

How to Use the WRF Registry

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Outline

- Examples
 - 1) Add output without recompiling
 - 2) Add a variable to the namelist
 - 3) Add an array
 - 4) Compute a diagnostic
 - 5) Add a physics package
 - 6) Tracer

Example 1: **ADD** output without recompiling

- Edit the namelist.input file, the time_control namelist record

```
iofields_filename = "myoutfields.txt" (MAXDOM)  
io_form_auxhist24 = 2 (choose an available stream)  
auxhist24_interval = 10 (MAXDOM, every 10 minutes)
```

- Place the fields that **you want** in the named text file `myoutfields.txt`

```
+ : h : 24 : RAINC , RAINNC
```

- Where “+” means ADD this variable to the output stream, “h” is the history stream, and “24” is the stream number

Example 1: **ZAP** output without recompiling

- Edit the namelist.input file, the time_control namelist record

```
iofields_filename = "myoutfields.txt" (MAXDOM)
```

- Place the fields **to remove** in the named text file `myoutfields.txt`

```
- : h : 0 : W , PB , P
```

- Where “-” means REMOVE this variable from the output stream, “**h**” is the history stream, and “**0**” is the stream number (standard WRF history file)

Example 1: What streams can I use?

- Generally history streams 10 – 24 are OK
- Avoid 21, 22, 23

HINT: Think of a stream as a separate file.
A history stream is a new output file

- Need LOTS more streams?
 - Edit WRF/arch/preamble

MAX_HISTORY = 25 <- - - *right now*

- ./clean -a, ./configure, ./compile, then re-run real and wrf

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Example 2: Add a variable to the namelist

- Use the examples for the **rconfig** section of the Registry
- Find a namelist variable similar to what you want
 - Integer *vs* real *vs* logical *vs* character
 - Single value *vs* value per domain
 - Select appropriate namelist record
- Insert your mods in all appropriate Registry files

Example 2: Add a variable to the namelist

- Remember that ALL Registry changes require that the WRF code be cleaned and rebuilt

```
./clean -a
```

```
./configure
```

```
./compile em_real
```

Example 2: Add a variable to the namelist

- Adding a variable to the namelist requires the inclusion of a new line in the Registry file:

```
rconfig integer my_option_1 namelist,time_control 1 0 - "my_option_1" "test namelist option"  
rconfig integer my_option_2 namelist,time_control max_domains 0
```

- Accessing the variable is through an automatically generated function:

```
USE module_configure  
INTEGER :: my_option_1 , my_option_2  
  
CALL nl_get_my_option_1( 1, my_option_1 )  
CALL nl_set_my_option_2( grid%id, my_option_2 )
```

Example 2: Add a variable to the namelist

- You also have access to the namelist variables from the grid structure ...

```
SUBROUTINE foo ( grid , ... )
```

```
USE module_domain
```

```
TYPE(domain) :: grid
```

```
print *,grid%my_option_1
```

Example 2: Add a variable to the namelist

- ... and you also have access to the namelist variables from config_flags

```
SUBROUTINE foo2 ( config_flags , ... )  
  
  USE module_configure  
  TYPE(grid_config_rec_type) :: config_flags  
  
  print *,config_flags%my_option_2
```

Example 2: Add a variable to the namelist

- What your variable looks like in the namelist.input file

```
&time_control  
run_days           = 0,  
run_hours          = 0,  
run_minutes        = 40,  
run_seconds        = 0,  
start_year         = 2006, 2006, 2006,  
my_option_1        = 17  
my_option_2        = 1, 2, 3
```

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Example 3: Add an Array

- Adding a state array to the solver, requires adding a single line in the Registry
- Use the Registry keyword information for a **state** or **I1** variable

Example 3: Add an Array

- Select a variable **similar** to one that you would like to add
 - 1d, 2d, or 3d
 - Staggered (X, Y, Z, or not “-”, *do not leave blank*)
 - Associated with a package
 - Part of a 4d array
 - Input (012), output, restart
 - Nesting, lateral forcing, feedback

