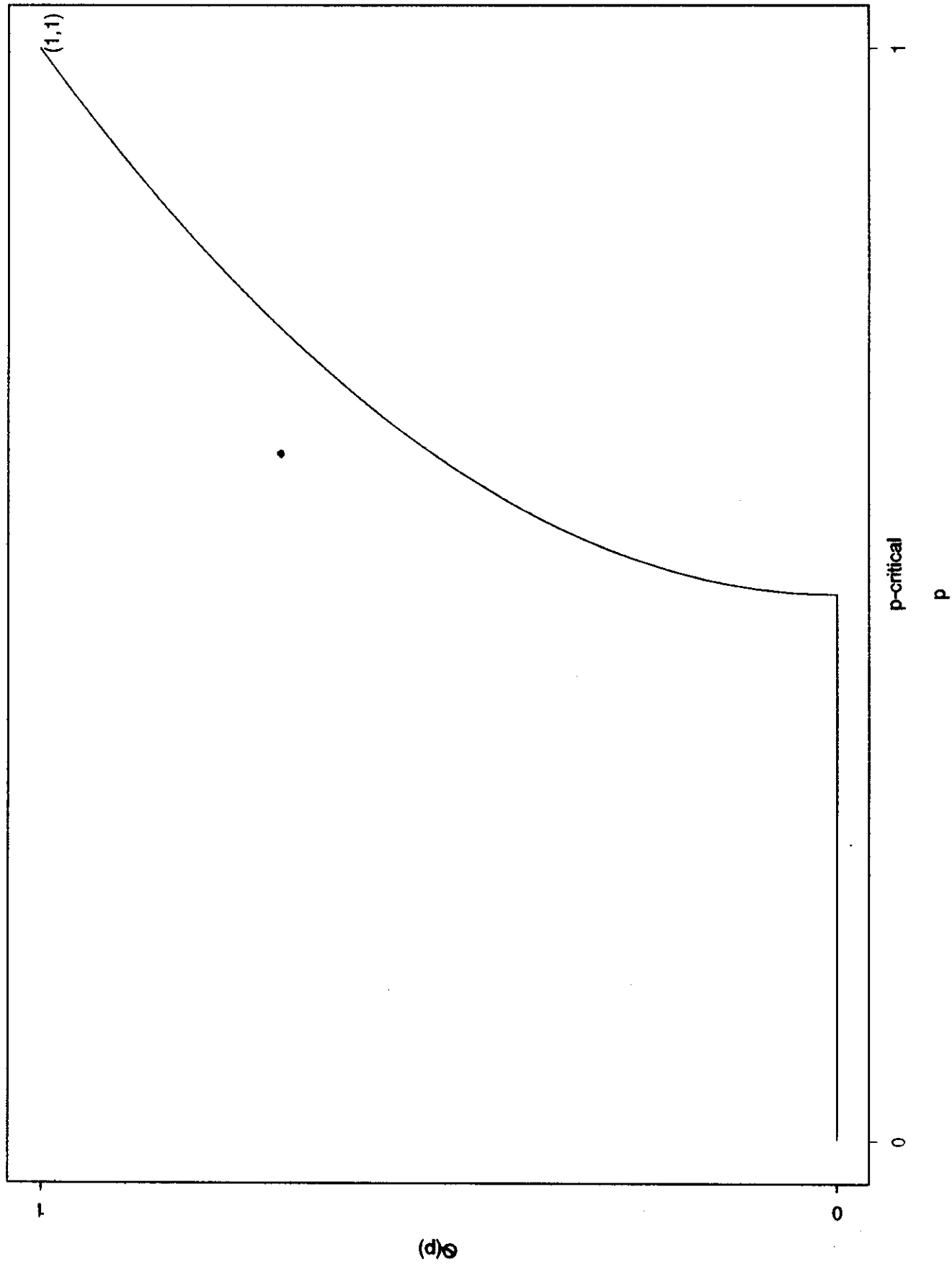
For a proof that the critical probability exists for  $d \geq 1$ , see Grimmett (p13). Grimmett asserts that the percolation probability  $\theta(p)$  behaves roughly as in Figure 2.

Figure 1 Modelling the Network of Fissures in the Chunk of Turf



Is the Center Grain of Dirt Wetted when the Chunk of Turf is Immersed in Water?

Figure 2 Percolation Probability V Probability



The critical probability  $p_c^1$  is clearly one, and it has been proved that  $p_c^2 = 1/2$  (Grimmett, Kesten). For  $d > 2$ , the value of the critical probability is unknown. In this paper I develop a method to estimate the critical probability in two and three dimensions. The method involves studying the percolation of electrical current across square and cubic lattice resistor networks.

## 2 Setup of the Square and Cubic Resistor Networks

For  $d = 2$ , let  $G_n^2 = \{1, 2, \dots, n\} \times \{1, 2, \dots, n\}$  be the interior nodes defining the square network, and let  $G_{n0}^2 = \{(0, m) : 1 \leq m \leq n\}$  and  $G_{n1}^2 = \{(n+1, m) : 1 \leq m \leq n\}$  be the left and right side nodes of  $G_n^2$ , respectively. Assign a potential of zero to each of the nodes on  $G_{n0}^2$  and a potential of one to each of the nodes on  $G_{n1}^2$ . Insulate the top and bottom sides of the network. Place silver bars along the left and right sides of the network and connect a battery, creating a potential difference of one across the resistor network. The two-dimensional network  $G_n^2$  connected to a battery is shown in Figure 3.

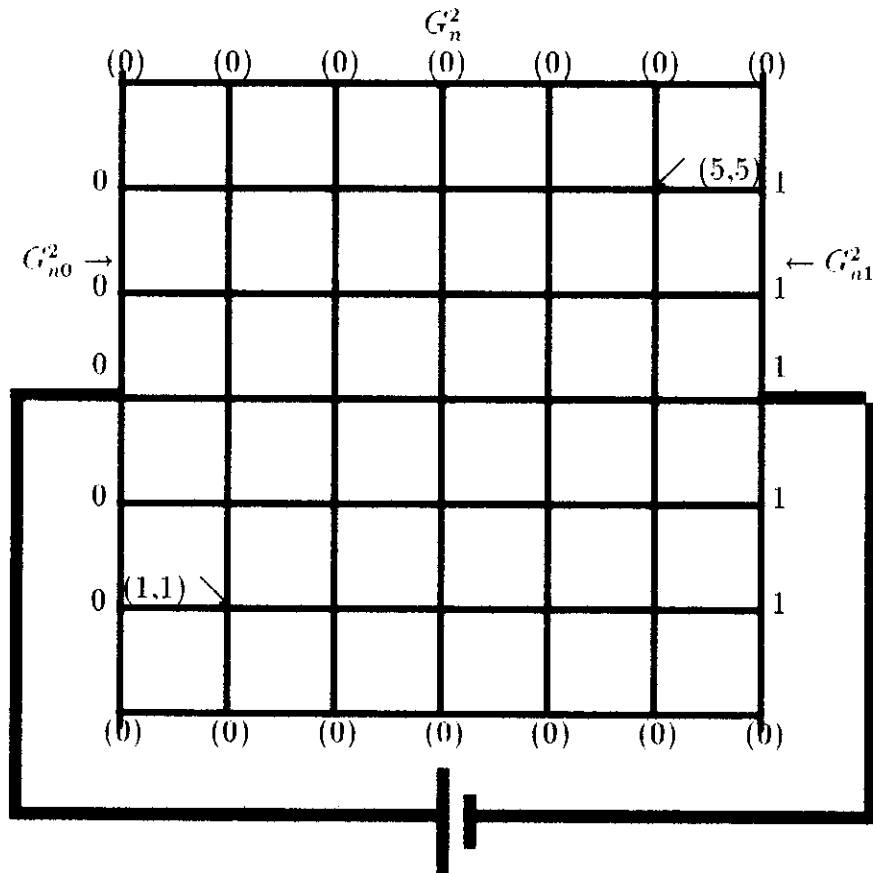


Figure 3:  $G_n^2$  connected to a battery

For  $d = 3$  the situation is similar. Here  $G_n^3 = \{1, 2, \dots, n\} \times \{1, 2, \dots, n\} \times \{1, 2, \dots, n\}$  are the interior nodes of the cubic network, and the nodes on the left and right faces are  $G_{n0}^3 = \{(0, l, m): 1 \leq l, m \leq n\}$  and  $G_{n1}^3 = \{(n, l, m): 1 \leq l, m \leq n\}$ . Assign zero potentials to the nodes on the left face of the cubic network and one potentials to the nodes on the right face of the network. Insulate the top, bottom, front, and back faces of the network. Place silver plates along the left and right faces of the cubic network and connect a battery, creating a potential difference of one across the resistor network.

Given a square or cubic network  $G_n^d$  and a probability  $p$  satisfying  $0 \leq p \leq 1$ , assign each edge, independently of all other edges, a resistance of  $1$  with probability  $p$  and a resistance of  $\beta$  with probability  $1 - p$ . What is the

effective resistance of the network? Given an object composed of a random mixture of two materials each having constant resistance, this problem is useful for determining the constitution of the mixture from measurements of effective resistance.

Suppose that  $\alpha = 1$  and  $\beta = \infty$ . This characterizes edges as being either open or closed; open if current can flow along it (corresponding to finite resistance 1) and closed if current cannot flow along it (corresponding to infinite resistance). This specialized problem lies in the domain of percolation theory, and is the object of this paper. I study the behavior of  $E_p(R_n^d)$ , the expected value of the resistance of the network  $G_n^d$ , as a function of the probability  $p$ . Of primary interest is determining for which values of  $p$  the expected value of the resistance  $E_p(R_n^d)$  is infinite. It is difficult to make accurate measurements of very large resistances, so the reciprocal expected conductance values  $E_p(C_n^d) = \frac{1}{E_p(R_n^d)}$  are studied.

### 3 The Map of Probabilities to Expected Conductance Values

Let  $V = 1$  be the voltage drop across the resistor network  $G_n^d$ , and let  $I_n$  be the total current between the left and right sides or faces of the network  $G_n^d$ . By Ohm's Law,  $C_n = I_n$ , so that  $C_n = 0$  is equivalent to no current passing from the right to the left side or face of the network. Let  $h_n^d : P \mapsto E_p(C_n^d)$  be the map of probabilities  $p \in [0, 1]$  to expected values of conductance across the network. It is very difficult to explicitly describe the map  $h$ ; I could only manage to express  $h_n^d(p)$  as a linear combination of edge conductance values and expected values of Kirchoff matrices. One evident property of  $h$  is that it is monotonic increasing, since  $p_1 < p_2$  implies that there is a greater chance that the network  $G_n^d(p_1)$  has more internal resistance than the network  $G_n^d(p_2)$ , which implies that more current is expected to cross  $G_n^d(p_2)$  than  $G_n^d(p_1)$ .

