



# *Installing the LMDZ model. Tutorial 1*

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# *First steps with LMDZ GCM: install\_lmdz.sh (1)*

```
wget http://www.lmd.jussieu.fr/~lmdz/pub/install_lmdz.sh  
chmod u+x install_lmdz.sh
```

```
./install_lmdz.sh -v 20171119.trunk -d 32x32x39
```

**1/ Get sources (netcdf, IOIPSL, LMDZ, ORCHIDEE)**

modipsl.20171119.trunk.tar.gz

bench\_lmdz\_32x32x39.tar.gz

**2/ Prepares makefiles and compiles the codes**

**3/ Launch a 1-day long test run at 32x32x39**



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# *First steps with LMDZ GCM: install\_lmdz.sh options (2)*

```
./install_lmdz.sh -h
```



# *First steps: What you need to run the LMDZ GCM (1)*

- Executable (LMDZ) file :

gcm.e

- Parameters files :

run.def, gcm.def, physiq.def, config.def, traceur.def

- Start files v :

- start.nc, startphy.nc

These files are created by the ce0l.e program or may be the result of previous runs

- Boundary conditions file v :

- limit.nc

Created by ce0l.e

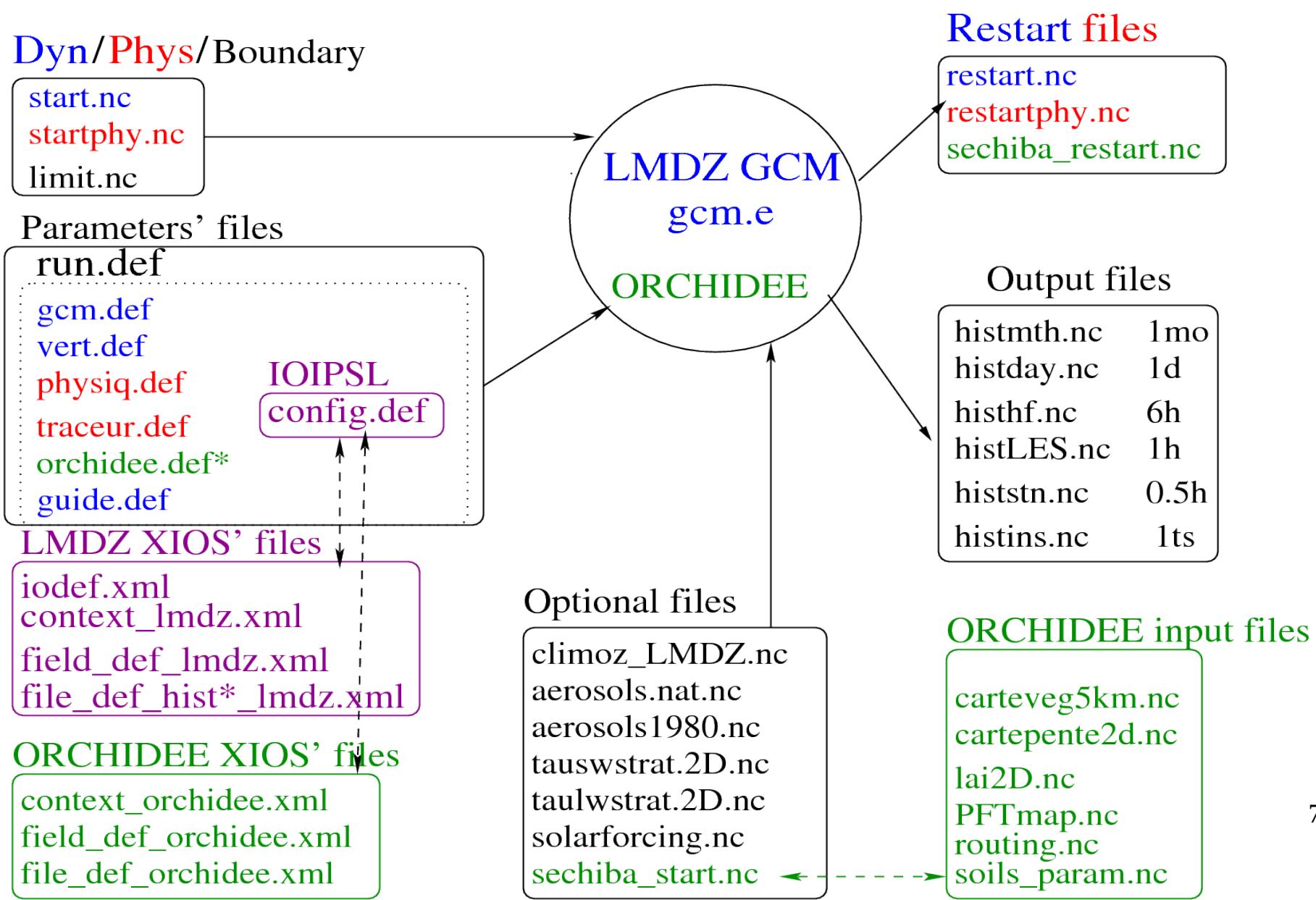
- Some optional input files v (depending on the simulation) :

- aerosols.nc, climoz\_LMDZ.nc, nudging input files (u.nc, v.nc,...), etc

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

# *First steps: What you need to run the LMDZ GCM (2)*

## I/O files for a LMDZ run





# *Installing the model LMDZ*

There are 3 ways to install LMDZ ; right choice depends on the machine you are using and the type of simulation (long, test or development) you run.

