

FYS4460:
UNORDERED SYSTEMS AND PERCOLATION
PERCOLATION TOPICS

ALEXANDER HAROLD SEXTON

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1. TOPIC 9: ALGORITHMS FOR PERCOLATION SYSTEMS

How do we generate a percolation system for simulations? How to analyze and visualize the systems? How to find spanning clusters and measure the percolation probability?

- Percolation systems are generated by defining a 2-d grid of random numbers.

- Define sites to be set if site is below some value p

```
1 L = 4
2 p = 0.5
3 z = np.random.random((L, L))
4 system = z < p
```

- Visualize system using `plt.imshow()` from `matplotlib`

- Example of this in figure 1.1.

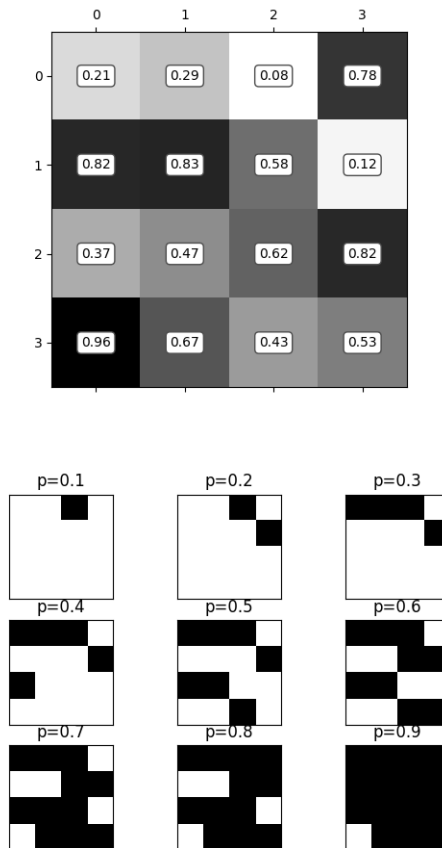


FIGURE 1.1. Percolation systems with varying p .

- Characterize clusters by which sites are connected
- Next-neighbour connectivity
- Percolating systems
- Use python libraries for analyzing imagery

```

1 import scipy.ndimage as sp
2 from skimage import measure
3 labels , n_features = sp.measurements.label(system)

```

- Returns a matrix of connected regions which are labeled
- Can be visualized easily as in figure 1.2

```

1 L = 10
2 p = 0.5
3 z = rand(L,L)
4 system = z<p
5 labels, n_features = measurements.label(system)
6 clusterArea = measurements.sum(system, labels, index=arange(labels.max() + 1))
7 areaImg = clusterArea[labels]
8 cmap = plt.cm.viridis
9 cmap.set_under('black')
10 imshow(areaImg, origin='lower', cmap=cmap, vmin=0.1, interpolation='none')
11 colorbar()
12 show()

```

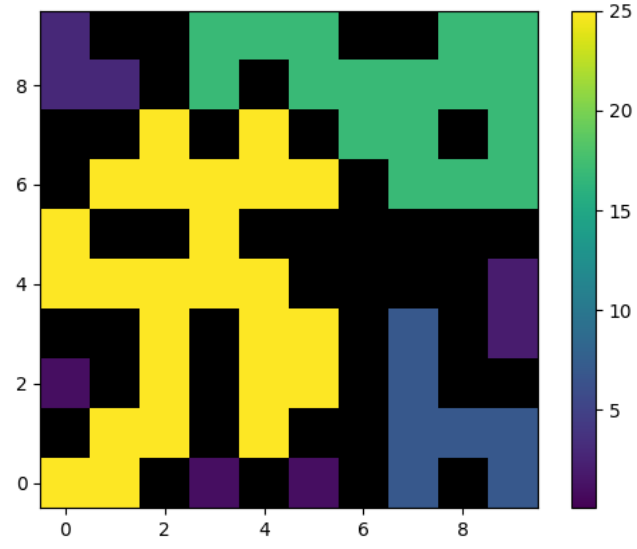


FIGURE 1.2. Visualization of the clusters in the system

- System of size $L = 100$ for p approaching p_c in figure 1.3

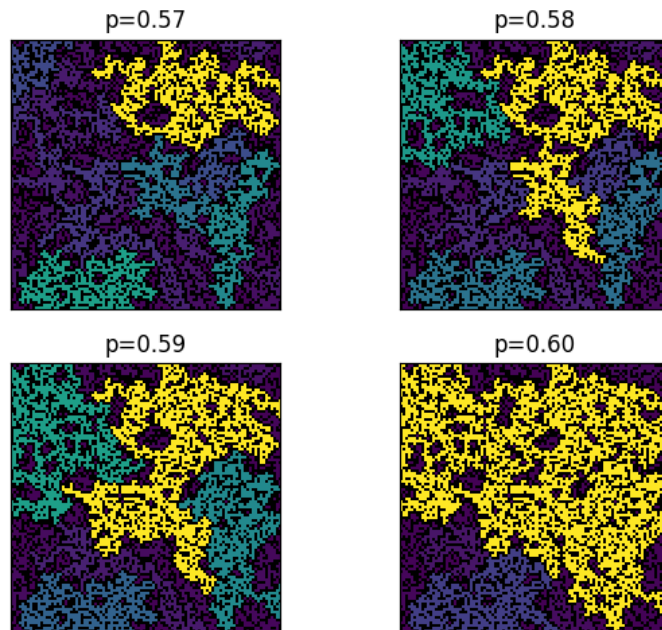


FIGURE 1.3. 100×100 systems with p approaching p_c .

- Use the bounding box of clusters to determine if system is percolating

```

1 def spanning_cluster_density(m, single_perc = False):
2     nx, ny = m.shape
3     labels, n_features = sp.measurements.label(m)
4     regions = measure.regionprops(labels)
5     density = 0
6     for region in regions:
7
8         x_start, y_start, x_stop, y_stop = region.bbox
9         dx = x_stop - x_start
10        dy = y_stop - y_start
11
12        if dx == nx or dy == ny:
13            density += region.extent
14            if single_perc:
15                break
16
17    return density

```

- Measure percolation probability by Monte Carlo simulations

```

1 fig, (ax1, ax2) = plt.subplots(2)
2 for L in L_arr:
3     Pi_arr = np.zeros(len(p_arr))
4     density_arr = np.zeros(len(p_arr))
5     for p, prob in enumerate(p_arr):

```

```

6     for i in range(n_samples):
7         z = np.random.random((L,L))
8         m = z < prob
9         density = spanning_cluster_density(m)
10        density_arr[p] += density
11        if density > 0:
12            Pi_arr[p] += 1
13
14    Pi_arr /= n_samples
15    density_arr /= n_samples
16    ax1.plot(p_arr, Pi_arr, label=f'L={L}')
17    ax2.plot(p_arr, density_arr, label=f'L={L}')

```

- Bonus: this code also computes the density of the spanning cluster. Example in figure 1.4

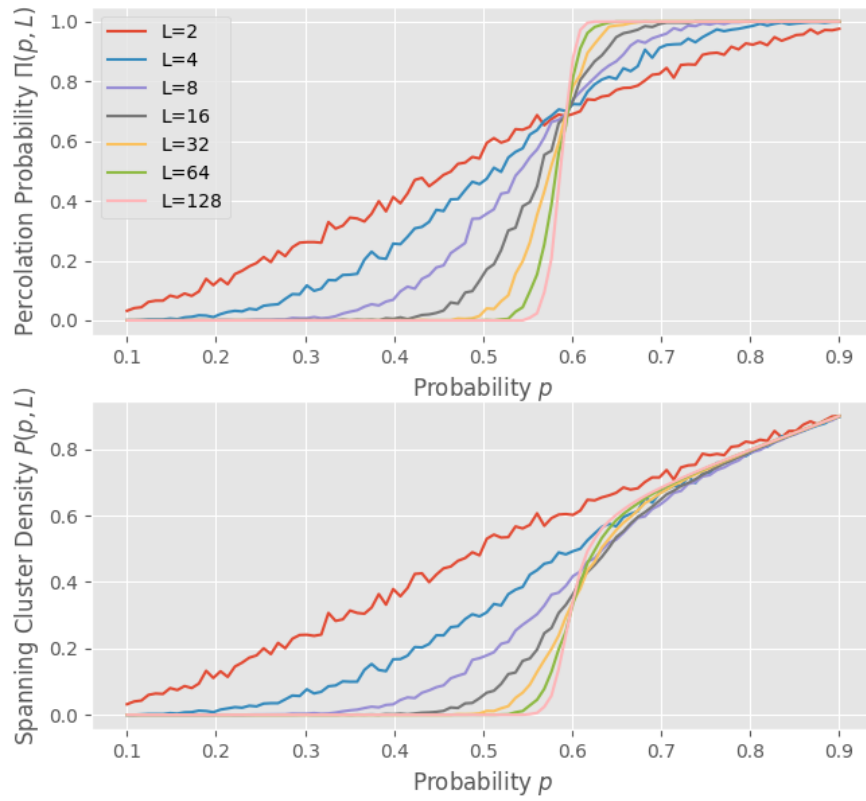


FIGURE 1.4. Spanning cluster density and percolation probability for different system sizes, $n = 1000$ samples.

2. TOPIC 10: PERCOLATION ON SMALL LATTICES

Discuss the percolation problem on a 2×2 lattice. Sketch $P(p, L)$ and $\Pi(p, L)$ for small L . Relate to your simulations. How do you calculate these quantities and how do you measure them in simulations?

- The 2×2 percolation problem is simple enough for us to list the possible outcomes by hand.

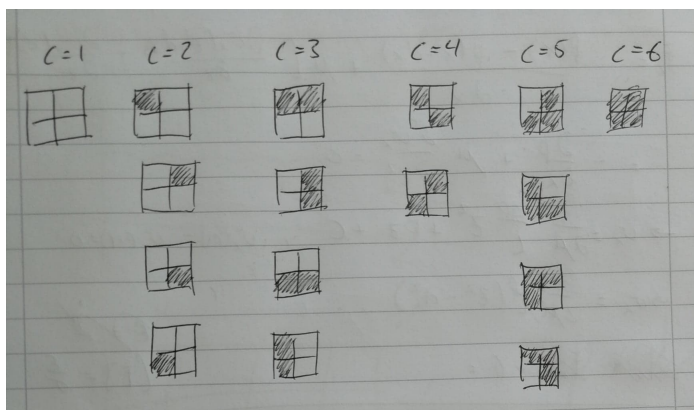


FIGURE 2.1. The 16 configurations of the 2×2 lattice.