Another property of the function  $h_n^d$  is that it can be explicitly calculated for  $p = 1$ . When  $p = 1$ , of course each edge of the resistor network has conductance one. Since the nodes of  $G_{n0}^d$  have zero potential and the nodes of  $G_{n1}^d$  have one potential, there is a voltage drop of  $\frac{1}{n+1}$  across each horizontal

edge. Kirchoff's Law gives, for each interior node  $x \in G_n^d \setminus G_{n0}^d \cup G_{n1}^d$ ,

$$\sum_{y \in G_n^d, y \text{ adjacent to } x} | \text{Potential}(x) - \text{Potential}(y) | = 0$$

By this law, since there are  $n + 1$  horizontal edges between the left and right sides or faces of the network, the following is deduced: the nodes directly to the left of the nodes on  $G_{n1}^d$  have potential  $\frac{n}{n+1}$ , the nodes two horizontal edges away from  $G_{n1}^d$  have potential  $\frac{n-1}{n+1}$ , and finally the nodes adjacent to  $G_{n0}^d$  have potential  $\frac{1}{n+1}$ . Now notice that there are  $n^{d-1}$  nodes on each vertical cross-section (a vertical line if  $d=2$ , a vertical plane into the page for  $d=3$ ) of the network. By Ohm's Law, the total current leaving the side or face  $G_{n0}^d$  is  $\frac{n^{d-1}}{n+1}$ , so that

$$E_1(C_n^d) = \frac{n^{d-1}}{n+1}.$$

## 4 The Forward Problem

Given a probability  $p \in [0, 1]$  and a resistor network  $G_n^2$  or  $G_n^3$ , assign to each edge of the network a conductance  $\mu$  with probability  $p$  and conductance  $\gamma$  with probability  $1 - p$ . In two dimensions connect a battery and set up the network as shown in Figure 3, and in three dimensions set up the network in the analogous way. The forward problem is to estimate the value of  $h_n^d(p)$ . In section 8 I describe an algorithm which calculates the sample mean conductance  $\bar{C}_n^d(p)$  corresponding to the probability  $p$ . I wrote a simple function in Splus titled *forward*, which takes the sample mean, the sample size, and the sample standard deviation as parameters and returns a 95% confidence interval about the true mean conductance  $h_n^d(p)$ . For given  $p$ , the confidence interval is calculated as

$$\left( \bar{C}_n^d(p) - 1.96 * \frac{\sigma}{\text{sqrt}(n)}, \bar{C}_n^d(p) + 1.96 * \frac{\sigma}{\text{sqrt}(n)} \right),$$

where  $\sigma$  is the sample standard deviation and 1.96 is the Z statistic corresponding to  $\alpha = .05$  in the standard normal distribution. In Table 1 there is a list of confidence intervals for several probabilities calculated on the two dimensional network with 20 nodes on a side  $G_{20}^2$ .



<i>Probability</i>	<i>Lower Limit for Confidence Bar</i>	<i>Upper Limit for Confidence Bar</i>
0.00	$9.52 e^{-11}$	$9.52 e^{-11}$
0.10	$1.19 e^{-11}$	$1.21 e^{-11}$
0.20	$1.60 e^{-11}$	$1.64 e^{-11}$
0.30	$2.56 e^{-11}$	$2.74 e^{-11}$
0.40	$5.66 e^{-11}$	$6.89 e^{-11}$
0.45	0.0023	0.0075
0.46	0.0058	0.013
0.47	0.0059	0.014
0.48	0.0098	0.019
0.49	0.015	0.026
0.50	0.022	0.035
0.51	0.027	0.041
0.52	0.043	0.058
0.53	0.049	0.068
0.54	0.060	0.078
0.55	0.066	0.085
0.60	0.162	0.182
0.70	0.349	0.369
0.80	0.548	0.564
0.90	0.752	0.764
1.00	0.952	0.952

Table 1: Confidence Intervals about values of  $h_{20}^2(p)$

## 5 The Inverse Problem

Let  $O$  be an object which is known to be a random mixture of two materials having constant conductance  $\mu$  and  $\gamma$  respectively. Approximate the object  $O$  by the appropriate square or cubic resistor network  $G_n^2$  or  $G_n^3$ . Connect a battery and set up the network as described in section 2. Measure the effective conductance of the network  $m$  times, and compute the sample mean  $C_n^d(p)$ . The inverse problem is to recover the constitution of the object  $O$  by estimating the probability  $p$ , which is the proportion of materials composing the object  $O$ . I wrote an Splus function titled *inverse* which takes a range of

probabilities, the corresponding sample means  $\bar{C}_n^d$ , the corresponding sample standard deviations, the sample size  $m$  and the dimension  $d$  as parameters and returns a 95% confidence interval about the true probability  $p$ . See the Appendix for information on how the function calculates the confidence interval. In Table 2 there is a list of confidence intervals about the true mean  $p$  corresponding to 100 effective conductance measurements taken on the two dimensional network with 20 nodes on a side  $G_{20}^2$ .

<i>Sample Mean</i>	<i>Lower Limit for Confidence Bar</i>	<i>Upper Limit for Confidence Bar</i>
0.001	0.0090	0.1710
0.01	0.0090	0.1900
0.0289	0.4510	0.5100
0.05	0.4910	0.5430
0.10	0.5530	0.5790
0.20	0.6140	0.6210
0.30	0.6610	0.6720
0.40	0.7110	0.7210
0.50	0.7630	0.7710
0.60	0.8130	0.8260
0.70	0.8620	0.8760
0.80	0.9100	0.9250
0.90	0.9750	0.9780

Table 2 Confidence bars about true probabilities on  $G_{20}^2$  with  $m=100$

## 5.1 The Critical Phenomenon of the Function $h$

Intuitively, for large  $n$  the function  $h_n^d$  is similar to the percolation probability function  $\theta$ . Given  $p$ ,  $\theta(p)$  is the probability that the origin of the infinite lattice  $Z^d$  is a member of an infinite open cluster. For a large finite lattice, the existence of an infinite open cluster is intuitively a similar condition to the existence of an open path from one side or face of the network to the other. Hence  $\theta^d(p) = 0$  corresponds to no open path existing from one side or face of the network to the other. Since it is known that a critical probability  $p_c^d$  exists for the function  $\theta^d(p)$ , it is reasonable to guess that the probability of an open path existing from one side or face of the network to the other

exhibits critical behavior. Moreover, I claim that the probability function describing for which probabilities there exists an open path from one side or face of the network to the other is equivalent to the probability function describing for which probabilities positive current flows. To see this, recall that by Ohm's Law  $C_n = I_n$ . The total current  $I_n$  is calculated by first solving for all of the interior potentials of the network. Then

$$C_n^d = I_n^d = \sum_{v \text{ a node adjacent to } G_{n0}^d} \text{Voltage}(v) \text{Conductance}(\text{edge to left of } v).$$

From this expression it is clear that  $C_n = 0$  if and only if there is an open path from  $G_{n0}^d$  to  $G_{n1}^d$ . Hence studying the probability of an open path existing from one side or face of the network to the other is equivalent to studying the probability of positive effective conductance.

Suppose that there is a critical probability  $\kappa^d$  for which no current flows across the network for probabilities less than  $\kappa^d$  and current flows across the network for probabilities greater than  $\kappa^d$ . Then, if  $\kappa^d$  exists, for large  $n$  the function  $h_n^d$  is of the form:

$$\begin{aligned} h_n^d(p) &= 0 \text{ if } p < \kappa^d \\ &> 0 \text{ if } p > \kappa^d. \end{aligned}$$

For  $d = 2$  I prove that  $\kappa^2$  exists, and is equal to  $p_c^2 = \frac{1}{2}$ . For  $d = 3$  I argue that  $\kappa^3$  exists. For  $d = 2$  and  $d = 3$  I describe an algorithm which maps probabilities to sample mean conductances of the network  $G_n^d$ . With this algorithm I approximate  $h_n^2$ , and use the approximation to empirically validate the theoretical result that  $\kappa^2 = \frac{1}{2}$ . I then approximate  $h_n^3$  and use the approximation to estimate  $\kappa^3$ . I conjecture that  $\kappa^3 = p_c^3$ , so that through this algorithm the unknown critical probability  $p_c^3$  is estimated.

## 6 Proof That $\kappa^2$ Exists and Equals $\frac{1}{2}$

**Theorem:**

Let  $G_n^2$  be a square resistor network, where each edge has conductance 1 with probability  $p$  and conductance 0 with probability

$1 - p$ . Let  $E_p(C_n^2)$  be the expected value of the conductance between  $G_{n0}^2$  and  $G_{n1}^2$ . Then there exists a positive integer  $N$  such that  $n \geq N \Rightarrow$

$$P\{E_p(C_n^2) = 0\} = 1 \text{ if } p < \frac{1}{2}, \quad (1)$$

$$P\{E_p(C_n^2) > 0\} = 1 \text{ if } p > \frac{1}{2}. \quad (2)$$

Equation (1) shows that for large enough square networks, no current will flow across the network if  $p < \frac{1}{2}$ , and equation (2) shows that positive current will flow if  $p > \frac{1}{2}$ . Hence  $\kappa^2$  exists and equals  $\frac{1}{2}$ . I prove this Theorem, leaving out some in-depth percolation theory arguments which digress from the focus of this paper.

## 6.1 Proof of (1)

Let  $p \in [0, \frac{1}{2})$ . To begin the proof of (1), recall that in section 5.1 I showed that the effective conductance  $C_n^2$  of the network is zero if and only if there is no conducting path from  $G_{n0}^2$  to  $G_{n1}^2$ . Hence

$$P\{E(C_n^2) = 0\} = P\{\text{There does not exist an open path from } G_{n0} \text{ to } G_{n1}\}.$$

Select an interior node near the center of the resistor network  $G_n^2$  to be the origin. Let  $W$  be the open cluster centered at the origin, and let  $\#W$  denote the number of edges in  $W$ . Finally let  $E_p(\#W)$  be the expected value of  $\#W$  on the network  $G_n^2$ .

Next, define a network  $H_n^2$  by stacking three networks  $G_n^2$  on top of each other. Then  $H_n^2$  is described by its interior nodes  $\{1, 2, \dots, n\} \cup \{1, 2, \dots, 3n\}$  and its left and right side nodes  $H_{n0}^2 = \{(0, m) : 1 \leq m \leq 3n\}$  and  $H_{n1}^2 = \{(n+1, m) : 1 \leq m \leq 3n\}$  (see Figure 4).

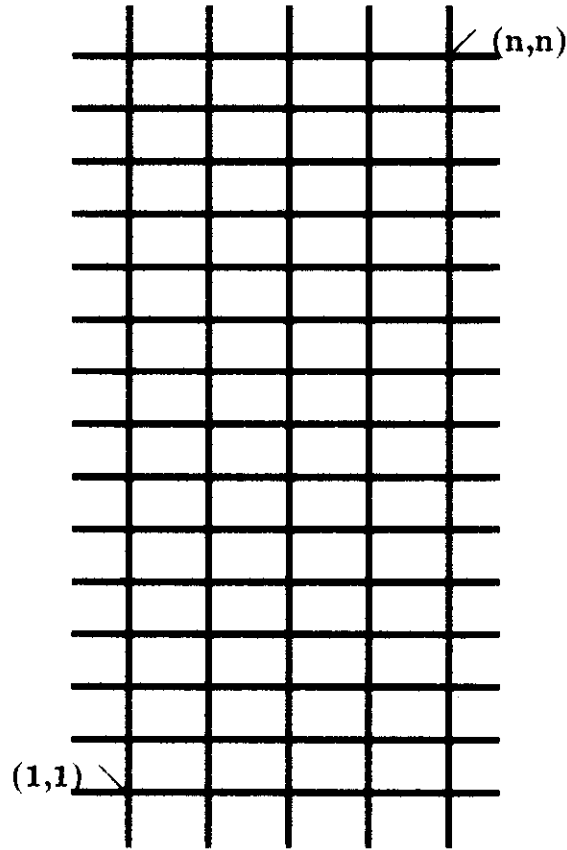


Figure 4:  $H_n^2$

Let  $\Psi(p)$  be the probability that there exists a conducting path from  $H_{n_0}^2$  to  $H_{n_1}^2$ . Harry Kesten, in Percolation Theory for Mathematicians (p 83), proved that if  $p \in [0, \frac{1}{2})$ , and  $E_p(\#W) < \infty$ , then  $\Psi(p) \rightarrow 0$  as  $n \rightarrow \infty$ . The proof is long and digresses from this paper's focus, so it is not included. Call Kesten's result Lemma 1:

Lemma 1:

If  $p \in [0, \frac{1}{2})$  and  $E_p(\#W) < \infty$ , then  $\Psi(p) \rightarrow 0$  as  $n \rightarrow \infty$ .

