

## Useful links

1. <https://userinfo.surfsara.nl/systems/lisa/user-guide/creating-and-running-jobs>

## Downloads

1. [UvA VPN](#)
  - a. Accessing the UvA network
2. [WinSCP:](#)
  - a. This enables the transfer of files between your local computer and the remote cluster using a GUI (Jorge: this is the cp command that I was referring to, you can use either this one or the sftp protocol).
  - b. Host name: `csn-h0.science.uva.nl`
  - c. Username & password: UvANetID & normal password
3. **Note:** Alternative to WinSCP: use `sftp` to access the cluster in a second command window:
  - easier way to copy and move files
  - a. `put file.pdf [local → remote]`
  - b. `get file.pdf [remote → local]`

## Entering the cluster

1. Open Windows Powershell
2. `ssh {uvanetid}@csn-h0.science.uva.nl`
3. Same password
4. h0 is the entrance node

## Navigating the cluster

1. Useful [unix commands](#):

<code>man ln</code>	List of commands
<code>cd</code>	Change directory
<code>cd ..</code>	Go up one directory
<code>cd ~</code>	Return to home directory
<code>ls</code>	List of files in current working directory

ls-lh	Long list of files with readable file size
ls-l	List with long format - show permissions
mkdir	Make a new directory
cp	Copy file (file_name new_directory)
mv	Move (file_name new_directory)

2. `cd /data`  
`mkdir {name}`  
`cd {name}`  
`mkdir {data} / {scripts} / {results}`

3. Use WinSCP/ sftp to copy the data over

## Sending a job to the cluster

1. Clone the [repository](#) into your scripts folder on the cluster
2. Be sure to use the command "git pull" to retrieve any changes committed to the GitHub
3. `squeue` - see list of current jobs that are in the queue (not running yet on SLURM)
  - a. Alternative `top -u {user}` (these jobs are currently running on the cluster)
4. `sbatch {slurm file}` → start job
5. `scancel {jobnumber}` → cancel if needed
6. Using the Vim editor:

Esc	Switch to command mode
i	Switch to edit mode
:	Start writing a command
:wq	Save and exit file

7. Example slurm

```
#!/bin/sh
#SBATCH -J Ferret_stuff
#SBATCH --output=/home/cbosman1/slurm/log/pupilMaster.out
#SBATCH --error=/home/cbosman1/slurm/log/pupilMaster.err
#SBATCH --mail-user conrado.bosman@me.com
#SBATCH --mail-type=ALL
#SBATCH --mem=64G

module load matlab
matlab -nodisplay -r "addpath /home/cbosman1/matlab/scripts/pupilProject/; cb_master_get_pupil_LFP; quit"
```

### Template:

```
#!/bin/sh
#SBATCH -J {identifier}
#SBATCH --output=/home/{name}/slurm/log/pupilMaster.out {where error
file should be saved}
#SBATCH --error=/home/{name}/slurm/log/pupilMaster.err {where error
file should be saved}
#SBATCH --mail-user {email address}
#SBATCH --mail-type=ALL
#SBATCH --mem=64G

module load matlab {version 2017a}

matlab -nodisplay -r "addpath /home/{path}/; {scriptname}; quit"
```

### Working with different input parameters:

1. sbatch is not interactive
2. Call multiple main files in the slurm file

```
matlab -nodisplay -r "addpath ~/scripts/; main(1); main(2); main(3);
quit"
```

3. Alternative:

```
matlab -nodisplay -r "addpath ~/scripts/; main; quit"
```

- The main.m file can call a "super\_main.m" file: "super\_main.m" will have a list of random variables that are output and stored somewhere as a variable.
- The main.m file then uses these stored variable as input

### Setting up MVGC

1. Reference paper: [The MVGC multivariate Granger causality toolbox: a new approach to Granger-causal inference](#)

2. Add MVGC toolbox to personal folder in the cluster. Add to path and use start up in matlab script
3. This message may come up in your .err file:

```
[mvgc startup] WARNING: Control System Toolbox(TM) does not seem to be present.  
[mvgc startup] Will use slower scripted routines (see utils/control directory).
```

- a.
- b. “The Matlab Control System Toolbox function `dlyap` solves equations of the form (A.5); if the Control System Toolbox is not available, then the MVGC toolbox function `utils/dlyap_aitr` implements an efficient iterative solver”
- c. This is however a lot slower:
- d. One can add the Matlab Control System Toolbox to your folder on the cluster. Make sure you add it to path in the main script (see “`gc_parfor_cluster`” for an example).

In case of trouble with using ‘startup’

```
origDir = pwd;  
cd ~/matlab/MVGC  
startup  
cd(origDir)
```

## Debugging in the cluster

It is possible to set up a testing bash environment to run the script. This can be set up to create an interactive matlab environment with a display.

Run the following commands to set up the testing environment:

```
srun --pty bash  
module load matlab  
matlab -r "addpath {path to script}; {scriptname}; quit"
```

*Note the final command can be copied from the .slurm file. Be sure to remove the “-nodisplay” flag to be able to see your outputs*