

SLURM @BIANCA

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More Slurm and other advanced UPPMAX techniques

S VERITA

- A closer look at Slurm
- Using the GPUs
- Job efficiency with the jobstats tool
- Advanced job submission



The Slurm Workload Manager

- Free, popular, lightweight
- Open source:
 - https://slurm.schedmd.com
- available at al SNIC centra
- UPPMAX Slurm userguide:
 - https://www.uppmax.uu.se/support/
 - user-guides/slurm-user-guide/





More on sbatch



• Recap:

sbatch	-A sens2023598	-t 10:00	-p core	-n 10	my_job.sh
Slurm batch	Project name	Maximum runtime	Partition ("job type")	#cores	job script



More on time limits



-t dd-hh:mm:ss

- \bullet 0-00:10:00 = 00:10:00 = 10:00 = 10
- \bullet 0-12:00:00 = 12:00:00
- 3-00:00:00 = 3-0
- 3-12:10:15



Job wall time

- Q: When you have no idea how long a program will take to run, what should you book?
 - A: very long time, e.g. 10-00:00:00
- Q: When you have an idea of how long a program would take to run, what should you book?
 - A: overbook by 50%



More on partitions



-p core

- "core" is the default partition
- ≤ 16 cores on Bianca
- a script or program written without any thought on parallelism will use 1 core

-p node

- if you wish to book full node(s)



- Quick testing
- The "devel" partition
 - max 2 nodes per job
 - up to 1 hour in length
 - only 1 at a time
 - -p devcore, -p devel
 - Any free nodes in the devel partition? Check status with
 - sinfo -p devel
 - jobinfo -p devel



Quick testing



- Any free nodes in the devel partition? Check status with
 - sinfo -p devel
 - jobinfo -p devel
 - more on these tools later
- High priority queue for short jobs
 - 4 nodes
 - up to 15 minutes
 - --qos=short



Debugging or complicated workflows

- Interactive jobs
 - handy for debugging a code or a script by executing it line by line or for using programs with a graphical user interface
 - salloc -n 80 -t 03:00:00 -A sens2023598
 - interactive -n 80 -t 03:00:00 -A sens2023598
 - up to 12 hours
 - useful together with the --begin=<time> flag
 - salloc -A sens2023598 --begin=2022-02-17T08:00:00

 asks for an interactive job that will start earliest tomorrow

 at 08:00



Parameters in the job script or the command line?

- Command line parameters override script parameters
- A typical script may be:

```
#!/bin/bash

#SBATCH -A sens2023598

#SBATCH -p core

#SBATCH -n 1

#SBATCH -t 24:00:00
```

Just a quick test:

```
sbatch -p devcore -t 00:15:00 jobscript.sh
```



Hands-on #1: sbatch/jobinfo

- login to Bianca
- find out which projects you're a member of using projinfo
- submit a short (10 min) test job; note the job ID
- find out if there are any free nodes in the devel partition
- submit a new job to use the devel partition
- write in the HackMD when you're done



Memory in core or devcore jobs

- -n X
- Bianca: 8GB per core
- Slurm reports the available memory in the prompt at the start of an interactive job



More flags



- J <jobname>
- email:
 - --mail-type=BEGIN, END, FAIL, TIME_LIMIT_80
 - --mail-user Don't use. Set your email correctly in SUPR instead.
- out/err redirection:
 - --output=slurm-%j.out and -error=slurm-%j.err by default, where %j will be replaced by the job ID
 - -- output=my.output.file
 - --error=my.error.file



More flags



- -C thin / -C 128GB
- -C fat / -C 256GB / -C 512GB
- Dependencies: --dependency
- Job array: --array
- More at https://slurm.schedmd.com/sbatch.html
 - or just man sbatch
 - not all options work on all systems!



- Monitoring jobs
- jobinfo a wrapper around squeue
 - lists running and pending jobs
 - jobinfo -u username
 - jobinfo -A sens2023598
 - jobinfo -u username --state=running
 - jobinfo -u username --state=pending
- One may also use the squeue command.
- bianca_combined_jobinfo (queued jobs of all projection)

Monitoring and modifying jobs



• scontrol

- scontrol show job jobid
- possible to modify the job details after the job has been submitted;
 some options, like maximum runtime, may be modified (=shortened)
 even after the job started
 - scontrol update JobID=jobid QOS=short
 - scontrol update JobID=jobid TimeLimit=1-00:00:00
 - scontrol update JobID=jobid NumNodes=10
 - scontrol update JobID=jobid Features=mem1TB



When a job goes wrong



scancel

- jobid
- -u username to cancel all your jobs
- -t state cancel pending or running jobs
- n name cancel jobs with a given name
- -i ask for confirmation



