

Getting started with LMDZ

LMDZ Training course December 9-10-11 2013

The LMDZ Team

How to install the LMDZ GCM : Get the sources and compile

Choosing which version of LMDZ to work with

Choose between the different available versions on the LMDZ web site:

<http://lmdz.lmd.jussieu.fr/utilisateurs/distribution-du-modele>
(in French)

<http://lmdz.lmd.jussieu.fr/utilisateurs/distribution-du-modele-en>
(in English)

NB: when on a French page, try clicking on the  flag to get English version

Ask the LMDZ team for more information on which versions are actually used :
lmdz-svp@lmd.jussieu.fr

Choosing which LMDZ version to work with

1 – « production » or reference versions :

- LMDZ4_AR5

svn co http://svn.lmd.jussieu.fr/LMDZ/LMDZ4/branches/LMDZ4_AR5

- the most tested/validated version
- production version used for CMIP5
- runs with « old physics » package (« ancienne physique »)

- LMDZ5_AR5

svn co http://svn.lmd.jussieu.fr/LMDZ/LMDZ4/branches/LMDZ5_AR5

- « new physics » version used for CMIP5

2 – « testing » version :

svn co http://svn.lmd.jussieu.fr/LMDZ/LMDZ5/branches/testing_LMDZ5

- less tested than reference version but more tested than development version
- runs with « old or new physics » packages (« ancienne et nouvelle physique »)

3 - « development » version :

svn co <http://svn.lmd.jussieu.fr/LMDZ/LMDZ5/trunk>

- Beware! This is often updated and may not be thoroughly tested!
- Some revisions are more tested than others; these development versions eventually become « testing » versions.

How to install LMDZ : getting the source codes & compiling

There are different ways to install LMDZ ; your choice depends on the machine you will be using and the type of simulation (long, test or development) you will run.

1) Installing LMDZ « by hand »

- get each source code one by one (IOIPSL, ORCHIDEE, LMDZ) and link them with the netcdf library installed on your machine.

2) Installing LMDZ using the install.sh script

- see afternoon tutorial
- the script will download the source codes needed and will compile them
- this method is recommended for Linux PC and used for developments and tests.

3) Installing LMDZ using modipsl and libIGCM

- you will need to install one of the configuration defined by modipsl (for example LMDZOR_v5)
⇒ follow the IPSL Training course.
- this is recommended for IDRIS and CCRT/TGCC and for long simulations as it provides tested reference versions and scripts for launching and following long simulations.

1) Installing LMDZ « by hand »

More information on the LMDZ site :

<http://lmdz.lmd.jussieu.fr/utilisateurs/guides/lmdz-pas-a-pas> (in French)

To do once and for all :

Install netcdf

Install IOIPSL

```
svn co http://forge.ipsl.jussieu.fr/igcmg/svn/modipsl/trunk modipsl
cd modipsl/util
./model IOIPSL
Modify AA_make.gdef
./ins_make [-t g95 ou -t egi]
cd ../modeles/IOIPSL/src
gmake
```

Get LMDZ

```
svn co http://svn.lmd.jussieu.fr/LMDZ/LMDZ4/branches/LMDZ4_AR5 LMDZ4
```

Compile LMDZ

Fill in the paths for netcdf, IOIPSL et ORCHIDEE libraries in the **makegcm** file or in the arch-XXX.path file if you use **makelmdz_fcm** or **makelmdz** scripts

```
./makelmdz_fcm -d « grid_resolution » -arch « XXX » gcm
./makelmdz -d « grid_resolution » -arch « XXX » gcm
./makegcm -d « grid_resolution » gcm
```

2) Installing LMDZ using the install.sh script (preliminary remarks)

Further details here : <http://lmdz.lmd.jussieu.fr/utilisateurs/script-install.sh-en> (in English)

- The install.sh script **compiles the model and the associated libraries** using any of the **g95**, **gfortran** , **ifort** or **pgf90** compilers and standard shell commands and utilities (gcc, wget, gunzip, tar, ...).
- This script will download the « testing » branch of LMDZ. The main modifications between different versions and the link between the proposed versions and the SVN revision numbers are summarized in the LISMOI (README) file.
- To modify some of the installation options, for example the directory where the model will be put or to inhibit the compilation of IOIPSL and/or NETCDF libraries you will have to modify variables defined in the install.sh script.
- Any of the **makegcm**, **makelmdz_fcm** or **makelmdz** scripts can be used to compile the model and may be used afterwards to recompile the model.
 - If **compile_with_fcm=1**, makelmdz_fcm will be used to compile the model.
 - If **compile_with_fcm=0**, makelmdz will be used.

