

# UL HPC School 2017

## PS2: HPC workflow with sequential jobs (test cases on GROMACS, Java and Python)

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<http://hpc.uni.lu>

## Latest versions available on **Github**:



UL HPC tutorials:

<https://github.com/ULHPC/tutorials>

UL HPC School:

<http://hpc.uni.lu/hpc-school/>

PS2 tutorial sources:

[http://ulhpc-tutorials.readthedocs.io/en/latest/basic/sequential\\_jobs/](http://ulhpc-tutorials.readthedocs.io/en/latest/basic/sequential_jobs/)





# Summary

- 1 Introduction**
- 2 Pre-requisites
- 3 Exercise 1: Parametric execution of Gromacs
- 4 Exercise 2: Watermarking images in Python
- 5 Exercise 3: Advanced use case, using a Java program: "JCell"
- 6 Conclusion



## Main Objectives of this Session

- Run sequential, parametric programs on the clusters
- Learn how-to use our set of launcher scripts
- Submit jobs
- use the cluster monitoring tools
  - ↳ Ganglia
  - ↳ Monika & Drawgantt

### Read the full subject of this PS here

- <http://git.io/5cYmPw>



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# Getting started

## 1 Connect to the cluster(s)

```
(laptop)$> ssh {iris,gaia,chaos}-cluster
```

## 2 Send files

```
(laptop)$> rsync -avz local_directory {iris,gaia,chaos}-cluster:
```

## 3 Retrieve files

```
(laptop)$> rsync -avz {iris,gaia,chaos}-cluster:path/to/files  
local_directory
```

## 4 Submit jobs

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OAR on Chaos/Gaia

Slurm on Iris

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oarsub -l  
oarsub ./program

srun -p interactive -qos qos-interactive -pty bash  
sbatch program

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## Tutorial link

**This tutorial is available on github !**

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- <https://git.io/vHyh3>



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## Gromacs

### **GROMACS:** GROningen MAchine for Chemical Simulations

versatile package for molecular dynamics, primarily designed for biochemical molecules

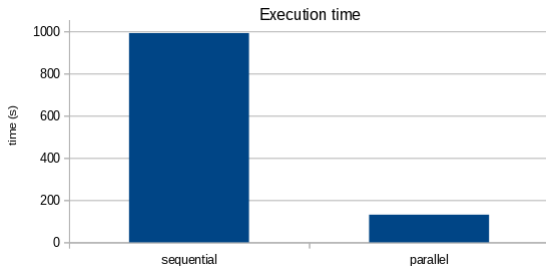
- very large codebase: 1.836.917 SLOC
- many applications in the package, several parallelization modes
- **mdrun**: computational chemistry engine, performing:
  - ↪ molecular dynamics simulations
  - ↪ Brownian Dynamics, Langevin Dynamics
  - ↪ Conjugate Gradient
  - ↪ L-BFGS
  - ↪ Steepest Descents energy minimization
  - ↪ Normal Mode Analysis
- **mdrun** - parallelized using MPI, OpenMP, pthreads and with support for GPU acceleration



## Comparison

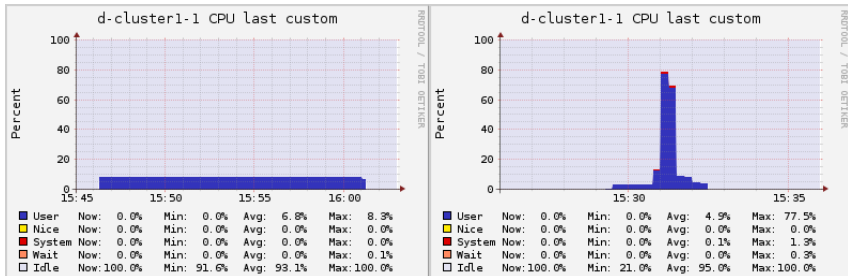
### 2 approaches

- Sequential (loop)
- Parallized (with GNU parallel)





# Comparison - Ganglia





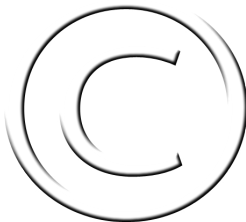
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## Watermark Application

- **Objective:** Apply a watermark to a given set of pictures
  - ↪ Simple Python script
  - ↪ Generic parallel launcher
  - ↪ Distribute the work on several nodes





## Source image





## Watermarked image





## Summary

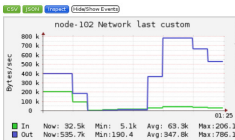
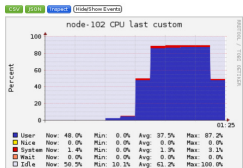
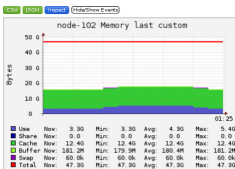
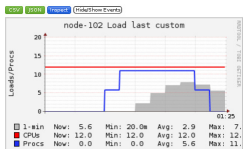
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## Jcell & cGAs

- **JCell**: a Java framework for working with genetic algorithms
  - ↳ Ex: Generational algorithm for the Combinatorial ECC problem
- Test the variations of these parameters:
  - ↳ *Mutation probability* and *Crossover probability*





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## Conclusion

- We have covered one of the most common workflow:
  - ↪ **parametric jobs**
- Our launchers can be improved!

### Perspectives

- Array jobs
- Best effort jobs
- Checkpoint/Restart mechanism



Thank you for your attention...

## Questions?

<http://hpc.uni.lu>

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