WRF: Set up and Run

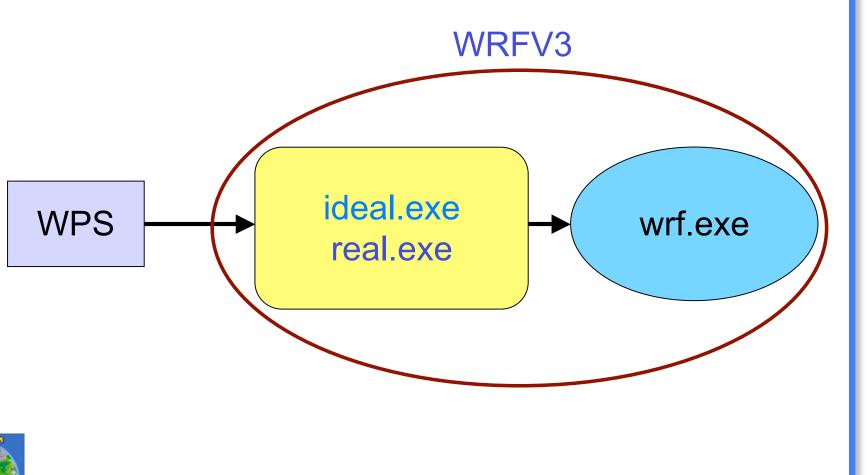
## Set Up and Run WRF

(ARW-Ideal and ARW-real)

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## WRF System Flowchart





#### **Outline**

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

This talk is complementary to 'Nesting' talk 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
- If you are running a real-data case, be sure that files from WPS are correctly generated:
  - met em.d01.\*, for ARW or
- Prepare namelist.input for runtime options.



### WRF test case directories

You have these choices in wrfv3/test/

```
(made at compile time):
```

```
├ 3d real-data
  em real
em quarter ss
em b wave
                      3d ideal
em les
em heldsuarez
em hill2d x
                                        ARW
em squall2d x
em squall2d y
                       2d ideal
em grav2d x
em seabreeze2d x
em scm xy
                       1d ideal
```

## Steps to Run

- 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*)
- 5. Run model executable, wrf.exe



## WRFV3/run directory

#### README namelist

LANDUSE . TBL ETAMPNEW DATA GENPARM. TBL RRTM DATA RRTMG SW DATA RRTMG LW DATA SOILPARM. TBL **VEGPARM. TBL** URBAN PARAM. TBL tr49t67 tr49t85 tr67t85 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
ETAMPNEW DATA
GENPARM TRI
RRTM DATA
RRTMG SW DATA
                            An example after
RRTMG LW DATA
SOILPARM. TBL
                            ARW real case
VEGPARM. TBL
                            compile
URBAN PARAM. TBL
tr49t67
tr49t85
tr67t85
gribmap.txt
grib2map.tbl
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 ARW Real-Data Case



## Running ARW Real-Data Case

 If you have compiled the em\_real 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
ETAMPNEW DATA -> ../../run/ETAMPNEW DATA
GENPARM.TBL -> ../../run/GENPARM.TBL
RRTM DATA -> ../../run/RRTM DATA
RRTMG SW DATA -> ../../run/RRTMG SW DATA
RRTMG LW DATA -> ../../run/RRTMG LW DATA
SOILPARM.TBL -> ../../run/SOILPARM.TBL
VEGPARM.TBL -> ../../run/VEGPARM.TBL
URBAN PARAM.TBL -> ../../run/URBAN PARAM.TBL
tr49t67 -> ../../run/tr49t67
tr49t85 -> ../../run/tr49t85
tr67t85 -> ../../run/tr67t85
gribmap.txt -> ../../run/gribmap.txt
grib2map.tbl -> ../../run/grib2map.tbl
namelist.input - require editing
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.* .
```

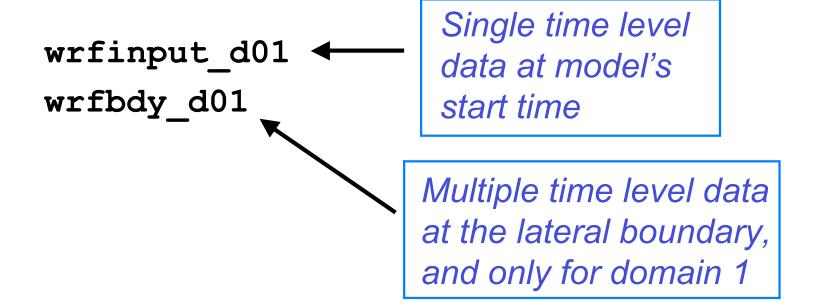


- Edit namelist.input file for runtime options (at mininum, one must edit &time\_control for start, end and integration times, and &domains for grid dimensions)
- Run the real-data initialization program:

```
./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:





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:

```
wrfout_d01_2005-08-28_00:00:00
```

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



# 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
```

→ 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 <u>input\_sounding</u>), or a pre-defined 3D input (e.g. <u>input\_jet</u> 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



- 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

# **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
run minutes
                        0,
run seconds
                        2000,
start year
start month
                        01,
                                     24,
                        24,
start day
                                     12,
                        12,
start hour
start minute
                        00,
                               00,
                                     00,
                        00,
                              00,
                                     00,
start second
                        2000
                              2000,
                                     2000
end year
                        01,
end month
                               01,
                                     01,
end day
                        25,
                               25,
                                     25,
                        12,
                               12,
                                     12,
end hour
                        00,
                               00,
                                     00,
end minute
                               00,
                        00,
                                     00,
end second
interval seconds
                        21600
history interval
                        180
                                             nest options
                      = 1000,
frame per outfile
                      = 360
restart interval
```



## 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.



## Notes on &time\_control

- Interval seconds:
  - Time interval between WPS output times, and LBC update frequency
- history\_interval:
  - Time interval in minutes 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



## &time control

```
io_form_history = 2,
io_form_restart = 2,
io_form_input = 2,
io_form_boundary = 2,
debug_level = 0,
```

IO format options:

= 1, binary

= 2, netcdf (most common)

= 4, PHDF5

= 5, Grib 1

=10, Grib 2

io\_form\_restart = 102 :
write output in patch
sizes: fast for large grids
and useful for restart file

Debug print control: Increasing values give more prints.



#### &domains

```
time step
                         = 180
time_step_fract_num
                         = 0,
time_step_fract_den
                         = 1,
max dom
                         = 1,
                         = 74,
e we
                         = 61,
e sn
                                    28, nest<sub>28,</sub>
                         = 28,
e vert
                                     options
num metgrid levels
                         = 21
num_metgrid_soil_levels
                         = 30000, 1000, 33
dx
                         = 30000, 1000, 3333
dy
                         = 1.0, 0.996, 0.\overline{99}, 0.98, \dots 0.0
eta levels
                         = 5000,
p_top_requested
```



#### 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; (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
- *dx, dy*.
  - grid distances in meters for ARW



#### Notes on &domains

- p\_top\_requested:
  - Pressure value at the model top.
  - Constrained by the available data from WPS.
  - Default is 5000 hPa
- eta\_levels:
  - Specify your own model levels from 1.0 to 0.0.
  - If not specified, program *real* will calculate a set of levels for you



#### Where do I start?

- Always start with a namelist template provided in a test case directory, whether it is a ideal case, or 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
  - 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), 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 and 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.error.0002 rsl.error.0002 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_12:12:00 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.
- 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 User's Guide, Chapter 5
- Also see 'Nesting Setup and Run' talk.

