



Numerical Galaxy Formation and Cosmology

Lecture/Exercise 6 - 1st Example: A cosmological simulation with Gadget-2

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Performing a cosmological simulation with Gadget-2

- main steps for creating initial conditions:
 - compile initial condition code (e.g. N-GenIC)
 - set parameters for initial condition code
 - create initial condition files
- main steps for running the simulations:
 - compile simulation code (e.g. Gadget-2)
 - set parameters for the simulation code
 - run simulation
- read simulation output and perform analysis (in two *weeks*)

Creating initial conditions

- download N-GenIC (by Volker Springel):

- ➔ `wget http://www.mpa-garching.mpg.de/gadget/n-genic.tar.gz`

- ➔ `tar -xvf n-genic.tar.gz`

- compile N-GenIC:

- requires MPI, FFTW and GSL libraries

- ➔ `module load openmpi/intel/11.1/1.4.2`

- ➔ `module load fftw/2.1.5_double`

- ➔ `module load gsl/1.14`

- ➔ `module load intel/11.1`

should work on IoA machines,
you can put this in your .profile file

- edit Makefile: `SYSTYPE="OpenSuse"` and `MPICHLIB = -lmpi`

- compile with `make`

Creating initial conditions

- parameters for N-GenIC, change e.g.:

Nmesh	32	% This is the size of the FFT grid used to
Nsample	32	% sets the maximum k that the code uses,
		← controls resolution used for initial displacements and velocities
TileFac	2	% Number of times the glass file is
		← controls particle number (for 16 ³ “glass” file -> 32 ² particles)
Box	60.0	% Periodic box size of simulation [Mpc/h]
UnitLength_in_cm	3.085678e24	% defines length unit of output (in cm/h), for Mpc/h
		← set box size and length units
Redshift	99.0	% Starting redshift
OmegaBaryon	0.045	% Baryon density (at z=0)
Sigma8	0.83	% power spectrum normalization
WhichSpectrum	1	% "1" selects Eisenstein & Hu spectrum,
		← set matter power spectrum
NumFilesWrittenInParallel	1	% limits the number of files that are

Creating initial conditions

- running N-GenIC:

- ➡ `mpiexec -np 2 ./N-GenIC ics.param` to run in parallel on two cores, or

- ➡ `./N-GenIC ics.param` to run as a serial code

- this creates the ICs file(s) in the output directory

Setting up the simulation

- download Gadget-2 (by Volker Springel):

- ➡ `wget http://www.mpa-garching.mpg.de/gadget/gadget-2.0.7.tar.gz`

- ➡ `tar -xvf gadget-2.0.7.tar.gz`

Setting up the simulation

- compile Gadget-2:

- edit Makefile, e.g:

```
< OPT    +=    -DUNEQUALSOFTENINGS
> #OPT    +=    -DUNEQUALSOFTENINGS

< OPT    +=    -DPMGRID=128
> OPT    +=    -DPMGRID=64

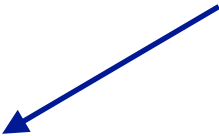
< #OPT    +=    -DDOUBLEPRECISION
< #OPT    +=    -DDOUBLEPRECISION_FFTW
> OPT    +=    -DDOUBLEPRECISION
> OPT    +=    -DDOUBLEPRECISION_FFTW

< OPT    +=    -DHAVE_HDF5
> #OPT    +=    -DHAVE_HDF5

< MPICHLIB =    -lmpich
> MPICHLIB =    -lmpi

< SYSTYPE="MPA"
> #SYSTYPE="MPA"
```

described in the Gadget2 user guide



same softening for all particles

size of PM grid (for real applications
should not be smaller than 256)

e.g. for computing with 64 bit floats
and when linking to double precision
FFTW

when not using HDF5 outputs

when using OpenMPI

turn off specific options for MPA
machines

- compile with `make`

Setting up the simulation

- setting the parameters for the simulation:
 - create a parameter file (changes compared to lcdm_gas.param):

```
InitCondFile      ../../N-GenIC/ICs/ics
OutputDir         ../sim/
SnapshotFileBase  snap
OutputListFilename parameterfiles/outputs.txt
```