- Can find exact expressions for $\Pi(p, L)$ and $P(p, L)$
- $\Pi(p, L) = \sum_c g_c \Pi(p, L|c) P(c)$
- Probabilities listed in table 2.1

TABLE 2.1. List of probabilities of the configurations of the 2×2 lattice, and the percolation probability.

c	g	$P(c)$	$\Pi(p, L c)$
1	1	$p^0(1-p)^4$	0
2	4	$p^1(1-p)^3$	0
3	4	$p^2(1-p)^2$	1
4	2	$p^2(1-p)^2$	0
5	4	$p^3(1-p)^1$	1
6	1	$p^4(1-p)^0$	1

- This gives $\Pi(p, L = 2) = 4p^2(1-p)^2 + 4p^3(1-p) + p^4 = p^2(p-2)^2$
- Density of the spanning cluster can be found from $p = \sum_{s=1}^{\infty} sn(s, p) + P(p, L)$

- Use same method as for the density of the spanning cluster:

$$\begin{aligned}
 (1) \quad P(p, L = 2) &= p - \sum_{s=1}^4 \sum_{c=1}^6 sn(s, p|c)P(c) \\
 (2) \quad &= p - [p(1-p)^3 + p^2(1-p)^2] \\
 (3) \quad &= p^2(2-p) \\
 (4) \quad &\text{(a bit unsure about this result, as we will see)}
 \end{aligned}$$

- Compare exact expressions to numerical results
- Monte Carlo simulations with $M = 1000$ samples
- Measurement of spanning cluster density: $P(p, L) = \frac{M_s}{L^d}$
- Calculated by the following algorithm

```

1 def spanning_cluster_density(m, single_perc = False):
2     nx, ny = m.shape
3     labels, n_features = sp.measurements.label(m)
4     regions = measure.regionprops(labels)
5     mass = 0
6     for region in regions:
7
8         x_start, y_start, x_stop, y_stop = region.bbox
9         dx = x_stop - x_start
10        dy = y_stop - y_start
11
12        if dx == nx or dy == ny:
13            mass += region.extent
14            if single_perc:
15                break
16    density = mass/(nx*ny)
17    return density

```

- Plot is shown in figure 2.2.
- Result from larger systems in figure 2.3

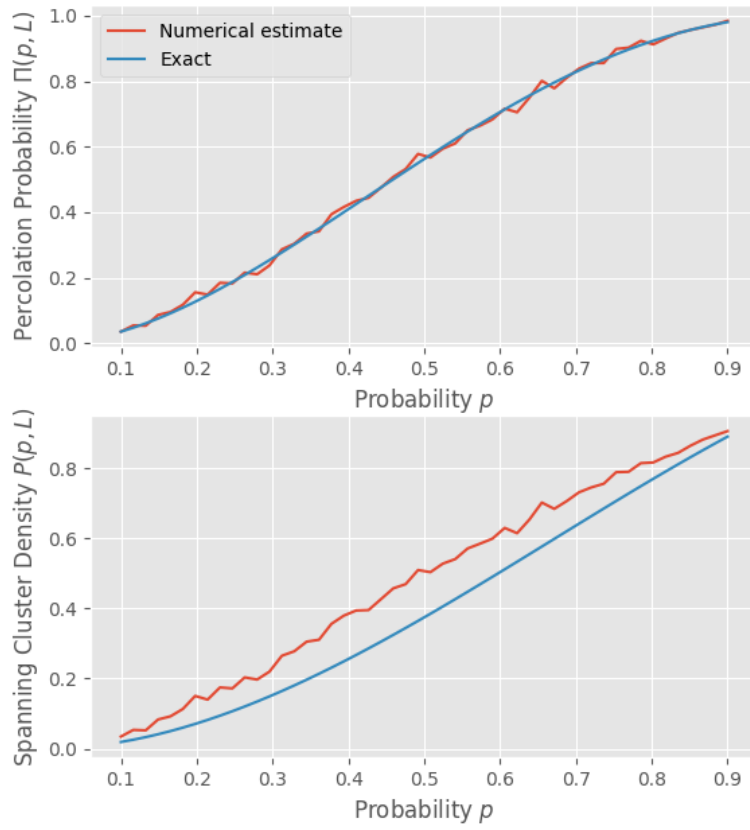


FIGURE 2.2. Density of the spanning cluster and percolation probability of the 2×2 lattice.

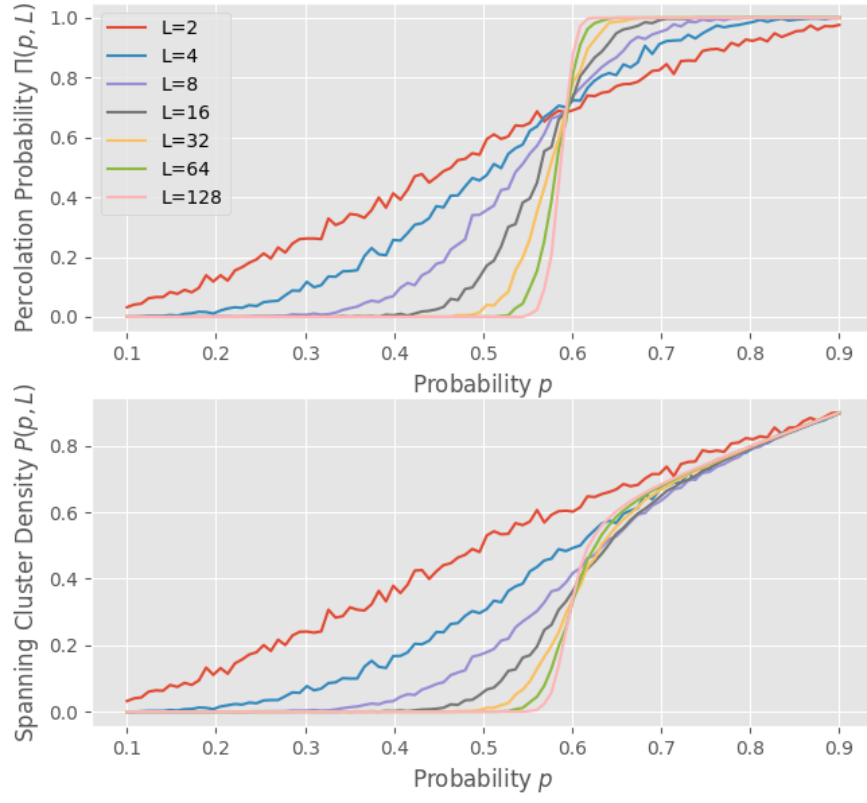


FIGURE 2.3. Density of the spanning cluster and percolation probability of different system sizes.

3. TOPIC 11: CLUSTER NUMBER DENSITY IN 1-D PERCOLATION

Define the cluster number density for 1-d percolation, and show how it can be measured. Discuss the behavior when $p \rightarrow p_c$. How does it relate to your simulations in two-dimensional systems?

- Definition of the cluster number density $n(s, p)$
- Definition of $P = sn(s, p)$
- Useful for characterizing clusters in a system
- 1-d percolation is a simple system, but also a powerful tool
- Cluster number density in 1-d: $n(s, p) = p^s(1 - p)^2$

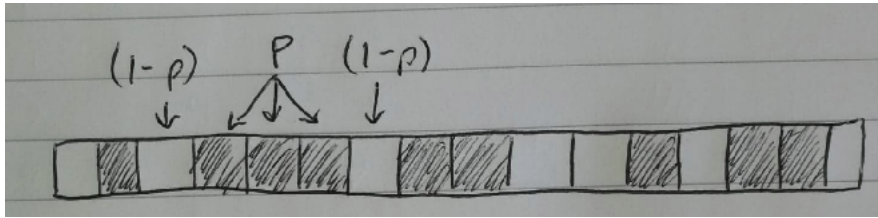


FIGURE 3.1. Percolation system in 1d.

- Check that it is normalized by calculating $p = \sum_{s=1}^{\infty} sn(s, p) + P$
- Numerical estimates of the cluster number density
- Can also estimate the probability by this method
- Using the above example: $\bar{p} = \sum_s \overline{sn(s, p)} = \sum_s \frac{sN_s}{L^d} = \frac{9}{16}$

TABLE 3.1. Table of values for the cluster number density from the example in figure 3.1.

s	N_s	$n(s, p)$
1	2	2/16
2	2	2/16
3	1	1/16

- Use of Monte Carlo methods for measuring good estimates: $\overline{n(s, p)} = \frac{N_s(M)}{ML^d}$
- Logarithmic binning

- Algorithm:

```

1 def log_bin(z, n_bins=10, base=10, log_max=None):
2     if log_max is None:
3         log_max = np.ceil(np.max(np.log(z))/np.log(base))
4     log_bins = np.logspace(0, log_max, n_bins, base=base)
5     log_hist, _ = np.histogram(z, bins=log_bins)
6     dz = np.diff(log_bins)
7     bin_mid = 0.5*(log_bins[1:] + log_bins[:-1])
8     log_hist_normed = log_hist / dz
9
10    return bin_mid, log_hist_normed

```

- Code for measuring $n(s, p)$

```

1 def cluster_number_density(L, p, n_samples, n_bins, logbase=10, log_max=None,
2     remove_zeros=False):
3     areas = []
4     for i in tqdm(range(n_samples)):
5         z = np.random.random((L, L))
6         m = z < p
7         labels, n_features = sp.measurements.label(m)
8         regions = measure.regionprops(labels)
9         for region in regions:
10            x_start, y_start, x_stop, y_stop = region.bbox
11            dx = x_stop - x_start
12            dy = y_stop - y_start
13            if dx != L and dy != L:
14                areas.append(region.area)
15
16    s, N_s = log_bin(areas, n_bins=n_bins, base=logbase, log_max=log_max)
17    nsp = N_s/(n_samples*L**2)
18    if remove_zeros:
19        idx = np.where(nsp > 1e-15)[0]
20        return s[idx], nsp[idx]
21    else:
22        return s, nsp

