

# Fractal Structure of Rainforests

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## 1 INTRODUCTION

Recent research in forest ecology has increasingly focused on the concept of self-organization and criticality in forest ecosystems, particularly in the context of gap dynamics in rainforests [1–4]. This research has revealed that non-linear systems, operating far from equilibrium and possessing extended spatial degrees of freedom, frequently transition spontaneously into a state characterized as “self-organized critical”. The pioneering work by Solé and Manrubia has been instrumental in this area. In this work, a paradigm to study fractal properties of low-canopy gaps in rainforests is established. The study focuses on two parts. Firstly, the examination of the Barro Colorado Island (BCI) rainforest in Panama. Then they formulate a *cellular automaton* that allows for a more general study of a rainforest model. They propose that the appearing patterns are indicative of a self-organized critical state. Moreover, they argue that this self-organized critical state should exhibit fractal behaviour because, otherwise, if it were a non-fractal distribution, we would be referring to a state that follows a non-arbitrary order. This distinction arises because they posit that the non-linear dynamical process of gap formation in forest ecosystems can give rise to the formation of fractal structures, treefall disturbances are important for the persistence of the ecosystem. Therefore, fractality would result from the fact that the distribution of gaps and rainforest does not adhere to a traditional geometry. Alternatively, it could stem from the jungle exhibiting self-similarity across different areas at a fixed scale.

In order to characterize this patterns, one can measure the fractal dimension. For a *self-similar* fractal composed of  $m$  copies of itself scaled by a factor  $s$ , the

fractal dimension  $D$  is defined as [5, p. 413]:

$$D = \frac{\log m}{\log s}. \quad (1)$$

However, most fractals are not self-similar and generalizations of this definition are needed. In fact, many definitions have been proposed (see [6] for a further details). In this work, we will consider the *Box Dimension* definition [5, p. 416]:

$$D = \lim_{\epsilon \rightarrow 0} \frac{\log N(\epsilon)}{\log(1/\epsilon)}, \quad (2)$$

where  $N(\epsilon)$  is the number of boxes of size  $\epsilon$  needed to cover the fractal, and if the value of  $D$  were a non-integer number, we would be dealing with a geometry that cannot be analyzed using “traditional” tools. More generally, we consider the **Correlation Dimension of Order  $q$**  defined by<sup>1</sup> [1, p. 32]:

$$D_q := \lim_{\epsilon \rightarrow 0} \frac{1}{(q-1)} \frac{\log[X(q)]}{\log(\epsilon)}, \quad q \in \mathbb{R}, \quad (3)$$

where:

$$X(q) := \sum_{i=1}^{N(\epsilon)} p_i^q. \quad (4)$$

Here,  $p_i$  is the probability that a box of size  $\epsilon$  is populated. This dimension provides us with a wealth of information about the geometric structure of the fractal set.

On the other hand, since real fractals are in fact multifractals [1, p. 34], we further consider the **Fractal Spectrum Dimension**  $f(\alpha)$  [7]:

$$f(\alpha(q)) = q\alpha(q) + D_q(1-q), \quad (5)$$

where  $\alpha(q)$  is the diverging exponent defined by  $p_k \approx \epsilon_k^\alpha$ , which can be determined through [7]:

$$\alpha(q) = \frac{d}{dq} [(q-1)D_q]. \quad (6)$$

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<sup>1</sup>Notice how for  $q = 0$  we re-obtain eq. (2).

This report will delve into their findings, specifically analysing the forest game model they introduced and its implications for understanding rainforest ecology.

## 2 FOREST GAME

The *Forest Game*, a cellular automaton model developed by Solé and Manrubia, serves as a crucial tool for simulating the dynamics of forest gaps and their macroscopic spatial regularities. This model, based on simple rules of tree growth, competition, and death, illustrates how complex patterns emerge from straightforward processes.