set input and output file paths

```
SnapFormat        2
```

set format of snapshot files

```
TimeBegin         0.01      % z=99, Begin of the simulation
```

```
OmegaBaryon       0.045
```

```
BoxSize           60.0
```

should match values used in N-GenIC

```
CpuTimeBetRestartFile 14400.0 ; here in seconds
```

```
MaxSizeTimestep      0.01
```

```
UnitLength_in_cm      3.085678e24 ; 1.0 Mpc/h
```

```
SofteningGas         0.05
```

```
SofteningHalo        0.05
```

```
SofteningGasMaxPhys  0.05
```

```
SofteningHaloMaxPhys 0.05
```

gravitational softening

here: 50 comoving kpc/h (~2.7%
of mean inter particle distance)

- create a file with the desired output scale factors (filename as chosen in parameter file)

Running the simulation

- running Gadget-2:

➔ `mpiexec -np 2 ./Gadget2 parameterfiles/param.txt` **uses 2 cores**

- or usually using a job script which is submitted to the scheduler (e.g. for Darwin)

```
#!/bin/bash
#SBATCH -J large
#SBATCH -A DIRAC-DP044
#SBATCH -p sandybridge
#SBATCH --nodes=32
#SBATCH --ntasks=512
#SBATCH --time=36:00:00
#SBATCH --mail-type=ALL
#SBATCH --no-requeue

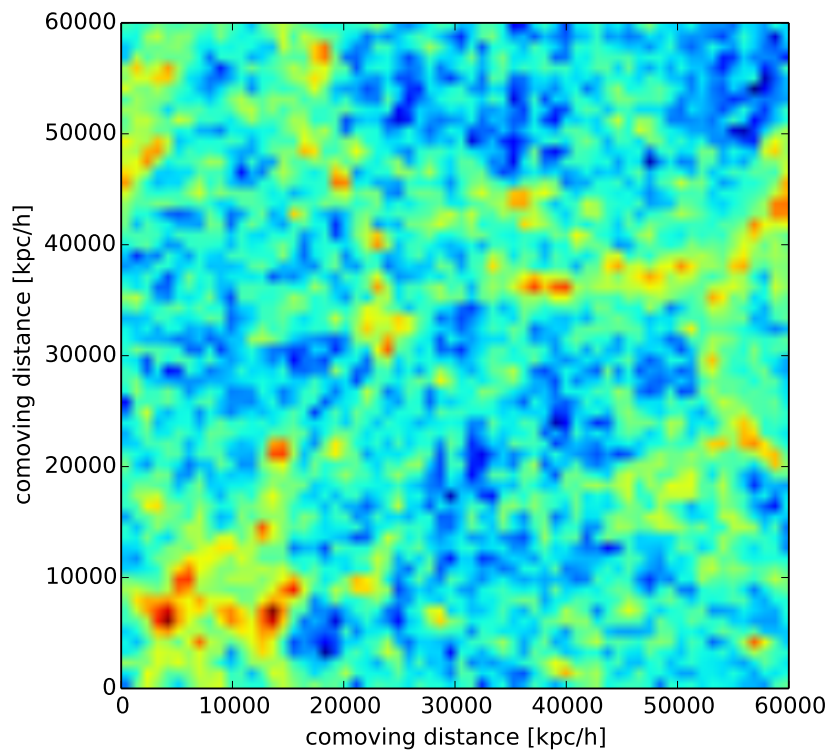
numnodes=$SLURM_JOB_NUM_NODES
numtasks=$SLURM_NTASKS
mpi_tasks_per_node=$(echo "$SLURM_TASKS_PER_NODE" | sed -e 's/^\([0-9][0-9]*\).*$/\1/')
np=$((numnodes*mpi_tasks_per_node))
workdir="$SLURM_SUBMIT_DIR"

# ... usually some more stuff here

cd $workdir
mpiexec -np $np ./Gadget2 parameterfiles/param.txt
```