An important result of percolation theory is that if the conclusion of Lemma 1 holds, then as  $n$  increases, the probability that  $\#W$  is greater than or equal to  $n$  decays exponentially. Call this result Lemma 2 (see Kesten, p 83):

**Lemma 2:**

If  $\Psi(p) \rightarrow 0$  as  $n \rightarrow \infty$ , then there exist constants  $0 < c, k < \infty$  such that

$$P_p\{\#W \geq n\} \leq ce^{-kn}.$$

By the definition of the percolation probability function  $\theta^2(p)$  given in section 1.1, evidently  $E_p(\#W) = \theta^2(p)$ . For  $p \in [0, \frac{1}{2})$ ,  $p < p_c^2$ , so  $E_p(\#W) = \theta^2(p) < \infty$ . Hence the hypothesis of Lemma 1 holds, and its conclusion is that  $\Psi(p) \rightarrow 0$  as  $n \rightarrow \infty$ . Then by Lemma 2, for constants  $0 < c, k < \infty$ ,

$$P_p\{\#W \geq n\} \leq ce^{-kn}. \quad (3)$$

Since current has to travel through at least  $n + 2$  horizontal edges to flow from  $G_{n0}^2$  to  $G_{n1}^2$ , clearly

$$P_p\{E(C_n^2) = 0\} \geq P\{\#W < n\}. \quad (4)$$

Since  $P_p\{\#W < n\} = 1 - P\{\#W \geq n\}$ , (3) and (4) imply that

$$1 - P_p\{E(C_n^2) = 0\} \leq P\{\#W \geq n\}. \quad (5)$$

Then by (3),  $1 - P_p\{E(C_n^2) = 0\} \leq ce^{-kn}$ , which implies that

$$P_p\{E(C_n^2) = 0\} \geq 1 - ce^{-kn}. \quad (6)$$

Clearly  $1 - ce^{-kn} \rightarrow 1$  as  $n \rightarrow \infty$ . Since (6) holds for  $p < \frac{1}{2}$ , (1) is proved.

## 6.2 Proof of (2)

Let  $p > \frac{1}{2}$ . Some definitions are required to prove (2). Define  $\Phi(p, hor)$  to be the probability that there exists an open path from  $G_{n0}^2$  to  $G_{n1}^2$ , and  $\Phi^*(p, hor)$  the probability that there exists a closed path (all edges have zero conductance) from  $G_{n0}^2$  to  $G_{n1}^2$ . In addition define  $\Phi(p, ver)$  to be the probability that there exists an open path from the top to the bottom of the square network, and  $\Phi^*(p, ver)$  to

be the probability that there exists a closed path from the top to the bottom of the square network. Notice that since there are the same number of horizontal edges as vertical edges in the network  $G_n^2$ ,  $\Phi(p, hor) = \Phi(p, ver)$  and  $\Phi^*(p, hor) = \Phi^*(p, ver)$ . Kesten (p 172) proves that

$$\Phi(p, hor) + \Phi^*(p, ver) \geq 1. \quad (7)$$

The proof requires detailed manipulation of many results of percolation theory, so I will not explain it. Recall that in section 5.1 I showed that  $E_p(C_n^2) > 0$  as soon as  $p$  is large enough so that the existence of an open path from  $G_{n0}^2$  to  $G_{n1}^2$  is assured. Hence

$$P_p\{E(C_n^2) > 0\} = \Phi(p, hor) \quad (8)$$

From (7),

$$\Phi(p, hor) \geq 1 - \Phi^*(p, ver) \quad (9)$$

Now, of course  $P_p\{\text{a given edge has conductance one}\} = p$  and  $P_p\{\text{a given edge has conductance zero}\} = 1 - p$ , so evidently  $P_p\{\text{a given edge has conductance zero}\} = P_{1-p}\{\text{a given edge has conductance one}\}$ . Extending this to sequences of edges,

$$\Phi^*(p, ver) = \Phi(1 - p, ver). \quad (10)$$

For  $p > \frac{1}{2}$ , obviously  $1 - p < \frac{1}{2}$ . Hence  $E_{1-p}(\#W) < \infty$  and Lemma 1 holds, so that  $\Psi(1 - p) \rightarrow 0$  as  $n \rightarrow \infty$ . Then by Lemma 2, for constants  $0 < c, k < \infty$ ,

$$P_{1-p}\{\#W \geq n\} \leq ce^{-kn}. \quad (11)$$

Now, since an open vertical path must consist of at least  $n + 1$  edges (since the network is  $n + 1$  edges high),

$$P_{1-p}\{E(C_n^2) = 0\} \geq P_{1-p}(\#W < n). \quad (12)$$

Since  $P_{1-p}\{E(C_n^2) = 0\} = 1 - P_{1-p}\{E(C_n^2) > 0\} = 1 - \Phi(1 - p, hor) = 1 - \Phi(1 - p, ver)$ ,

$$1 - \Phi(1 - p, ver) \geq P_{1-p}\{\#W < n\}. \quad (13)$$

Evidently  $P_{1-p}(\#W < n) = 1 - P_{1-p}(\#W \geq n)$ , so that (11) becomes

$$1 - P_{1-p}(\#W < n) \leq ce^{-kn}, \text{ or} \quad (14)$$

$$P_{1-p}(\#W < n) \geq 1 - ce^{-kn}. \quad (15)$$

Then by (13) and (15),

$$1 - \Phi(1 - p, ver) \geq 1 - ce^{-kn}. \quad (16)$$

From (9) and (10),

$$\Phi(p, hor) \geq 1 - \Phi(1 - p, ver). \quad (17)$$

Then by (16) and (17),  $\Phi(p, hor) \geq 1 - ce^{-kn}$ , and by the characterization (8),

$$P_p\{E(C_n^2) > 0\} \geq 1 - ce^{-kn}. \quad (18)$$

Clearly  $1 - ce^{-kn} \rightarrow 1$  as  $n \rightarrow \infty$ . Since (18) holds for  $p > \frac{1}{2}$ , (2) is proved.

## 7 Argument for the Existence of $\kappa^3$

At the end of section 5.1 I conjectured that there exists a critical probability  $\kappa^3$  such that for  $p < \kappa^3$  the expected conductance value of the network  $G_n^3$  is zero and for  $p > \kappa^3$  the expected conductance value is positive. Define  $H_n^3$  to be the network resulting from stacking three cubic networks  $G_n^3$  on top of each other;  $H_n^3$  is the exact analog of  $H_n^2$ . Define  $\Gamma_c^3$  as:

$$\Gamma_c^3 = \sup\{p : \lim P_p\{\Upsilon(p) = 0\}\}, \quad (19)$$

where  $\Upsilon(p) = P_p\{\text{There exists an open path between the left and right faces of the network } H_n^3\}$ .

Kesten (p333) proved that for  $p < \Gamma_c^3$ ,  $\Upsilon(p) = 0$ , and for  $p > \Gamma_c^3$ ,  $\Upsilon(p)$  is uniformly bounded from zero for all  $n$ , at least for a subsequence of  $n$ 's. Hence there is a critical probability describing the probability of an open path stretching across the network  $H_n^3$ . Since the probability of nonzero expected current flow is equivalent to the probability of the existence of an open path from the left to the right face of the network, the existence of  $\kappa^3$  is implied.



## 8 The Approximating Graphs of the Functions $h_n^2$

I wrote two Fortran programs titled SquareConductances.exe and CubicConductances.exe which create values approximating the functions  $h_n^2$  and  $h_n^3$  respectively. For each probability value  $p = 0.00, 0.01, 0.02, 0.03, \dots, 0.98, 0.99, 1.00$ , SquareConductances.exe calculates a sample mean  $\bar{C}_n^2(p)$ , where  $m$  is the sample size. For given  $p$ , the following algorithm is performed  $m$  times to calculate a sample effective conductance: Using a random number generator, assign to each edge, independently of all other edges, a conductance of 1 with probability  $p$  and a conductance of zero with probability  $1 - p$ . By Ohm's Law, since the potential difference between  $G_{n1}^2$  and  $G_{n0}^2$  is one, the effective conductance is equal to the total current flowing from  $G_{n1}^2$  to  $G_{n0}^2$ . The total current is calculated by first computing the potentials of the interior nodes of the network. These interior potentials are computed from the boundary information of zero potentials on left side nodes, one potentials on right side nodes, and zero current on top and bottom side nodes.

To solve for the interior potentials, Kirchoff's Law is used to set up a system of equations  $Ax = b$ , where  $x$  is the vector of interior potentials to be solved for,  $A$  is the  $n^2$  by  $n^2$  Kirchoff matrix containing information supplied by Kirchoff's Law, and  $b$  is a vector of length  $n^2$  containing boundary information supplied by Kirchoff's Law. There are two techniques that I used to solve the linear system  $Ax = b$ . The first is a Linpackd routine which directly solves the linear system  $Ax = b$ . This method produces an accurate solution, but computation time is long and storage space must be allocated to declare the  $n^2$  by  $n^2$  Kirchoff matrix  $A$ . In the two dimensional case, these problems can be handled, but in the three dimensional case they are intolerable. The Kirchoff matrix in three dimensions is  $n^3$  by  $n^3$ , and for  $n > 6$  it becomes impossible to declare the Kirchoff matrix. Hence in addition to writing programs that utilize Linpackd routines to explicitly solve the system  $Ax = b$ , I wrote programs using the Gauss-Seidel iteration scheme. This

method only requires enough storage space to hold the solution vector  $x$ , and runs faster.

## 8.1 The Gauss-Seidel Iteration Scheme

Let  $L$  be the lower diagonal matrix consisting of the lower diagonal elements of  $A$ , and  $U$  the upper diagonal matrix consisting of the upper diagonal elements of  $A$ . Then clearly  $A = L + U$ , and the linear system  $Ax = b$  becomes  $(L + U)x = b$ . To construct an iteration scheme, suppose that  $j$  indicates the  $j$ th iteration. The goal is to find a convergent sequence of vectors  $x^j$  to the solution vector  $x$ . The idea behind Gauss-Seidel iteration is to estimate  $x^{j+1}$  and  $x^j$  for each iteration and stop the iterations when  $x^{j+1}$  and  $x^j$  are sufficiently close to each other, i.e.  $\|x^{j+1} - x^j\| \leq \epsilon$  for a prescribed  $\epsilon > 0$ . This of course does not guarantee convergence, but if the terms  $x^{j+1}$  and  $x^j$  are near enough, it is very likely that the final term  $x^{j+1}$  closely approximates the solution.