```

- Behaviour of $n(s, p)$ when $p \rightarrow p_c$ in 1-d. Figure 3.2

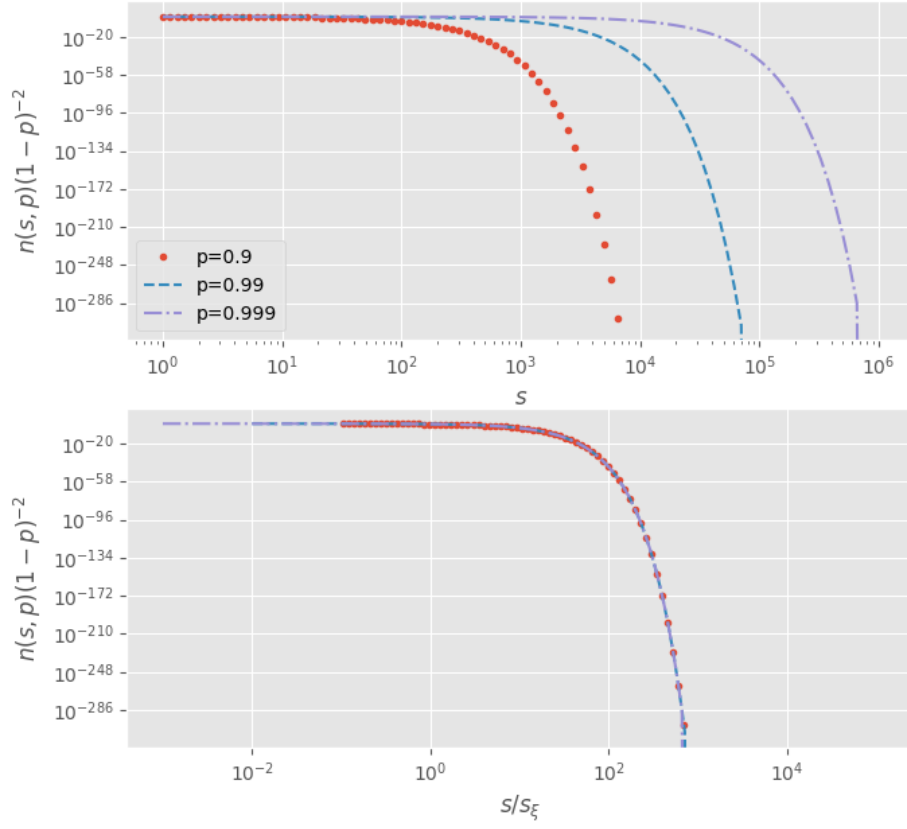


FIGURE 3.2. Plot of $n(s, p)(1-p)^{-2}$ as a function of s and s/s_{ξ} for a 1d system of length $L = 10^6$.

- Similar behaviour in 2 dimensions. Figure 3.3

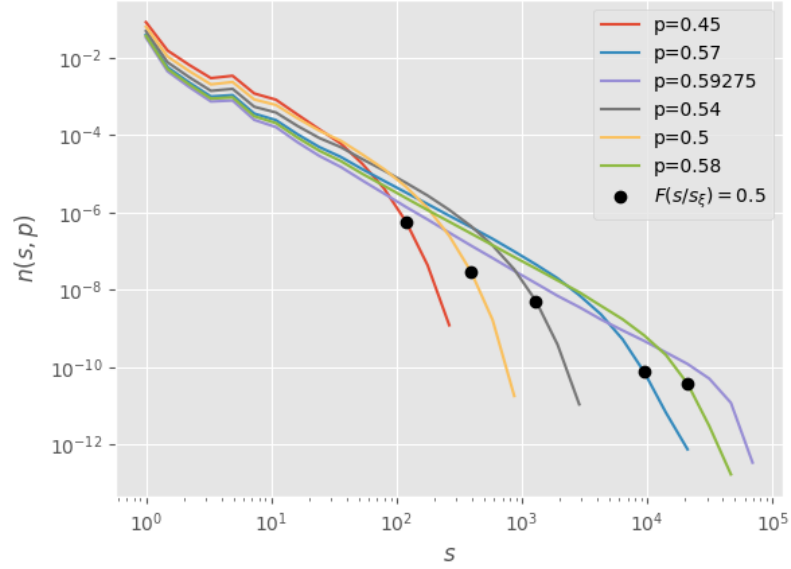


FIGURE 3.3. Cluster number density of a 2d system of size $L = 512$ for values of p approaching p_c .

- Understanding the behaviour by rewriting:

$$(5) \quad n(s, p) = (1 - p)^2 p^s = (1 - p)^2 e^{s \ln p} = (1 - p)^2 e^{-s/s_\xi},$$

where

$$(6) \quad s_\xi = -\frac{1}{\ln p}.$$

- Exponential power-law behaviour can be seen in the above loglog plot: $n(s, p) \propto s^{-\tau}$
- Estimating the scaling exponent τ
- Results from 2d: $\tau = 1.957$
- "True" value $\tau = 187/91 \approx 2.055$

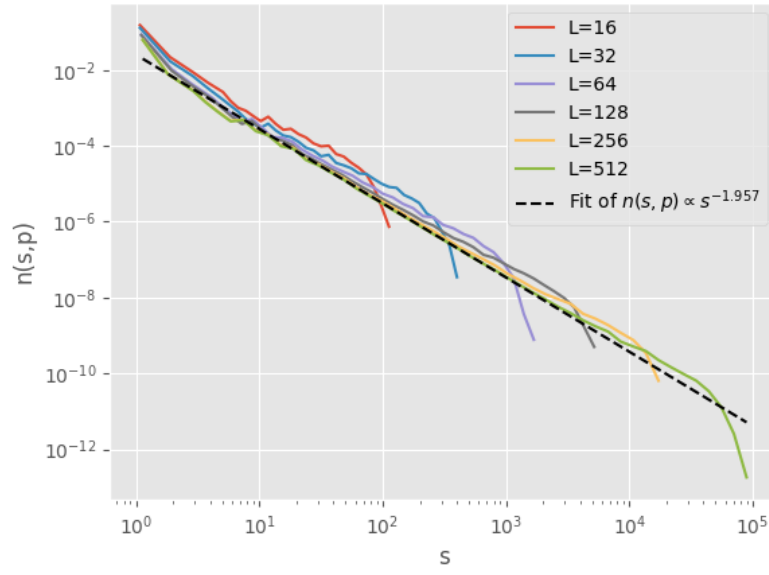


FIGURE 3.4. Cluster number density for varying system sizes at $p = p_c$. We also show the fitted exponent τ of the power law behaviour.

4. TOPIC 12: CLUSTER SIZE IN 1-D PERCOLATION

Introduce the characteristic cluster size for the 1-d percolation problem, and discuss their behavior when $p \rightarrow p_c$. Relate to your simulations on two-dimensional percolation

- Cluster number density in 1d: $n(s, p) = (1 - p)^2 p^s$
- How does it behave in relation to s when $p \rightarrow p_c$? Figure 4.1

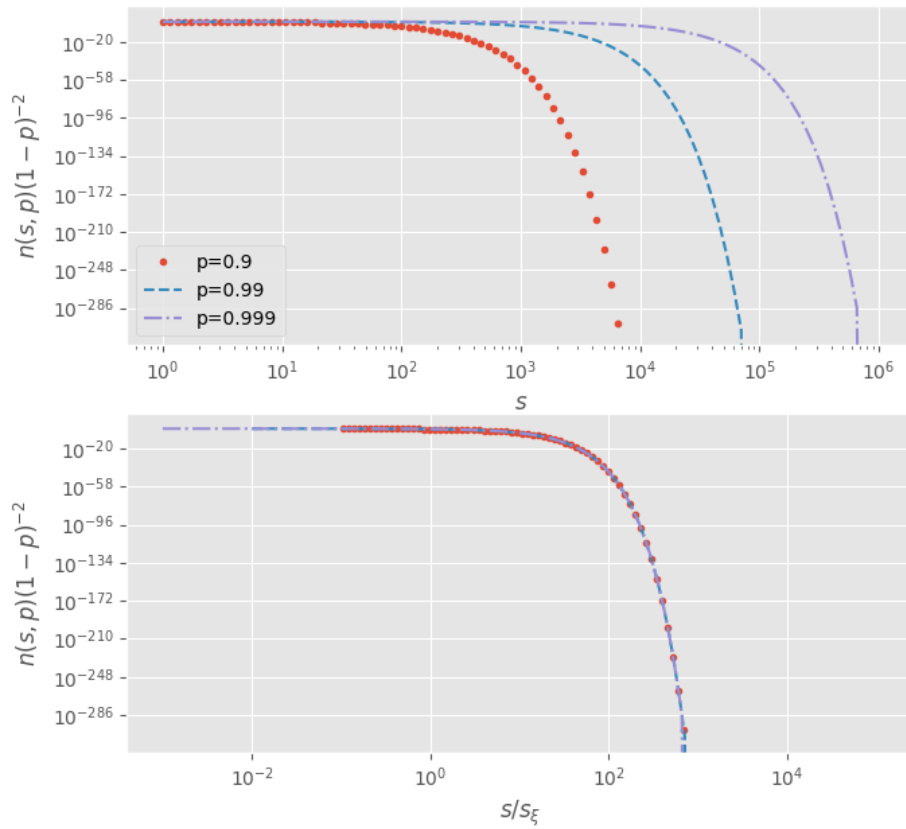


FIGURE 4.1. Plot of $n(s, p)(1 - p)^{-2}$ as a function of s and s/s_ξ in a 1d system of size $L = 10^6$.

- Definition of the characteristic cluster size s_ξ
- Diverges as $p \rightarrow p_c$

- Can understand this by rewriting the cluster number density:

$$(7) \quad n(s, p) = (1 - p)^2 p^s = (1 - p)^2 e^{s \ln p} = (1 - p)^2 e^{-s/s_\xi},$$

where

$$(8) \quad s_\xi = -\frac{1}{\ln p}.$$

- s_ξ causes exponential fall-off of $n(s, p)$
- Also the case in 2d, figure 4.2

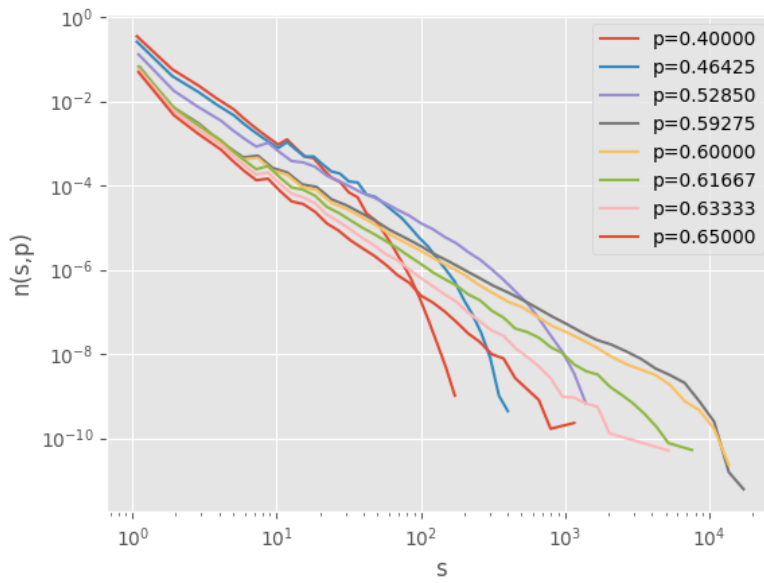


FIGURE 4.2. Plot of $n(s, p)$ for varying p in a 2d system of size $L = 200$.

