



## HPC cluster Initiation





### **i-Trop Presentation**



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### i-Trop Presentation





Requests form

https://bioinfo.ird.fr/index.php/cluster-fr/

- o Account
- o Software
- O Projects
- Incidents: contact <u>bioinfo@ird.fr</u>

### • Howtos:

https://bioinfo.ird.fr/index.php/tutorials-fr/howtosfor-hpc-cluster-itrop/

### • Slurm Tutorials:

https://bioinfo.ird.fr/index.php/tutorials-fr/slurm/

### • FAQ:

https://bioinfo.ird.fr/index.php/faq-fr/



### ARCHITECTURE



- A logical unit composed of several servers
- A powerful unique machine
- Allow to obtain high computing performance
- A bigger capacity storage
- More reliable
- A better ressources availability



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### **Cluster components**



### • Master Node

Handle ressources and jobs priorities

Computing nodes
 Resources (CPU or RAM memory)



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Handle ressources and jobs priorities

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 Resources (CPU or RAM memory)

# STORAGE



San
 Storage



### Architecture: components role

### • 1 Master node



bioinfo-master1.ird.fr

Role :

- Launch and prioritize jobs on computing nodes
- Accessible from the Internet
- Connection :

ssh login@bioinfo-master1.ird.fr



### Architecture: components role

### 1 Master node



bioinfo-master1.ird.fr

### • 32 nodes



nodeX X : 0..31 Role :

- Launch and prioritize jobs on computing nodes
- Accessible from the Internet
- Connection :

ssh login@bioinfo-master1.ird.fr

Role :

- Used by the master to execute jobs
- Not accessible from the Internet
- node0 to node31





### Step 1: Connection, srun





### Analyses step of the cluster

Connection to bioinfomaster1.ird.f r and resources reservation



Step 1 srun ou sbatch



### **Partitions**

Partitions	Use	Nodes RAM features	Nodes Core features
short	Short jobs < 1 jour	48 to 64 GB	12 cores
normal	Short jobs max 14 days	64 to 512 GB	12 to 112 cores
long	45 days > long jobs > 14 days	48 GB	12 to 24 cores
highmem	Jobs with memory needs	512GB	112 cores
supermem	Jobs with big memory needs	1TB	40 cores
gpu	Need for analyses on GPU cores	192GB	24 cpu et 8 GPU cores



- Partition to work on GPUs processors : basecalling, MiniOn etc..
- Restricted access to gpu\_account group
- Request access with arguments to do here:

https://itrop-glpi.ird.fr/plugins/formcreator/front/formlist.php



YES



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Rules	Partition	Tools example	comments
basecalling, demultiplexing, correction	gpu	medaka, guppy, machine learning tools	Restricted access
assembling >100G RAM	supermem	miniasm, flye, raven, smartdenovo	Target genome> 400 Mb (Rice genome doesn't need 100GB)
genomicsbd (gatk) > 100G RAM	supermem	GATK genomicsDB	Target genome of more than 400 Mb (>10 samples)
assemblings => 35G et < 100G RAM	highmem	miniasm, flye, raven, smartdenovo	Target genome between 100 and 400 Mb
Pops structure	long		
simulations	long		
metagenomic	normal	quiime2, frogs	
mapping	normal	bwa, minimap2, hisat2	Need a lot of cores not too many RAM Tool number of cores = number of cores to reserve
genotypage	normal	GATK haplotypecaller, samtools mpileup, bcftools	Need a lot of cores not too many RAM Tool number of cores = number of cores to reserve
stats	normal	R	
scripts test	short	bash, python, R	



### Architecture: elements role

### I Master node



bioinfomaster1.ird.fr

#### Role :

- Launch and prioritize jobs on compting nodes
- Accessible from the Internet

### • 32 Nodes



#### Rôle :

- Used by the master to execute jobs
- Not accessible from the Internet

### • 1 san



bioinfo-san.ird.fr (san)

#### Role :

- Store users data
- Accessible from the Internet
- To transfer data : via filezilla or scp



#### san server



### node servers









### node servers san server virtual link /users /users /projects /projects virtual link /scratch /scratch /scratch











### Analyses steps of the cluster



mkdir





Step 2:srun, partition



2



### Data transfer on the i-Trop Cluster





### Data transfer on i-Trop cluster

/users, /projects/medium/, /projects/large, /projects/xl, projects/xxl ou /share





### Analyses steps of the cluster





Copy your data from your PC/MAC to the san if they are not on the cluster





Step 3: filezilla

### Go to the **<u>Practice3</u>** of the github



• Copy between 2 remote servers :

scp -r source destination

• Syntax if the source is remote:

scp -r server\_name:/path/file\_to\_copy local\_folder

• Syntaxe si la destination est distante :

scp -r /path/file\_to\_copy nameserver:/path/remote\_server

Ex: scp -r san:/home/tando/data/folder /scratch/tando/



### Analyses step on the cluster

Connection to bioinfomaster1.ird. fr and resources reservation

Step 1



scp





### **Step 4: scp to nodes**

### Go to the Practice4 of the github



- Allow to choose the version of software you want to use
- 2 types of softwares : bioinfo : includes all the bioinformatics softwares
   ( example BEAST) system : includes all the system softwares(example JAVA)
- Overcome the environment variables