## **1) Installing LMDZ by hand**

- get source code for each component you need (IOIPSL, ORCHIDEE, LMDZ) and link them with the netcdf library installed on your machine.

## **2) 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, TGCC, CINES and for long simulations, as it provides tested reference versions and scripts for launching and monitor long simulations.

## **3) Installing LMDZ using the `install_Lmdz.sh` script**

- this is what you're about to do.
- the script will download the source codes needed and will compile them
- recommended method for Linux PC ; used for developments and tests.



# *Using MODIPSL to install the model*

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- this is what you're about to do.
- the script will download the source codes needed and will compile them
- recommended method for Linux PC ; used for developments and tests.



# *Using MODIPSL to install the model*

- Recommended for IDRIS, TGCC and CINES for long runs. See :  
[https://forge.ipsl.jussieu.fr/igcmg\\_doc/wiki/Doc](https://forge.ipsl.jussieu.fr/igcmg_doc/wiki/Doc)

## Main configurations defined in modipsl :

- **LMDZOR\_v5.2** : LMDZ coupled with ORCHIDEE (**MPI-OMP**)
- **LMDZOR\_v5** : LMDZ coupled with ORCHIDEE
- **LMDZINCA** : LMDZ coupled with INCA
- CMIP5 ocean-atmosphere : **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 Plateforme group.

► Next courses : **To Be Determined**

[https://forge.ipsl.jussieu.fr/igcmg\\_doc/wiki/Train](https://forge.ipsl.jussieu.fr/igcmg_doc/wiki/Train)



# Using MODIPSL to install the model

More information on the LMDZ site (in French) :

<http://lmdz.lmd.jussieu.fr/utilisateurs/guides/lmdz-pas-a-pas>

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/LMDZ5/branches/testing LMDZ5
```

Compile LMDZ

Fill in the paths for netcdf, IOIPSL et ORCHIDEE libraries in the arch-XXX.path used by **makelmdz\_fcm** or **makelmdz** scripts.

```
./makelmdz_fcm -d <grid_resolution> -arch <XXX> gcm
./makelmdz      -d <grid_resolution> -arch <XXX> gcm
```



# *Getting the source codes and compiling*

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# Using *install\_lmdz.sh* – Contents

Further details (in French), in particular the main modifications between versions:

<http://www.lmd.jussieu.fr/~lmdz/Distrib/LISMOI.trunk>

It will do most of the work for you, using standard shell tools and commands (gcc, wget, gunzip, tar, ...):

- Download the required codes archives
- **Install** ancillary **libraries** (netcdf, modipsl, ioipsl)
- Install land surface model ORCHIDEE if needed
- Choose adequate compiler options and build a Makefile
- **Install LMDZ** using any of **makelmdz\_fcm** and **makelmdz** scripts
- **Run a test bench**

Some parameters can be modified using command line options (short manual if you launch the script with **-h** option):

- |                              |  |
|------------------------------|--|
| • <b>-v version:</b>         | choose a version/branch [YYMMDD.]trunk/testing |
| • <b>-r release_nb:</b>      | choose a particular svn release                |
| • <b>-parallel mode:</b>     | sequential/mixed parallelism                   |
| • <b>-d grid_resolution:</b> | choose model grid resolution                   |
| • <b>-bench:</b>             | launch or not a test bench                     |
| • <b>-name MODEL:</b>        | choose model folder name                       |
| • ...                        | LMDZversion.release                            |
- mpi\_omp or none**
- nlonxnlatxnlev**
- 1/0**



# Using *install\_lmdz.sh* – Options

Some other parameters are defined in the script itself only ; you have to edit them:

• <b>compilo</b>	choose compiler	<b>gfortran / mpif90 / pgf90 / ifort</b>
• <b>getlmdzor</b>	get a LMDZ archive	<b>0/1</b>
• <b>netcdf</b>	install NetCDF library	<b>0/1</b>
• <b>check_linux</b>	check basic tools availability	<b>0/1</b>
• <b>ioipsl</b>	install IOIPSL I/O library	<b>0/1</b>
• <b>veget</b>	install ORCHIDEE land surface model	<b>0/1</b>
• <b>compile_with_fcm</b>	use makelmdz_fcm or makelmdz	<b>0/1</b>
• ...		

- If you need to recompile later, use makelmdz[\_fcm].

This script is in constant evolution. Currently: OK for ADA (IDRIS) and a linux PC.

- Later: more machines, with XIOS.



# *Using install.sh – Download and launch it*

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

➡ **TP1**



# *Choosing which LMDZ version 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)

**NB:** use the flag to switch from a French to an English page, for existing translated pages.

Ask the LMDZ team for more information on which versions are actually used :

[lmdz-svp@lmd.jussieu.fr](mailto: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](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/LMDZ5/branches/LMDZ5\\_AR5](http://svn.lmd.jussieu.fr/LMDZ/LMDZ5/branches/LMDZ5_AR5)

- ***new physics*** version used for CMIP5

## 2 – *testing version :*

svn co [http://svn.lmd.jussieu.fr/LMDZ/LMDZ5/branches/testing\\_LMDZ5](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.



# *Running the model*

