

## Cellular Automata Assignment: Diffusion-Limited Aggregation

Diffusion-limited aggregation (DLA) is the process whereby particles undergoing a random walk due to Brownian motion cluster together to form aggregates of such particles. This model was proposed to simulate certain types of aggregation, for instance metal ions diffusing through a fluid and sticking to a charged electrode. “Diffusion” because the particles forming the structure, also called *Brownian tree* or *cluster*, wander around randomly before attaching themselves (“aggregating”) to the structure. “Diffusion-limited” because the particles are considered to be in low concentrations and therefore don’t interact together. Other examples can be found in non-living and living nature, e.g. mineral deposition, snowflake growth, lightning paths or corals growth.

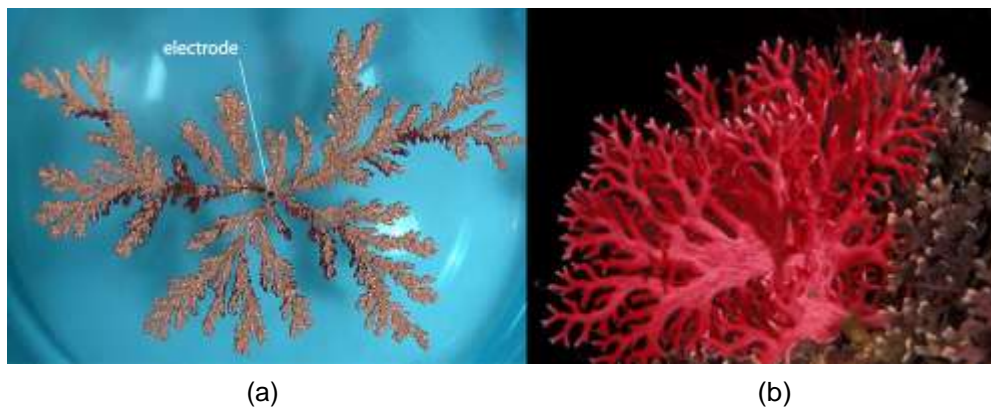


Figure 1: (a) DLA structure grown from a copper sulfate solution in an electrodeposition cell. (b) Red coral *Errina novaezealandiae* in the Te Awaatu Marine Reserve in Fiordland.

## Cellular Automaton

The goal of this exercise is to implement a two-dimensional CA that mimics a DLA process. Launch Octave (or Matlab), open *ca.m* and run it. Have a look at the content of the file to become familiar with the main variables (top part). The automaton space is formed by an array of cells, which size is given by the integer parameters *nx* and *ny*. *The x-axis represents the vertical axis and the y-axis the horizontal axis*. Motionless, non-interacting particles (blue cells) are initially present in the CA space. Their density can be changed through the variable *particlesDensity*.

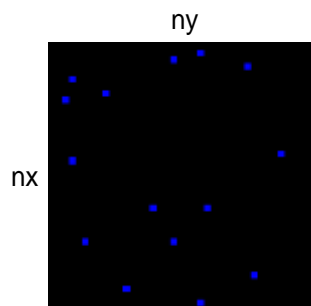


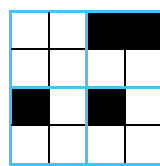
Figure 2: First run of *ca.m*. The above automaton space is defined by an array of 40x40 cells. Motionless particles (in blue) are initially present in the environment.

## Implementation of a Pseudo Brownian Motion

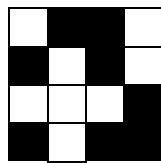
This section describes how to obtain particles that undergo a random walk. These particles are assumed to be in low concentrations, so no interaction between them are modeled.

Start by setting in the code the number of time steps to  $T=7000$ . Run `ca.m` and observe the motion of the particles. Note that you can speed up or slow down the walk of the particles by adjusting the parameter `delay`. You can also change the size of the automaton space (`nx` and `ny` must be divisible by 2) or resize the window to get a better visibility. Try to qualify the movement of the particles. What kind of neighborhood is implemented? Does the observed motion feature some randomness? If it's possible, find the deterministic and random components of the motion.

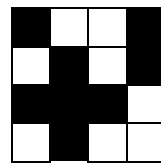
Try to understand what does the code between the tags "Pseudo Brownian motion" and "End". The automaton space is divided into blocks of  $2 \times 2$  cells inside which the position of the particles is updated. At each time step, two random, complementary matrices containing "0" and "1" elements are generated (`cw` and `ccw`). With your understanding of the code, find and draw the updated position of the particles at time  $t+1$  for the initial configuration of the following  $4 \times 4$  CA. Consider vectors `xind = yind = [1, 3]` which define the indexes of the upper-left cell of each block, and black cells below to represent particles and/or "1" matrix elements. (Hint: only the elements in `cw` that are defined by `xind` and `yind` are used, that are here `cw(1, 1)`, `cw(3, 1)`, `cw(1,3)`, and `cw(3,3)`.)