Example 3: Add an Array

- Copy the “similar” field’s line and make a few edits
- Remember, no Registry change takes effect until a “clean -a” and rebuild

```
state  real  h_diabatic  ikj  misc  1  -      r      \
      "h_diabatic"    "PREVIOUS TIMESTEP CONDENSATIONAL HEATING"

state  real  msft      ij  misc  1  -      i012rhdu=(copy_fcnm)  \
      "MAPFAC_M"      "Map scale factor on mass grid"

state  real  ht      ij  misc  1  -      i012rhdu      \
      "HGT"           "Terrain Height"

state  real  ht_input  ij  misc  1  -      -      \
      "HGT_INPUT"     "Terrain Height from FG Input File"

state  real  TSK_SAVE  ij  misc  1  -      -      \
      "TSK_SAVE"      "SURFACE SKIN TEMPERATURE"      "K"
```

Example 3: Add an Array

- Usually modify Registry.*core_name*_COMMON or Registry.*core_name*, where *core_name* might be EM

```
state  real  h_diabatic  ikj  misc  1  -      r      \
      "h_diabatic"      "PREVIOUS TIMESTEP CONDENSATIONAL HEATING"

state  real  msft      ij  misc  1  -      i012rhdu=(copy_fcnm) \
      "MAPFAC_M"      "Map scale factor on mass grid"

state  real  ht      ij  misc  1  -      i012rhdu      \
      "HGT"      "Terrain Height"

state  real  ht_input  ij  misc  1  -      -      \
      "HGT_INPUT"      "Terrain Height from FG Input File"

state  real  TSK_SAVE  ij  misc  1  -      -      \
      "TSK_SAVE"      "SURFACE SKIN TEMPERATURE"      "K"
```

Example 3: Add an Array

- Add a new 3D array that is sum of all moisture species, called **all_moist**, in the Registry.EM_COMMON
 - Type: real
 - Dimensions: 3D and ikj ordering, not staggered
 - Supposed to be output only (history): h
 - Name in netCDF file: ALL_MOIST

```
state      real    all_moist      ikj      \  
misc              1      -      h      \  
"ALL_MOIST"      \  
"sum of all of moisture species"      \  
"kg kg-1"
```

Example 3: Add an Array

- Registry **state** variables become part of the derived data structure usually called **grid** inside of the WRF model.
- WRF → WRF model top → integrate → solve_interface → solve
- Each step, the **grid** construct is carried along for the ride
- No source changes for new output variables required until below the solver routine when dereferenced by first_rk_step_part1 for the physics drivers

Example 3: Add an Array

- Top of solve_em.F
- **grid** is passed in
- No need to declare any new variables, such as all_moist

```
!WRF:MEDIATION_LAYER:SOLVER
```

```
SUBROUTINE solve_em ( grid , &  
  
config_flags , &
```

Example 3: Add an Array

- In **solve_em**, add the new array to the call for the microphysics driver
- Syntax for **variable=local_variable** is an association convenience
- All state arrays are contained within grid, and must be **de-referenced**

```
CALL microphysics_driver(                                &
    QV_CURR=moist(ims,kms,jms,P_QV), &
    QC_CURR=moist(ims,kms,jms,P_QC), &
    QR_CURR=moist(ims,kms,jms,P_QR), &
    QI_CURR=moist(ims,kms,jms,P_QI), &
    QS_CURR=moist(ims,kms,jms,P_QS), &
    QG_CURR=moist(ims,kms,jms,P_QG), &
    QH_CURR=moist(ims,kms,jms,P_QH), &
    all_moist=grid%all_moist                        , &
```

Example 3: Add an Array

- After the array is re-referenced from grid and we are **inside the microphysics_driver** routine, we need to
 - Pass the variable through the argument list
 - Declare our passed in 3D array

```
,all_moist &
```

```
REAL,  DIMENSION(ims:ime ,kms:kme ,jms:jme ), &  
        INTENT(OUT)  ::  all_moist
```

