

# Set Up and Run WRF (ARW-Ideal, ARW-real, and NMM-real)

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#### WRF System Flowchart



## Outline

- Running WRF code
  - Before you run..
  - Running idealized case
  - Running ARW real-data case
  - Running NMM real-data case
- Basic runtime options for a *single* domain run (*namelist*)
- Check output
- Simple trouble shooting



• Running a nested case: later

#### Before You Run ..

• Check and make sure appropriate executables are created in WRFV3/main/ directory:

For ARW:

- ideal.exe
- real.exe
- wrf.exe
- ndown.exe
- tc.exe

For NMM:

- real\_nmm.exe
- wrf.exe

- If you are running a real-data case, be sure that files from WPS are correctly generated:
  - met\_em.d01.\*, for ARW or
  - met\_nmm.d01.\* for NMM



Prepare **namelist**.input for runtime options.

#### WRF test case directories

You have these choices in WRFV3/test/



5

## Steps to Run

- 1. cd to *run/* or one of the *test case* directories
- 2. Link or copy WPS output files to the directory for real-data cases
- 3. Edit *namelist.input* file for the appropriate grid and times of the case
- 4. Run initialization program (*ideal.exe, real.exe*, or *real\_nmm.exe*)
- 5. Run model executable, wrf.exe



#### WRFV3/run directory

README.namelist LANDUSE. TBL GENPARM. TBL SOILPARM. TBL **VEGPARM**. TBL URBAN PARAM. TBL RRTM DATA RRTMG SW DATA RRTMG LW DATA CAM ABS DATA CAM AEROPT DATA ozone.formatted ozone lat.formatted ozone plev.formatted ETAMPNEW DATA tr49t67 tr49t85 tr67t85

WRF

gribmap.txt grib2map.tbl (a few more) these files are model physics data files: they are used to either initialize physics variables, or make physics computation more efficient

## WRFV3/run directory after compile

LANDUSE.TBL SOILPARM.TBL VEGPARM.TBL GENPARM.TBL URBAN\_PARAM.TBL RRTM\_DATA RRTMG\_SW\_DATA RRTMG\_LW\_DATA ETAMPNEW\_DATA tr49t67 tr49t85 tr67t85

An example after ARW real case compile

namelist.input -> ../test/em\_real/namelist.input real.exe -> ../main/real.exe wrf.exe -> ../main/wrf.exe ndown.exe -> ../main/ndown.exe ...(a few more)



## Running an Idealized Case ARW only



- If you have compiled an ideal case, you should have:
   ideal.exe ideal case initialization program
   wrf.exe model executable
- These executables are linked to: WRFV3/run

and

WRFV3/test/em\_test-case

 $\rightarrow$  One can go to either directory to run.



Go to the desired *ideal* test case directory: e.g. cd test/em\_quarter\_ss

If there is `run\_me\_first.csh' in the directory, run it first - this links physics data files to the currect directory:

./run\_me\_first.csh



Then run the ideal initialization program:

./ideal.exe

The input to this program is typically a sounding file (file named *input\_sounding*), or a pre-defined 2D input (e.g. *input\_jet* in **em\_b\_wave** case).

Running ideal.exe creates WRF initial condition file: wrfinput\_d01

Note that wrfbdy file is not needed for idealized cases.



• To run the model interactively, type

./wrf.exe >& wrf.out &

for single processor (serial) or SMP run. Or

```
mpirun -np N ./wrf.exe &
```

for a MPI run (where N is the number of processors requested)

 Successful running of the model executable will create a model history file called wrfout\_d01\_<date>

```
e.g. wrfout_d01_0001-01-01_00:00:00
```



Based on start date defined in namelist

- Edit **namelist**.input file to change options.
- For your own case, you may provide a different sounding.
- You may also edit dyn\_em/ module\_initialize\_<case>.F to change other aspects of the initialization.

#### Note:

- For 2D cases and baroclinic wave case, ideal.exe must be run serially
- For all 2D cases, wrf.exe must be run serially or with SMP



For the 1D case, compile and run serially

## Running **ARW** Real-Data Case



## Running **ARW** Real-Data Case

 If you have compiled the <u>em\_real</u> case, you should have:

real.exe - real data initialization program
wrf.exe - model executable
ndown.exe - program for doing one-way nesting
tc.exe - program for TC bogusing

These executables are linked to:

WRFV3/run

and

WRFV3/test/em\_real



• One can go to either directory to run.

