

## Working on ciclad

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ciclad is an IPSL computing server located on the Jussieu campus in Paris, France.

## Documentation

■ <http://ciclad-web.ipsl.jussieu.fr>

■ <http://ciclad-web.ipsl.jussieu.fr/ciclad-utilisation.pdf>

hotline : svp-ciclad@...

## The machines and file systems

The front-end machine can be accessed via the `ciclad.jussieu.ipsl.fr` IP.

Data files must be placed in `/data/` or in the filesystem dedicated to your project.

## Shared account

cf. [Repository for shared files and shared tools](#) / `ipslfs/igcmg/IGCM` belonging to the account `igcmg` (`${HOME}=/home/igcmg`)

```
$ id -a igcmg
uid=31575(igcmg) gid=31575(igcmg) groups=31575(igcmg)
```

## Individual account

You must belong to the `igcmg` users' group.

cf. to check the result of the command

```
id -a
```

## How to define your environment

Add the following line in your login file (e.g. `/home/igcmg/.bashrc`) :

```
. /home/igcmg/.atlas_env_ciclad_ksh
```

for the FORTRAN compiler, the NetCDF library, ferret,...

## To find out the selected implementation in the MPI library

```
mpi-selector --query
```

If the answer is different from :

```
default:openmpi-1.4.2-gfortran-x86_64
level:system
```

or if you are not using the FORTRAN compiler `gfortran` you must change this environment with `mpi-selector --set` by choosing among the possibilities returned by the command `mpi-selector --list`.

## Job manager commands

torque/maui tool

## End-of-job messages

To receive the end-of-job messages returned by the job itself (e.g. end of simulation, error,...) you must specify your email address in the file `${HOME}/.forward`.

**How to choose the number of processes?**

**Example of job for a MPI executable**

**libIGCM specificities on ciclad**