Example 3: Add an Array

- After the array is re-referenced from grid and we are **inside the microphysics_driver** routine, we need to
 - Zero out the array at each time step

! Zero out moisture sum.

```
DO j = jts,MIN(jde-1,jte)  
DO k = kts,kte  
DO i = its,MIN(ide-1,ite)  
    all_moist(i,k,j) = 0.0  
END DO  
END DO  
END DO
```

Example 3: Add an Array

- After the array is re-referenced from grid and we are **inside the microphysics_driver** routine, we need to
 - At the end of the routine, for each of the **moist species that exists**, add that component to **all_moist**

```
DO j = jts,MIN(jde-1,jte)
  DO k = kts,kte
    IF ( f_QV ) THEN
      DO i = its,MIN(ide-1,ite)
        all_moist(i,k,j) = all_moist(i,k,j) + &
                               qv_curr(i,k,j)
      END DO
    END IF
  END IF
```

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Example 4: Compute a Diagnostic

- Problem: Output global average and global maximum and lat/lon location of maximum for 10 meter wind speed in WRF
- Steps:
 - Modify solve to compute wind-speed and then compute the local sum and maxima at the end of each time step
 - Use reduction operations built-in to WRF software to compute the global qualities
 - Output these on one process (process zero, the “monitor” process)

Example 4: Compute a Diagnostic

Compute local sum and local max and the local indices of the local maximum

```
--- File: dyn_em/solve_em.F (near the end) ---  
  
! Compute local maximum and sum of 10m wind-speed  
  sum_ws = 0.  
  max_ws = 0.  
  DO j = jps, jpe  
    DO i = ips, ipe  
      wind_vel = sqrt( grid%u10(i,j)**2+ grid%v10(i,j)**2 )  
      IF ( wind_vel .GT. max_ws ) THEN  
        max_ws = wind_vel  
        idex = i  
        jdex = j  
      ENDIF  
      sum_ws = sum_ws + wind_vel  
    ENDDO  
  ENDDO
```

Example 4: Compute a Diagnostic

- Compute global sum, global max, and indices of the global max (WRF intrinsics)

```
! Compute global sum
  sum_ws = wrf_dm_sum_real ( sum_ws )

! Compute global maximum and associated i,j point
  CALL wrf_dm_maxval_real ( max_ws, idex, jdex )
```

Example 4: Compute a Diagnostic

- On the process that contains the maximum value, obtain the latitude and longitude of that point; on other processes set to an artificially low value.
- The use parallel reduction to store that result on every process

```
IF ( ips .LE. idex .AND. idex .LE. ipe .AND. &  
    jps .LE. jdex .AND. jdex .LE. jpe ) THEN  
    glat = grid%xlat(idex,jdex)  
    glon = grid%xlong(idex,jdex)  
ELSE  
    glat = -99999.  
    glon = -99999.  
ENDIF
```

```
! Compute global maximum to find glat and glon  
glat = wrf_dm_max_real ( glat )  
glon = wrf_dm_max_real ( glon )
```

Example 4: Compute a Diagnostic

- Output the value on process zero, the “monitor”

```
! Print out the result on the monitor process
IF ( wrf_dm_on_monitor() ) THEN
  WRITE(outstring,*) 'Avg. ',sum_ws/((ide-ids+1)*(jde-jds+1))
  CALL wrf_message ( TRIM(outstring) )
  WRITE(outstring,*) 'Max. ',max_ws,' Lat. ',glat,&
                    ' Lon. ',glon
  CALL wrf_message ( TRIM(outstring) )
ENDIF
```