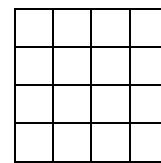
CA space (t)



cw



ccw



CA space (t+1)

You observe that the motion of a particle is actually limited to the block it originally belongs to. Which simple modification could be adopted to allow the particles to move across the entire automaton space?

What happens if you replace in the code `s=0` by `s=mod(t,2)`? Run the modified m-file, observe and comment the new motion of the particles.

## Growth of Brownian Trees

This section aims to simulate the growth of *Brownian trees* (or "*sticky*" clusters) using the cellular automaton developed at the end of the previous section. Beforehand, operate the following modifications where required. Then run `ca.m`.

- $nx = ny = 200$
- $particlesDensity = 0.1$
- $T = 7000$
- $delay = 0$

- `enableBrownianTree = 1`

At time  $t=0$ , the Brownian tree is composed of a single, yellow cell placed at the center of the automaton space. Observe how the tree grows from moving particles that stick to it. What is the condition required for a blue particle to stick and become part of the tree? Figure 3 illustrates the growth of two trees resulting from two successive runs of *ca.m*.

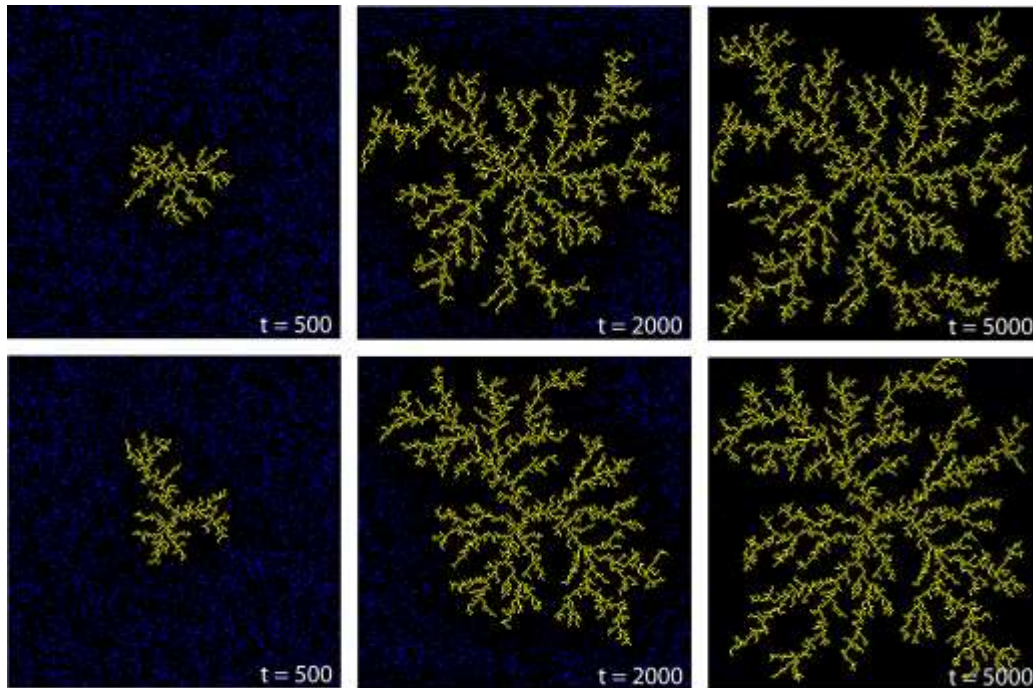
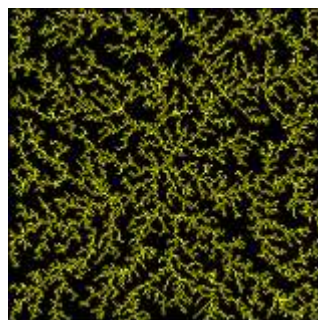


Figure 3: Growth of Brownian trees using a cellular automaton that mimics a diffusion-limited aggregation (DLA) process. The two different trees are the result of two successive runs of *ca.m* (with identical parameter values). Snapshots are taken at time  $t=500$ , 2000 and 5000 iterations.

Have a look at the structure of the following tree. What distinct it from the two ones displayed in Figure 7 ( $t=5000$ )? Try different values for the parameter responsible of this change and observe how it affects the structure of the tree.



At which part of the tree does the growth occur mainly? How do you expect the size of a Brownian tree (number of cells that compose it) to vary in function of the time?