



**Barcelona
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Centro Nacional de Supercomputación

OmpSs-2 Hands-on

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Extract and configure oss-2-ee package

(1) Extracting the sources

```
$ tar -xvzf oss-2-ee.tar.gz
<list of extracted files>
```

(2) Main directory

```
$ cd oss-2-ee
$ ls
02-examples
03-fundamentals
04-mpi+ompss-2
05-cuda+ompss-2
06-openacc+ompss-2
local
configure
configure-dc
paraver-cfgs
README.md
```

(3) Configure

```
$ source configure
Basic configuration...
  Mercurium compiler at /apps/ompss/bin
  Extrae library at /apps/extrae/bin
  Paraver utility at /apps/wxparaver/bin
  ...
Additional libraries...
  MPI library at /opt/mpi/bullxmpi/.../lib
  MKL library at /opt/.../mkl/lib/intel64/
  ATLAS library at /opt/.../ATLAS/3.9.51/lib
```

(0) Remainder: exercises are in /apps/PM/training/PUMPS-2022

```
$ cp /apps/PM/training/PUMPS-2022/oss-2-ee.tar.gz $HOME
```

(1) nn-session / exercise

```
$ cd 02-examples/cholesky-openblas  
$
```

(2) Directory contents

```
$ ls  
cholesky.c  
cholesky.h  
Makefile  
README.rst
```

(3) README.rst file (script)

```
$ vi README.rst  
$
```

(4) Run “make” will create...

```
$ make  
Building: program  
Creating: mutirun.sh, run-once.sh, trace.sh and  
extrae.xml
```

- » an executable
- » different scripts to run your programs
- » an extrae.xml file (as default to get your Paraver traces)

```
$ ls  
cholesky.c      cholesky.h      cholesky  
extrae.xml      Makefile        multirun.sh  
README.rst      run-once.sh     trace.sh
```

Running oss-2-ee package (CTE Power 9) [1]



Checking script configuration

```
$ vi run-once.sh  
$
```

```
#!/bin/bash  
#SBATCH --job-name=oss-2-ee  
#SBATCH --ntasks=1  
#SBATCH --cpus-per-task=12  
#SBATCH --qos=debug  
#SBATCH --gres=gpu:1  
PROGRAM=cholesky  
#INSTRUMENT=./xxxxx.sh  
$INSTRUMENT ./$PROGRAM 4096 512 1
```

run-once.sh

Submitting the script

```
$ sbatch run-once.sh  
Submitted batch job 1234567
```

Check submitted jobs with:
\$ squeue

Exercise directory

```
$ ls  
cholesky.c      cholesky.h      cholesky  
extrae.xml      Makefile        multirun.sh  
README.rst      run-once.sh     trace.sh
```

Directory after execution

```
$ ls  
cholesky.c      cholesky.h      cholesky  
extrae.xml      Makefile        multirun.sh  
oss-2-ee_nnn.err oss-2-ee_nnn.out README  
run-once.sh     trace.sh
```

Running oss-2-ee package (CTE Power 9) [2]

Checking script configuration

```
$ vi run-once.sh
$
```

```
#!/bin/bash
#SBATCH --job-name=oss-2-ee
#SBATCH --ntasks=1
#SBATCH --cpus-per-task=12
#SBATCH --qos=debug
#SBATCH --gpus=gpu:1
PROGRAM=cholesky
#INSTRUMENTATION=cholesky
##SBATCH --qos=debug
$INSTRUMENTATION #SBATCH --reservation = PUMPS+AI22
```

Submitting the script

```
$ sbatch run-once.sh
Submitted batch job 1234567
```

Exercise directory

```
$ ls
cholesky.c      cholesky.h      cholesky
extrae.xml      Makefile        multirun.sh
README.rst      run-once.sh     trace.sh
```

Directory after execution

```
$ ls
cholesky.c      cholesky.h      cholesky
extrae.xml      Makefile        multirun.sh
oss-2-ee_nnn.err oss-2-ee_nnn.out README
run-once.sh     trace.sh
```