- Generalizing s_ξ when $p \rightarrow p_c$, Taylor expansion of $\ln p \approx -(1 - p)$:

$$(9) \quad s_\xi \approx \frac{1}{1 - p} = \frac{1}{p_c - p} = |p - p_c|^{-1/\sigma},$$

- s_ξ marks the cross-over between the two behaviours of $n(s, p)$.

- Characterize s_ξ by measurements of $n(s, p)$

- Can be defined as

$$(10) \quad \frac{n(s, p)}{n(s, p_c)} = F(s/s_\xi) = 0.5.$$

- Use this in simulations of $n(s, p)$ to find σ in 2d, figure 4.3

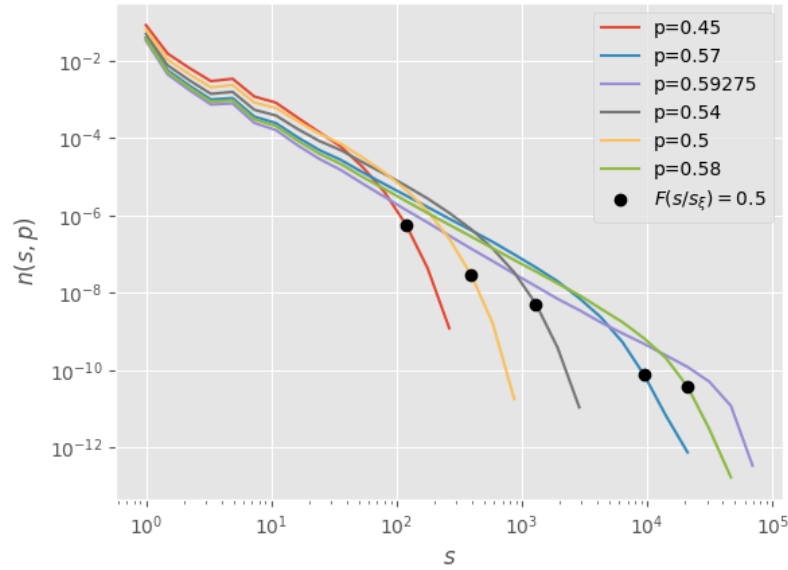


FIGURE 4.3. Cluster number density of a 2d system of size $L = 512$ for values of p approaching p_c , with corresponding values of s_ξ .

```

1 def log_bin(z, n_bins=10, base=10, log_max=None):
2     if log_max is None:
3         log_max = np.ceil(np.max(np.log(z))/np.log(base))
4     log_bins = np.logspace(0, log_max, n_bins, base=base)
5     log_hist, _ = np.histogram(z, bins=log_bins)
6     dz = np.diff(log_bins)
7     log_bins = 0.5*(log_bins[1:] + log_bins[:-1])
8     log_hist_normed = log_hist / dz
9
10    return log_bins, log_hist_normed
11
12 def cluster_number_density(L, p, n_samples, n_bins, logbase=10, log_max=None,
13    remove_zeros=False):
14    areas = []
15    for i in tqdm(range(n_samples)):
16        z = np.random.random((L, L))
17        m = z < p
18        labels, n_features = sp.measurements.label(m)
19        regions = measure.regionprops(labels)
20        for region in regions:
21            x_start, y_start, x_stop, y_stop = region.bbox
22            dx = x_stop - x_start
23            dy = y_stop - y_start
24            if dx != L and dy != L:
25                areas.append(region.area)
26
27    s, N_s = log_bin(areas, n_bins=n_bins, base=logbase, log_max=log_max)
28    nsp = N_s/(n_samples*L**2)
29    if remove_zeros:

```

```

29     idx = np.where(nsp > 1e-15)[0]
30     return s[idx], nsp[idx]
31 else:
32     return s, nsp
33
34 def estimate_s_xi(L, n_samples, n_bins):
35     pc = 0.59275
36     eps = 1e-14
37     p_vals = [0.45, 0.50, 0.54, 0.57, 0.58]
38     logbase = 1.3
39     remove_zeros = False
40     s, nc = cluster_number_density(L, pc, n_samples, n_bins, logbase,
41     remove_zeros=remove_zeros)
42     log_max = np.log(s[-1])/np.log(logbase)
43     nonzero = np.where(nc > eps)[0]
44     sxi_vals = []
45     F_vals = []
46     p_array = []
47     for i, p in tqdm(enumerate(p_vals)):
48         s, n = cluster_number_density(L, p, n_samples, n_bins, logbase, log_max
49         , remove_zeros)
50         nonzero = np.where(n > eps)[0]
51
52         temp_sxi = []
53         temp_n = []
54         for i in range(len(n)):
55             if n[i] > eps and nc[i] > eps:
56                 if n[i]/nc[i] <= 0.5:
57                     temp_sxi.append(s[i])
58                     temp_n.append(n[i])
59
60         if len(temp_n) > 0:
61             idx = np.argmax(temp_n)
62             F = temp_n[idx]
63             sxi = temp_sxi[idx]
64             sxi_vals.append(sxi)
65             F_vals.append(F)
66             p_array.append(p)
67
68     return sxi_vals, p_array

```

- Linear fit $\rightarrow \sigma = 0.462$
- "True" value: $\sigma = 36/91 \approx 0.395$.
- Plot of s_ξ as function of p in figure 4.4

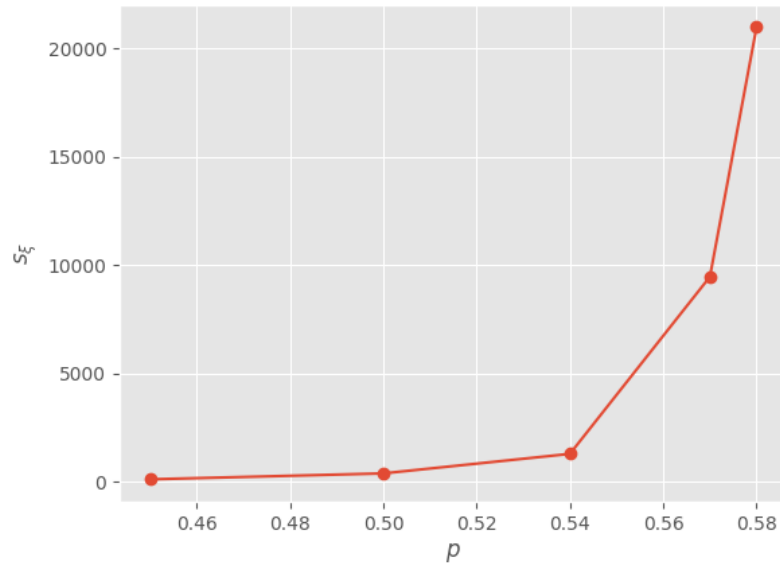


FIGURE 4.4. s_ξ as a function of p for a system of size $L = 512$.

5. TOPIC 13: MEASUREMENT AND BEHAVIOR OF $P(p, L)$ AND $\Pi(p, L)$

Discuss the behavior of $P(p, L)$ and $\Pi(p, L)$ in a system with a finite system size L . How do you measure these quantities?

- Definition of $\Pi(p, L)$

- Definition of $P(p, L)$

$$P(p, L) = \frac{M_S}{L^d}$$

- Expected behaviour of these quantities
- Small vs large systems

- Measurements of $\Pi(p, L)$:

$$(11) \quad \Pi(p, L) = \frac{\sum_{i=1}^n \pi_i}{n},$$

where

$$(12) \quad \pi_i = \begin{cases} 1 & \text{if percolating} \\ 0 & \text{else.} \end{cases}$$

- Measurements of $P(p, L)$:

$$P(p, L) = \frac{\sum_{i=1}^n M_{S,i}}{nL^2}$$

- Algorithm for finding spanning clusters

```

1 def spanning_cluster_density(m, single_perc = False):
2     nx, ny = m.shape
3     labels, n_features = sp.measurements.label(m)
4     regions = measure.regionprops(labels)
5     density = 0
6     for region in regions:
7
8         x_start, y_start, x_stop, y_stop = region.bbox
9         dx = x_stop - x_start
10        dy = y_stop - y_start
11
12        if dx == nx or dy == ny:
13            density += region.extent
14            if single_perc:
15                break
16
17    return density

```

- Algorithm for estimating Π and P :

```

1 def estimate_Pi_P(L_arr, p_arr, n_samples):
2     for L in L_arr:
3         Pi_arr = np.zeros(len(p_arr))
4         density_arr = np.zeros(len(p_arr))
5         for p, prob in enumerate(p_arr):
6             for i in range(n_samples):
7                 z = np.random.random((L,L))
8                 m = z < prob

```

```

9         density = spanning_cluster_density(m)
10        density_arr[p] += density
11        if density > 0:
12            Pi_arr[p] += 1
13
14        Pi_arr /= n_samples
15        density_arr /= n_samples
16    return Pi_arr, density_arr

```

- Results for varying system sizes with $M = 1000$ samples, figure 5.1

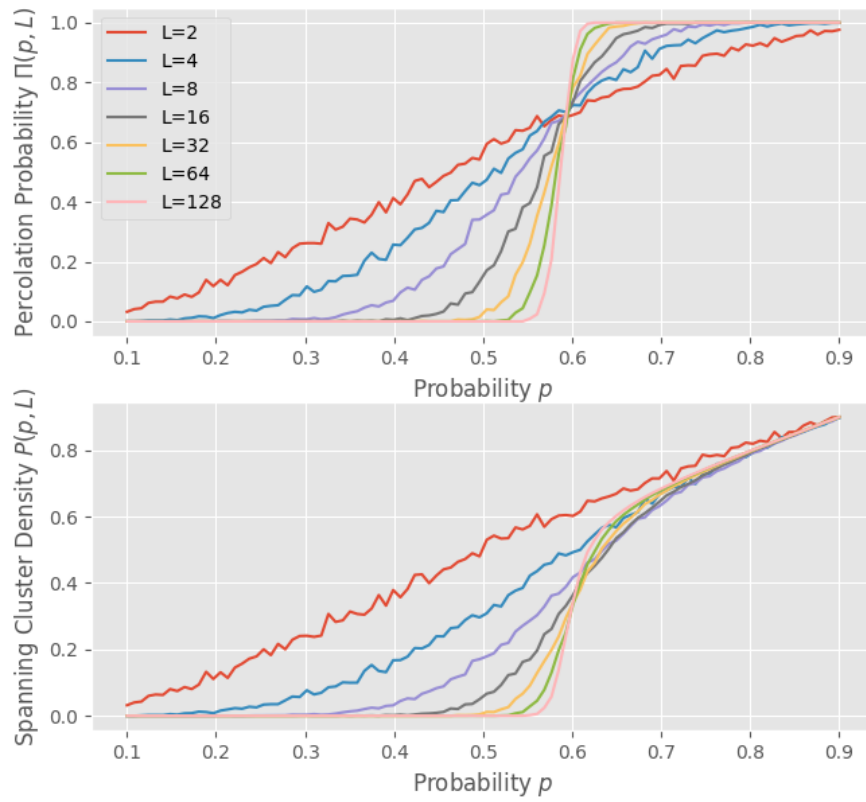


FIGURE 5.1. Spanning cluster density and percolation probability for different system sizes, $n = 1000$ samples.