- See the available modules : module avail
- Obtain infos on a particular module: module whatis + module name
- Load a module : module load + modulename
- List the loaded module : module list
- Unload a module : module unload + modulename
- Unload all the modules : module purge



### Analyses steps of the cluster





Step 5 module





### **Step 5: module environment**





### Analyses steps of the cluster

Connection to bioinfomaster1.ird. fr and resources reservation

Step 1







- Load the software version to launch
- Launch the data analysis

*\$~ command <options> <arguments>* 

With *command*: the command to launch





### **Step 6: launch the analysis**





• Copy between 2 remote servers :

scp -r source destination

• Syntax if the source is remote:

scp -r server\_name:/path/file\_to\_copy local\_folder

• Syntax if the destination is remote :

scp -r /path/file\_to\_copy nameserver:/path/remote\_server



### Analyses steps of the cluster







**Step 7: Retrieve results** 



### Go to the <u>Practice7</u> of the github



- Scratch = temporary spaces
- Verify that the copy is OK
- Use rm command

cd /scratch rm -rf nom\_rep



### Analyses steps of the cluster







**Step 8: data deletion** 





### Scripts to visualize/delete temporary data

- Scripts location: /opt/scripts/scratch-scripts
- Visualize data on scratchs: scratch\_use.sh

sh /opt/scripts/scratch-scripts/scratch\_use.sh

• Delete data on scratch: clean\_scratch.sh

sh /opt/scripts/scratch-scripts/clean\_scratch.sh



### Main Slurm commands

Commande	Description	Example
sruntime=0X:00pty bash -i	Interactive way to connect to a node for X minutes	sruntime=02:00:00pty bash -i Connection for 2 hours
sbatch	Launch a analyses in background via a script	sbatch script.sh
sinfo	Informations on partitions	sinfo
scancel	Deletion of jobs <job_id></job_id>	scancel 1029
squeue	Infos onjobs	squeue -u tando
scontrol show job <job_id></job_id>	Infos on active job <job_id></job_id>	scontrol show job 1029
sacct -j <job_id></job_id>	Infos on finished job <job_id></job_id>	sacct -j 1029

More Slum infos here: <a href="https://bioinfo.ird.fr/index.php/tutorials-fr/slurm/#part-2">https://bioinfo.ird.fr/index.php/tutorials-fr/slurm/#part-2</a>



### **Options of sbatch, srun, salloc**

Options	Description	Example
job-name= <name></name>	Name the job	sbatchjob-name=tando_blast
-p <partition></partition>	Choose a partition	sbatch -p highmem
nodelist= <nodex></nodex>	Choose a particular node	sbatch -p normalnodelist=node14
<pre>-n <nbre_taches></nbre_taches></pre>	Launch several instance of a command	srun -n 4 hostname
-c <nb_cpu_par_tache></nb_cpu_par_tache>	Allocate the number of cpus per task	srun -n 4 -c 2 hostname
mail-user= <emailaddress></emailaddress>	Send a email	sbatchmail- user=ndomassi.tando@ird.fr
mail-type= <event></event>	Send a email when: END: end of the job FAIL: abort BEGIN: beginning ALL: all	sbatchmail-type=BEGIN



### LAUNCH A JOB



- Scheduler choose resources automatically
- Use up to 24 cores at the same time
- Possibility to configure this choice
- Jobs launch in background
  - $\rightarrow$  possibility to turn off your PC/MAC
  - $\rightarrow$  automatic results retrieving



- Execute a script via Slurm
- Use:

\$~ **sbatch** script.sh

With script.sh : the name of the script



### First part of the script (in green): sge execution options with the key word #SBATCH

#### #!/bin/bash



### bash scripts syntax

### In the 2nd part of the script: the command to execute

nom\_variable1="valeur\_variable1"
nom\_variable2="valeur\_variable2"

sleep 30 hostname





### Launch a script with sbatch





### **Satisfaction survey**

### It is mandatory for you to fill this form to have your account extend :

https://itrop-survey.ird.fr/index.php/432222?lang=fr



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URL: https://bioinfo.ird.fr/- http://www.southgreen.fr
"



- Include a budget for bioinformatics resources in your answer to projects funding
- A need in hard drives, renew machines etc...
- Available quotations

**Projects** 

• Contact : help, needs definition, quotations...



# Thank you for you attention !



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