Instrumenting oss-2-ee package

Checking configuration scripts

```
$ vi run-once.sh trace.sh
$
```

```
#!/bin/bash
#SBATCH --job-name=oss-2-ee
#SBATCH --cpus-per-task=1
#SBATCH --qos=debug
#SBATCH --gres=gpu:1
PROGRAM=cholesky
#INSTRUMENT= ./xxxxxx.sh
$INSTRUMENT./$PROGRAM 4096 512 1
```

run-once.sh

or

INSTRUMENT= ./graph.sh

```
#!/bin/bash
. . .
export EXTRAE_CONFIG_FILE=extrae.xml
export LD_PRELOAD=/path/to/extrae/lib?
$*
```

trace.sh

Exercise directory

```
$ ls
cholesky.c      cholesky.h      cholesky
extrae.xml      Makefile        multirun.sh
README.rst      run-once.sh     trace.sh
```

Submitting the script

```
$ sbatch run-once.sh
Creating: cholesky.prv, cholesky.pcf cholesky.raw
```

Compile and execute (examples)

Exercise's location: 02-examples

```
~oss-2-ee:$ ls
02-examples          03-fundamentals
04-mpi+ompss-2       05-cuda+ompss-2
06-openacc+ompss-2   ...
```

Compile and execute (guidelines)

- Code is completely annotated, you DON'T need to modify it. Review source code, check different directives and their clauses
- Check scalability → compute speed-up
- Obtain traces, explore them
- Check the README file (more goals)

Hands-on session's contents

~~Exercise 2.1: cholesky [-mkl][-openblas]~~

~~Exercise 2.2: matmul [] [-dep] [-red] [-wdep]~~

Remember to configure your system

```
$ source configure
Basic configuration...
```

Check running scripts (before submit)

```
$ vi run-once.sh
$
```

- Scripts **run-once.sh**, **multi-run.sh** or **trace.sh**
- Job scheduler configuration front-matter: queue, hardware resources, etc.
- Other runtime's options used in the script

Fundamentals

Exercise's location: 03-fundamentals

```
~oss-2-ee:$ ls
02-examples          03-fundamentals
04-mpi+ompss-2       05-cuda+ompss-2
06-openacc+ompss-2   ...
```

OmpSs-2 Fundamentals (guidelines)

- Code are incomplete
 - » Lack of clauses (in some cases)
 - » Lack of directives (is some others)
- Check scalability → compute speed-up
- Obtain traces, explore them
- Check the README file (more goals)

Hands-on session's contents

Continue previous session's exercises

Exercise 3.1: dotproduct [-dep] [-red]

Exercise 3.2: axpy [-dep] [-wdep]

Remember to configure your system

```
$ source configure
Basic configuration...
```

Check running script (before submit)

```
$ vi run-once.sh
$
```

- Scripts **run-once.sh**, **multi-run.sh** or **trace.sh**
- Job scheduler configuration front-matter: queue, hardware resources, etc.
- Other runtime's options used in the script

OmpSs-2 + CUDA

Exercise's location: 05-cuda+ompss-2

```
oss-2-ee:~$ ls
02-examples          03-fundamentals
04-mpi+ompss-2       05-cuda+ompss-2
06-openacc+ompss-2   ...
```

OmpSs-2 + CUDA (guidelines)

- Review source code, check different directives and their clauses
- Check scalability → compute speed-up
- Obtain traces, explore them
- Check the README file (more goals)

Hands-on session's contents

Continue previous session's exercises

Exercise 5.1: SAXPY kernel

Exercise 5.2: Matmul kernel

Exercise 5.3: Nbody kernel

Remember to configure your system

```
$ source configure
Basic configuration...
```

Check running script (before submit)

```
$ vi run-once.sh
$
```

- Scripts `run-once.sh`, `multi-run.sh` or `trace.sh`
- Job scheduler configuration front-matter: queue, hardware resources, etc.
- Other runtime's options used in the script

OmpSs-2 + OpenACC

Exercise's location: 06-openacc+...

```
oss-2-ee:$ ls
02-examples          03-fundamentals
04-mpi+ompss-2       05-cuda+ompss-2
06-openacc+ompss-2   ...
```

OmpSs-2 + OpenACC (guidelines)

- Review source code, check different directives and their clauses
- Check scalability → compute speed-up
- Obtain traces, explore them
- Check the README file (more goals)

Hands-on session's contents

Continue previous session's exercises

Exercise 6.1: SAXPY kernel

Exercise 6.2: Matmul kernel

Exercise 6.3: Nbody kernel

Remember to configure your system

```
$ source configure
Basic configuration...
```

