IN-CLASS WORK: SCIENTIFIC TALKS & XFIG & DLA (CONTINUED)

Scientific Talks: Take a few minutes to think about very good scientific talks. Which? What made the talks so good?

Xfig Intro

I will guide you through the following main commands of xfig, which is drawing tool:

- To get started: Type on the command line: xfig & This will open a new window.
- drawing tools: background grid, circle, line, text, picture, grouping, scaling, copying, editing.
- To save an xfig session use File → SaveAs and give your xfig-file a name ending with .fig. You can get back to this session any time on the command line with xfig filename.fig & or within xfig with File → Open.
- To make an eps-file out of your figure use File → Export, make sure to choose "EPS (Encapsulated Postscript)" and choose the same filename but with the ending .eps. This eps-file can then be included in your latex file for the paper. (Later into the course I will also show you a variation of latex, latex beamer, which we will use to make talk-slides. You will be able to use the same eps-files for the paper and for the talk and therefore your work on the eps-files for your paper will be very handy for your talk preparation.)

Optional: Comment for Advanced xfig Users who like Latex:

Figure(s) for Model Section To practice a bit, start working on a figure for your talk, which is helpful to explain your model.

DLA (continued)

You find the solutions to last's class in-class work in

~kvollmay/share.dir/inclass2023.dir/classfractal*.py

I will walk us through ~kvollmay/share.dir/inclass2023.dir/classfractal5b.py

6. Stick to Cluster

Next we will work on rule V, the sticking of a random walker particle to the cluster, if the random walker is next to a cluster cell. We use "von Neumann neighbors", which means a neighbor cell up,down,left or right.

6a. Next add to your program of 5b. that whenever the random walker is next (left, right, top, bottom) to a particle of the cluster then the random walk stops (walkstop update). Use the flow chart to decide where to add the necessary lines. If you have kept the print(x,y,LATMID,r) from 5a (and the same seed) then you can check that your program is working right.

6b. Now add to your program of 6a that you also have an integer variable npart which is initialized to be npart=1 and gets increased by one whenever a particle sticks to the cluster. Also update lattice whenever a particle sticks. Whenever a particle sticks to the cluster, you also need to check if RMAX has grown and if so, then you need to update RMAX. Add this to your program.

7. Finish Program: Loop Over Particles

Now you are ready to finish your DLA program! Add to your program a while loop over particles. Condition for this while loop are both that the wanted number of cluster particles NPARTMAX has not yet been reached and that the radius for the stopping of a too far out random walker fits into the lattice. Use the flow chart of class to decide where to add this while-loop. Use the constants LATSIZE=100, NPARTMAX=50. Comment out the printing of (x,y,LATMID,r), but print the resulting lattice at the end (so after the particle-loop). In case you would like not to print the complete lattice, you may use the following commands

```
plt.imshow(lattice[int(LATMID-RMAX-2):int(LATMID+RMAX+2),\
    int(LATMID-RMAX-2):int(LATMID+RMAX+2)],interpolation='nearest')
plt.savefig('frame7.pdf')
```