Example 4: Compute a Diagnostic

- Output from process zero of a multi-process run

```
--- Output file: rsl.out.0000 ---  
. . .  
Avg.      5.159380  
Max.      15.09370      Lat.      37.25022      Lon.      -67.44571  
Timing for main: time 2000-01-24_12:03:00 on domain 1:      8.96500 elapsed secs.  
Avg.      5.166167  
Max.      14.97418      Lat.      37.25022      Lon.      -67.44571  
Timing for main: time 2000-01-24_12:06:00 on domain 1:      4.89460 elapsed secs.  
Avg.      5.205693  
Max.      14.92687      Lat.      37.25022      Lon.      -67.44571  
Timing for main: time 2000-01-24_12:09:00 on domain 1:      4.83500 elapsed secs.  
. . .
```

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Example 5: Input periodic SSTs

- Add a new physics package with time varying input source to the model
- This is how we could supply a time varying value to the model for a field that is traditionally fixed
- Example is sea surface temperature

Example 5: Input periodic SSTs

- Problem: adapt WRF to input a time-varying lower boundary condition, e.g. SSTs, from an input file for a new surface scheme
- Given: Input file in WRF I/O format containing 12-hourly SST's
- Modify WRF model to read these into a new state array and make available to WRF surface physics

Example 5: Input periodic SSTs

- Steps
 - Add a new state variable and definition of a new surface layer package (that will use the variable) to the Registry
 - Add to variable stream for an unused Auxiliary Input stream
 - Adapt physics interface to pass new state variable to physics
 - Setup namelist to input the file at desired interval

Example 5: Input periodic SSTs

- Add a new state variable to Registry/Registry.EM_COMMON and put it in the variable set for input on Auxiliary Input Stream #4

#	type	symbol	dims	use	tl	stag	io	dname	description	units
state	real	nsst	ij	misc	1	-	i4h	"NEW_SST"	"Time Varying SST"	"K"

- Also added to History and Restart
- Result:
 - 2-D variable named grid%**nsst** defined and available in solve_em
 - Dimensions: ims:ime, jms:jme
 - Input and output on the AuxInput #4 stream will include the variable under the name NEW_SST

Example 5: Input periodic SSTs

- Pass new state variable to surface physics

```
--- File: dyn_em/module_first_rk_step_part1.F ---
```

```
CALL surface_driver(                                     &
    . . .
! Optional
&      ,QV_CURR=moist(ims,kms,jms,P_QV) , F_QV=F_QV      &
&      ,QC_CURR=moist(ims,kms,jms,P_QC) , F_QC=F_QC      &
&      ,QR_CURR=moist(ims,kms,jms,P_QR) , F_QR=F_QR      &
&      ,QI_CURR=moist(ims,kms,jms,P_QI) , F_QI=F_QI      &
&      ,QS_CURR=moist(ims,kms,jms,P_QS) , F_QS=F_QS      &
&      ,QG_CURR=moist(ims,kms,jms,P_QG) , F_QG=F_QG      &
&      ,NSST=grid%nsst                                     & ! new
&      ,CAPG=grid%capg, EMISS=grid%emiss, HOL=hol,MOL=grid%mol &
&      ,RAINBL=grid%rainbl,SR=grid%em_sr                  &
&      ,RAINNCV=grid%rainncv,REGIME=regime,T2=grid%t2,THC=grid%thc &
    . . .
```

Example 5: Input periodic SSTs

- Add new variable nsst to Physics Driver in Mediation Layer

```
--- File: phys/module_surface_driver.F ---

SUBROUTINE surface_driver(                                     &
    . . .
    ! Other optionals (more or less em specific)
    &      nsst                                               &
    &      ,capg,emiss,hol,mol                                &
    &      ,rainncv,rainbl,regime,t2,thc                      &
    &      ,qsg,qvg,qcg,soilt1,tsnav                          &
    &      ,smfr3d,keepfr3dflag                               &
    . . .
                                                    ))

    . . .
REAL, DIMENSION( ims:ime, jms:jme ), OPTIONAL, INTENT(INOUT) :: nsst
```

- By making this an “Optional” argument, we preserve the driver’s compatibility with other cores and with versions of WRF where this variable hasn’t been added.