The Forest Game is defined by a two-dimensional square lattice of  $L \times L$  points together with periodic boundary conditions [3, p. 533]. Each cell  $(i, j)$  represents a tree of height  $S_n(i, j)$  at the time step  $n$ . The tree height is bounded by  $S_0 \leq S_n(i, j) \leq S_c$ , where  $S_0$  is the minimum and  $S_c$  is the maximum tree size. The game's rules are as follows:

1. **Birth rule:** A tree of size  $S_0$  is born with probability  $p_b$  in each empty cell.
2. **Growth rule:** A tree of size  $S_n(i, j)$  will grow to:

$$S_{n+1}(i, j) = S_n(i, j) + \Delta n(i, j), \quad (7)$$

where  $\Delta n(i, j)$  is the Heaviside function defined as:

$$\Delta n(i, j) := \Theta \left[ \mu - \frac{\gamma}{8} \sum_{\langle r, s \rangle} S_n(r, s) \right]. \quad (8)$$

Here,  $\mu$  is a parameter and  $\gamma$  is the interaction strength. The sum is over the 8 nearest neighbours of the cell  $(i, j)$ , that is, we consider the [Moore neighbourhood](#).

3. **Death rule:** A tree of size  $S(i, j)$  dies with probability  $p_d$  or if  $S(i, j) \geq S_c$ .
4. **Gap formation rule:** A dying tree will form a gap of radius  $R$  if the size of neighbours (not necessarily the nearest ones) is less than or equal to the size of the tree. That is [2, p. 51]:

$$\sum_{B(R)} S_n(r, s) \leq S_n(i, j) \quad (9)$$

where  $B(R)$  is the neighbourhood of radius  $R$  around the cell  $(i, j)$ .

The significance of the Forest Game lies in its ability to replicate observed natural phenomena using a minimal set of rules. By adjusting parameters such as tree growth rate, death probability, and competition, the model can exhibit a variety of patterns that resemble real-world forest dynamics. The game's results support the idea that rainforest ecosystems might be operating near a critical point, where a small change could lead to significant transformations in the forest structure. This insight has profound implications for understanding the stability and biodiversity of rainforests, providing a theoretical framework for future ecological studies and conservation strategies.

### 2.1 Numerical Implementation

Regarding the computational implementation, we consider a random forest as the initial state of the system. We have taken the following parameters by default  $S_0 = 0.1$ ,  $S_c = 30$ ,  $\mu = \gamma = 1$  and  $p_b = 0.3$ .

In order to compute the fractal measures, we have followed the paper's methodology calculating  $X(q)$  by [2, p. 51]:

$$X_q(\epsilon) = \sum_{j=1}^{j_{\max}} N(j) \left( \frac{j}{N_P} \right)^q. \quad (10)$$

where  $N(j)$  is the number of boxes containing  $j$  gap points,  $N_P$  is the total number of gap points and  $j_{\max}$  is the maximum number of gap points in a box. In this manner, we have taken  $j_{\max} = 16$  and  $\epsilon = 1/20$ . The code is available in [appendix A](#).