#### WRFV3/test/em\_real directory

LANDUSE.TBL -> ../../run/LANDUSE.TBL GENPARM.TBL -> ../../run/GENPARM.TBL SOILPARM.TBL -> ../../run/SOILPARM.TBL VEGPARM.TBL -> ../../run/VEGPARM.TBL URBAN\_PARAM.TBL -> ../../run/URBAN\_PARAM.TBL RRTM\_DATA -> ../../run/RRTM\_DATA RRTMG\_SW\_DATA -> ../../run/RRTMG\_SW\_DATA RRTMG\_LW\_DATA -> ../../run/RRTMG\_LW\_DATA ETAMPNEW\_DATA -> ../../run/ETAMPNEW\_DATA tr49t67 -> ../../run/tr49t67 tr49t85 -> ../../run/tr67t85

```
namelist.input - editing required
real.exe -> ../../main/real.exe
wrf.exe -> ../../main/wrf.exe
ndown.exe -> ../../main/ndown.exe
...(a few more)
```



- One must successfully run WPS, and create met\_em.\* file for more than one time period
- Link or copy WPS output files to the run directory:

cd test/em\_real

ln -s ../../WPS/met\_em.d01.\*



- Edit namelist.input file for runtime options (at mininum, one must edit <u>&time\_control</u> for start, end and integration times, and <u>&domains</u> for grid dimensions)
- Run the real-data initialization program:

   /real.exe, if compiled serially / SMP, or
   mpirun -np N ./real.exe, or
   mpirun -machinefile file -np N ./real.exe
   for a MPI job
   where N is the number of processors requested, and
   file has a list of CPUs for the MPI job



 Successfully running this program will create model initial and boundary files:





N: the number of time periods processed

• Run the model executable by typing:

```
./wrf.exe >& wrf.out &
Or
```

mpirun -np N ./wrf.exe &

• Successfully running the model will a create model *history* file:

Based on start date defined in namelist

and a *restart* file if **restart\_interval** is set to a time within the range of the forecast time:



wrfrst\_d01\_2008-08-28\_12:00:00

## Running NMM Real-Data Case



## Running NMM Real-Data Case

- If you have compiled the *nmm\_real*, you should have:
   real\_nmm.exe NMM real date initialization
   program
   wrf.exe NMM model executable
- These executables are linked to: WRFV3/run

and

WRFV3/test/nmm\_real



 $\blacktriangleright$  One can go to either directory to run.

#### WRFV3/test/nmm\_real directory

LANDUSE.TBL -> ../../run/LANDUSE.TBL GENPARM.TBL -> ../../run/GENPARM.TBL SOILPARM.TBL -> ../../run/SOILPARM.TBL VEGPARM.TBL -> ../../run/VEGPARM.TBL URBAN\_PARAM.TBL -> ../../run/URBAN\_PARAM.TBL RRTM\_DATA -> ../../run/RRTM\_DATA RRTMG\_SW\_DATA -> ../../run/RRTMG\_SW\_DATA RRTMG\_LW\_DATA -> ../../run/RRTMG\_LW\_DATA ETAMPNEW\_DATA -> ../../run/ETAMPNEW\_DATA tr49t67 -> ../../run/tr49t67 tr49t85 -> ../../run/tr49t85 tr67t85 -> ../../run/tr67t85

namelist.input - editing required real\_nmm.exe -> ../../main/real\_nmm.exe wrf.exe -> ../../main/wrf.exe



...

- One must successfully run WPS, and create met\_nmm. \* file for more than one time period
- Link or copy WPS output files to the run directory:

cd test/nmm real

ln -s ../../WPS/met\_nmm.d01.\*



- Edit namelist.input file for runtime options (at minimum, one must edit &time\_control for start, end and integration time, and &domains for grid dimensions)
- Run the real-data initialization program in MPI:

mpirun -np N ./real\_nmm.exe

Or serially: ./real\_nmm.exe >& real\_nmm.out

• Successfully running this program will create model initial and boundary files:

wrfinput\_d01



wrfbdy\_d01

Run the model executable by typing in MPI:
 mpirun -np N ./wrf.exe

Or serially: ./wrf.exe >& wrf.out

• Successfully running the model will create a model *history* file:

wrfout\_d01\_2005-08-28\_00:00:00

and a *restart* file if **restart\_interval** is set to a time within the range of the forecast time:

wrfrst\_d01\_2008-08-28\_12:00:00



#### **Basic namelist Options**



## What is a namelist?

- A Fortran namelist contains a list of *runtime* options for the code to read in during its execution. Use of a namelist allows one to change runtime configuration without the need to recompile the source code.
- Fortran 90 namelist has very specific format, so edit with care:

```
&namelist-record - start
/ - end
```