- Π Linear for small systems, approaches step function for large systems.
- Expected behaviour for an infinite system:

$$(13) \quad \Pi(p, \infty) = \begin{cases} 0 & p < p_c \\ 1 & p \geq p_c \end{cases}$$

- Similar behaviour for P below the percolation threshold.
- Linear behaviour for $p > p_c$.
- Behaviour of $P(p_c, L)$ as the system size is increased.

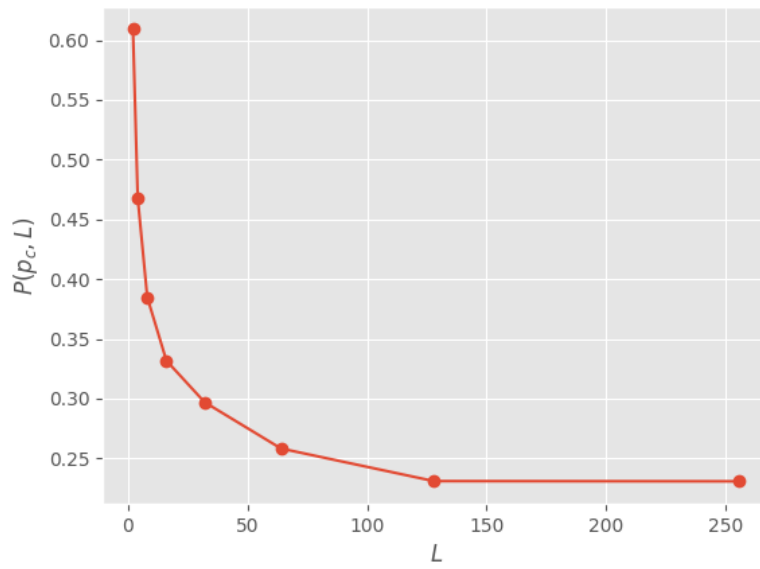


FIGURE 5.2. Spanning cluster density for different system sizes at $p = p_c$.

6. TOPIC 14: THE CLUSTER NUMBER DENSITY

Introduce the cluster number density and its applications: Definition, measurement, scaling and data-collapse.

- Definition of $n(s, p)$
- $P(\text{site is part of cluster of size } s) = sn(s, p)$
- Characterizes the size of clusters, basis for scaling theory in percolation
- Can be studied exactly in 1d, restricted to numerical estimates in 2d:

$$\overline{n(s, p)} = \frac{N_s}{ML^2}.$$
- Measured by Monte Carlo simulations
- Logarithmic binning
- Effectively measures $\overline{n(s, p)}\Delta s_i$

```

1 def log_bin(z, n_bins=10, base=10, log_max=None):
2     if log_max is None:
3         log_max = np.ceil(np.max(np.log(z))/np.log(base))
4     log_bins = np.logspace(0, log_max, n_bins, base=base)
5     log_hist, _ = np.histogram(z, bins=log_bins)
6     dz = np.diff(log_bins)
7     bin_mid = 0.5*(log_bins[1:] + log_bins[:-1])
8     log_hist_normed = log_hist / dz
9
10    return bin_mid, log_hist_normed

```

- Algorithm for measuring $n(s, p)$

```

1 def cluster_number_density(L, p, n_samples, n_bins, logbase=10, log_max=None,
2     remove_zeros=False):
3     areas = []
4     for i in tqdm(range(n_samples)):
5         z = np.random.random((L, L))
6         m = z < p
7         labels, n_features = sp.measurements.label(m)
8         regions = measure.regionprops(labels)
9         for region in regions:
10             x_start, y_start, x_stop, y_stop = region.bbox
11             dx = x_stop - x_start
12             dy = y_stop - y_start
13             if dx != L and dy != L:
14                 areas.append(region.area)
15
16    s, N_s = log_bin(areas, n_bins=n_bins, base=logbase, log_max=log_max)
17    nsp = N_s/(n_samples*L**2)
18    if remove_zeros:

```



```

18     idx = np.where(nsp > 1e-15)[0]
19     return s[idx], nsp[idx]
20 else:
21     return s, nsp

```

- $n(s, p)$ as a function of s in figure 6.1
- Can propose the following scaling form: $n(s, p) \propto s^{-\tau}$

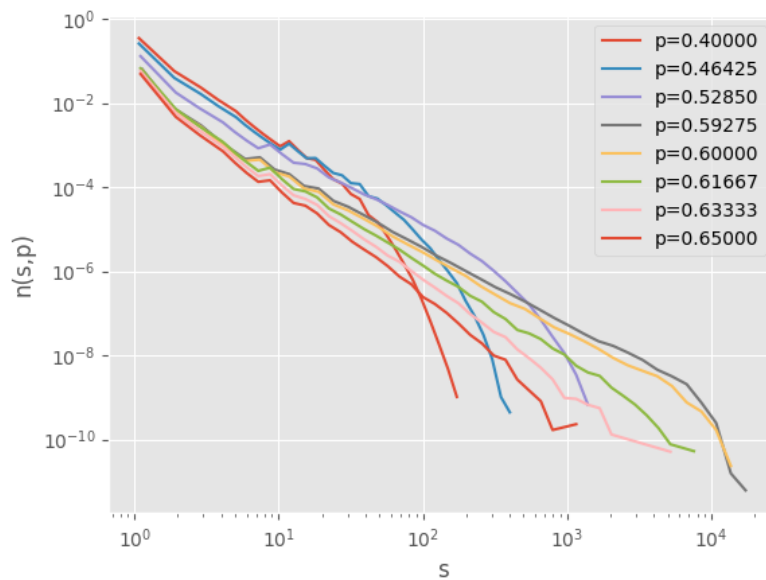


FIGURE 6.1. Plot of $n(s, p)$ for varying p in a 2d system of size $L = 200$.

- Estimating τ : $\log n(s, p) = -\tau \log s$
- Result in figure 6.2
- Estimated value: $\tau = 1.957$, "true" value: $\tau = 187/91 \approx 2.055$

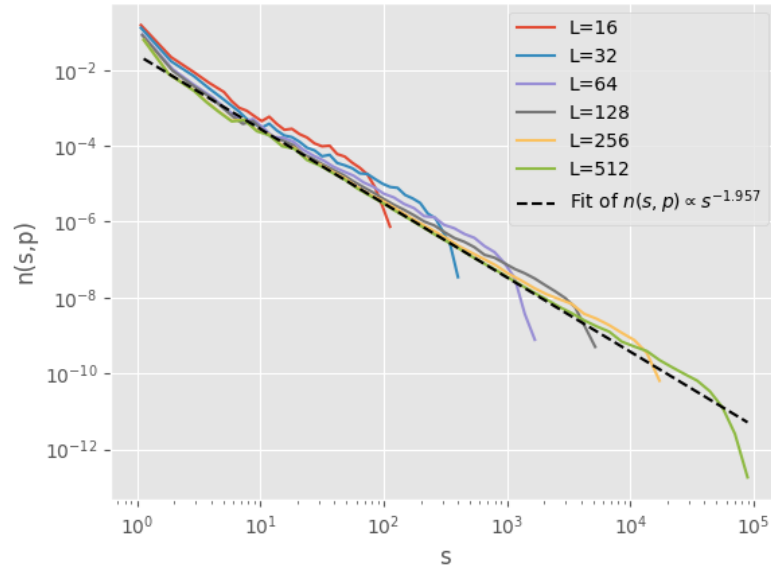


FIGURE 6.2. Cluster number density for varying system sizes at $p = p_c$. We also show the fitted exponent τ of the power law behaviour.

- Can also find s_ξ from these simulations
- Use result from 1d percolation to develop scaling theory of $n(s, p)$:
 $s_\xi \propto |p - p_c|^{-1/\sigma}$
- Propose that $n(s, p)$ has the form:

$$(14) \quad n(s, p) = s^{-\tau} F(s/s_\xi) = s^{-\tau} F\left((p - p_c)^{1/\sigma} s\right),$$

- Define s_ξ to be at the point $\frac{n(s, p)}{n(s, p_c)} = F(s/s_\xi) = 0.5$
- Algorithm for finding s_ξ :

```

1 def estimate_s_xi(L, n_samples, n_bins):
2     pc = 0.59275
3     eps = 1e-14
4     p_vals = [0.45, 0.50, 0.54, 0.57, 0.58]
5     logbase = 1.3
6     remove_zeros = False
7     s, nc = cluster_number_density(L, pc, n_samples, n_bins, logbase,
8     remove_zeros=remove_zeros)
9     log_max = np.log(s[-1])/np.log(logbase)
10    nonzero = np.where(nc > eps)[0]
11    sxi_vals = []
12    F_vals = []

```

```

12 p_array = []
13 for i, p in tqdm(enumerate(p_vals)):
14     s, n = cluster_number_density(L, p, n_samples, n_bins, logbase, log_max
15     , remove_zeros)
16     nonzero = np.where(n > eps)[0]
17
18     temp_sxi = []
19     temp_n = []
20     for i in range(len(n)):
21         if n[i] > eps and nc[i] > eps:
22             if n[i]/nc[i] <= 0.5:
23                 temp_sxi.append(s[i])
24                 temp_n.append(n[i])
25
26     if len(temp_n) > 0:
27         idx = np.argmax(temp_n)
28         F = temp_n[idx]
29         sxi = temp_sxi[idx]
30         sxi_vals.append(sxi)
31         F_vals.append(F)
32         p_array.append(p)
33
34 return sxi_vals, p_array

```

- Perform line fit to estimate σ :

$$\log s_\xi = -\frac{1}{\sigma} \log |p - p_c|$$

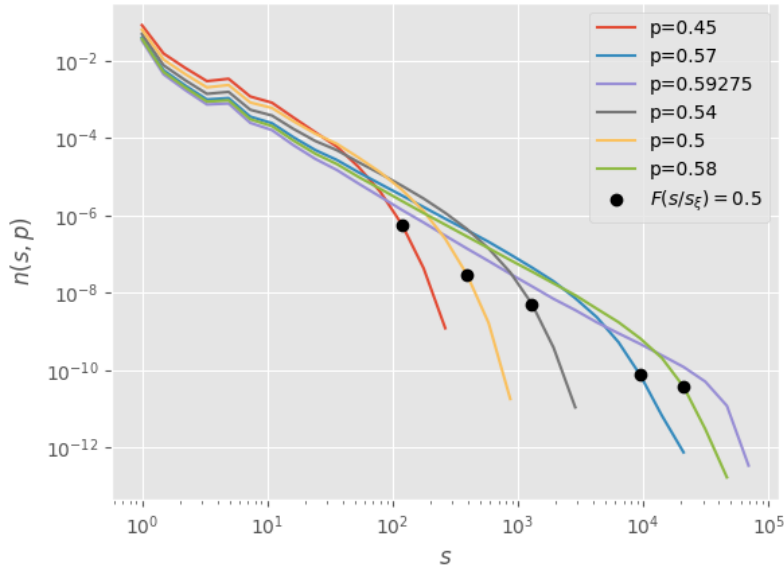


FIGURE 6.3. Cluster number density of a 2d system of size $L = 512$ for values of p approaching p_c .