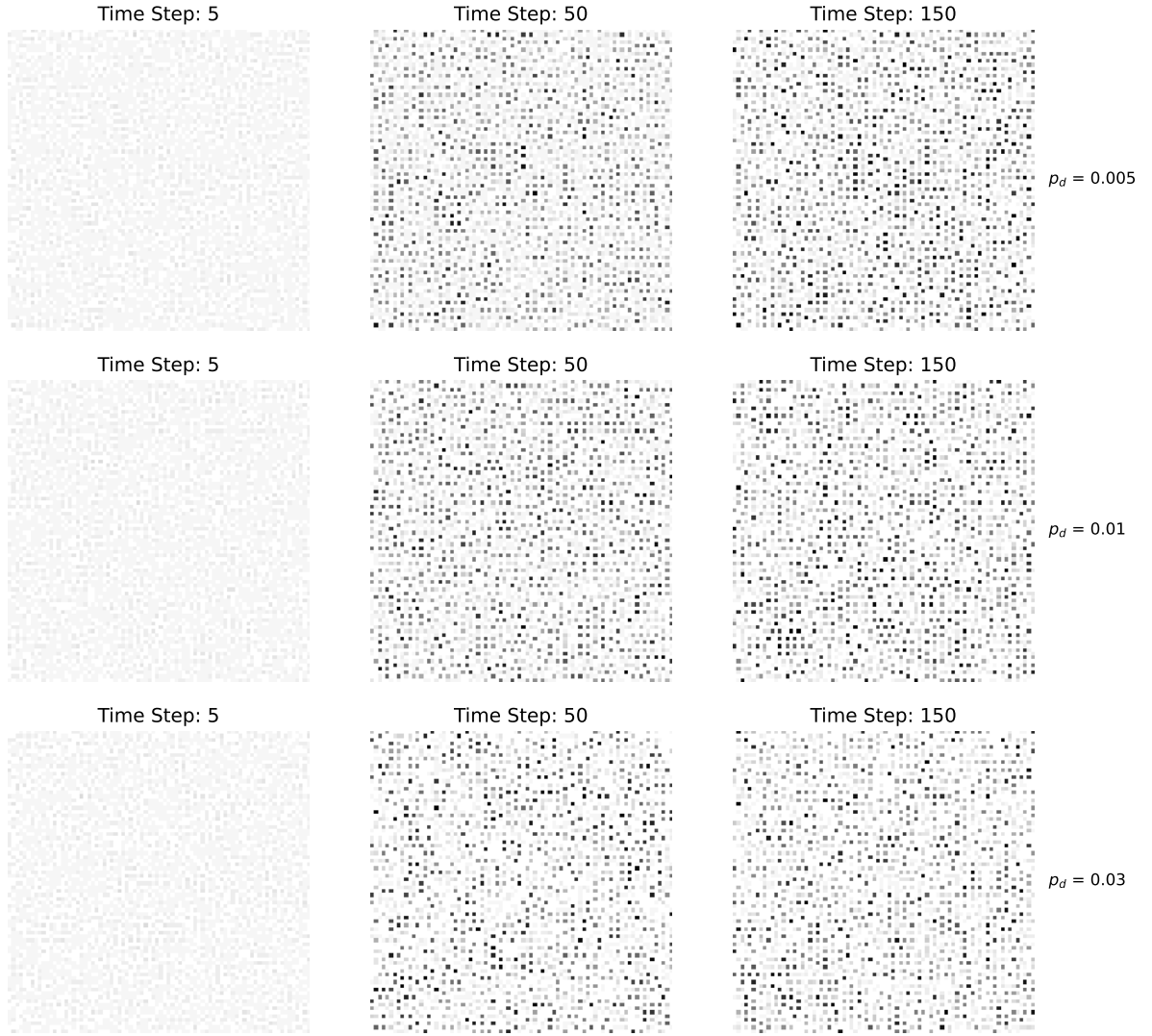


Figure 1: Images of the forest simulation according to the automaton presented earlier for different times and different death probabilities ( $p_d = 0.005, 0.01, 0.03$ ) are provided. As can be observed, they all start from a more or less empty forest with low density, and over time, they form the distribution of trees and gaps that we can see. It can be observed that with the increase in the death probability, for the last time, there are fewer darker black points, indicating lower density, and more white gaps. This is directly related to the fractal behavior of this system and will be discussed later. However, we can see how this is a easily recognizable phenomenon in the images of the forest. A lattice of  $80 \times 80$  was used, and the birth probability was  $p_b = 0.3$

### 3 RESULTS

#### 3.1 Biomass Evolution

In this section, we will present the results obtained from the Forest Game. We will start by measuring the *biomass* evolution in time<sup>2</sup>. Then, we will compute the fractal dimension and spectrum dimension for an equilibrium state.

We define the biomass at time step  $n$  by:

$$B_n = \sum_{i,j}^L S_n(i,j). \quad (11)$$

That is, we sum the size of all the trees in the lattice at time step  $n$ . In fig. 2 we show the evolution of the biomass for  $L = 40$  in  $t \in \{1, \dots, 700\}$ . We can see that initially, the biomass increases significantly, but after  $n \approx 50$  time steps, it reaches an equilibrium state. At this point, the rainforest attains an attractor of the self-organized critical state, that is, it fluctuates around a

<sup>2</sup>Time here is discrete, representing either years or seasons, depending on study of interest.

constant value. These fluctuations arise from the non-linearity of the gap formation mechanism.