Now I motivate why Gauss-Seidel should work. If  $x^{j+1}$  and  $x^j$  are very near, then  $(Lx^{j+1} + Ux^j) = b$  approximates the linear system  $Ax = b$ . Under the hypothesis that  $L$  is invertible, this approximation becomes

$$x^{j+1} = L^{-1}(b - Ux^j) = F(x^j) \quad (\text{say}). \quad (20)$$

So for each  $j > 1$ ,  $x^{j+1}$  is expressed as a function of  $x^j$ , and  $x^{j+1}$  converges to a solution  $x$  if  $F(x) = x$ , for in this case  $Lx = b - Ux$  by (20), so that  $(L + U)x = b$  and hence  $Ax = b$ . Thus it is enough to show that  $F$  has a fixed point, which is true if  $F$  is a contraction mapping. Now,  $F$  is a contraction mapping if

$$\|x^{j+2} - x^{j+1}\| \leq C \|x^{j+1} - x^j\| \text{ for some } C < 1 \text{ and for all } j \geq 1. \quad (21)$$

From (20),  $x^{j+2} = L^{-1}(b - Ux^{j+1})$ . Hence  $\|x^{j+2} - x^{j+1}\| = \|L^{-1}(b - Ux^{j+1}) - L^{-1}(b - Ux^j)\| = \|L^{-1}b - L^{-1}Ux^{j+1} - L^{-1}b + L^{-1}Ux^j\| = \|L^{-1}U(x^j - x^{j+1})\| \leq \|L^{-1}U\| \|x^j - x^{j+1}\|$ . Hence  $F$  is a contraction mapping if  $\|L^{-1}U\| < 1$ . It is very difficult to show this for a given

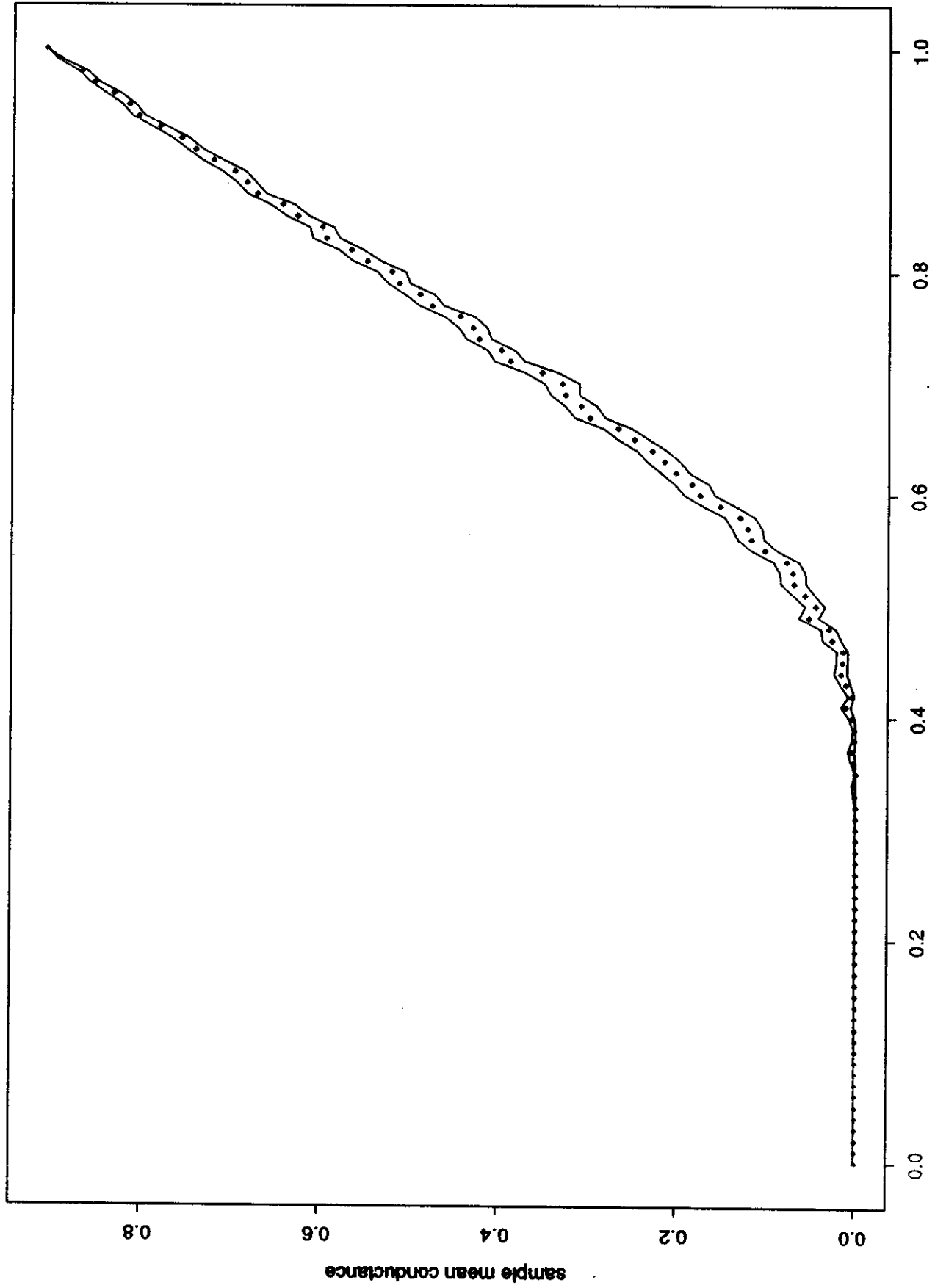
matrix  $A$ , but since the Kirchoff matrix  $A$  is symmetric positive definite, this holds. In addition, the necessary hypothesis that the lower diagonal matrix  $L$  of the Kirchoff matrix  $A$  be invertible is satisfied. Hence the Gauss-Seidel method should produce a reliable solution vector  $x$  of interior potentials.

Now that the interior potentials are known, the total current  $I_n^2$  can be calculated using Ohm's Law:

$$I_n^2 = \sum_{v \text{ a node adjacent to } G_{n0}^2} \text{Potential}(v) \text{Conductance}(\text{edge to left of } v). \quad (22)$$

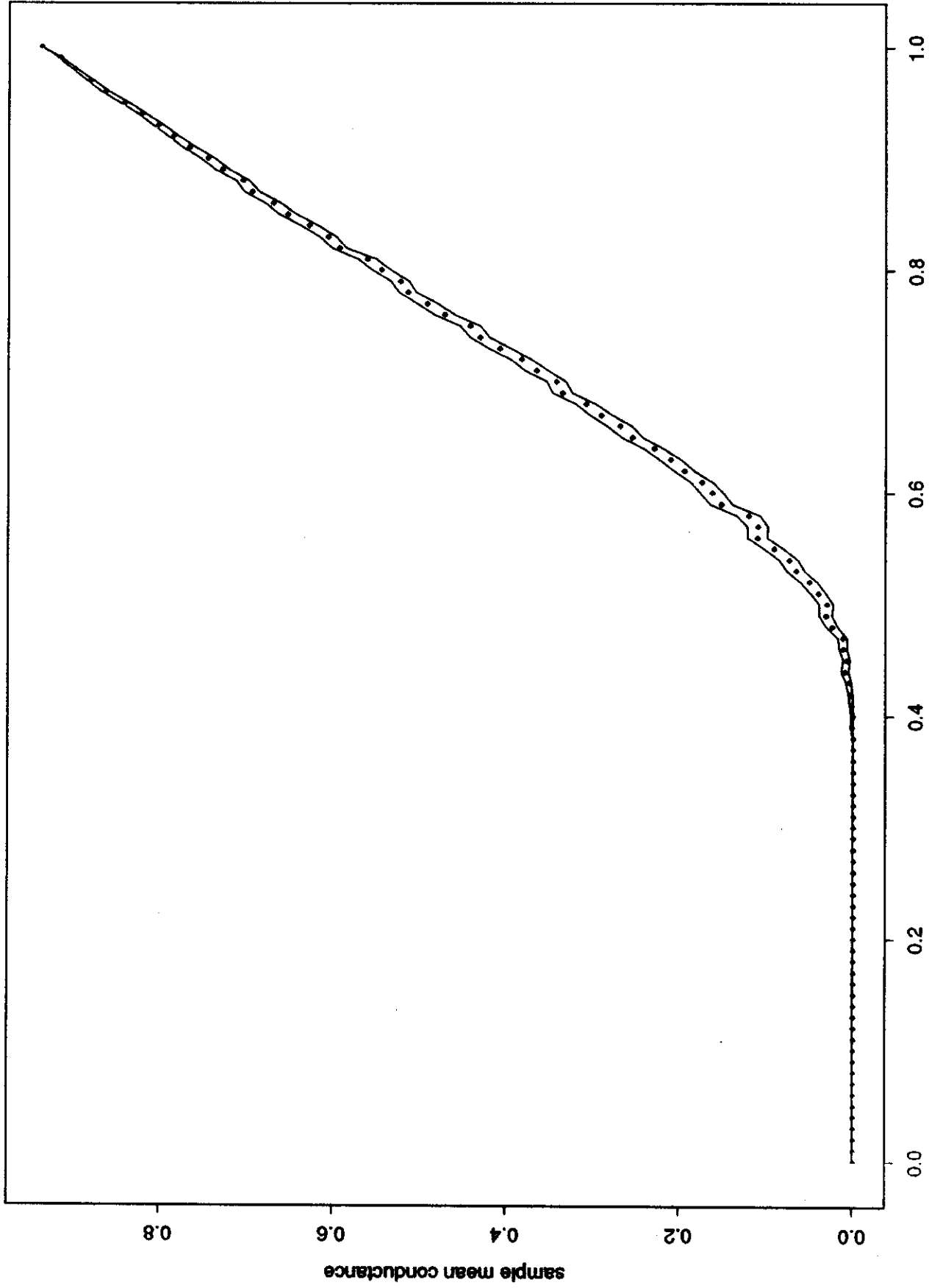
In addition to calculating sample mean conductances, the program SquareConductances.exe also computes sample standard deviations. For each  $p$  these sample standard deviations are used to calculate 95% confidence intervals about the true mean conductance. Figures 5, 6, and 7 are plots of  $\bar{C}_n^2(p)$  versus  $p$  for square networks with 10, 15, and 20 nodes on a side respectively. In the plots,  $d$  is the dimension of the network  $G_n^2$ ,  $n$  is the number of nodes on a side of the network and  $m$  is the sample size of computed conductances used to calculate the sample mean for each  $p$ . The value *test* is the small positive number which is used as a stopping criteria for the Gauss-Seidel iteration scheme- when  $\|x^{j+1} - x^j\| \leq \text{test} * \|x^j\|$  the iteration scheme halts. In general, if  $\text{test} = 10^{-z}$ , where  $z$  is a positive integer, then  $C_n^2$  is computed with an accuracy of about  $z$  significant digits. Finally, *seed* is a negative integer which is used to initiate the sequence of random numbers produced by the random number generator. For a given negative integer, the sequence of random numbers will always be the same. This is convenient because a given experiment can be exactly replicated. For each graph, 95% confidence intervals about the true mean conductances are included, which are calculated by the Splus function *forward*, which I described in section 4.

Figure 5 Sample Mean Conductance V Probability  $d=2, n=10, m=100, test=.0001, seed=-1$



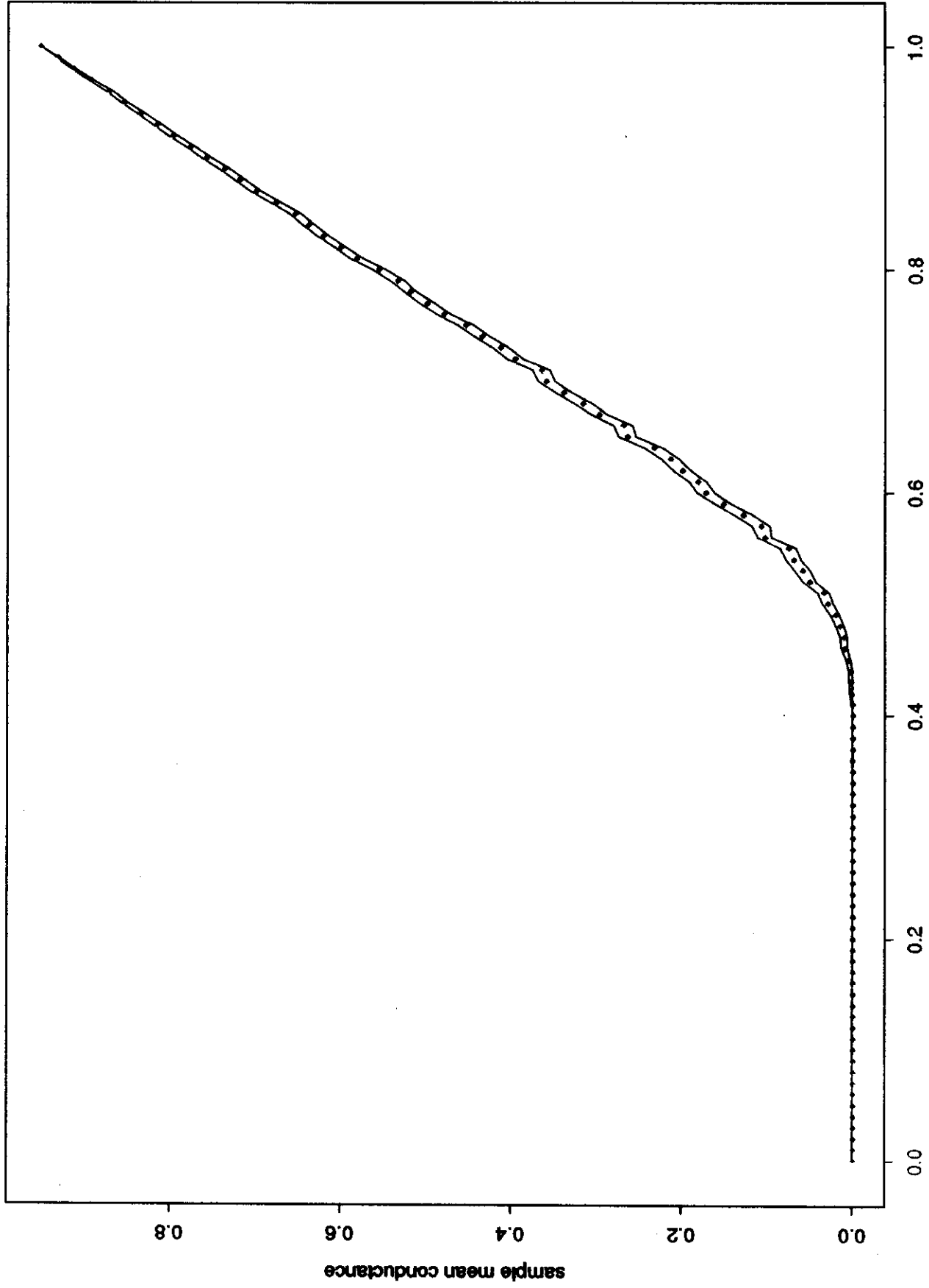
For any particular p-value, the lines give a 95% confidence interval for true mean conductance

