



CSCS

Centro Svizzero di Calcolo Scientifico
Swiss National Supercomputing Centre

ETH zürich



Summer School 2016

Programming Environment

<http://github.com/eth-cscs/SummerSchool2016/wiki> - DAY1

18th July 2016

Summary

- Accessing CSCS
- Remote display
- Compiling my code
- Running my code
- Editing my code
- Transferring files from/to CSCS
- Debugging tools



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Accessing CSCS

Accessing CSCS: SSH

Secure shell: 2 steps

- Piz Daint is hidden: ela / daint10x / daint0x / `nidxxx`
- frontend first `Ela`: `ssh -X courseNN@ela.cscs.ch`
- then login node `Piz Daint`: `ssh -X daint`

```
ssh -X course51@ela.cscs.ch
```

```
=====
                IMPORTANT REMINDER FOR USERS of CSCS facilities
                help@cscs.ch - +41 91 610 82 10 - http://user.cscs.ch
=====

course51@ela1:~
```

```
ssh -X daint
```

```
course51@ela1:~ ssh -X daint
course51@daint101:~ xclock
```



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Compiling the code

Compiling the code: setup your PE

Cray Programming Environment

- 4 compilers available: `CCE`*, `GNU`, `INTEL`, `PGI`
- 4 predefined Programming Environment:
 - `PrgEnv-cray`, `PrgEnv-gnu`, `PrgEnv-intel`, `PrgEnv-pgi`
 - `echo $PE_ENV` to get the current PrgEnv
- 3 wrappers available: `ftn` (Fortran), `cc` (C), `CC` (C++)
 - You `must` use the wrappers to compile for the compute node,
 - The wrappers support serial, OpenMP, MPI and Cuda codes.
- The wrappers are based on the `module` command
 - dynamic modification of the `programming environment` via modulefiles,
 - the wrappers will detect the loaded `modulefiles`,
 - and will automatically add the needed flags and libraries.

Compiling the code: module list

module list

```
*course51@daint103:~ $ module list
Currently Loaded Modulefiles:
  1) modules/3.2.10.3
  2) eswrap/1.1.0-1.020200.1231.0
  3) switch/1.0-1.0502.60522.1.961.ari
  4) craype-network-aries
* 5) cce/8.3.12
  6) craype/2.4.0
  7) totalview-support/1.1.4
  8) totalview/8.11.0
* 9) cray-libsci/13.0.4
 10) udreg/2.3.2-1.0502.10518.2.17.ari
 11) ugni/6.0-1.0502.10863.8.29.ari
 12) pmi/5.0.10-1.0000.11050.0.0.ari
 13) dmapp/7.0.1-1.0502.11080.8.76.ari
 14) gni-headers/4.0-1.0502.10859.7.8.ari
 15) xpmem/0.1-2.0502.64982.5.3.ari
 16) dvs/2.5_0.9.0-1.0502.2188.1.116.ari
 17) alps/5.2.4-2.0502.9822.32.1.ari
 18) rca/1.0.0-2.0502.60530.1.62.ari
 19) atp/1.8.2
*20) PrgEnv-cray/5.2.82
*21) craype-sandybridge
 22) slurm
*23) cray-mpich/7.2.2
 24) ddt/6.0
```

Compiling the code: module swap

module swap from CCE to GNU

```
course51@daint103:~ module swap PrgEnv-cray PrgEnv-gnu
Currently Loaded Modulefiles:
 4) craype-network-aries
*5) gcc/4.8.2
 6) craype/2.4.0
 7) craype-sandybridge
 9) cray-mpich/7.2.2
13) cray-libsci/13.0.4
*24) PrgEnv-gnu/5.2.82

course51@daint103:~ ftn --version
GNU Fortran (GCC) 4.8.2 20131016 (Cray Inc.)
```

module swap from GNU to CCE

```
course51@daint103:~ module swap PrgEnv-gnu PrgEnv-cray
Currently Loaded Modulefiles:
 4) craype-network-aries
 5) craype-sandybridge
 7) cray-mpich/7.2.2
*9) cce/8.3.12
10) craype/2.4.0
13) cray-libsci/13.0.4
*24) PrgEnv-cray/5.2.82

course51@daint103:~ ftn -V
Cray Fortran : Version 8.3.12
```