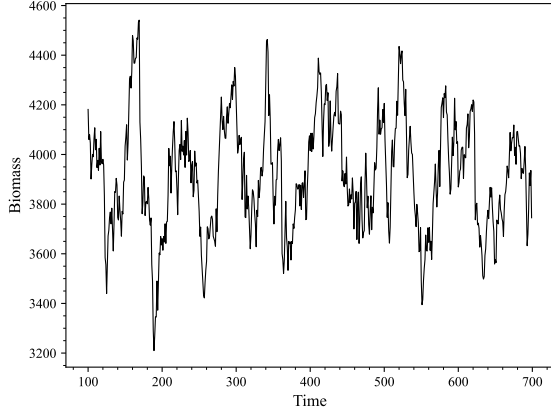


Figure 2: Biomass Dynamics. We compute the biomass for  $L = 40$  with  $p_d = 0.01$  in  $t \in \{1, \dots, 700\}$  via eq. (11).

Furthermore, we have analysed the fourier spectrum from  $n = 100$  to ensure the equilibrium state. We observe in fig. 3 that the behaviour of the spectrum scales like  $f^{-\beta}$ , with a  $\beta$  value falling within the range [missing value for us].

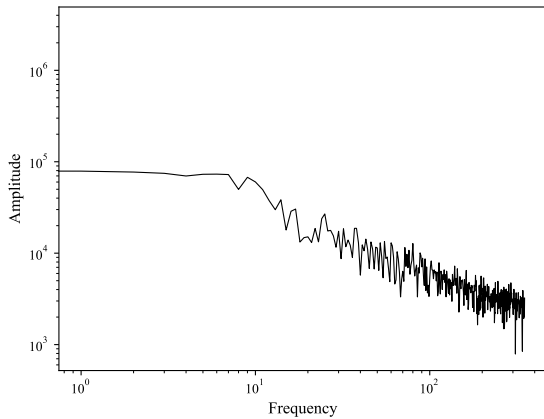


Figure 3: Fourier Transform of the Biomass. For the same values as in fig. 2. The obtained  $\beta$  value is 0.833. Near to the value of the reference [3]

### 3.2 Fractal Structure

We now study the fractal structure of the rainforest. We have computed the fractal correlation dimension and the fractal spectrum dimension for<sup>3</sup>  $p_d \in \{0.005, 0.01, 0.03\}$  and  $L = 80$ . In fig. 4 we show  $D_q$  for  $q \in [-8, 8]$  (RHS plot). We observe that as the mortality rate increases, the fractal dimension tends towards

2. This implies that as the likelihood of mortality and gap formation rises, the critical state of the system becomes more random. All these values are close to those obtained from a similar study on a real rainforest. In the case of BCI [1], the results indicate a fractal dimension of around 1.86. In our case, we observe a certain dependence of the fractal dimension on the death probability, which increases as the probability of death increases. The average results obtained for the fractal dimension of the cases simulated by the automaton are as follows: for  $p_d = 0.005$ ,  $D_0 = 1.81 \pm 0.02$ , for  $p_d = 0.01$ ,  $D_0 = 1.90 \pm 0.01$ , and, for  $p_d = 0.03$ ,  $D_0 = 1.96 \pm 0.02$ . In all cases, we obtain values less than 2, indicating a fractal structure. The dependence on death probability suggests that as the probability of a tree dying and forming a gap increases, the distribution is affected and approaches an integer value of 2. We could hypothesize that this is the limit when the death probability is maximum, effectively killing the forest and completely losing any fractality. Furthermore, with the figure 4, we can obtain values for other important magnitudes in the study of fractals, such as the information dimension ( $D_1$ ) or correlation dimension ( $D_2$ ). Therefore, once this relationship is detected, it is interesting to view this model as a tool for the calculation and control of forest health.

Therefore, we observe that our automaton can predict the distribution of a real rainforest and its fractal properties.

<sup>3</sup>These are typical rates for BCI rainforest [8]