Figure 6 Sample Mean Conductance V Probability  $d=2, n=15, m=100, test=.0001, seed=-1$



For any particular p-value, the lines give a 95% confidence interval for true mean conductance

Figure 7 Sample Mean Conductance V Probability  $d=2, n=20, m=100, test=.0001, seed=-1$



For any particular p-value, the lines give a 95% confidence interval for true mean conductance

## 8.2 Analysis of the Approximating Graphs of the Functions $h_n^2$

The most striking feature of the graphs in Figures 5, 6 and 7 is that the average conductances are all zero or very near zero until a region where suddenly nonzero average conductances appear. The critical probability  $\kappa^2$  must lie somewhere in this region. A comparison of Figures 5, 6 and 7 reveals that as  $n$  increases, the region of change from zero to nonzero average conductances narrows. This suggests that the program SquareConductances.exe can be used to approximate  $\kappa^2$  to an accuracy as great as desired by studying larger networks. The problem of course is its enormous running time (on the MSCC computers SquareConductances.exe took about 10,000 minutes to complete for a network with  $n = 20$ ).

Note that as  $n$  increases from 10 to 15 to 20, the region of change from zero to nonzero average conductances moves to the right. Nonzero average conductances begin to appear at about  $p = .4$  for the case  $n = 10$ , at about  $p = .44$  for the case  $n = 15$  and at about  $p = .47$  for the case  $n = 20$ . This suggests that as  $n$  gets large, nonzero average conductances will begin to appear at probabilities closer and closer to  $\frac{1}{2}$ . Hence this empirical approximation of the critical probability  $\kappa^2$  agrees with its theoretical value of  $\frac{1}{2}$ . Since  $\kappa^2$  is equal to the critical probability  $p_c^2$ ,  $p_c^2$  is estimated by this algorithm.

Another feature of the graphs in Figures 5, 6 and 7 is that they are approximately linear for values of  $p$  greater than the approximated critical probability. Comparing Figures 5 and 6 with Figure 7 shows that as  $n$  increases, the graph gets more linear. Moreover, in section 3 I showed that for given  $n$ ,  $h_n^2(1) = \frac{n}{n+1}$ , so that this linear portion has slope approximately equal to  $\frac{2n}{n+1}$ . This leads to my conjecture that as  $n \rightarrow \infty$ ,  $h_\infty^2$  is of the following form:

$$h_{\infty}^2 = \begin{cases} 0 & \text{if } p < \kappa^2 = \frac{1}{2} \\ 2p - 1 & \text{if } p \geq \frac{1}{2}. \end{cases}$$

The conjectured graph of the function  $h_{\infty}^2$  is depicted in Figure 8.

## 9 The Approximating Graphs of the Functions $h_n^3$

I wrote a Fortran program titled `CubicConductances.exe` which creates values of the function  $h_n^3$ . The program is exactly analogous to `SquareConductances.exe` except that sample mean conductances across the network  $G_n^3$  are calculated for the 31 probabilities  $p = 0.01, 0.04, 0.07, 0.10, 0.13, 0.16, \dots, 0.94, 0.97, 1.00$  instead of the 101 probabilities  $p = 0.00, 0.01, 0.02, 0.03, \dots, 0.98, 0.99, 1.00$  as in `SquareConductances.exe`. The reason for this is that the Gauss-Seidel algorithm used to solve for the interior potentials of a cubic network takes  $n$  times as long as the Gauss-Seidel algorithm used to solve for the interior potentials of a square network. Ohm's Law, Kirchoff's Law, and the Gauss-Seidel iteration scheme are essentially unchanged. The only difference is that each interior node of a three dimensional lattice network has six neighboring edges whereas each interior node of a two dimensional lattice network has four neighboring edges. Figures 9 and 10 depict plots of  $\bar{C}_n^3(p)$  versus  $p$  for three dimensional networks with 3 and 5 nodes on a side, respectively.



Figure 8 Conjectured Graph of the Two Dimensional Function  $h$  as  $n$  approaches infinity

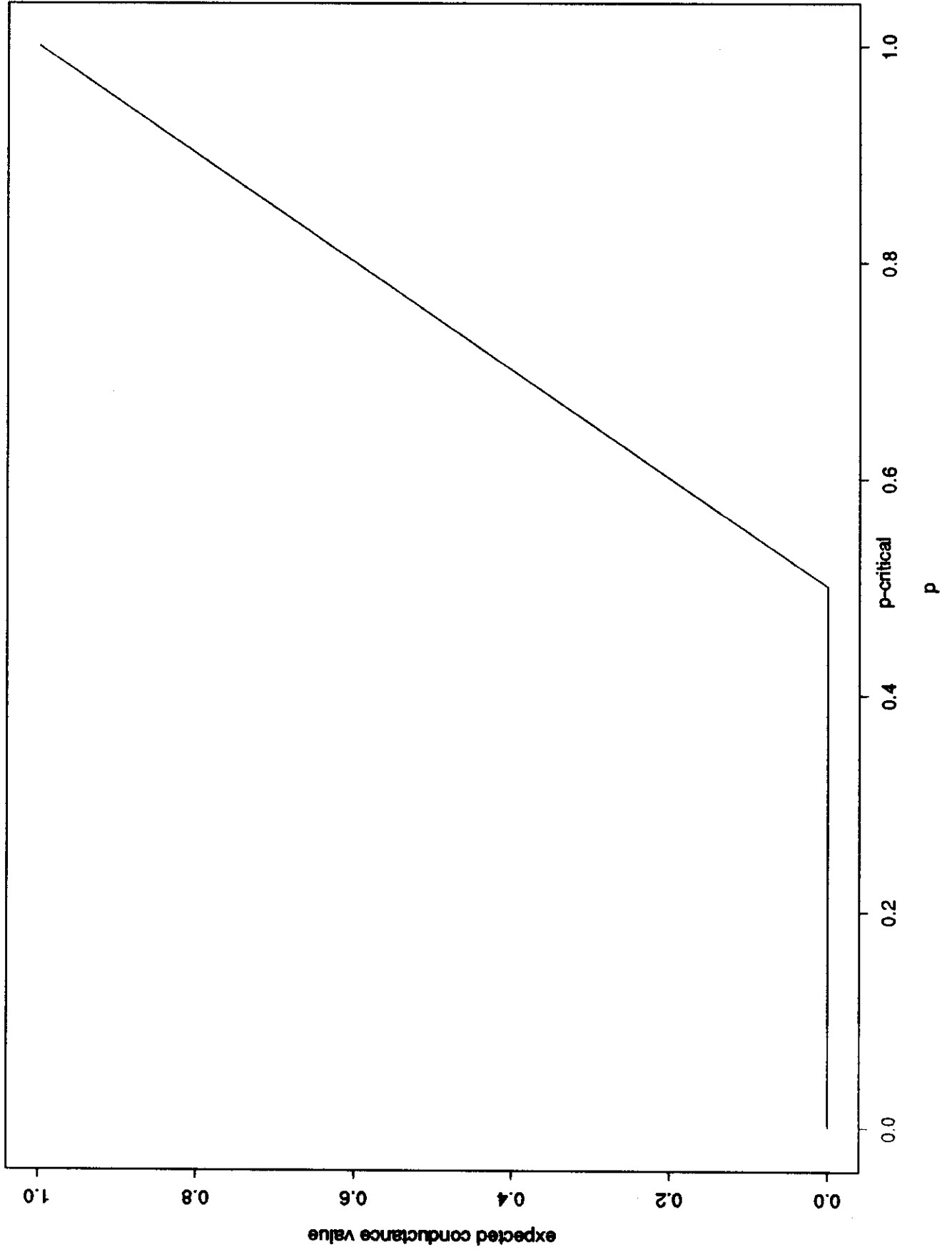


Figure 9 Sample Mean Conductance V Probability  $d=3, n=3, m=50, test=.0001, seed=-1$