Compiling the code: module avail

module avail

```
course51@daint103:~ module avail
```

```
# --- COMPILERS ---
```

```
PrgEnv-cray/5.2.82 PrgEnv-gnu/5.2.82  
PrgEnv-intel/5.2.82 PrgEnv-pgi/5.2.82  
cce/8.3.12 gcc/4.8.2 intel/15.0.1.133 pgi/15.3.0  
cray-mpich/7.2.2 Python/3.5.1-CrayGNU-2016.03
```

```
# --- TOOLS ---
```

```
ddt/6.0  
perftools/6.2.4  
scalasca/1.4.2 scorep/1.4.2 vampir/9.0  
visit/2.10 paraview/5.0
```

```
# --- LIBS ---
```

```
cray-libsci/13.0.4  
cray-hdf5-parallel/1.8.14 cray-netcdf-hdf5parallel/4.3.3.1  
cray-petsc-64/3.5.3.1 cray-tpsl-64/1.5.0 cray-trilinos/11.12.1.3  
fftw/3.3.4.3
```

```
# --- GPU ---
```

```
craype-accel-nvidia35  
cudatoolkit/6.5.14-1.0502.9836.8.1  
cray-libsci_acc/3.1.1
```

```
# --- APPS ---
```

```
cp2k/2.6 espresso/5.1.2 gromacs/5.0.6  
lammps/10Feb15 namd/2.9 nwchem/6.3r2 vasp/5.3
```

Compiling the code: module show/help

module avail cray-hdf5-parallel

```
course51@daint103:~ module avail cray-hdf5

----- /opt/cray/modulefiles -----
cray-hdf5/1.8.11          cray-hdf5-parallel/1.8.11
cray-hdf5/1.8.12          cray-hdf5-parallel/1.8.12
cray-hdf5/1.8.13          cray-hdf5-parallel/1.8.13
cray-hdf5/1.8.14(default) cray-hdf5-parallel/1.8.14(default)
cray-hdf5/1.8.16          cray-hdf5-parallel/1.8.16
cray-hdf5/1.8.9           cray-hdf5-parallel/1.8.9
```

module show cray-hdf5-parallel

```
module show cray-hdf5-parallel # CCE
      setenv  HDF5_DIR /opt/cray/hdf5-parallel/1.8.14/CRAY/83

module swap PrgEnv-cray PrgEnv-gnu # GNU
module show cray-hdf5-parallel # GNU
      setenv  HDF5_DIR /opt/cray/hdf5-parallel/1.8.14/GNU/4.9
```

module help cray-hdf5-parallel

```
course51@daint103:~ module help cray-hdf5-parallel
# Module Specific Help for 'cray-hdf5-parallel/1.8.14'
# Doc: http://www.hdfgroup.org/HDF5/doc/index.html
# ...
```

Compiling the code: module load/rm

module load cray-hdf5-parallel

```
course51@daint103:~ module load cray-hdf5-parallel
course51@daint103:~ module list
course51@daint103:~ which h5dump
# /opt/cray/hdf5/1.8.14/bin/h5dump
```

module rm cray-hdf5-parallel

```
course51@daint103:~ module rm cray-hdf5-parallel
course51@daint103:~ which h5dump
# not set
```

Compiling the code: mini-app

Get the src

```
course51@daint103:~ git clone \  
https://github.com/eth-cscs/SummerSchool2016.git \  
SummerSchool2016.git  
# Cloning into 'SummerSchool2016.git'...
```

Compile the Fortran version

```
course51@daint103:~ cd SummerSchool2016.git/miniapp/serial/fortran/  
course51@daint103:~ make clean  
# rm -f main *.o *.i *.mod output.*  
  
course51@daint103:~ make  
# ftn -O3 -fopenmp -c stats.f90 -o stats.o  
# ftn -O3 -fopenmp -c data.f90 -o data.o  
# ftn -O3 -fopenmp -c operators.f90 -o operators.o  
# ftn -O3 -fopenmp -c linalg.f90 -o linalg.o  
# ftn -O3 -fopenmp *.o main.f90 -o main
```