• As a general rule:

Multiple columns: domain dependent Single column: value valid for all domains



#### &time\_control

run days = domain 1 option run hours 24 = run minutes = 0,run seconds Ο, = = 2000,start year 2 00 01, start month 01 24, 24 start day 12, 12, start hour 12, start minute 00, 00, 00, 00, 00, start second 00, 2000 2000, 2000 end year end month 01, 01, 01, 25, 25, 25, end day 12, 12, 12, end hour 00, 00, 00, end minute 00, 00, end second 00 interval seconds = 21600history interval = 180 60 nest options frame per outfile = 1000, 1000, 1000 restart interval = 360, restart .true., =



#### Notes on &time\_control

- *run\_*\* time variables:
  - Model simulation length: wrf.exe and domain 1 only
- *start*\_\* and *end*\_\* time variables:
  - Program *real* will use WPS output between these times to produce lateral (and lower) boundary file
  - They can also be used to specify the start and end of simulation times for the coarse grid if *run\_\** variables are not set (or set to 0).



#### Notes on &time\_control

- *interval\_seconds*:
  - Time interval between WPS output times, and lateral BC (and lower BC) update frequency
- *history\_interval*:
  - Time interval in <u>minutes</u> when a history output is written
  - The time stamp in a history file name is the time when the history file is first written, and multiple time periods may be written in one file. e.g. a history file for domain 1 that is first written for 1200 UTC Jan 24 2000 is



wrfout\_d01\_2000-01-24\_12:00:00

#### Notes on &time\_control

- *frame\_per\_outfile*:
  - Number of history times written to one file.
- restart\_interval:
  - Time interval in minutes when a restart file is written.
  - By default, restart file is not written at hour 0.
  - A restart file contains only one time level data, and its valid time is in its file name, e.g. a restart file for domain 1 that is valid for 0000 UTC Jan 25 2000 is
     wrfrst d01 2000-01-25 00:00:00
- restart:



– whether this is a restart run

#### Notes on *restart*

- What is a *restart* run?
  - A restart run is a continuation of a model run.
- How to do a *restart* run:
  - In the first run, set <u>restart\_interval</u> to a value that is within the model integration time.
  - A restart file will be created. e.g.

wrfrst\_d01\_2000-01-25\_00:00:00

- When doing a restart run:
  - Set restart = .true.,
  - Set start times to restart times in namelist



#### &time\_control





#### &domains




### Notes on &domains

- *time\_step, time\_step\_fract\_num, time\_step\_frac\_den*:
  - Time step for model integration in seconds.
  - Fractional time step specified in separate integers of numerator and denominator.
  - ARW: 6xDX; NMM: 2.25xDX (DX is grid distance in km)
- e\_we, e\_sn, e\_vert:
  - Model grid dimensions (staggered) in X, Y and Z directions.
- *num\_metgrid\_levels*:
  - Number of *metgrid* (input) data levels.
- *num\_metgrid\_soil\_levels*:
  - Number of soil data levels in the input data
  - Found by typing ncdump -h met\_\*.d01.<date> | more
- *dx, dy*:
- WRF
- grid distances: in meters for ARW; in degrees for NMM.

