# Embarrassingly Parallel workflows and how to execute them



# **Embarrassingly parallel workflows**

### What?

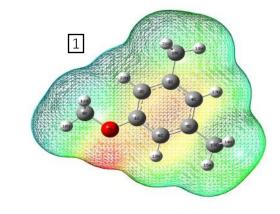
• Executing many independent small jobs simultaneously to use the hardware resource efficiently.

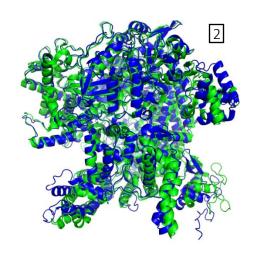
### When?

- It is not possible to combine multiple runs/tasks into a single application.
- Large parameter sweep based on the same application or workflow with dependent pipeline steps.
- Workflow involves runs that are depending on several independent software which do not scale beyond a certain number of cores.

### Why?

- Nodes are getting denser with respect to resources such as number of cores/threads,
   GPU cards, Memory.
- Applications do not necessarily scale even within a node: shared memory (OpenMP, MPI SHM, pthreads)







<sup>1.</sup> Mathiyalagan, A., Manimaran, K., Muthu, K., Rajakantham, M., Density functional theory study on the electronic structures and spectral properties of 3,5-Dimethylanisole dye sensitizer for solar cell applications, Results in Chemistry, Volume 3, 2021

### How?

#### **Tools**

- Linux background processes (within a node) along with numactl
  - ► Supported on Snellius and requires prior knowledge of bash scripting and process placement using **numactl**.
  - Gives most manual control if one knows what they are doing.
- SLURM arrays (to perform similar jobs on multiple allocations)
  - Supported on Snellius and requires prior knowledge of SLURM array options.
  - ► Is limited to cgroups supported by SLURM which means each allocation cannot be below the minimum allocation within SLURM. (Snellius -> 16 cores and 28 GiB of memory on "rome" partition)
  - ► One has to wait for each task in the array to get allocated and finish.
- STOPOS
  - ▶ Not supported in Snellius anymore. A tool to execute embarrassingly parallel jobs, requires prior knowledge of bash scripting.
- QCG-PilotJob (https://qcg-pilotjob.readthedocs.io/en/develop/)
  - Supported on Snellius and requires prior knowledge of Python scripting.







### There are many more tools















### **QCGPilot-Job**

### Job manager inside SLURM:

- Submits jobs within a single large allocation across multiple nodes.
- Does not have to wait for small individual allocations like in SLURM arrays.
- Tries to optimize resources for each job based on requirements such that the queue is always full.
- Possibility of restarting failed/timed out jobs.
- One can also specify complex result-based job dependencies.

### Some caveats and drawbacks:

- One needs to estimate, the total time of all the jobs combined.
- Job/system state files get produced in the submission directory.
- Might bring some complexity into execution of a Singularity containers with embedded MPI-based codes.
- In the case of problems, the log files may not be very expressive.



### Workflow

### SLURM job script

Describes resources, loads the QCG-PilotJob module and runs the QCG-PilotJob python script

# QCG-PilotJob python script

Invokes the QCG-PilotJob
Manager that aggregates all
your jobs along with
resources and conditions
based on which jobs need to
be invoked

### **Executable**

This is the main application that one wants to execute

### Output

Job output with QCG-PilotJob metrics



# QCGPilot-Job: Set up the environment

- The QCG-PilotJob package is available as a module in the 2023 environment
- Note that the environment from the login node is **not transferred** to the compute nodes, unless defined in, e.g., the **.bashrc** file. Thus, you have to re-load the QCG-PilotJob package in your job script.

```
$ module purge
$ module load 2023
$ module load QCG-PilotJob/0.14.1-foss-2023a
```



### Workflow

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# QCGPilot-Job: SLURM job script

```
run.sh
#SBATCH -n 16
module load 2023
module load QCG-PilotJob/0.14.1-foss-2023a
python test.py
```