2) Installing LMDZ using the install.sh script (how to download and launch it)

```
> wget http://www.lmd.jussieu.fr/~lmdz/Distrib/install.sh
> chmod +x install.sh
( eventually modify some variables in the script )
> ./install.sh
```

The variables to modify :

version	see LISMOI file (to eventually know the available versions)
MODEL	PATH to install the model
veget=1	to compile with a « frozen » ORCHIDEE version
netcdf=1	to install netcdf library
bench=1	to launch a test run at the end of compiling stage
compilo	compiler to use : g95 / ifort / pgf90 / gfortran
compile_with_fcm	=1 use of makelmdz_fcm (available since revision 1578) =0 use of makelmdz (makegcm before)
grid_resolution	possible choices for the included bench : 32x24x11, 48x36x19, 96x71x19

3) Installing LMDZ using modipsl and libIGCM

- Recommended for IDRIS and CCRT/TGCC for long runs

See :

<https://forge.ipsl.jussieu.fr/igcmg/wiki/platform/documentation> (in French)

Main configurations defined in modipsl :

- [LMDZOR v5.2 configuration](#) : LMDZ coupled with ORCHIDEE (**MPI-OMP**) !

- [LMDZOR v5 configuration](#) : LMDZ coupled with ORCHIDEE

- [LMDZINCA configuration](#) : LMDZ coupled with INCA

- CMIP5 ocean-atmosphere configurations : [IPSLCM5A](#), [IPSLCM5B](#),
[IPSLCM5_v5](#)

Main points :

- Download modipsl
- Choose a given configuration (containing the different models)
- Compile all models using one main Makefile
- Launch the run (simulation and post-treatment) using libIGCM

• **modipsl and libIGCM courses** - « highly recommended » also for false beginners - are given at “Maison de la simulation” by the IPSL Plate-forme group.

► Next courses : **Spring 2014**

<https://forge.ipsl.jussieu.fr/igcmg/wiki/platform/en/training>

What you need to run the LMDZ model

- **Parameters files** :
 - run.def, gcm.def, physiq.def, config.def, traceur.def
- **Start files** : ♣
start.nc, startphy.nc
These files are created by the ce0l.e program or may be the result of previous runs
- **Boundary conditions file** : ♣
limit.nc
Created by ce0l.e
- **Executable (LMDZ) file** :
gcm.e
- **Some optional input files (depending on the simulation)** : ♣
aerosols.nc, climoz_LMDZ.nc, nudging input files (u.nc, v.nc,..)

♣ : these files have to be interpolated on the horizontal grid of the model

How to run the model

```
> ls  
start.nc startphy.nc limit.nc config.def gcm.def orchidee.def  
physiq.def run.def traceur.def gcm.e
```

```
> ./gcm.e
```

or

```
> ./gcm.e > out_lmdz.x 2>&1
```

To carry on a simulation that has been run, you have to copy the restart files obtained at the end of the previous run as new initial start files:

```
> mv restart.nc start.nc  
> mv restartphy.nc startphy.nc
```

```
> ./gcm.e
```

Has your run completed successfully ? (1)

YES

▶ you will then have a message saying « **Everything is cool** » on the standard output or in the output text file.

▶ The code will have created 2 restart files

`restart.nc` and `restartphy.nc`

needed to carry on your run

▶ and some output diagnostic files

`histday.nc`, `hismth.nc`, etc. ...

to explore/view using `ferret`, `grads`, ...