- Our results: $\sigma = 0.462$, "true" value $\sigma = 36/91 \approx 0.395$

- Divergent behaviour of s_ξ shown in figure 6.4

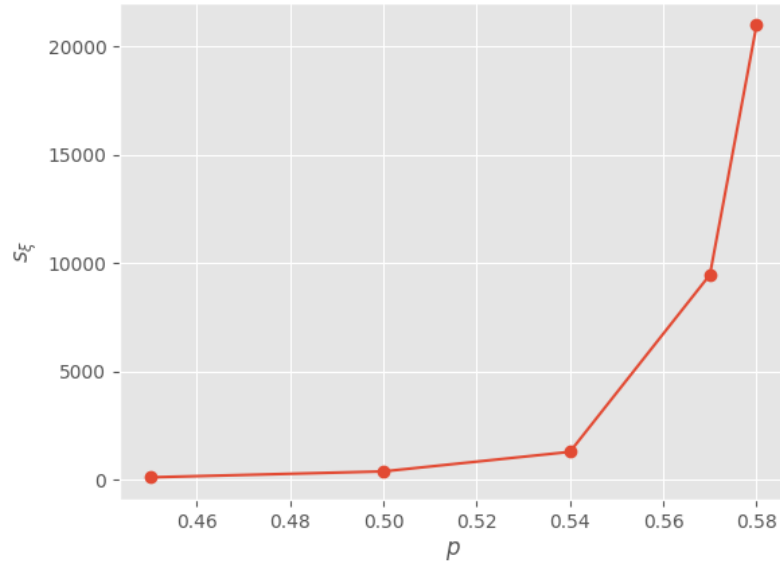


FIGURE 6.4. s_ξ as a function of p for a system of size $L = 512$.

- Data collapse: plot $s^\tau n(s, p)$ as a function of $s|p - p_c|^{1/\sigma}$, figure 6.5

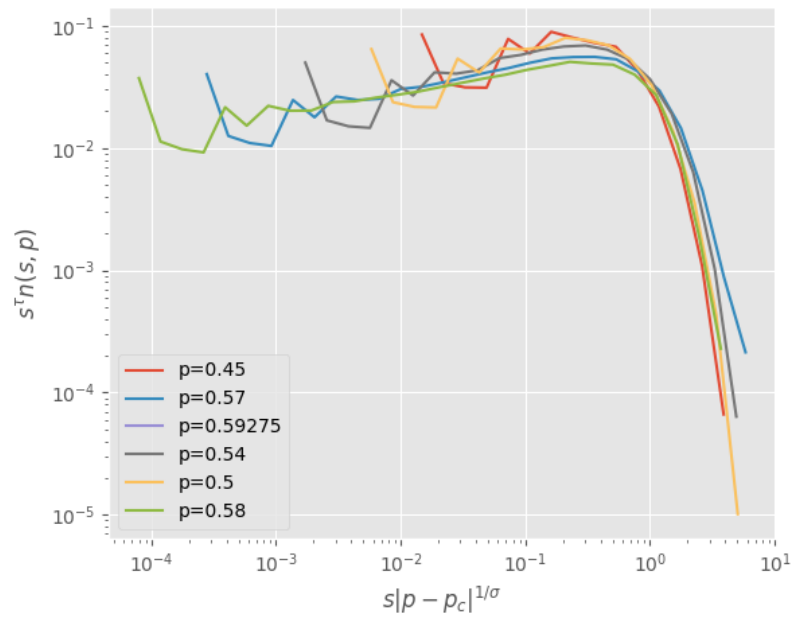


FIGURE 6.5. Data-collapse of $n(s, p)$.

7. TOPIC 15: FINITE SIZE SCALING OF $\Pi(p, L)$

Discuss the behavior of $\Pi(p, L)$ in a system with a finite system size L . How can we use this to find the scaling exponent ν , and the percolation threshold, p_c ?

- Definition of $\Pi(p, L)$
- Expectations to it's behaviour
- Can be studied exactly in 1d and small 2d systems
- Restricted to numerical estimates in higher dimensions
- Use this as basis for a scaling theory
- Measurements by Monte Carlo simulations

$$(15) \quad \Pi(p, L) = \frac{\sum_{i=1}^n \pi_i}{n},$$

where

$$(16) \quad \pi_i = \begin{cases} 1 & \text{if percolating} \\ 0 & \text{else.} \end{cases}$$

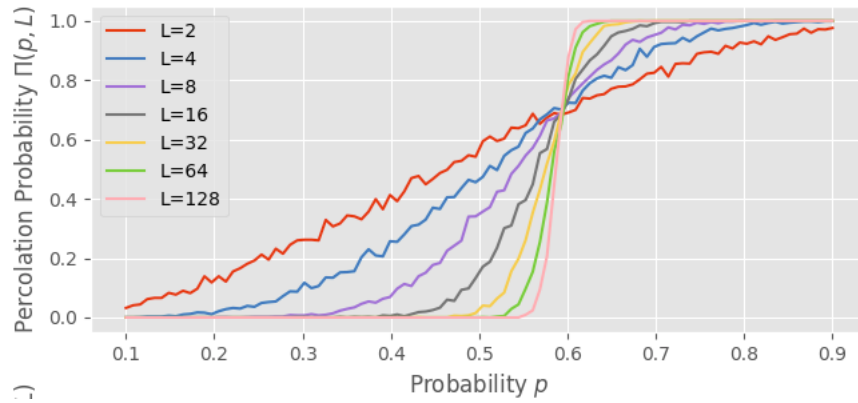


FIGURE 7.1. Percolation probability as a function of p for varying L .

- Behaviour of the different system sizes
- Estimating the percolation threshold

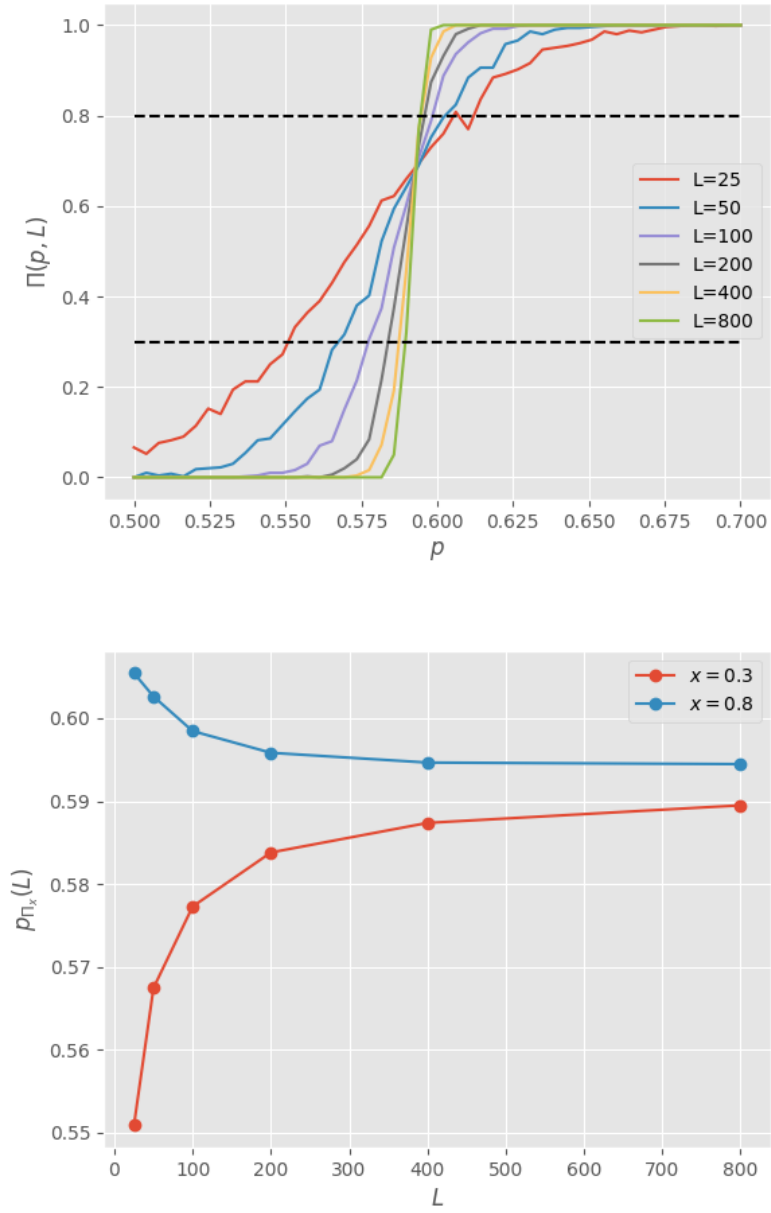


FIGURE 7.2. Plot of $\Pi(p, L)$ and p_{Π_x} at $x = 0.3, 0.5$ for varying system sizes.