Check running script (before submit)

```
$ vi run-once.sh
$
```

- Scripts **run-once.sh**, **multi-run.sh**
- Job scheduler configuration front-matter: queue, hardware resources, etc.
- Other runtime's options used in the script



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Visualizing paraver traces [1/5]

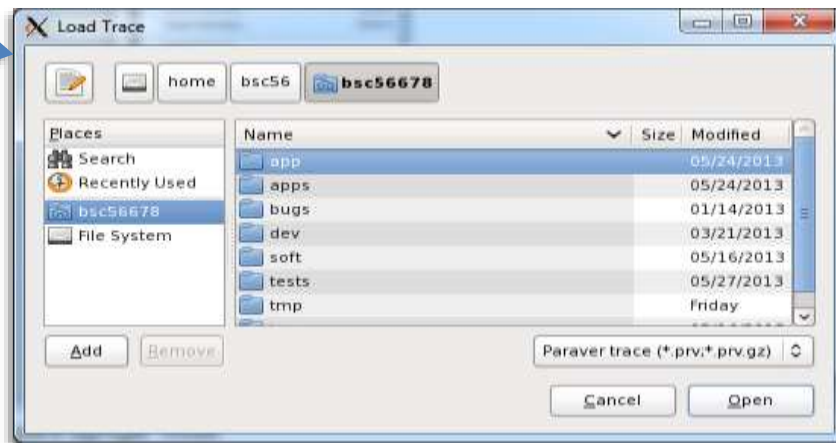
Running paraver tool

```
$ paraver
$
```

— Load a Paraver trace

Suite exercise directory

```
$ ls
cholesky.c      cholesky-d
cholesky.h      cholesky-i
cholesky-i.pcf  cholesky-i.prv
cholesky-i.raw  cholesky-p
extrae.xml      Makefile
multirun.sh     README.rst
run-once.sh     trace.sh
```



Visualizing paraver traces [2/5]

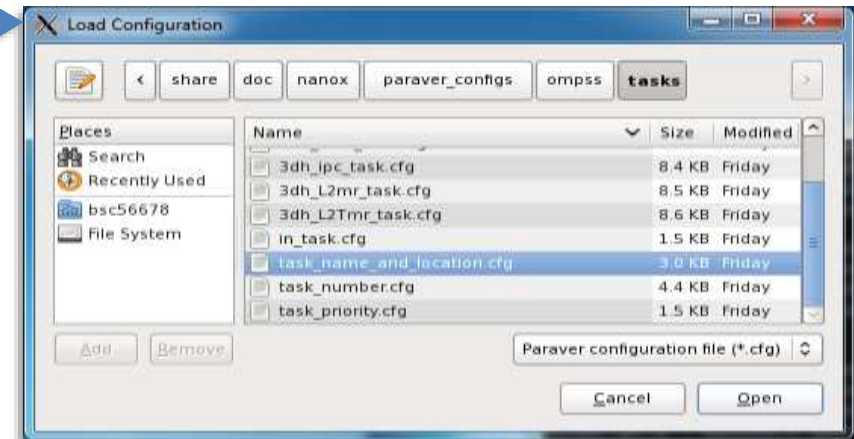
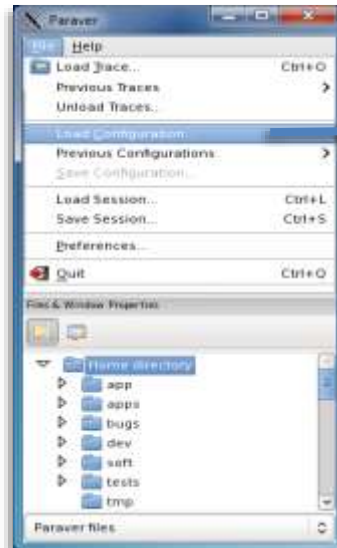
Running paraver tool

```
$ paraver
$
```

- Load a Paraver trace
- Load a configuration file

Suite root directory

```
$ ls
01-examples          02-beginners
03-gpu-devices       04-mpi+ompss
common-files         configure.sh
paraver-cfgs         README.rst
```



Visualizing paraver traces [3/5]