Priority



- Roughly:
- The first job of the day has elevated priority
- Other normal jobs run in the order of submission (subject to scheduling)
- Projects exceeding their allocation get successively into the lower priority category
- Bonus jobs run after the jobs in the higher priority categories



Priority



- In practice:
 - submit early = run early
 - bonus jobs always run eventually, but may need to wait until the night or weekend
- In detail:

https://www.uppmax.uu.se/support/faq/running-jobs-faq/your-priority-in-the-waiting-job-queue/



Hands-on #2:

sbatch/squeue/scancel/scontrol/jobinfo

- submit a new job; note the job ID
- check all your running jobs
- what is the priority or your recently-submitted job?
- submit a new job to run for 24h; note the job ID
- modify the name of the job to "wrongjob" and the maximum runtime to 7days, for example
- cancel your job with name "wrongjob"





Determining job efficiency

• jostats - custom-made UPPMAX tool



Job efficiency



- jobstats a tool in the fight for productivity
 - it works only for jobs longer than 5-15 minutes
 - r jobid check running jobs
 - A project check all recent jobs of a given project
 - -p jobid produce a CPU and memory usage plot
 - https://www.uppmax.uu.se/support/user-guides/ jobstats-user-guide/



Hands-on #3: jobstats

- Generate jobstats plots for your jobs
 - Firstly, find some job IDs from this month
 - finishedjobinfo -m username
 - Write down the IDs from some interesting jobs.
 - Generate the images:

```
$ jobstats -p ID1 ID2 ID3
```

- Look at the images
 - \$ eog *png &



Hands-on #3: jobstats

- Which of the plots
 - Show good CPU or memory usage?
 - Indicate that the job requires a fat node?

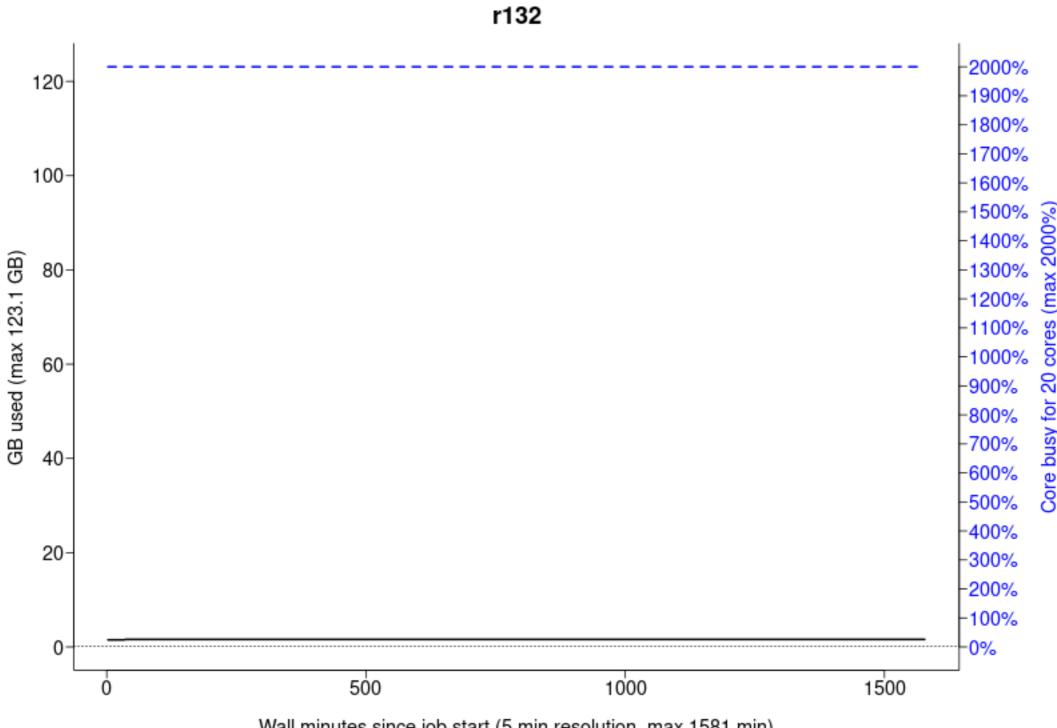


11217530 COMPLETED on rackham end: 2019-12-17T12:35:44 runtime: 1-02:16:12

User:

Proj: snic2019-1-12 Jobname: fesn_x

Flags: none



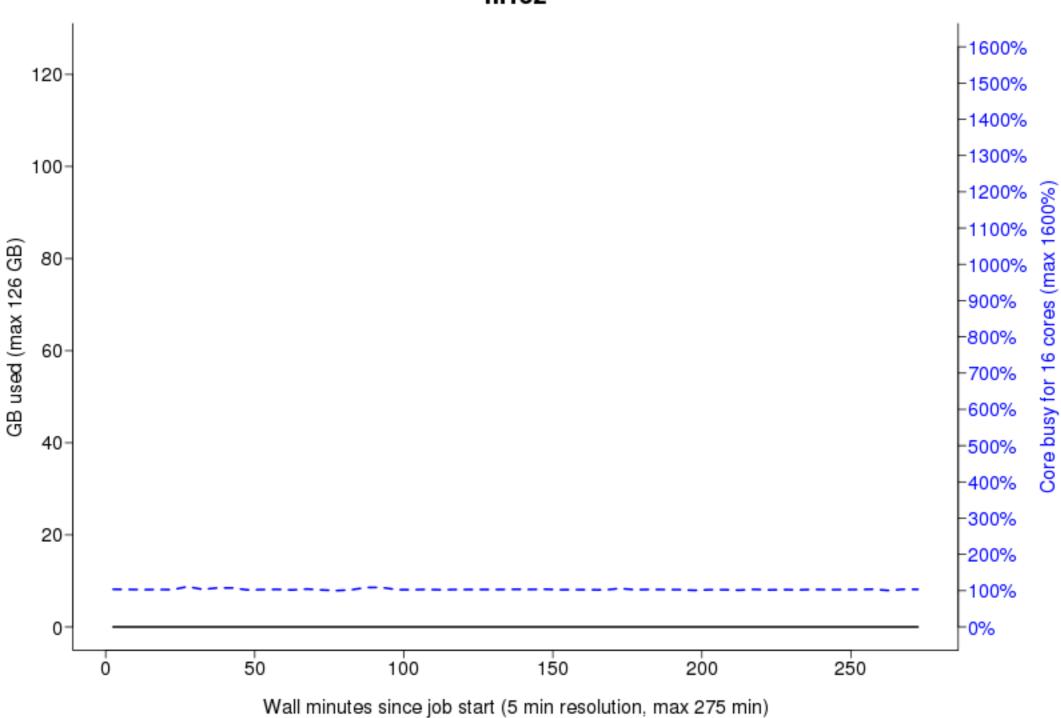


8804061 COMPLETED on milou end: 2016-10-10T20:30:26 runtime: 04:36:33

User:

Proj: b2015110 overbooked:12%, !!half_overbooked, !!severely_overbooked, Jobname: nf-MergeBam_(2) cores_overbooked:16:2, mem_overbooked:126:0, core_mem_overbooked:15.8:0





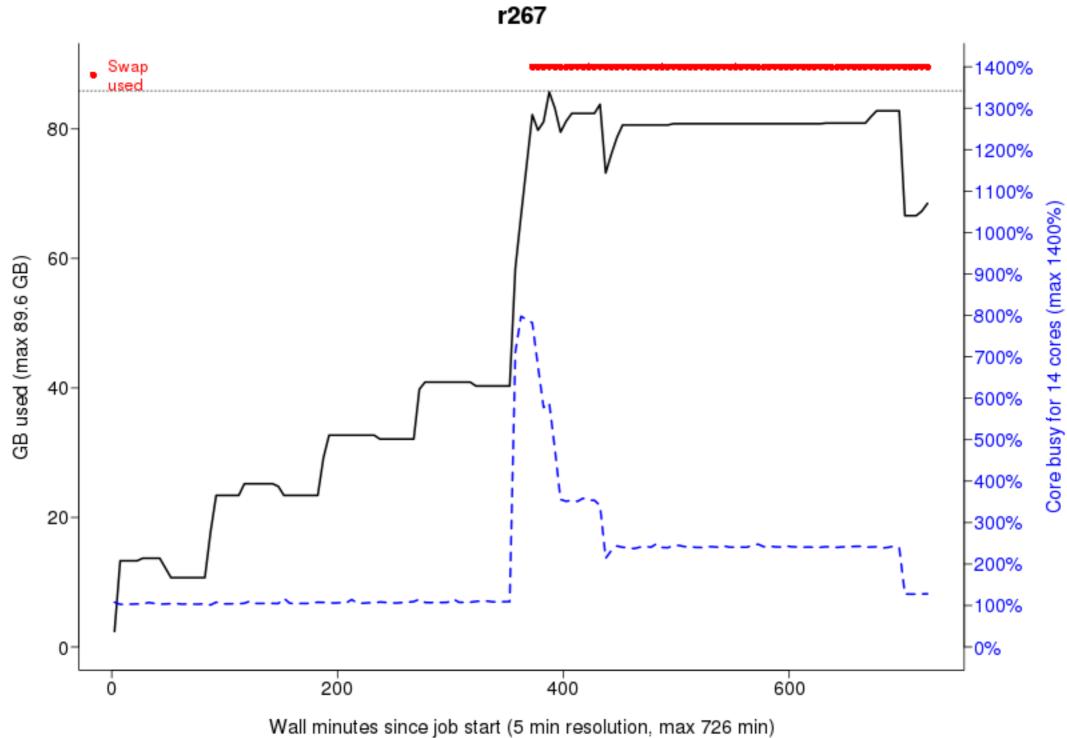


607031 OUT_OF_MEMORY on rackham end: 2018-08-25T07:29:39 runtime: 12:08:5

User:

Proj: snic2017-1-355 Jobname: pgd.apps

!!swap_used, cores_overbooked:14:11







Different flavours of Slurm

Job script examples and workflows



Simple workflow

```
#!/bin/bash
#SBATCH -J jobname
#SBATCH -A sens2023598
#SBATCH -p core
#SBATCH -n 10
#SBATCH -t 10:00:00
module load software/version
module load python/3.9.5
./my-script.sh
```

./another-script.sh

./myprogram.exe



Job dependencies

- sbatch jobscript.sh submitted job with jobid1
- sbatch anotherjobscript.sh submitted job with jobid2
- --dependency=afterok: jobid1: jobid2 job will only start running after the successful end of jobs jobid1: jobid2
- very handy for clearly defined workflows
- One may also use --dependency=afternotok: jobid in case you'd like to resubmit a failed job, OOM for example, to a node with a higher memory: -C mem215GB or -C mem512



I/O intensive jobs: \$SNIC TMP

```
#!/bin/bash
#SBATCH -J jobname
#SBATCH -A sens2023598
#SBATCH -p core
#SBATCH -n 1
#SBATCH -t 10:00:00
module load bioinfotools
module load bwa/0.7.17 samtools/1.14
export SRCDIR=$HOME/path-to-input
cp $SRCDIR/foo.pl $SRCDIR/bar.txt $SNIC TMP/.
cd $SNIC TMP
./foo.pl bar.txt
cp *.out $SRCDIR/path-to-output/.
```



OpenMP or multi-threaded job

```
#!/bin/bash
#SBATCH -A sens2023598
#SBATCH --exclusive
#SBATCH -p node
#SBATCH --ntasks-per-node=1
#SBATCH --cpus-per-task=20
#SBATCH -t 01:00:00
module load uppasd
export OMP NUM THREADS=20
```

sd > out.log



GPU nodes on Bianca

- Nodes with Nvidia A100 40 GB
- All GPU nodes have at least 256 GB RAM (fat nodes) with 16
 CPU cores
- 1 or 2 GPUs per node
- SBATCH options:

```
#SBATCH -C gpu
#SBATCH --gres=gpu:1
#SBATCH --gpus-per-node=1
```





Running on several nodes: MPI jobs

```
#!/bin/bash -1
#SBATCH -J rsptjob
#SBATCH -mail-type=FAIL
#SBATCH -A sens2023598
#SBATCH -t 00-07:00:00
#SBATCH -p node
#SBATCH -N 4
### for jobs shorter than 15 min (max 4 nodes):
###SBATCH --qos=short
module load RSPt/2021-10-04
export RSPT SCRATCH=$SNIC TMP
```

srun -n 80 rspt

rm -f apts dmft_lock_file e_entropy efgArray.dat.0 efgData.out.0
energy_matrices eparm_last interstitialenergy jacob1 jacob2 locust.toppsala
out last pot last rspt fft wisdom.* runs.a symcof new
UNIVERSITET

Job arrays

- VERITA VERITA VUR
- Submit many jobs at once with the same or similar parameters
- Use \$SLURM ARRAY TASK ID in the script in order to find the correct path

```
#!/bin/bash
#SBATCH -A sens2023598
#SBATCH -p node
#SBATCH -N 2
#SBATCH -t 01:00:00
#SBATCH -J jobarray
#SBATCH --array=0-19
#SBATCH --mail-type=ALL, ARRAY TASKS
 SLURM ARRAY TASK ID tells the script which iteration to run
echo $SLURM ARRAY TASK ID
cd /pathtomydirectory/dir $SLURM ARRAY TASK ID/
srun -n 40 my-program
env
```

• You may use scontrol to modify some of the job arrays.



Snakemake and Nextflow

- Conceptually similar, but with different flavours
- First define steps, each with an input, an output, and a command that transforms the input into output
- Then just ask for the desired output and the system will handle the rest



Hands-on #4: make it your own

- use 2 or 3 of the sample job scripts as a starting point for your own job script
- tweak them so that you run something closer to your research; or just feel free to experiment
- paste at least one of the examples in the HackMD
- great if you could add a comment what the job script is about





Feedback on Slurm

- what did you find useful?
- not so useful?
- what is most challenging while editing your job script / worksflow?
- something that was not covered that you'd like to know about?
- please provide your feedback in the HackMD





Where to go from here?

Code documentation

SNIC training newsletter - software-specific training events included

https://coderefinery.org/workshops/upcoming/

https://nbis.se/training/events.html (bio)

email support@uppmax.uu.se or https://supr.naiss.se/support/



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