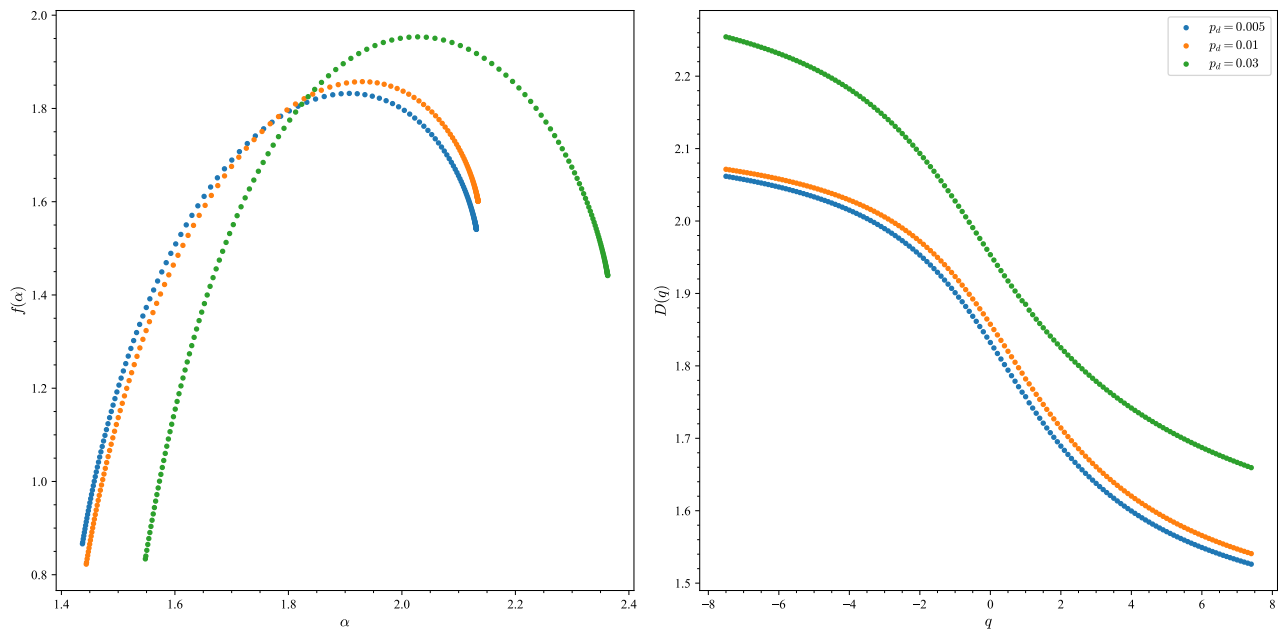


Figure 4: a) Multifractal spectrum and b) Fractal dimension of the forest for different values of the death probability. A lattice of 80x80 was used, and the birth probability was  $p_b = 0.3$

#### 4 CONCLUSIONS

The present report has served to rework the results obtained in previous works with the intention of highlighting the relevance of such approaches, ranging from the study of fractals to applications in ecological scenarios. Through the formulation of an automaton with non-linear rules, a critical state has been achieved, manifesting behaviors such as self-organization or multifractality in the distribution of gaps. However, not only does the gap distribution exhibit distinct behavior, but also the biomass, showcasing temporal self-similarity. Nevertheless, this doesn't necessarily represent the limit of the automaton, as many of the conditions that constitute it are limited or simplistic.

In the current context, concerns about climate change and its impact on the biosphere could prompt a revival of such approaches. Exploring the relationship between an increase in mortality and its potential effects on the fractal geometry of clearings in a forest becomes pertinent. However, with the rules established so far, we have achieved a system that self-organizes to reach a critical state, displaying spatial patterns that are self-similar.