submit job with: `sbatch job.sh`

Results - matter density field



our example: 32^3 particles,
~4 core minutes

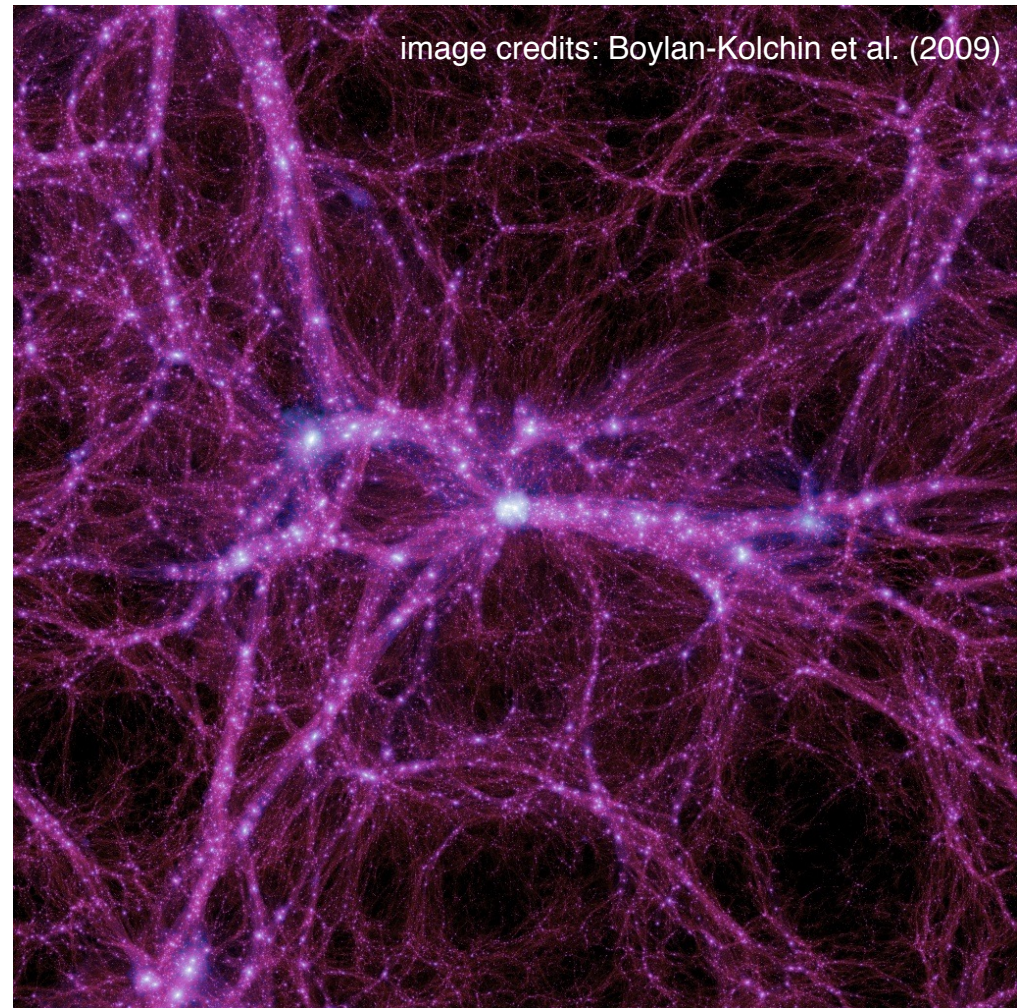


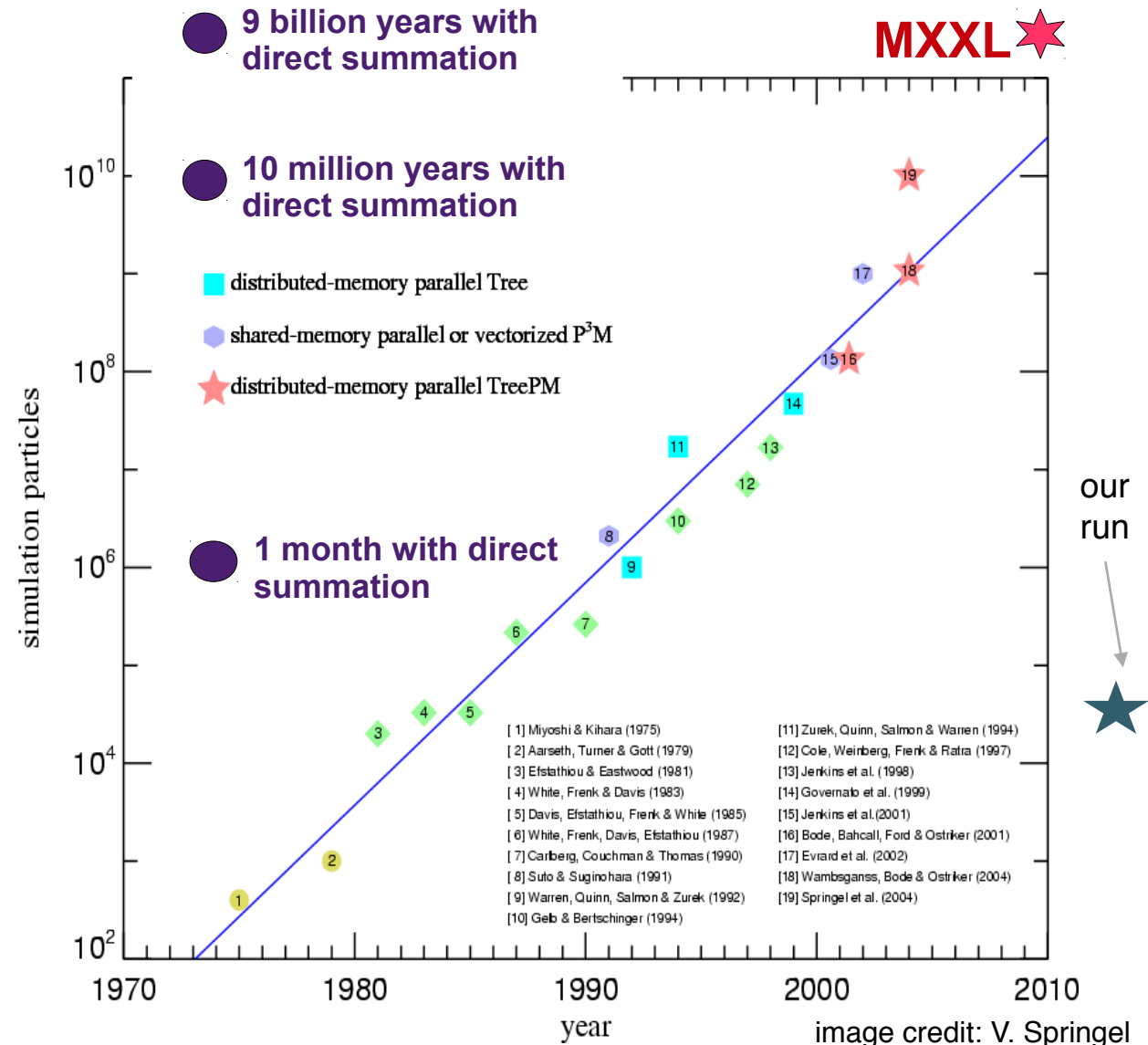
image credits: Boylan-Kolchin et al. (2009)

100 Mpc / h

Millennium II simulation: 2160^3 particles, 1.4 million core hours

Size of cosmological simulations over time

- computers double speed every 18 months (Moore's law)
- particle number in simulations doubles every 16-17 months
- only possible with algorithms that scale close to $\sim N$ (or $N \log(N)$)



Literature & Outlook

- “The cosmological simulation code GADGET-2”, V. Springel, 2005, MNRAS, 364, 1105, arXiv:astro-ph/0505010
- Gadget-2 user guide in “Documentation” sub-folder
- Next lecture:
 - ***Simulating a galaxy merger***
- In two weeks:
 - ***Analysing the simulation outputs of the cosmological box***
- ***Bonus exercise:***
 - ***Do a higher resolution run, e.g. 64^3 or 128^3***
(Do not forget to increase Nmesh, Nsample, TileFac when creating ICs, as well as PMGRID when compiling Gadget2. The softening should be reduced accordingly.)