- For an infinite system:

$$(17) \quad \Pi(p, \infty) = \begin{cases} 0 & p < p_c \\ 1 & p \geq p_c, \end{cases}$$

- Not a viable strategy for estimating p_c

- Instead develop a scaling theory
- From above results: Π is dependent on both ξ and L
- Proposed scaling form: $\Pi(p, L) = \xi f\left(\frac{L}{\xi}\right)$
- Use $\xi = \xi_0 |p - p_c|^{-\nu}$:
- $\Pi(p, L) = f(L \xi_0 |p - p_c|^\nu) = f(\xi_0 (L^{1/\nu} (p - p_c))^\nu)$
- Introduce $\phi(u) = f(\xi_0 u^{1/\nu})$
- Finite size scaling ansatz: $\Pi(p, L) = \phi(L^{1/\nu} (p - p_c))$
- Goal: use this result to estimate p_c and ν
- Problem if both are unknown
- From above results and figure:

$$(18) \quad x = \Pi(p_x, L) = \phi(L^{1/\nu} (p_x(L) - p_c))$$

$$(19) \quad \phi^{-1}(x) = (p_x - p_c) L^{1/\nu} = C_x.$$

- Subtract equation from itself:
- $dp = p_{x_1} - p_{x_2} = (C_{x_1} - C_{x_2}) L^{-1/\nu}$
- Estimate ν by line fit:
- $\log dp = \log(C_{x_1} - C_{x_2}) - \frac{1}{\nu} \log(L)$
- Estimated: $\nu = 1.405$, "true" value $\nu = 4/3 \approx 1.333$
- Can now use this to estimate p_c :
- $p_x(L) = p_c + C_x L^{-1/\nu}$
- Results in figure 7.3

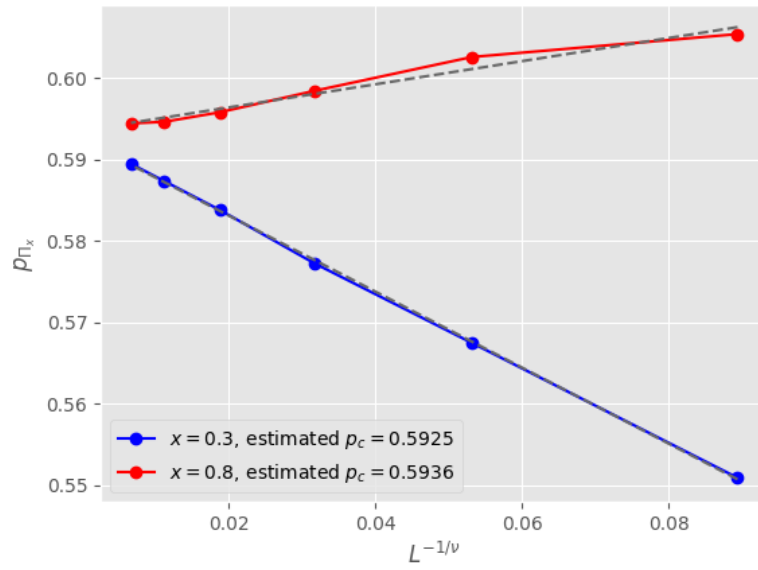


FIGURE 7.3. Plot of p_x as a function of $L^{-1/\nu}$ for $x = 0.3$ and $x = 0.8$. The y -intercept is the estimated value for p_c .

- Shape of the scaling function $\phi(x)$:
- Plot $\Pi(p, L)$ as a function of $(p - p_c)L^{-1/\nu}$

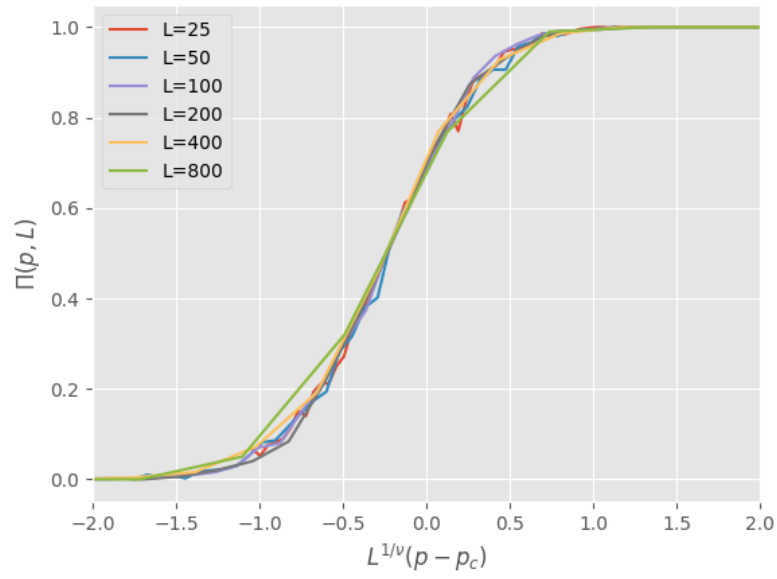


FIGURE 7.4. Data-collapse of $\Pi(p, L)$.

8. TOPIC 16: EFFECTIVE PERCOLATION THRESHOLD

Define and discuss the effective percolation threshold in a system with a finite system size L . How can we measure it, and what consequences does it have that the effective percolation threshold is different from the actual percolation threshold?

- System of finite size L will have a probability p which on average gives rise to a spanning cluster
- Dependent on system size, $p_x(L)$
- Approaches p_c as $L \rightarrow \infty$
- Definition of effective percolation threshold: $\Pi(p_x, L) = x$
- Plot of $\Pi(p, L)$, figure 8.1

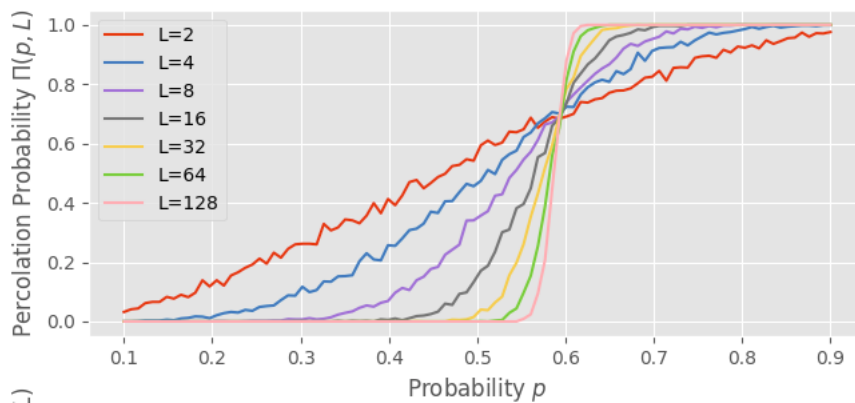


FIGURE 8.1. Percolation probability as a function of p for different L .

- Simplest approach: define p_x where $\Pi(p, L) = 1/2$.
- Can measure $p_x(L)$ by interpolation:

```

1 x_vals = [0.3, 0.8]
2 L_vals = [25, 50, 100, 200, 400, 800]
3 for j, x in enumerate(x_vals):
4     for i, L in enumerate(L_vals):
5         Pi = data[i,:] #data contains Pi_p_L loaded from file
6         idx = np.argmax(Pi > x)
7         pc = p[idx-1] + (x-Pi[idx-1])*(p[idx] - p[idx-1])/(Pi[idx] - Pi[idx-1])
8         p_pi_vals[j,i] = pc

```

- Results from calculations of Π with $M = 500$ samples, figure 8.2
- Convergence towards a finite value

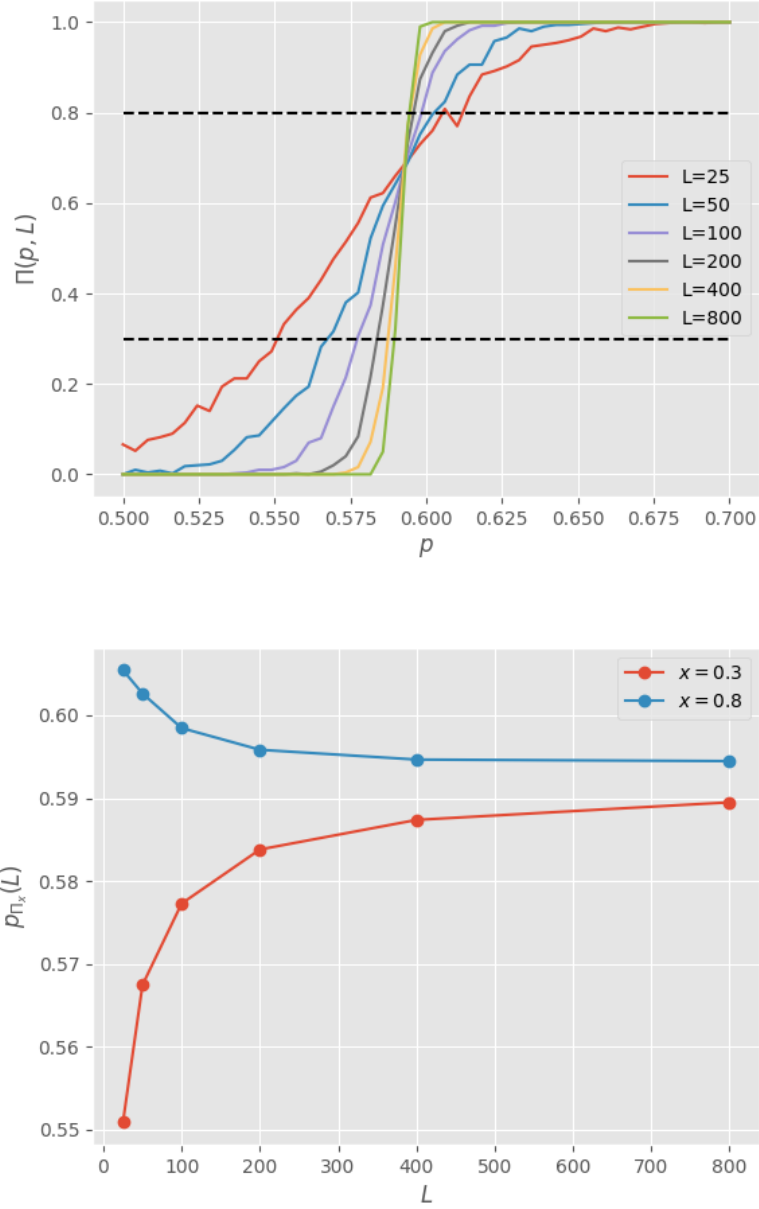


FIGURE 8.2. Plot of $\Pi(p, L)$ and p_{Π_x} at $x = 0.3, 0.5$ for varying system sizes.