Compile the C++ version

```
course51@daint103:~ cd SummerSchool2016.git/miniapp/serial/cxx/  
course51@daint103:~ make clean; make  
# CC -O3 -fopenmp -c stats.cpp -o stats.o  
# CC -O3 -fopenmp -c data.cpp -o data.o  
# CC -O3 -fopenmp -c operators.cpp -o operators.o  
# CC -O3 -fopenmp -c linalg.cpp -o linalg.o  
# CC -O3 -fopenmp *.o main.cpp -o main
```



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Running the code

Running the code: interactive session (1)

salloc

- The job submission system used at CSCS is `Native SLURM`.
- `salloc` allows to connect to the compute node (`man salloc`)
- `salloc --res=summerschool`: our reservation for the course
- `salloc -N`: number of compute nodes (8 cores max per node)
- `salloc -t`: duration of the session in minutes (default is 1h)
- http://user.cscs.ch/getting_started/running_jobs/

salloc

```
course51@daint102:~ salloc --res=summerschool -t120 -N2
salloc: Pending job allocation 1145157
salloc: job 1145157 queued and waiting for resources
salloc: job 1145157 has been allocated resources
salloc: Granted job allocation 1145157
course51@daint102:~ echo "Hourra :-)"
```

Running the code: interactive session (2)

Other useful Slurm commands

- `squeue -u $USER` : what is the status of my salloc session ?
- `scontrol show job $JOBID` : more details about my session ?
- `scancel $JOBID` : cancel my session

squeue/scontrol/scancel

```
course51@daint102:~ squeue -u course51
      JOBID      USER      ACCOUNT    ST          REASON  NODES    PRIORITY
      1145201   course51  courses    PD           None         2       15044

course51@daint102:~ scontrol show job 1145201
JobId=1145201 Name=bash
UserId=course51(22854) GroupId=courses(30340)
Priority=15044 Nice=0 Account=courses QOS=normal
JobState=RUNNING Reason=None Dependency=(null)
...
NodeList=nid000[68-69]
BatchHost=daint03
NumNodes=2 NumCPUs=16 CPUs/Task=1 ReqB:S:C:T=0:0:*:*
Socks/Node=* NtasksPerN:B:S:C=0:0:*:*1 CoreSpec=0
MinCPUsNode=1 MinMemoryCPU=2G MinTmpDiskNode=0
WorkDir=/users/course51

course51@daint102:~ scancel 1145201
```

Running the code: interactive session (3)

```
srun [options] myexecutable
```

- To run an application on the compute nodes, use the srun command.
- Adapt the command to your needs:
 - `-n`: Total number of MPI tasks
 - `--ntasks-per-node`: Number of MPI tasks per compute node (8 max per node)
 - `-c`: Number of OpenMP threads per task

Running the miniapp (serial version)

```
course51@daint102:~ srun -n 1 ./main 256 256 200 0.01
=====
                        Welcome to mini-stencil!
version  :: Fortran90 serial
mesh     :: 256 * 256      dx = 3.9215688593685627E-3
time     :: 200 time steps from 0 .. 1.00000000000000002E-2
=====
simulation took 5.2622885461896658 seconds
13761 conjugate gradient iterations
2615.0219394495398 CG iterations per second
1328 nonlinear newton iterations
-----
Goodbye!
```




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Editing the code

Editing the code

Many text editors are available

- `vim` filename (X version: gvim)
- `emacs` -nw filename (X version: emacs)
- X only:
 - `gedit`
 - `/apps/pilatus/nedit/5.5/Linux-x86/bin/nedit`



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Moving data from/to CSCS

Moving data: scp

SCP

- Getting a file: `scp USER@FROM:remotefile localfile`
- Sending a file: `scp localfile USER@TO:remotefile`
- Add the `-rp` flag to `scp` to copy an entire directory

Getting 1 (or more) file from CSCS

```
jg@mylaptop:
scp course51@ela.cscs.ch:~/SummerSchool2016.git/miniapp/*.pdf .
# miniapp.pdf          30%  576KB 140.9KB/s   00:09 ETA
# miniapp.pdf          52%  992KB 128.1KB/s   00:06 ETA
# miniapp.pdf          85% 1600KB 116.1KB/s   00:02 ETA
# miniapp.pdf         100% 1875KB 104.2KB/s   00:18

jg@mylaptop: evince miniapp.pdf &
```

Sending 1 (or more) file to CSCS

```
jg@mylaptop: scp          mycode.c          course51@ela.cscs.ch:~
```