```
ls
```

```
start.nc startphy.nc limit.nc config.def gcm.def orchidee.def  
physiq.def run.def traceur.def gcm.e L???.def
```

```
./gcm.e
```

or

```
./gcm.e > lmdz.out 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 ?*

**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](#), [histmth.nc](#), etc. ...

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



# *Has your run completed successfully ?*

**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**, **integrd: 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.**

Instability in the physics are likely to be detected by **hgardfou**, which checks the model temperature has realistic values.

Instability in the dynamics most often end up the run 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**



# *Some 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 for the last “robust” version :

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

To get a particular revision, for example :

```
svn co -r 2692 http://svn.lmd.jussieu.fr/LMDZ/LMDZ5/trunk LMDZ5
```



# *Some basic svn commands* (2)

To know what revision and branches you've downloaded:

```
svn info
```

```
Path: .
URL: http://svn.lmd.jussieu.fr/LMDZ/LMDZ5/trunk/libf
Repository Root: http://svn.lmd.jussieu.fr/LMDZ
Repository UUID: e51f81be-29bc-408f-98e3-ee85b5628ff9
Revision: 2699
  2699
Node Kind: directory
Schedule: normal
Last Changed Author: oboucher
Last Changed Rev: 2699
Last Changed Date: 2016-11-07 22:23:46 +0100 (Mon, 07 Nov 2016)
```



# *Some basic svn commands* (3)

To know which files have been modified with respect to the extracted version:

```
svn stat
```

```
[rdzt079@ada337: misc]$ svn stat
?      regr_lint_m.F90
!      regr1_lint_m.F90
!
?      regr3_lint_m.F90
?
?      regr_conserv_m.F90
!
?      regr1_conserv_m.F90
M      slopes_m.F90
```

To see modifications with respect to last version on svn server:

```
svn -u stat
```

```
[rdzt079@ada337: misc]$ svn -u stat
?                  regr_lint_m.F90
!          2699    regr1_lint_m.F90
!
!          2699    regr3_lint_m.F90
?
?          2699    regr_conserv_m.F90
!
?          2699    regr1_conserv_m.F90
M          2699    slopes_m.F90
Status against revision:  2723
```



# Some basic svn commands (4)

To check the differences between different versions:

`svn diff`

```
Index: regr_pr_int_m.F90
=====
--- regr_pr_int_m.F90 (revision 2699)
+++ regr_pr_int_m.F90 (working copy)
@@ -27,7 +27,7 @@
     use netcdf95, only: nf95_inq_varid, handle_err
     use netcdf, only: nf90_get_var
     use assert_m, only: assert
-    use regr1_lint_m, only: regr1_lint
+    use regr_lint_m, only: regr_lint
     use mod_phys_lmdz_mpi_data, only: is_mpi_root
     use mod_grid_phy_lmdz, only: nbp_lon, nbp_lat, nbp_lev
     use mod_phys_lmdz_transfert_para, only: scatter2d
@@ -96,7 +96,8 @@
 
     ! Regrid in pressure at each horizontal position:
     do i = 1, klon
-        v3(i, nbp_lev:1:-1) = regr1_lint(v2(i, :), (/0., plev/), pplay(i, nbp_lev:1:-1))
+        call regr_lint(1,v2(i,:),(/0.,plev/),pplay(i,nbp_lev:1:-1), &
+                      v3(i,nbp_lev:1:-1))
         ! (invert order of indices because "pplay" is in descending order)
     end do
```

To check the differences between revision x1 and the working version:

`svn diff -r x1`



# *Some basic svn commands* (5)

To check the differences between x1 and x2 revisions:

```
svn diff -r x1:x2
```

Use **HEAD** to point to the last revision

To update your version with respect to the last revision of the branch you initially downloaded from:

```
svn update
```

To update your version with respect to revision x1 of the branch you initially downloaded from:

```
svn update -r x1
```

*Note that this **update** can be towards an older revision.*