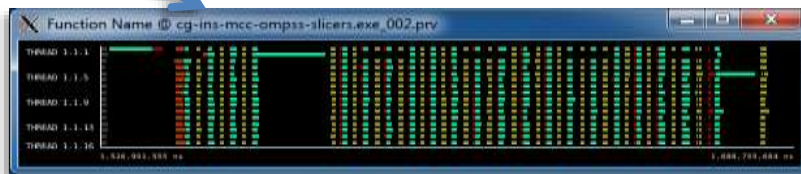
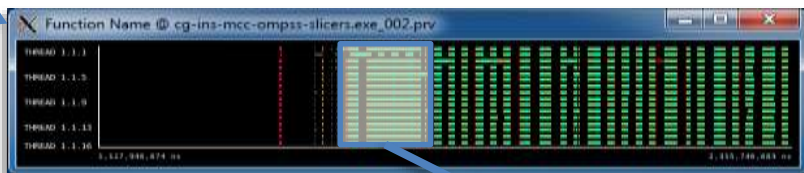
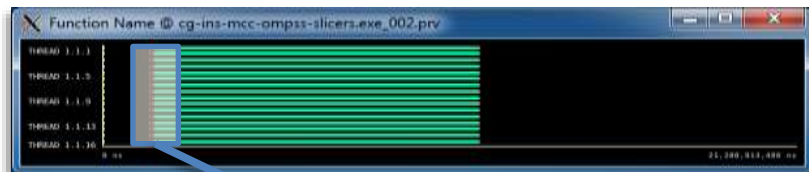
Running paraver tool

```
$ paraver
$
```

- Load a Paraver trace
- Load a configuration file
- Trace analysis (zoom in, details)

Suite root directory

```
$ ls
01-examples          02-beginners
03-gpu-devices       04-mpi+ompss
common-files         configure.sh
paraver-cfgs         README.rst
```



Visualizing paraver traces [4/5]

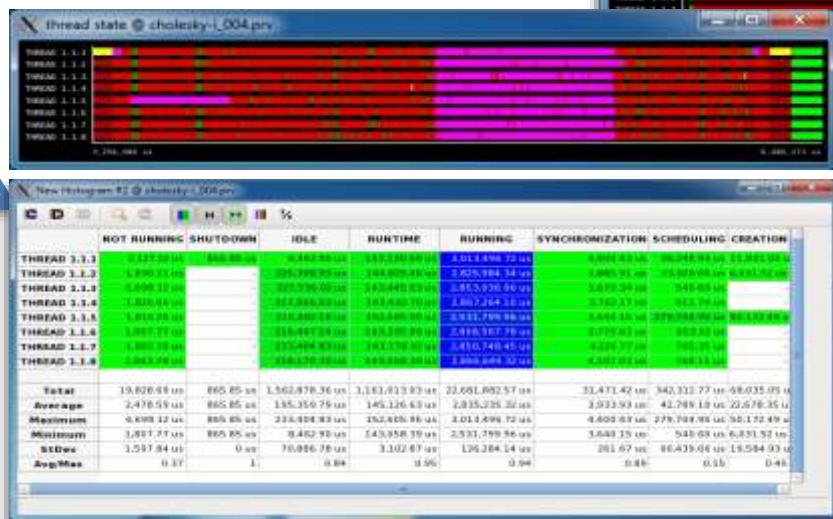
Running paraver tool

```
$ paraver
$
```

- Load a Paraver trace
- Load a configuration file
- Trace analysis (zoom in, details)
- Histograms to summarize traces

Suite root directory

```
$ ls
01-examples          02-beginners
03-gpu-devices       04-mpi+omps
common-files         configure.sh
paraver-cfgs        README.rst
```



Running paraver tool

```
$ paraver  
$
```

- Load a Paraver trace
- Load a configuration file
- Trace analysis (zoom in, details)
- Histograms to summarize traces
- Other configuration files
 - » ompss / runtime / thread_state.cfg
 - » ompss / runtime / nanos_API.cfg
 - » ompss / tasks /
task_name_and_location.cfg
 - » ompss / cuda / ...
 - » hwc / papi / performance / ...

Suite root directory

```
$ ls  
01-examples          02-beginners  
03-gpu-devices       04-mpi+ompss  
common-files         configure.sh  
paraver-cfgs         README.rst
```

- more info about paraver instrumentation tool

<http://pm.bsc.es/ompss-docs/user-guide>