### Workflow

### SLURM job script

# python script

QCG-PilotJob

This is the main application that one wants to invoke

Executable

### Output

Describe resources, loads the QCG-PilotJob module and runs the QCG-PilotJob python script Invokes the QCG-PilotJob
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Job output with QCG-PilotJob metrics



- Work with the **QCG-PilotJob** manager consists of a few mandatory steps:
  - Creation of an instance of QCG-PilotJob Manager
    - Creation of a class object
  - Waiting for all submitted jobs to finish
    - will wait until some or all submitted to the QCG-PilotJob Manager jobs finish
  - Call for the **finish()** method at the end of the script
    - this method terminates the background thread in which the instance of a QCG-PilotJob Manager has been run



### test.py

```
manager = LocalManager()
print('available resources: ', manager.resources())
manager.wait4all()
manager.finish()
```

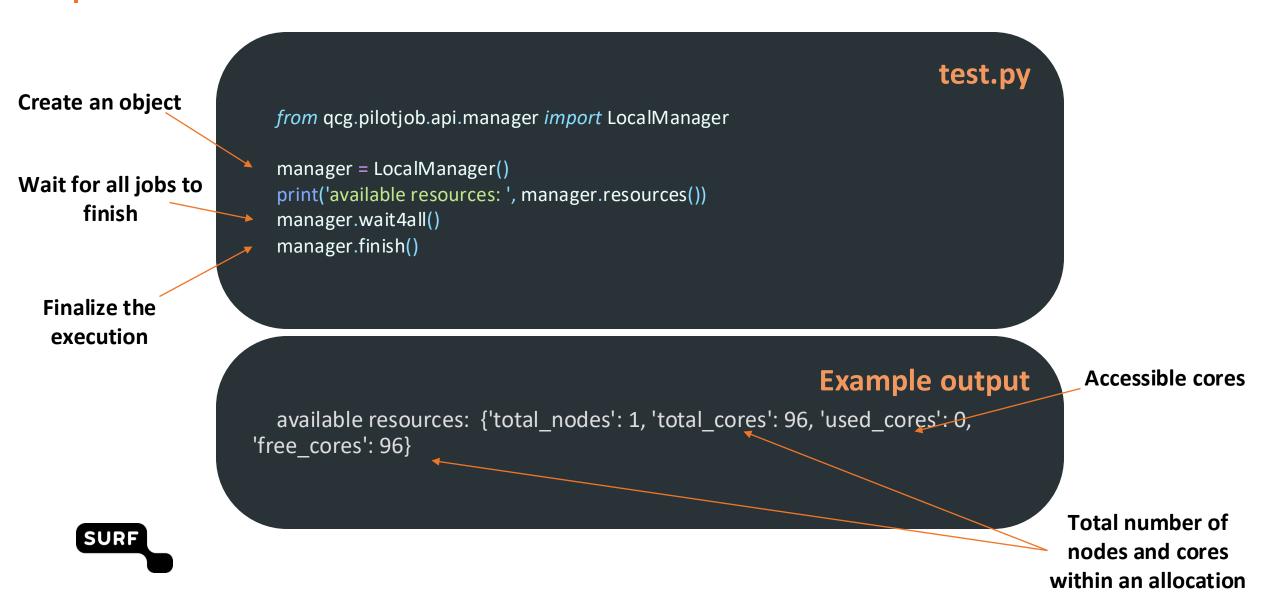
from qcg.pilotjob.api.manager import LocalManager

# **Example output**

```
available resources: {'total_nodes': 1, 'total_cores': 96, 'used_cores': 0, 'free_cores': 96}
```



test.py **Create an object** from qcg.pilotjob.api.manager import LocalManager manager = LocalManager() Wait for all jobs to print('available resources: ', manager.resources()) finish manager.wait4all() manager.finish() Finalize the execution **Example output** available resources: {'total\_nodes': 1, 'total\_cores': 96, 'used\_cores': 0, 'free\_cores': 96} SURF