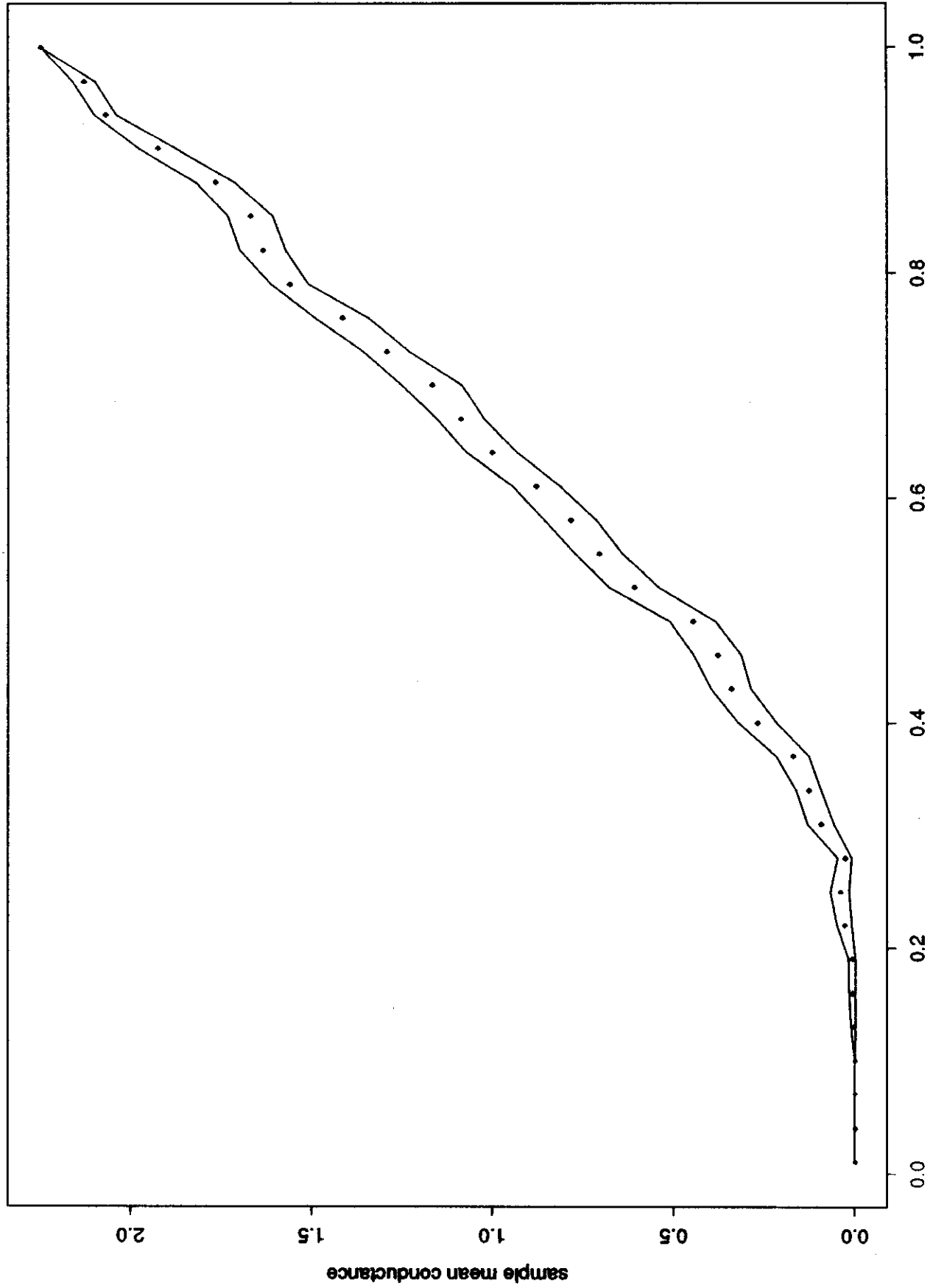
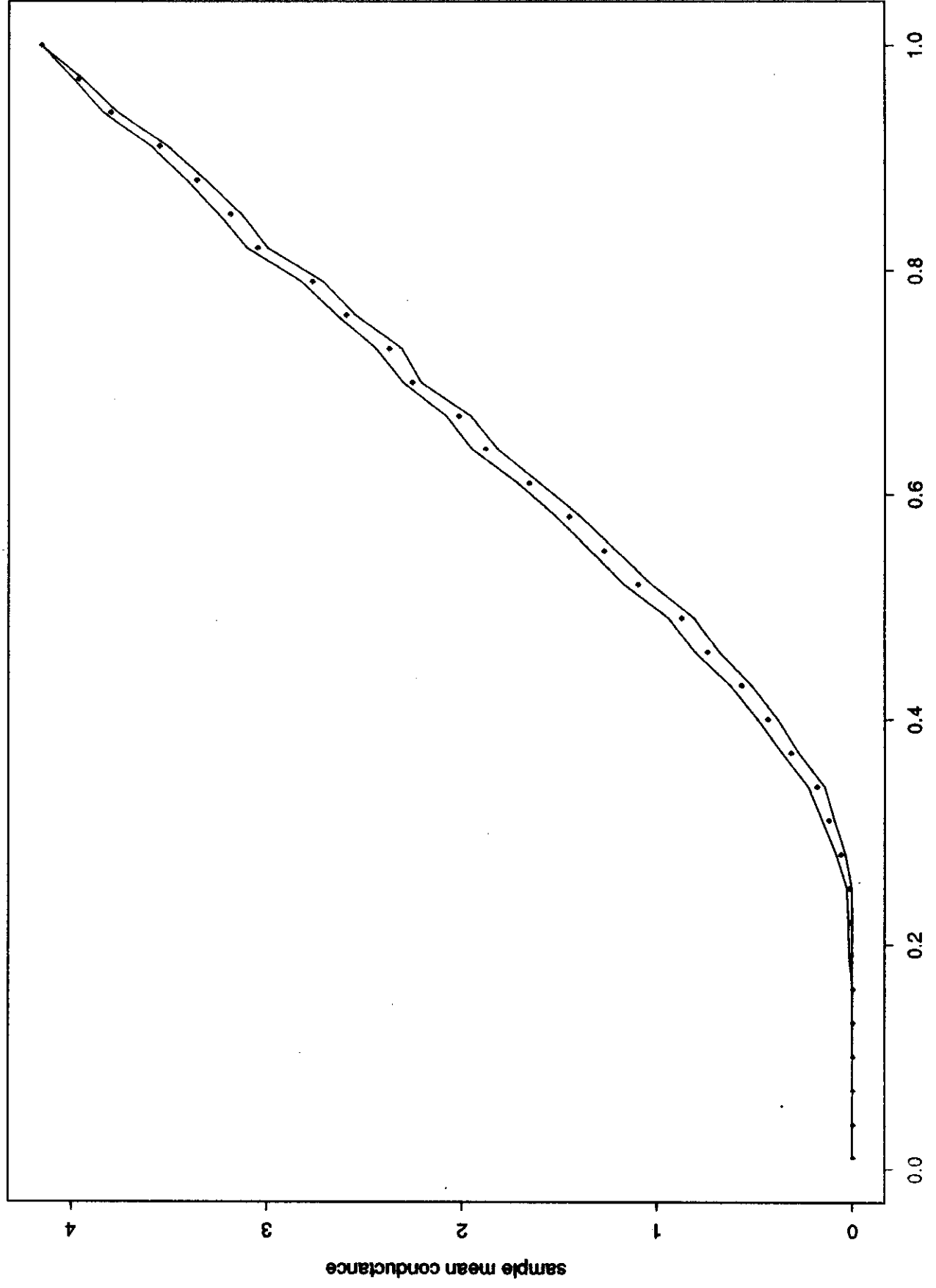


Figure 10 Sample Mean Conductance V Probability  $d=3, n=5, m=50, test=.0001, seed=-1$



For any particular p-value, the lines give a 95% confidence interval for true mean conductance

## 9.1 Analysis of the Approximating Graphs of the Functions $h_n^3$

Notice that in Figure 9 nonzero average conductances begin to appear approximately at about  $p = .20$ , and in Figure 10 nonzero average conductances begin to appear at about  $p = .25$ . I conjecture that just as in the two dimensional case, as  $n$  increases nonzero average conductances begin to appear at larger and larger probabilities. Grimmett (p13) proved that for  $d \geq 1$ ,  $p_c^{d+1} < p_c^d$ . Hence  $\kappa^{d+1} < \kappa^d$  for  $d \geq 1$ , and thus  $p_c^3 = \kappa^3 < p_c^2 = \frac{1}{2}$ . This gives the modest estimate that  $\kappa^3 \in (\frac{1}{4}, \frac{1}{2})$ . To obtain better estimates for the critical probability  $\kappa^3$ , the algorithm CubicConductances.exe needs to be optimized.

## 10 Improving the Algorithms

One simple way of quickening the programs SquareConductances.exe and CubicConductances.exe is to use a relaxation parameter in the Gauss-Seidel iteration scheme. The Gauss-Seidel iteration scheme could also be made faster by lowering the value of *test*. I used the value of .0001 for the algorithms. Lowering this value would drastically reduce the run time, but would sacrifice much accuracy. For example, on a three dimensional network with 5 nodes on a side, it is known that  $h_5^3(1) = \frac{25}{6} \doteq 4.167$ . When CubicConductances.exe is run on the network  $G_5^3$  and *test* = .0001, the sample mean computed for  $p = 1$  is 4.158, .10 away from the true mean 4.167. On the other hand, when *test* = .001 the sample mean computed for  $p = 1$  is 4.080, .17 away from the true mean 4.167. Another way to decrease computer run time would be to modify SquareConductances.exe and CubicConductances.exe to calculate sample means for a small range of probabilities. For example, if the goal is to approximate  $\kappa^2$ , running SquareConductances.exe for probabilities  $p = .47, .48, .49, .50, .51, .52, .53$  would be sufficient. Another way of speeding up the routines is to use smaller sample sizes to compute the sample means. The 95% confidence intervals

are fairly narrow in all of the graphs, so it appears the sample size could be substantially lowered before a substantial amount of accuracy is lost.

## 11 Conclusion

In this paper, I studied the behavior of electrical current flowing across disordered square and cubic resistor networks composed of random resistors. With the construction of forward and inverse problems, I empirically studied the map  $h : P \rightarrow E_p(C_n^d)$  and its inverse. Considering that the 95% confidence intervals in Figures 5, 6, 7 are fairly narrow, I conclude that the algorithm SquareConductances.exe accurately solves the forward and inverse problems in two dimensions. In addition, the graphs created by the algorithm SquareConductances.exe clearly validate that the critical probability  $\kappa^2$  exists. Figure 7 shows that with 95% confidence, the critical probability  $\kappa^2$  lies somewhere between .47 and .53 . Since I proved in section 6 that  $\kappa^2 = \frac{1}{2}$  , this is an accurate approximation.

Since I could not run the algorithm CubicConductances.exe on large networks due to computer time restraints, CubicConductances.exe only modestly solves the forward and inverse problems. As well, it only modestly estimates the unknown critical probability  $\kappa^3$  . At best, Figure 10 shows that with 95% confidence  $\kappa^3$  is greater than .25 . In conclusion, the algorithm SquareConductances.exe accurately solves the forward and inverse problems and estimates  $\kappa^2$  , but the algorithm CubicConductances.exe must be optimized and run on larger networks if it is to produce accurate solutions to the forward and inverse problems and accurately estimate  $\kappa^3$  .

## 12 References

1 Grimmett, Geoffrey. Percolation . New York: Springer-Verlag, 1989.

2 Kesten, Harry. Percolation Theory for Mathematicians . Boston: Birkhauser, 1982.

## 13 Appendices

```
*      PROGRAM SquareConductances.exe
*
*   This program deals with square resistor networks. Potentials of
*   zero are assigned to the boundary nodes on the left side of the square
*   network and potentials of one are assigned to the boundary nodes on
*   the right side of the square network. The boundary nodes on the
*   top and bottom sides of the network are insulated.
*   This program has a controlling loop over probabilities valued 0,.01
*   .02,.03 . . . .98,.99,1.00. For each probability p, a random
*   generator (the function RANDOM, in the file RANDOM.F) assigns each edge
*   of the network a conductance value MU with probability p and a conductance
*   value GAMMA with probability 1 - p. MU and GAMMA are prompted for, and can
*   be given nonnegative values. Then the program calculates the effective
*   conductance between the left and right sides of the square. This procedure
*   is repeated M times for each probability p, and the average of the
*   effective conductances is computed. As well, for each p the sample
*   standard deviation of the conductances is calculated. The probabilities,
*   the corresponding average conductances, and the sample standard deviations
*   are written to a file called 2Dexps.dat. This program utilizes the
*   Gauss-Seidel iteration scheme to calculate the interior potentials
*   of the network. A value TEST is prompted for; the smaller the value
*   of TEST the more accurate the interior potentials will be. In general
*   if TEST is of order  $10^{-z}$ , where z is a positive integer, then the
*   effective conductance computed will have z significant digits.
*   The Gauss-Seidel iteration scheme is faster than the alternative
```

```

* method of calculating interior potentials, via setting up the
* matrix equation  $Ax=b$  and using Lipackd routines. However, the
* Gauss-Seidel scheme used in this program does not include a relaxation
* parameter, which would speed the program up much more.

*
* INDEXING
*
* The indexing that I used was the following: the leftmost boundary
* node on on the top face is the first node, and from there proceed from
* left to right and then top to bottom. Hence for a network with N nodes
* on a side there are  $4N*N**2$  total nodes in the network, and the last node
* is the rightmost node on the bottom face.
*
* The interior nodes have there own indexing: the leftmost interior node
* adjacent to the top side of the network is the first node, and from
* there proceed from left to right then top to bottom.
*
* Edges are indexed in the following manner: Horizontal edges are
* numbered from top to bottom, left to right, and vertical edges are
* numbered from left to right, top to bottom.

*
* VARIABLES
* N      The number of nodes on a side
* M      The number of iterations of the random generator
* NUMEGS The number of edges of the network
* LIMIT  The maximum number of edges for a network
* MAXDIM The maximum number of nodes on a side
* HORCDS Stores the conductances of the horizontal edges
*        moving left to right then top to bottom
* VERCDS Stores the conductances of the vertical edges
*        moving top to bottom then left to right
* CURNTS Stores the currents of the edges adjacent to West face
*        boundary nodes
* A      The solution vector of interior node potentials
* ALAST  Used to tell when the solution vector A has converged to