#### References

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## A CODE

```
1 def rainforest(L, S0, p):
2     """Defines a random initial state of the forest. That is a square
3         ↳ lattice of size LxL with a probability p of having a tree of size
4         ↳ S0 in each cell.
5
6     Args:
7         L (int): Lattice size
8         S0 (float): Minimum tree size
9         p (float): Probability of having a tree of size S0 in each cell
10
11     Returns:
12         np.array((L,L)): Lattice with the initial state of the forest
13     """
14     lattice = np.zeros((L, L))
15
16     # Create a mask where each cell has a probability (p) of being True
17     growth_mask = np.random.random((L, L)) <= p
18
19     # Set cells in the lattice to S0 where the growth mask is True
20     lattice[growth_mask] = S0
21
22     return lattice
23
24 def birth(lattice, p_birth, S0):
25     """Defines the birth rule for the forest. That is, a tree of size S0 is
26         ↳ born with probability p_birth in each empty cell.
27
28     Args:
29         lattice (np.matrix): forest lattice
30         p_birth (float): probability of birth
31         S0 (float): minimum tree size
32
33     Returns:
34         np.matrix: forest lattice after applying the birth rule
35     """
36     # Create a mask for empty cells
37     empty_mask = (lattice == 0)
38
39     # Generate a random array of the same shape as the lattice
40     random_values = np.random.random(lattice.shape)
41
42     # Create a birth mask where cells are empty and the random value is
43         ↳ below the birth threshold
44     birth_mask = empty_mask & (random_values <= p_birth)
45
46     # Set cells in the lattice to S0 where the birth mask is True
47     lattice[birth_mask] = S0
```

```

44
45     return lattice
46
47 def heaviside_vectorized(mu, gamma, lattice):
48     """heaviside_vectorized calculates the Heaviside function values for
49         ↳ each cell in the lattice. This accounts for  $\Delta n_{ij}$  in the
50         ↳ growth rule.
51
52     Args:
53         mu (float): parameter
54         gamma (float): interaction strength
55         lattice (np.matrix): forest lattice
56
57     Returns:
58         np.matrix: Heaviside function values for each cell in the lattice
59     """
60     # Define the kernel for neighbor sum
61     kernel = np.array([
62         [1, 1, 1],
63         [1, 0, 1],
64         [1, 1, 1]
65     ])
66     neighbors_sum = convolve2d(lattice, kernel, mode='same', boundary='wrap'
67         ↳ ) # https://docs.scipy.org/doc/scipy/reference/generated/scipy.
68         ↳ signal.convolve2d.html
69
70     # Calculate the Heaviside function values
71     heaviside_values = mu - (gamma / 8) * neighbors_sum
72     heaviside_values[heaviside_values < 0] = 0 # Set negative values to zero
73
74     return heaviside_values
75
76 def growth(lattice, S0, Sc, mu, gamma):
77     """Defines the growth rule for the forest.
78
79     Args:
80         lattice (np.matrix): forest lattice
81         S0 (float): minimum tree size
82         Sc (float): maximum tree size
83         mu (float): parameter
84         gamma (float): interaction strength
85
86     Returns:
87         np.matrix: forest lattice after applying the growth rule
88     """
89
90     # Create a mask for cells that are eligible for growth
91     growth_mask = (lattice <= Sc) & (lattice >= S0)

```



```

89     # Calculate the Heaviside values for the entire lattice
90     heaviside_values = heaviside_vectorized(mu, gamma, lattice)
91
92     # Apply growth only to eligible cells
93     lattice[growth_mask] += heaviside_values[growth_mask]
94
95     return lattice
96
97 def find_nth_moore_layer(lattice, i, j, n):
98     """Find the nth layer of Moore neighbors in a square lattice, i.e, a
99         ↳ square  $(2n+1) \times (2n+1)$  excluding the inner  $(2n-1) \times (2n-1)$  square. (
100         ↳ see https://mathworld.wolfram.com/MooreNeighborhood.html for more
101         ↳ details).
102
103     Args:
104         i (int): x coordinate of the cell.
105         j (int): y coordinate of the cell.
106         n (int): n-th Moore neighborhood.
107         grid_size (int): The size of the grid.
108
109     Returns:
110         list: list of tuples with the coordinates of the nth Moore neighbors
111             ↳ .
112     """
113     L = lattice.shape[0]
114     neighbors = []
115     for dx in range(-n, n+1):
116         for dy in range(-n, n+1):
117             if max(abs(dx), abs(dy)) == n: # This condition excludes inner
118                 ↳ squares
119                 # Wrap around the grid using modulo operator to maintain
120                 ↳ periodic boundaries
121                 neighbors.append(((i + dx) % L, (j + dy) % L))
122
123     return neighbors
124
125 def find_neighbors_at_R(lattice, i, j, R):
126     """Finds the neighbors of a cell exactly at a radius R. Here we consider
127         ↳ euclidian distance.
128
129     Args:
130         lattice (np.matrix): forest lattice
131         i (int): x coordinate of the cell
132         j (int): y coordinate of the cell
133         R (float): radius
134
135     Returns:
136         list: list of tuples with the coordinates of the neighbors
137     """