#### Notes on &domains

- p\_top\_requested:
  - Pressure value at the model top.
  - Constrained by the available data from WPS.
  - Default is 5000 Pa
- eta\_levels:
  - Specify your own model levels from 1.0 to 0.0.
  - If not specified, program *real* will calculate a set of levels
- *ptsgm* (NMM only):
  - Pressure level (Pa) at which the WRF-NMM hybrid coordinate transitions from sigma to pressure (default: 42000 Pa)



## Where do I start?

- Always start with a *namelist* template provided in a test case directory, whether it is an ideal case, ARW or NMM real data case.
  - A number of namelist templates are provided in test/test\_<case>/ directories

For example: in *test/em\_real/*, there are namelist.input.4km ~ 4 km grid size namelist.input.jun01 ~ 10 km grid size namelist.input.jan00 ~ 30 km grid size



## Where do I start?

- Use document to guide the modification of the namelist values:
  - run/README.namelist
  - test/em\_real/examples.namelist
  - User's Guide, Chapter 5 (online version has the latest)
  - Full list of namelists and their default values can be found in Registry files: Registry.EM (ARW), Registry.NMM and registry.io\_boilerplate (IO options, shared)



# To run a job in a different directory..

- Directories *run*/ and test\_<case>/ are convenient places to run, but it does not have to be.
- Copy or link the content of these directories to another directory, including physics data files, wrf input and boundary files, wrf namelist and executables, and you should be able to run a job anywhere on your system.



### **Check Output**



### **Output After a Model Run**

- Standard out/error files:
   wrf.out, or rsl.\* files
- Model history file(s):

wrfout\_d01\_<date>

Model restart file(s), optional

wrfrst\_d01\_<date>



#### Output from a multi-processor run

The standard out and error will go to the following files for a MPI run: mpirun -np 4 ./wrf.exe →

rsl.out.0000	rsl.error.0000
rsl.out.0001	rsl.error.0001
rsl.out.0002	rsl.error.0002
rsl.out.0003	rsl.error.0003

There is one pair of files for each processor requested



### What to Look for in a standard out File?

Check run log file by typing tail wrf.out, or tail rsl.out.0000

You should see the following if the job is successfully completed: wrf: SUCCESS COMPLETE WRF



### How to Check Model History File?

• Use ncdump:

ncdump -v Times wrfout\_d01\_<date>
to check output times. Or
 ncdump -v U wrfout\_d01\_<date>
to check a particular variable (U)

- Use **ncview** or **ncBrowse** (great tools!)
- Use post-processing tools (see talks later)



#### What is in a *wrf.out* or *rsl* file?

- A print of namelist options
- Time taken to compute one model step:

 Timing for main: time 2000-01-24\_12:03:00 on domain
 1:
 3.25000 elapsed seconds.

 Timing for main: time 2000-01-24\_12:06:00 on domain
 1:
 1.50000 elapsed seconds.

 Timing for main: time 2000-01-24\_12:09:00 on domain
 1:
 1.50000 elapsed seconds.

 Timing for main: time 2000-01-24\_12:09:00 on domain
 1:
 1.50000 elapsed seconds.

 Timing for main: time 2000-01-24\_12:12:00 on domain
 1:
 1.55000 elapsed seconds.

• Time taken to write history and restart file:

Timing for Writing wrfout\_d01\_2000-01-24\_18:00:00 for domain 1: 0.14000 elapsed seconds.

• Any model error prints: (example from ARW run)

5 points exceeded cfl=2 in domain 1 at time 4.200000 MAX AT i,j,k: 123 48 3
 cfl,w,d(eta) = 4.165821

An indication the model has become numerically unstable



## Simple Trouble Shooting



### Often-seen runtime problems

- module\_configure: initial\_config: error reading
  namelist: &dynamics
  - > Typos or erroneous namelist variables exist in namelist record &dynamics in namelist.input file
- input\_wrf.F: SIZE MISMATCH: namelist ide,jde,num\_metgrid\_levels= 70 61 27 ; input data ide,jde,num\_metgrid\_levels= 74 61 27
  - > Grid dimensions in error



#### Often-seen runtime problems

- Segmentation fault (core dumped)
  - Often typing `unlimit' or `ulimit -s unlimited' or equivalent can help when this happens quickly in a run.
- If you do: grep cfl rsl.error.\* and see 121 points exceeded cfl=2 in domain 1 at time 4.200000 MAX AT i,j,k: 123 48 3 cfl,w,d(eta)= 4.165821
  - Model becomes unstable due to various reasons. If it happens soon after the start time, check input data, and/or reduce time step.



#### References

- Information on compiling and running WRF, and a more extensive list of namelist options and their definition / explanations can be found in the ARW and NMM User's Guide, Chapter 5
- Also see 'Nesting Setup and Run' and 'Other Runtime Options' talks.