```

\*           within TEST of the true potentials  
 \* IXINTS An array containing the indices of the network corresponding  
 \*           to the interior nodes  
 \* INVIKS The inverse of IXINTS- an array containing the indices of the  
 \*           interior nodes corresponding to the network indices  
 \* IPVT    Used for the DGEFA and DGESL routines  
 \* ( )BND S Stores the indices of the (N)orth, (E)ast  
 \*           (S)outh, and (W)est boundary nodes  
 \* LIXHOR LIXHOR and RIXHOR store the indices of the nodes straddling  
 \* RIXHOR the horizontal edges  
 \* TIXVER TIXVER and BIXVER store the indices of the nodes straddling  
 \* BIXVER the vertical edges  
 \* RE       The resistance from the left to right face  
 \* CT       The total current flowing from left to right  
 \* CDTS     An array of conductances from the left to right face. Each  
 \*           element holds one conductance corresponding to one set of  
 \*           edge conductances  
 \* STDEVS Vector of length 101, which will be filled with the sample  
 \*           standard deviations  
 \* P        Vector of probabilities  
 \* ECD      The answer, which is the average of the elements of CD  
 \* EXPS     Store all the expectations  
 \* MU       One of the possible values for the conductance of an edge  
 \* GAMMA    The other possible conductance value  
 \* SEED     The value used as a parameter to RANDOM  
 \* TEST     The small value which determines how many Gauss-Seidel iterations  
 \*           take place.  
 \* ANORM    The infinity norm of the vector of interior potentials, ALAST  
 \* UNORM    The infinity norm of the difference between the most recently  
 \*           calculated vector of interior potentials A and the previously  
 \*           calculated vector of interior potentials ALAST

```

    INTEGER N, M, NUMEGS, LIMIT, MAXDIM, SEED,S,
    $ NBHORS(1:4),IXEGS(1:4)
    PARAMETER (LIMIT = 420)
    PARAMETER (MAXDIM=20)
    REAL A(MAXDIM**2),ALAST(MAXDIM**2),CDS(1:4),ECD,CT,
  
```



```

$ P(101),MU,GAMMA, EXPS(101), TEST, STDEVS(1:101),CDTS(1:101)
REAL HORCDS(1:LIMIT), VERCDS(1:LIMIT), ANORM,UNORM
INTEGER IXINTS (1:LIMIT), INVIXS(1:LIMIT)
INTEGER NBNDS(1:MAXDIM),EBNDS(1:MAXDIM)
INTEGER SBNDS(1:MAXDIM),WBNDS(1:MAXDIM)
INTEGER LIXHOR(1:LIMIT),RIXHOR(1:LIMIT)
INTEGER TIXVER(1:LIMIT),BIXVER(1:LIMIT)

*
* SUBROUTINES CALLED
*
*
* SUBROUTINES USED FOR INDEXING
*
* INDINTNODES    Fills IXINTS with the indices of the interior nodes
* INVINDINTNODES Fills INVIXS with the indices of the network
*                corresponding to the indices of the interior nodes
* INDBNDNODES    Fills NBND, EBND, WBND, and SBND with the indices
*                of the network.  Boundary nodes are ordered starting
*                in the leftmost position in the South face moving
*                counterclockwise about the network.
* HORMAP         Fills LIXHOR and RIXHOR
* VERMAP         Fills TIXVER and BIXVER
*
* OTHER SUBROUTINES
*
* FILLCDS        Gets the neighboring conductances for any interior node
* RANDOM         A random number generator, called from external.  The code
*                was found in NUMERICAL RECIPES.  A value SEED is required
*                to start RANDOM, and the same sequence of random numbers
*                is always produced for a given SEED.
*
*****
*** BEGIN
*
      INTRINSIC SQRT,ABS,REAL,MOD
      PRINT *, 'Enter the size of the square resistor (1-20)'
      READ *, N

```

```

NUMEGS = 2*N*(N+1)

PRINT *
PRINT *, 'Enter the number of conductances to compute'
READ *, M

PRINT *
PRINT *, 'Enter MU(>=0), the conductance value with probability p'
READ *, MU

PRINT *
PRINT *, 'Enter GAMMA(>=0), the conductance value with prob. 1-p'
READ *, GAMMA

PRINT *
PRINT *, 'Enter TEST, the approximation level for Gauss-Seidel'
READ *, TEST

DO 100 I = 0,100
  P(I+1) = REAL(I)/REAL(100)
100 CONTINUE
* Fill P with probabilities .00 to 1.00

CALL INDINTNODES (IXINTS, N)
* Fill IXINTS with the indices of the interior nodes

CALL INVINDINTNODES (INVIXS, N)

CALL INDBNDNODES (NBNDS, EBNDS, SBNDS, WBNDS, NMAXDIM)
* Fill NBNDS, EBNDS, SBNDS, and WBNDS with the indices of the boundary nodes

CALL HORMAP (LIXHOR, RIXHOR, N)
* Fill LIXHOR and RIXHOR with the indices of nodes straddling horizontal edges

CALL VERMAP (TIXVER, BIXVER, N)
* Fill TIXVER and BIXVER with the indices of nodes straddling vertical edges

```

```

* We now have the indices of the boundary and interior nodes

      DO 110 I = 1,101
        STDEVS(I) = 0
110   CONTINUE

* Initialize STDEVS to zero

*
*
**** CONTROLLING DOUBLE LOOP

      PRINT *
      PRINT *, 'Please enter a negative integer for a SEED'
      READ *, SEED
* Get a value for SEED to use in the random number generating function
* RANDOM

      DO 120 PP = 1,101
        ECD = 0
* Initialize the expected conductance to zero

      DO 130 J = 1, M

* Use the random number generator, found in NUMERICAL RECIPES,
* to determine the conductance of each edge.
      DO 6000 I = 1, NUMEGS/2
        IF (RANDOM(SEED) .LE. P(PP)) THEN
          HORCDS(I) = MU
        ELSE
          HORCDS(I) = GAMMA
        ENDIF
        IF (RANDOM(SEED) .LE. P(PP)) THEN
          VERCDS(I) = MU
        ELSE
          VERCDS(I) = GAMMA
        ENDIF

```

```

6000  CONTINUE

      DO 140 I = 1,N**2
        A(I) = 0
140    CONTINUE
* Initialize A with zeros

150    DO 160 I = 1,N**2
        ALAST(I) = A(I)
160    CONTINUE

      DO 170 S = 1, N**2

        CALL FILLCDS (CDS,HORCDS,VERCDS,IXINTS,LIXHOR,RIXHOR,TIXVER,
          $ BIXVER,NBHORS,IXEGS,N,NUMEGS, S, MAXDIM)
* Get the neighboring conductances surrounding the Sth interior node.
* CDS(1) is the left neighbor's conductance, CDS(2) is the right
* neighbor's conductance, CDS(3) is the above neighbor's conductance,
* and CDS(4) is the bottom neighbor's conductance.

* Take care of top, left corner
      IF (S .EQ. 1) THEN
        IF (CDS(1)+CDS(2)+CDS(4) .GT. .5) THEN
          A(S) = (A(S+1)*CDS(2)+A(S+N)*CDS(4))/
          $ (CDS(1)+CDS(2)+CDS(4))
        ENDIF

* Take care of top, right corner
      ELSEIF (S .EQ. N) THEN
        IF (CDS(1)+CDS(2)+CDS(4) .GT. .5) THEN
          A(S) = (A(S-1)*CDS(1)+CDS(2)+A(S+N)*CDS(4))/
          $ (CDS(1)+CDS(2)+CDS(4))
        ENDIF

* Take care of bottom, left corner
      ELSEIF (S .EQ. N**2 - N + 1) THEN
        IF (CDS(1)+CDS(2)+CDS(3) .GT. .5) THEN

```

```

          A(S) = (A(S+1)*CDS(2)+A(S-N)*CDS(3))/
$          (CDS(1)+CDS(2)+CDS(3))
      ENDIF

* Take care of bottom, right corner
      ELSEIF (S .EQ. N**2) THEN
          IF (CDS(1)+CDS(2)+CDS(3) .GT. .5) THEN
              A(S) = (A(S-1)*CDS(1)+CDS(2)+A(S-N)*CDS(3))/
$              (CDS(1)+CDS(2)+CDS(3))
          ENDIF

* Take care of left edge, excluding corners
      ELSEIF ((MOD(S-1,N) .EQ. 0) .AND.
$          (S .NE. 1) .AND. (S .NE. N**2 - N + 1)) THEN
          IF (CDS(1)+CDS(2)+CDS(3)+CDS(4) .GT. .5) THEN
              A(S) = (A(S+1)*CDS(2)+A(S-N)*CDS(3)+A(S+N)*CDS(4))/
$              (CDS(1)+CDS(2)+CDS(3)+CDS(4))
          ENDIF

* Take care of right edge, excluding corners
      ELSEIF ((MOD(S,N) .EQ. 0) .AND.
$          (S .NE. N) .AND. (S .NE. N**2)) THEN
          IF (CDS(1)+CDS(2)+CDS(3)+CDS(4) .GT. .5) THEN
              A(S) = (A(S-1)*CDS(1)+CDS(2)+A(S-N)*CDS(3)+A(S+N)*
$ CDS(4))/(CDS(1)+CDS(2)+CDS(3)+CDS(4))
          ENDIF

* Take care of top edge, excluding corners
      ELSEIF ((S .GT. 1) .AND. (S .LT. N)) THEN
          IF (CDS(1)+CDS(2)+CDS(4) .GT. .5) THEN
              A(S) = (A(S-1)*CDS(1)+A(S+1)*CDS(2)+A(S+N)*CDS(4))/
$              (CDS(1)+CDS(2)+CDS(4))
          ENDIF

* Take care of bottom edge, excluding corners
      ELSEIF ((S .GT. N**2-N+1) .AND. (S .LT. N**2)) THEN
          IF (CDS(1)+CDS(2)+CDS(3) .GT. .5) THEN

```

```

          A(S) = (A(S-1)*CDS(1)+A(S+1)*CDS(2)+A(S-N)*CDS(3))/
$          (CDS(1)+CDS(2)+CDS(3))
      ENDIF
    ELSE
      IF (CDS(1)+CDS(2)+CDS(3)+CDS(4) .GT. .5) THEN
        A(S) = (A(S-1)*CDS(1)+A(S+1)*CDS(2)+A(S-N)*CDS(3)+
$        A(S+N)*CDS(4))/(CDS(1)+CDS(2)+CDS(3)+CDS(4))
      ENDIF

    ENDIF

170  CONTINUE

      ANORM = 0
      UNORM = 0
      DO 500 I = 1,N**2
        IF (ABS(A(I)-ALAST(I)) .GT. UNORM) THEN
          UNORM = ABS(A(I)-ALAST(I))
        ENDIF
      * Fill UNORM with the maximum component value of ABS(A(I)-ALAST(I))

      IF (ABS(ALAST(I)) .GT. ANORM) THEN
        ANORM = ABS(ALAST(I))
      ENDIF
      * Fill ANORM with the maximum component value of ABS(ALAST(I))

500  CONTINUE

      IF (UNORM .GT. TEST*ANORM) GOTO 150
      * Test if another iteration is necessary

      *
      * Now calculate the current across the network CT

      CT = 0
      * Initialize the value of the CT to zero.