Has your run completed successfully ? (2)

NO

You must find out what the problem is !!!

Look for an error message in the output text file. Search for one of the following key words/phrase « **Houston, we have a problem** », « **STOP** », « **hgardfou** », «**integrdr: negative surface pressure** » etc.

Different typical errors :

- technical problem : a missing input file, an error in one of the *.def file
- **problem with the model's stability.**

If it is an instability in the physics, **hgardfou** will probably detect it as it tests if the model temperature has realistic values.

If it is an instability in the dynamics, you will most often end up with a « **negative surface pressure** » error message

> In any of these cases you will most probably have to adjust some flags in the .def files. See the talk tomorrow.

- **you have some source code modifications that might not have been thoroughly tested or validated.**

Some basic svn commands

Basic svn commands (1)

To get last revision of a directory with its sub-directories :

```
> svn co http://svn.lmd.jussieu.fr/LMDZ/LMDZ5/trunk LMDZ5
```

or

```
> svn co http://svn.lmd.jussieu.fr/LMDZ/LMDZ4/branches/LMDZ4\_AR5 LMDZ4
```

To get a particular revision, for example :

```
> svn co -r 1575 http://svn.lmd.jussieu.fr/LMDZ/LMDZ5/trunk LMDZ5
```

Basic svn commands (2)

> **svn info** : to know what revision and branches you've downloaded

```
[jghattas@woodyard trunk]$ svn info
Chemin : .
URL : http://svn.lmd.jussieu.fr/LMDZ/LMDZ5/trunk
Racine du dépôt : http://svn.lmd.jussieu.fr/LMDZ
UUID du dépôt : e51f81be-29bc-408f-98e3-ee85b5628ff9
Révision : 1572
Type de nœud : répertoire
Tâche programmée : normale
Auteur de la dernière modification : lguez
Révision de la dernière modification : 1572
Date de la dernière modification: 2011-09-13 10:58:59 +0200 (mar. 13 sept. 2011)
```


Basic svn commands (3)

> **svn stat** : to know which files have been modified with respect to the extracted version

```
[jghattas@woodyard libf]$ svn stat
?      grid/dimensions.h
?      grid/dimension/dimensions.32.24.11
M      phylmd/concvl.F
M      phylmd/phys_output_mod.F90
```

> **svn -u stat** : to see modifications with respect to the last version on the svn server.

```
[jghattas@woodyard libf]$ svn -u stat
*      1572    phylmd/mod_phys_lmdz_omp_data.F90
*      1572    phylmd/fisrtilp_tr.F
M      *      1572    phylmd/phys_output_mod.F90
*      1572    phylmd/mod_phys_lmdz_mpi_data.F90
*      1572    phylmd/soil.F90
M      *      1572    phylmd/concvl.F
*      1572    phylmd/conf_phys.F90
*      1572    phylmd/sw_aeroAR4.F90
?      grid/dimensions.h
?      grid/dimension/dimensions.32.24.11
État par rapport à la révision    1576
```

Basic svn commands (4)

> **svn diff** : to check the differences between different versions

```
[jgips1@dhcp-236 phylmd]$ svn diff phys_output_mod.F90
Index: phys_output_mod.F90
```

```
=====
--- phys_output_mod.F90      (révision 1572)
+++ phys_output_mod.F90      (copie de travail)
@@ -1695,10 +1695,11 @@
     IF (nqtot>=3) THEN
         DO iq=3,nqtot
             iiq=niadv(iq)
+            print*,'iq = ',iq
             o_trac(iq-2) = ctrl_out((/ 4, 5, 1, 1, 1, 10 /))

-            ENDDO
+            END
         ENENDIF
```

> **svn diff -r x1** : differences between the x1 revision and the working version
> **svn diff -r x1:x2** : differences between x1 and x2 revisions.
Use HEAD to point to the last revision

Basic svn commands (5)

> **svn update** : to update your version with respect to the last revision of the branch you initially downloaded from.

> **svn update -r x1** : to update your version with respect to revision x1 of the branch you initially downloaded from.

Note that this “update” can be towards an older revision.