- Consequences for $p_x \neq p_c$:

- Find scaling of Π and ξ , and predict p_c
- From scaling theory of $\Pi(p, L)$:

$$(20) \quad \Pi(p, L) = \xi f(L/\xi) = f\left(\xi_0 \left(L^{1/\nu}(p - p_c)\right)^\nu\right)$$

$$(21) \quad = \phi\left(L^{1/\nu}(p - p_c)\right),$$

where

$$(22) \quad \phi(u) = f\left(\xi_0 u^{1/\nu}\right).$$

- Estimating the scaling of ξ :
- Apply $\phi^{-1}(x)$ and use $\Pi(p_x, L) = x$:
- $(p_x(L) - p_c)L^{1/\nu} = \phi^{-1}(x) = C_x$
- Fit line to our data: $\log(p_x(L) - p_c) = \log(C_x) - \frac{1}{\nu} \log(L)$
- Requires that p_c is known. Can instead use:
- $dp = \log(C_{x_1} - C_{x_2}) - \frac{1}{\nu} \log(L)$
- Estimated: $\nu = 1.405$, "true" value $\nu = 4/3 \approx 1.333$.
- Can now estimate p_c if we wish, results in figure 8.3

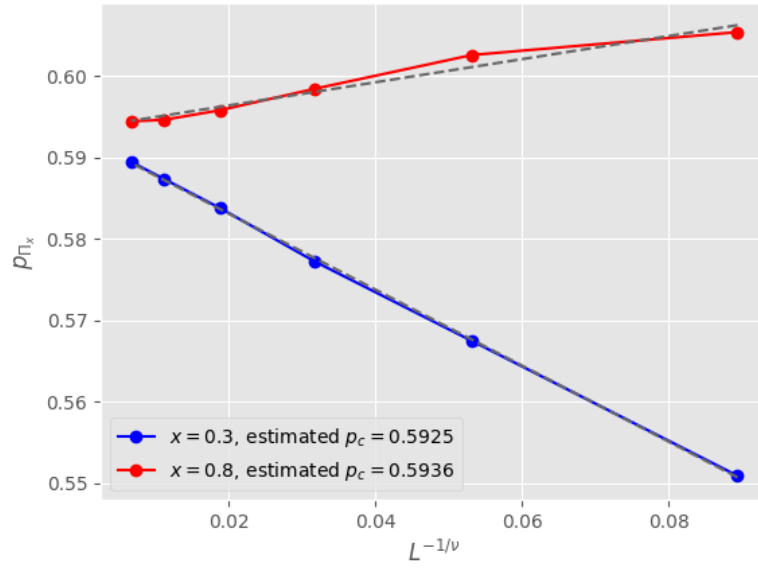
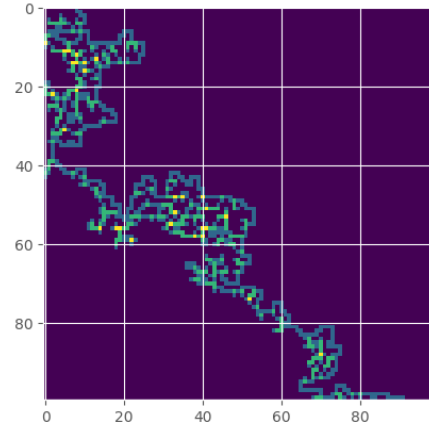


FIGURE 8.3. Plot of the effective percolation threshold for $x = 0.3$ and $x = 0.8$. The y -intercept is the estimated value for p_c .

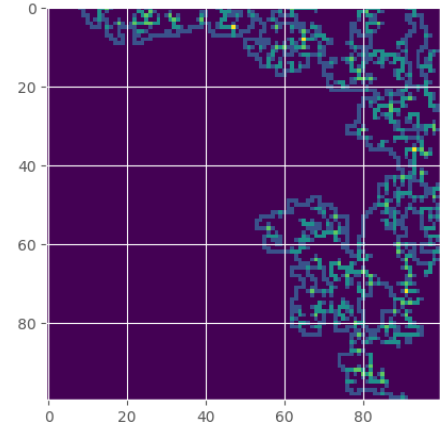
9. TOPIC 17: SUBSETS OF THE SPANNING CLUSTER

Introduce and discuss the scaling of subsets of the spanning cluster. How can we measure the singly-connected bonds, and how does it scale?

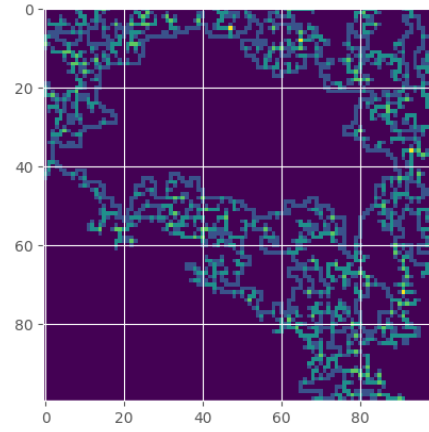
- Definition of the subsets of the spanning cluster:
- Singly connected bonds
- Backbone
- Dangling ends
- visualization of the subsets in figure 9.1



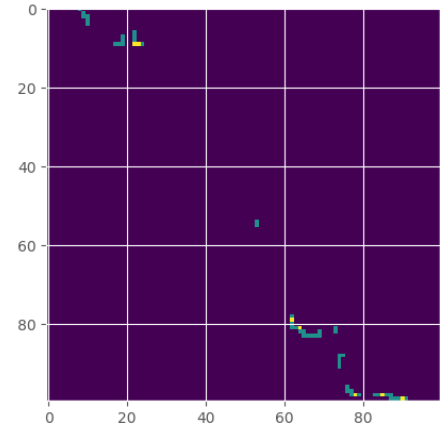
Left walker



Right walker



Backbone (left + right)



Singly connected bonds

FIGURE 9.1. Visualization of the paths of the right and left walkers.

- Scaling of the subset geometries
- Propose similar scaling form as $M \propto L^D$

$$(23) \quad M_{SC} \propto L^{D_{SC}}$$

$$(24) \quad M_{min} \propto L^{D_{min}}$$

$$(25) \quad M_{max} \propto L^{D_{max}}$$

$$(26) \quad M_{BB} \propto L^{D_{BB}}$$

$$(27) \quad M_{DE} \propto L^{D_{DE}}.$$

- Hierarchy of the exponents:

$$(28) \quad D_{SC} \leq 1 \leq D_{min} \leq D_{max} \leq D_{BB} \leq D \leq d.$$

- Scaling of dangling ends:

- $M_{DE} = M - M_{BB} \propto L^D - L_{BB}^D = L^{D-D_{BB}}$

- For an infinite system:

- $\frac{M_{DE}}{M} \propto 1 - \frac{L_{BB}^D}{L^D} = 1 - L^{D_{BB}-D}$

- Scaling of the singly connected bonds

- Use Monte Carlo simulation

- Measure M_{SC} for $L = 2^k$ for $k \in [4, 10]$.

- Linear fit of $\log M_{SC} \propto D_{SC} \log L$,. Figure 9.2

- Estimated: $D_{SC} = 0.7822$, "true" value: 0.75

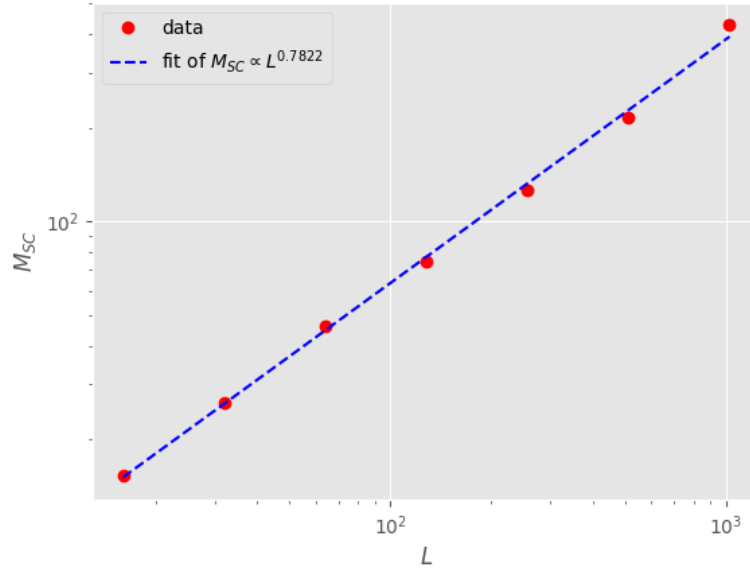


FIGURE 9.2. Mass of the singly connected bonds as a function of system size.

- Density of the singly connected bonds:
- $P_{SC} = \frac{M_{SC}}{L^d} \propto L^{D_{SC}-d}$
- Measure as a function of $(p - p_c)$
- Plot in figure 9.3

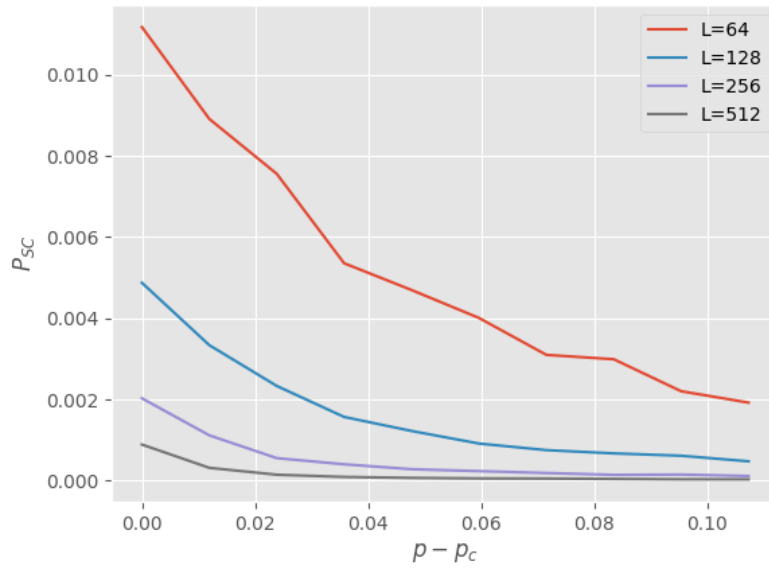


FIGURE 9.3. Density of the singly connected bonds as a function of $p - p_c$ for various L .

- Power-law behaviour of the density
- Propose $P_{SC} \propto (p - p_c)^{-x}$
- Fit to the data: $x = 1.06$. Plot in figure 9.4

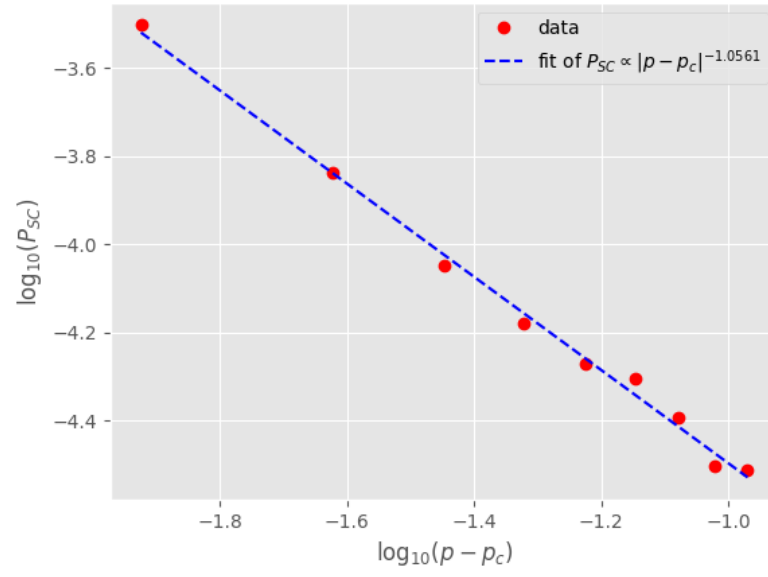


FIGURE 9.4. Loglog plot of P_{SC} as a function of $(p - p_c)$ for a sysstem of size $L = 512$