```

```
DO 180 K = 1, N
  CT = CT + HORCDS(1 + (N+1)*(K-1))*A(1 + N*(K-1))
180 CONTINUE
```

```
  CDTS(J) = CT
  ECD = ECD + CT
130 CONTINUE
```

```
  ECD = ECD/M
  DO 190 I = 1, M
    STDEVS(PP) = STDEVS(PP) + (CDTS(I)-ECD)**2
190 CONTINUE
```

```
  STDEVS(PP) = STDEVS(PP)/(M-1)
  STDEVS(PP) = SQRT(STDEVS(PP))
```

```
  EXPS(PP) = ECD
120 CONTINUE
```

```
  CALL OUTPUT(N,P,EXPS,M,MU,GAMMA,STDEVS)
```

```
*****
```

```
  END
```

```
* PROGRAM FORWARDGS.F completed
```

```
  SUBROUTINE FILLCDS(CDS,HORCDS,VERCDS,IXINTS,LIXHOR,RIXHOR,TIXVER,
$ BIXVER,NBHORS,IXEGS,N,NUMEGS, Q, MAXDIM)
```

```
  INTEGER NBHORS(1:4), IXEGS(1:4), MAXDIM,
$ N, Q, NUMEGS, IPTQ
  REAL HORCDS(NUMEGS), VERCDS(NUMEGS),CDS(1:4)
  INTEGER IXINTS(NUMEGS), LIXHOR(NUMEGS), RIXHOR(NUMEGS)
  INTEGER TIXVER(NUMEGS), BIXVER(NUMEGS),NBHORS(1:4)
```

```

      IPTQ = IXINTS (Q)
* IPTQ contains the node index of the Qth interior node.
* The IXINTS (Q)th node has neighboring nodes with indices
* IPTQ - 1 to the left, IPTQ + 1 to the right,
* IPTQ - (N+2) above, and IPTQ + (N+2) below.
* Now the subroutine builds the array NBHORS, containing the indice
* numbers of the nodes surrounding the Qth interior node.

* The first two positions in the array are occupied by the
* indices of the nodes forming horizontal edges, and the second
* two positions are occupied by the indices of the nodes forming
* vertical edges.

```

```

      NBHORS (1) = IPTQ - 1
      NBHORS (2) = IPTQ + 1
      IF (IPTQ .LE. 2*N + 1) THEN
        NBHORS (3) = IPTQ - N - 1
      ELSE
        NBHORS (3) = IPTQ - N - 2
      ENDIF

      IF (IPTQ .GE. N**2 + 2*N - 1) THEN
        NBHORS (4) = IPTQ + N + 1
      ELSE
        NBHORS (4) = IPTQ + N + 2
      ENDIF

```

```

* Now the subroutine will fill the CDS array with the conductances.
* CDS(1) contains the conductance of the left edge, CDS(2) the
* right, CDS(3) the above, and CDS(4) the below.

* This loop locates the indices of the edges which are straddled
* by two nodes with known indices.

```

```

      DO 200 J = 1, NUMEGS
        IF (LIXHOR(J) .EQ. NBHORS(1)) THEN
          IF (RIXHOR(J) .EQ. IPTQ) THEN

```



```

        IXEGS(1) = J
        IXEGS(2) = J + 1
    ENDIF
ENDIF

    IF (TIXVER(J) .EQ. NBHORS(3)) THEN
        IF (BIXVER(J) .EQ. IPTQ) THEN
            IXEGS(3) = J
            IXEGS(4) = J + 1
        ENDIF
    ENDIF
ENDIF

```

```

200    CONTINUE

```

```

        CDS(1) = HORCDS(IXEGS(1))
        CDS(2) = HORCDS(IXEGS(2))
        CDS(3) = VERCDS(IXEGS(3))
        CDS(4) = VERCDS(IXEGS(4))
    END

```

```

*

```

```

*     SUBROUTINES FOR INDEXING

```

```

*

```

```

*

```

```

* HORMAP This subroutine fills the inputed array B1 with the node indices
* which are to the left of each edge, i.e. the ith element of B1 contains
* the index of the left node which is straddled by the ith horizontal
* edge. Similarly the array C1 is filled with the node indices which are
* to the right of each edge, i.e. the ith element of C1 contains the index
* of the right node which is straddled by the ith horizontal edge.

```

```

SUBROUTINE HORMAP (B1, C1, N)

```

```

    INTEGER COUNTB, COUNTC, N
    INTEGER B1(2*N*(N+1)), C1(2*N*(N+1))
    COUNTB = -1
    COUNTC = -2

```

```

DO 210 I = 1, N
    COUNTB = COUNTB + 1
    COUNTC = COUNTC + 2
    DO 220 J = 1, N+1

        B1(N*(I-1) + COUNTB + J) = I*N + J + COUNTC
        C1(N*(I-1) + COUNTB + J) = I*N + J + COUNTC + 1

220     CONTINUE
210     CONTINUE

```

END

\*  
\* VERMAP This function is similar to HORMAP, it fills B1 and C1 with the  
\* indices of nodes which straddle vertical edges to the left and right  
\* respectively.

SUBROUTINE VERMAP (B1, C1, N)

INTEGER N, B1(2\*N\*(N+1)),C1(2\*N\*(N+1))

```

DO 230 I = 1, N
    DO 240 J = 1, N+1

```

```

        IF (J .EQ. 1) THEN
            B1((N+1)*(I-1) + J) = I
        ELSE IF (J .EQ. N+1) THEN
            B1((N+1)*(I-1) + J) = 2*N + N**2 - 1 + I
        ELSE IF (J .EQ. 2) THEN
            B1((N+1)*(I-1) + J) = I + N + 1
        ELSE
            B1((N+1)*(I-1) + J) = B1((N+1)*(I-1)
$           + J - 1) + N + 2
        END IF

```

```

            IF ((B1((N+1)*(I-1)+J) .GT. N**2 + 2*N - 1) .OR.

```

```

$           (B1((N+1)*(I-1)+J) .LE. N)) THEN
              C1((N+1)*(I-1) + J) = B1((N+1)*(I-1)+J)
$           + N + 1
              ELSE
              C1((N+1)*(I-1) + J) = N+2 + B1((N+1)*
$           (I-1) + J)
              ENDIF
240          CONTINUE
230          CONTINUE

          END

```

\* INDINTNODES takes an array of integers and an integer N as paramaters  
 \* The subroutine inputs the index of the ith interior node in the ith position  
 \* of the array.

```

          SUBROUTINE INDINTNODES (ARRAY, N)

          INTEGER N, ARRAY(N**2)
          DO 250 I = 1, N
              DO 260 J = 1, N

                  ARRAY (N*(I-1) + J) = I*N + 2*I + J - 1
260          CONTINUE
250          CONTINUE
          END

```

\*  
 \* INVINDINTNODES This subroutine fills the ith position of the  
 \* inputted ARRAY with the index of the node corresponding to the  
 \* ith interior node. It is the inverse routine to INDINTNODES.

```

          SUBROUTINE INVINDINTNODES(ARRAY, N)

          INTEGER N, ARRAY(N**2 + 4*N)

          DO 270 I = 1, N

```

```

        DO 280 J = 1,N
            ARRAY(I*N + 2*I + J - 1) = N*(I-1)+J
280     CONTINUE
270     CONTINUE
        END

```

\* INDBNDNODES takes four arrays of integers and an integer N as paramaters.  
 \* The four arrays correspond to the North, East, South, and West boundary node  
 \* The subroutine inputs the index of the ith boundary node in the ith position  
 \* of the appropriate array.

```

        SUBROUTINE INDBNDNODES (AN,AE,AS,AW,N,MAXDIM)

        INTEGER N, MAXDIM
        INTEGER AN(MAXDIM),AE(MAXDIM),AS(MAXDIM),AW(MAXDIM)

        DO 290 I = 1, N
            AN (N + 1 - I) = I
            AE (N + 1 - I) = N + I*N + 2*I
            AS (I) = N**2 + 3*N + I
            AW (I) = I*N + 2*I - 1
290     CONTINUE
        END

```

```

        SUBROUTINE OUTPUT(N,P,EXPS,M,MU,GAMMA,STDEVS)
        INTEGER N, M
        REAL EXPS(101),P(101),STDEVS(101),MU,GAMMA
        PRINT *
        PRINT *, 'The expectations of the conductances'
        PRINT *, ' will be written to the file '
        PRINT *, '2Dexps.dat.'
        PRINT *, 'in the following format:'
        PRINT *, ' the first row has the value of N'
        PRINT *, ' the second row has the value of M'
        PRINT *, ' the third row has the value of MU'
        PRINT *, ' the fourth row has the value of GAMMA'
        PRINT *, ' the next 101 rows contain the probabilities'

```

```

PRINT *,' 0.00-1.00'
PRINT *,' the next 101 rows contain the corresponding'
PRINT *,' average values of the computed conductances'
PRINT *,' the final 101 rows contain the standard deviations'
PRINT *,' for the calculated conductances'
PRINT *,' There is one standard deviation for each probability'
PRINT *

OPEN (1,FILE='2Dexps.dat',ERR=999)
WRITE(1,*) N
WRITE(1,*) M
WRITE(1,'(D16.8)') MU
WRITE(1,'(D16.8)') GAMMA
DO 300 I=1, 101
  WRITE (1,'(D16.8)') P(I)
300  CONTINUE
DO 310 I = 1, 101
  WRITE (1, '(D16.8)') EXPS(I)
310  CONTINUE
DO 320 I = 1, 101
  WRITE (1, '(D16.8)') STDEVS(I)
320  CONTINUE
RETURN
999  STOP 'Error in writing to 2Dexps.dat'
END

```

```

inverse _ function(x,y,smean,m,sigma,d) {

# x is the vector of probabilities
# y is the vector of sample mean conductances
# smean is the sample mean calculated by making m measurements of
# effective conductances and averaging them
# m is the sample size
# sigma is the sample standard deviation
# d is the dimension of the network

# Assuming that the data is normally distributed,
# this function returns a 95% confidence interval about the proportion p
# of materials, one having conductance 1 and the other having conductance 0.

# Assign lowerlimit and upperlimit the lower and upper bounds of the 95%
# confidence interval about the true mean conductance.

lowerlimit _ smean - (1.96*sigma)/sqrt(m)
upperlimit _ smean + (1.96*sigma)/sqrt(m)

l _ length(y)
if (d == 2) k _ 1001
else k _ 331
hors _ rep(0,k)
vers _ rep(0,k)

# Fill the vector hors with the probabilities .000,.001,.002,...,.998,
# .999,1.000 if d=2, and .01,.013,.016,.019,.022,...,.997,1.000 if d=3

if (d == 2) hors _ c(0:1000)/1000
else hors _ seq(.01:1,by=.003)

# Fill the vector vers with interpolated sample mean conductances; i.e.
# between any two values of y insert 9 values equidistant from each other
# Hence vers contains the values of y with added values

```

```

for (i in 1:(l-1)) {
for (j in 1:10) {
vers[j+(i-1)*10] _ y[i+1] - (j/10)*(y[i+1] - y[i])
}
}
vers[(l-1)*10+1] _ y[l]

# Now find the index upperindex which indicates the element of vers closest
# to upperlimit and find the index lowerindex which indicates the element
# of vers closest to lowerlimit

upperdists _ rep(0,k)
lowerdists _ rep(0,k)
upperminimum _ 10
lowerminimum _ 10

upperindex _ 0
lowerindex _ 0

for (i in 1:k) {

upperdists[i] _ abs(upperlimit-vers[i])
lowerdists[i] _ abs(lowerlimit-vers[i])

if (upperdists[i] < upperminimum) {
upperminimum _ upperdists[i]
upperindex _ i}

if (lowerdists[i] < lowerminimum) {
lowerminimum _ lowerdists[i]
lowerindex _ i}

}

# Now hors[lowerindex] is the lower confidence limit for the true probability

```

```
# and hors[upperindex] is the upper confidence limit for the true probability
return(hors[lowerindex],hors[upperindex])
}
```