# **QCGPilot-Job: Create jobs**

- To execute multiple jobs within a single SLURM allocation we need to:
  - Create an object of the **Jobs** class
  - Populate it with some parameters (describe jobs)
  - Submit the created object to **QCG-PilotJob Manager**

```
from qcg.pilotjob.api.job import Jobs

jobs = Jobs()
jobs.add(...)
jobs_ids = manager.submit(jobs)
print('submited jobs: ', str(job_ids))
...
```



# QCGPilot-Job: Job description (main parameters)

- **name** the job name
- **exec** path to the executable program
- **args** executable program arguments
- **script** bash script content
- **stdin** path to file which content should be passed to the standard input stream
- **stdout** path to the file where standard output stream should be saved
- **stderr** path to the file where standard error stream should be saved
- **modules** list of modules that should be loaded before job start
- **numCores** number of required cores specification
- **numNodes** number of required nodes specification
- **iteration** iterations definition
- model model of execution



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### **QCGPilot-Job: execution models**

- **default** only a single process is launched within the allocation
- threads is designed for running OpenMP tasks on a single node
- openmpi the processes are started with the mpirun command
- intelmpi same but for Intel MPI
- srunmpi the processes are started with the srun command

Read more about execution models in the official docs:

https://qcg-pilotjob.readthedocs.io/en/develop/execution\_models.html



### QCGPilot-Job: Calling for an executable

```
test.py
from qcg.pilotjob.api.manager import LocalManager
from qcg.pilotjob.api.job import Jobs
manager = LocalManager()
jobs = Jobs()
jobs.add(script='echo "job ${it} executed at `date` @ `hostname`"',
     name='test_job',
    stdout='job.out.${it}',
    iteration=4)
job names = manager.submit(jobs)
print('submited job names: ', str(job_names))
job status = manager.status(job names)
print('job status: ', job_status)
manager.wait4all()
manager.finish()
```



# Hands-on



### Hands-on

### **Clone the repository**

• <a href="https://github.com/sara-nl/course-module-qcg-pilotjob">https://github.com/sara-nl/course-module-qcg-pilotjob</a>

```
$ cd ~ $ git clone https://github.com/sara-nl/course-module-qcg-pilotjob
```

\$ cd course-module-qcg-pilotjob



### Hands-on #1

#### Task

• Execute **512 independent short jobs** on two nodes on the "rome" partition. Note that one node on the "rome" partition has 128 cores.

### Steps

- Go to hands-on/example\_01
- Take a look at qcg\_job.py and qcg\_job.sh
- Submit the SLURM job script, check the results

```
from qcg.pilotjob.api.job import Jobs
from qcg.pilotjob.api.manager import LocalManager
manager = LocalManager()
jobs = Jobs()
for job in range(512):
  jobs.add(name="job {}".format(job),
       exec='hostname',
       stdout='output/job.{}.out'.format(job),
       stderr='output/job.{}.err'.format(job),
       model='default',
       numCores={ "exact" : 1},
       iteration=1)
print("-- submit jobs --")
manager.submit(jobs)
print("-- wait for all jobs --")
manager.wait4all()
manager.finish()
print("-- finished")
```



### Hands-on #2

#### Task

• Compute the average of one column in 1000 CSV files. Aggregate the results in one CSV file.

### Steps

- Go to hands-on/example\_02
- Take a look at README.md, qcg\_job\_v1.py and qcg\_job.sh
- Extract the input data (input.tar.gz)
- Submit the SLURM job script, check the results

```
job names = []
for filename in glob(os.path.join("input/*.csv"), recursive = False):
  job names.append(os.path.basename(filename))
manager = LocalManager()
jobs = Jobs()
for jname in job names:
  print("submit {}".format(jname))
  jobs.add(name=jname,
       exec='python3',
       args=["average.py",
          os.path.join("input", jname)],
       stdout='average {}'.format(jname),
       stderr='job.{}.err'.format(jname),
       modules=[
      "2023", "Python/3.11.3-GCCcore-12.3.0"],
       iteration=1
jobs.add(name="aggregate",
     script='cat average *.csv | sort',
     stdout='result.csv',
     stderr='aggregate.err')
```