```

```

131 n_rows, n_cols = lattice.shape
132 max_offset = int(np.ceil(R)) # Maximum offset to consider
133
134 # Create arrays for row and column offsets
135 row_offsets = np.arange(-max_offset, max_offset + 1)
136 col_offsets = np.arange(-max_offset, max_offset + 1)
137
138 # Calculate the grid of distances considering periodic boundaries
139 row_distances = np.minimum(np.abs(row_offsets), n_rows - np.abs(
    ↪ row_offsets))**2
140 col_distances = np.minimum(np.abs(col_offsets), n_cols - np.abs(
    ↪ col_offsets))**2
141 grid_distances = np.sqrt(row_distances[:, np.newaxis] + col_distances)
142
143 # Find neighbors exactly at the radius R
144 neighbor_mask = grid_distances == R
145 neighbor_offsets = np.argwhere(neighbor_mask) - max_offset
146
147 # Calculate neighbor coordinates with periodic boundary conditions
148 neighbors = np.mod(np.array([i, j]) + neighbor_offsets, [n_rows, n_cols
    ↪ ])
149
150 return list(map(tuple, neighbors))
151
152 def generate_radius_list(max_radius):
153     """Generates a list of radii to consider from 1 to max_radius.
154
155     Args:
156         max_radius (float): maximum radius
157
158     Returns:
159         np.array: array of radii
160     """
161     radius_set = set()
162     for x in range(max_radius + 1):
163         for y in range(x + 1):
164             r2 = x**2 + y**2
165             if r2 <= max_radius**2:
166                 radius_set.add(r2)
167     return np.sqrt(sorted(radius_set))[1:] # Remove the first element (R =
    ↪ 0)
168
169 def gap_form(lattice, i, j, radius_list, Moore = False):
170     """Applies the gap formation rule to a cell. That is, if the sum of the
    ↪ trees in the neighborhood is less than or equal to the tree in the
    ↪ cell, then the trees in the neighborhood die. The radius of the
    ↪ neighborhood is defined by the inequality.
171
172     Args:

```

```

173     lattice (np.matrix): forest lattice
174     i (int): x coordinate of the cell
175     j (int): y coordinate of the cell
176     radius_list (np.array): array of radii to consider
177     Moore (bool, optional): Flag to choose which distance to use in gap
    ↪ formation. Defaults to False.

178
179     Returns:
180         np.matrix: forest lattice after applying the gap formation rule
181     """
182     S = lattice[i, j]
183     S_nn = 0
184     for R in radius_list:
185         neighbors = find_nth_moore_layer(lattice, i, j, R) if Moore else
    ↪ find_neighbors_at_R(lattice, i, j, R)
186         xs, ys = list(zip(*neighbors)) # From [(x1,y1), ...] to [x1, ...], [
    ↪ y1, ...]
187         S_nn += np.sum(lattice[xs, ys])
188         if S_nn <= S:
189             lattice[xs, ys] = 0
190         else:
191             break
192
193     return lattice
194
195 def death(lattice, pd, Sc, Moore = False):
196     """Defines the death rule for the forest. A given tree dies randomly
    ↪ with probability pd. A tree that reaches the maximum size Sc also
    ↪ dies.

197
198     Args:
199         lattice (np.matrix): forest lattice
200         pd (float): probability of death
201         Sc (float): maximum tree size
202         Moore (bool, optional): Flag to choose which distance to use in gap
    ↪ formation. Defaults to False.

203
204     Returns:
205         np.matrix: forest lattice after applying the death rule
206     """
207
208     L = lattice.shape[0]
209
210     # Create masks for cells above Sc and cells that die randomly
211     above_sc_mask = (lattice >= Sc)
212     death_mask = np.random.random(lattice.shape) <= pd
213     gap_mask = above_sc_mask | (lattice > 0) & death_mask # combined mask
214     gap_indices = np.argwhere(gap_mask) # array of indices of cells that
    ↪ will form gaps

```

---

```

215
216     # Generate a list of radii to consider
217     max_radius = L // 2 # Maximum radius is half the size of the grid
218     radius_list = [i for i in range(1, max_radius + 1)] if Moore else
        ↪ generate_radius_list(max_radius)
219
220     # Apply gap_form to the necessary cells
221     for i, j in gap_indices:
222         lattice = gap_form(lattice, i, j, radius_list, Moore)
223
224     # Set to zero the cells that died
225     lattice[above_sc_mask] = 0
226     lattice[(lattice > 0) & death_mask] = 0
227
228     return lattice

```