

# Exercises – Grid Workshop

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## 1. Installation of the ARC client

Install and configure ARC on your computer. Documentation on installing ARC can be found at <http://www.nordugrid.org/arc/>

When configured the command "arcproxy --version" should display:

```
arcproxy version 1.1.0
```

## 2. Installation of grid-certificate

*Terena Certificate:*

Use the information at

[http://snicdocs.nsc.liu.se/wiki/Grid\\_certificates#Exporting\\_Terena\\_certificate\\_for\\_use\\_with\\_Grid\\_tools](http://snicdocs.nsc.liu.se/wiki/Grid_certificates#Exporting_Terena_certificate_for_use_with_Grid_tools)

To export your grid-certificate and install it in the .globus directory in your home directory.

Make sure the file userkey.pem mode is set to 400 and usercert.pem to 644.

*NorduGrid Certificate:*

Make sure you install the usercert.pem received from the NorduGrid CA in you .globus directory in your home directory.

*All*

Verify your certificate using the command **grid-cert-info -subject**. This command should display your Distinguished Name (DN) in the terminal window.

## 3. Configure the ARC grid client.

Use the information in the slides to configure your ARC client for SweGrid. If configured correctly, the command **arcinfo** should display the following:

```
Execution Service: arc-ce01.pdc.kth.se
  URL: ARC0:ldap://arc-ce01.pdc.kth.se:2135/nordugrid-cluster-name=arc-
ce01.pdc.kth.se,Mds-Vo-name=local,o=Grid
  Queue: easy
  Health State: ok
```

```
Execution Service: svea.c3se.chalmers.se
  URL: ARC0:ldap://svea.c3se.chalmers.se:2135/nordugrid-cluster-
name=svea.c3se.chalmers.se,Mds-Vo-name=local,o=Grid
  Queue: svea
  Health State: ok
```

...

```
Execution Service: siri.lunarc.lu.se
URL: ARC0:ldap://siri.lunarc.lu.se:2135/nordugrid-cluster-
name=siri.lunarc.lu.se,Mds-Vo-name=local,o=Grid
Queue: arc
Health State: ok
```

#### 4. Querying the information system

Use the **arcinfo** command to query more information on specific resources. Use the **--cluster** switch to specify a resource explicitly. Use the **--long** switch to show more information on a resource. The following list shows the available resources at SweGrid:

- arc-ce.smokerings.nsc.liu.se - NSC
- siri.lunarc.lu.se - LUNARC
- grad.uppmax.uu.se - UPPMAX
- svea.c3se.chalmers.se - C3SE
- jeannedarc.hpc2n.umu.se - HPC2N
- arc-ce01.pdc.kth.se - PDC

#### 5. Submitting testjobs

Use the **arctest** command to send testjobs to the resources. Use the built-in help using the **--help** switch to submit a job that prints the environment on a node.

Use **arcstat** to monitor the status of the submitted jobs.

Use **arckill** to kill a running job.

Use **arcget** to retrieve a finished job.

Use **arcclean** to clean a job that has not been downloaded with **arcget**.

#### 6. Creating a simple job

Create a job description and needed files for a job that prints the current date and lists the files in the current working directory. The wall time of the job should be 5 minutes.

#### 7. Creating a simple job sweep

Modify the code from <http://www.scipy.org/Cookbook/CoupledSpringMassSystem> to be a job sweep modifying the m1 parameter in the range  $1.0 \geq m1 \leq 2.0$ .

Executing the following code on the grid with different parameters:

```
python two_springs_solver.py > two_springs.dat
```

The python-files for this exercise can be found at the following URLs:

- [dl.dropbox.com/u/2888586/grid\\_workshop/ex7/two\\_springs.py](https://dl.dropbox.com/u/2888586/grid_workshop/ex7/two_springs.py)
- [dl.dropbox.com/u/2888586/grid\\_workshop/ex7/two\\_springs\\_solver.py](https://dl.dropbox.com/u/2888586/grid_workshop/ex7/two_springs_solver.py)
- [dl.dropbox.com/u/2888586/grid\\_workshop/ex7/two\\_springs\\_plot.py](https://dl.dropbox.com/u/2888586/grid_workshop/ex7/two_springs_plot.py)

## 8. Creating a MPI Job

Submit the following MPI example to grid using 8 cores:

```
program hello_mpi

use mpi

call mpi_init(ierr)
call mpi_comm_size(MPI_COMM_WORLD,npes, ierr)
call mpi_comm_rank(MPI_COMM_WORLD,irank,ierr)

print *,'Hello MPI! I am ',irank,' of ',npes

call mpi_finalize(ierr)

end program
```

Use the environment ENV/MPI/OPENMPI-1.4.4/GCC