Example 5: Input periodic SSTs

- Add call to Model-Layer subroutine for new physics package to Surface Driver

```
--- File: phys/module_surface_driver ---

!$OMP PARALLEL DO  &
!$OMP PRIVATE ( ij, i, j, k )
  DO ij = 1 , num_tiles
    sfclay_select: SELECT CASE(sf_sfclay_physics)

      CASE (SFCLAYScheme)
        . . .

      CASE (NEWSFCSCHEME)  ! <- This is defined by the Registry "package" entry

        IF (PRESENT(nsst)) THEN
          CALL NEWSFCSCHEME(
                                &
                                nsst,                                &
                                ids,ide, jds,jde, kds,kde,          &
                                ims,ime, jms,jme, kms,kme,          &
                                i_start(ij),i_end(ij), j_start(ij),j_end(ij), kts,kte  )
        ELSE
          CALL wrf_error_fatal('Missing argument for NEWScheme in surface driver')
        ENDIF
        . . .
      END SELECT sfclay_select
    ENDDO
  !$OMP END PARALLEL DO
```

- Note the PRESENT test to make sure new optional variable nsst is available

Example 5: Input periodic SSTs

- Add definition for new physics package NEWSHEME as setting 4 for namelist variable `sf_sfclay_physics`

<code>rconfig</code>	<code>integer</code>	<code>sf_sfclay_physics</code>	<code>namelist,physics</code>	<code>max_domains</code>	<code>0</code>
<code>package</code>	<code>sfclayscheme</code>	<code>sf_sfclay_physics==1</code>	-	-	
<code>package</code>	<code>myjsfcscheme</code>	<code>sf_sfclay_physics==2</code>	-	-	
<code>package</code>	<code>gfssfcscheme</code>	<code>sf_sfclay_physics==3</code>	-	-	
<code>package</code>	<code>newsfcscheme</code>	<code>sf_sfclay_physics==4</code>	-	-	

- This creates a defined constant NEWSFCSCHEME and represents selection of the new scheme when the namelist variable `sf_sfclay_physics` is set to '4' in the namelist.input file
- **clean -a** and recompile so code and Registry changes take effect

Example 5: Input periodic SSTs

- Setup namelist to input SSTs from the file at desired interval

```
    --- File: namelist.input ---  
  
&time_control  
  . . .  
  auxinput4_inname      = "sst_input"  
  auxinput4_interval_h  = 12  
  . . .  
/  
  
  . . .  
&physics  
  sf_sfclay_physics     = 4, 4, 4  
  . . .  
/
```

- Run code with sst_input file in run-directory

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Tracer Example

1. Modify Registry for new fields.

Use the “tracer” array with a new 3D component

Use existing NML option

2. Initialize data in real.

Identify (i,j) location

Spread in “PBL”

3. Set values in solver.

“Release” per time step



Tracer Example

Registry/Registry.EM add our new field “PLUME” as part of “TRACER” array.

```
#      New tracer for example
state  real  plume  ikjftb  tracer \
      1  -  irhusdf=(bdy_interp:dt) \
      "PLUME"  "Fukushima Tracer"  " "

#      4D arrays need an associated package
package  tracer_test3  tracer_opt==3  - \
      tracer:plume
```

Tracer Example

Modify the real and WRF programs to initialize and continuously re-supply the
“PLUME” array

dyn_em/module_initialize_real.F (initial value from real.exe)

dyn_em/solve_em.F (continuous plume in wrf.exe)

```
! Add in the Fukushima initial venting.
```

```
IF ( ( its .LE. 50 ) .AND. ( ite .GE. 50 ) .AND. &  
    ( jts .LE. 50 ) .AND. ( jte .GE. 50 ) ) THEN  
    tracer(50,1:5,50,P_plume) = 1.
```

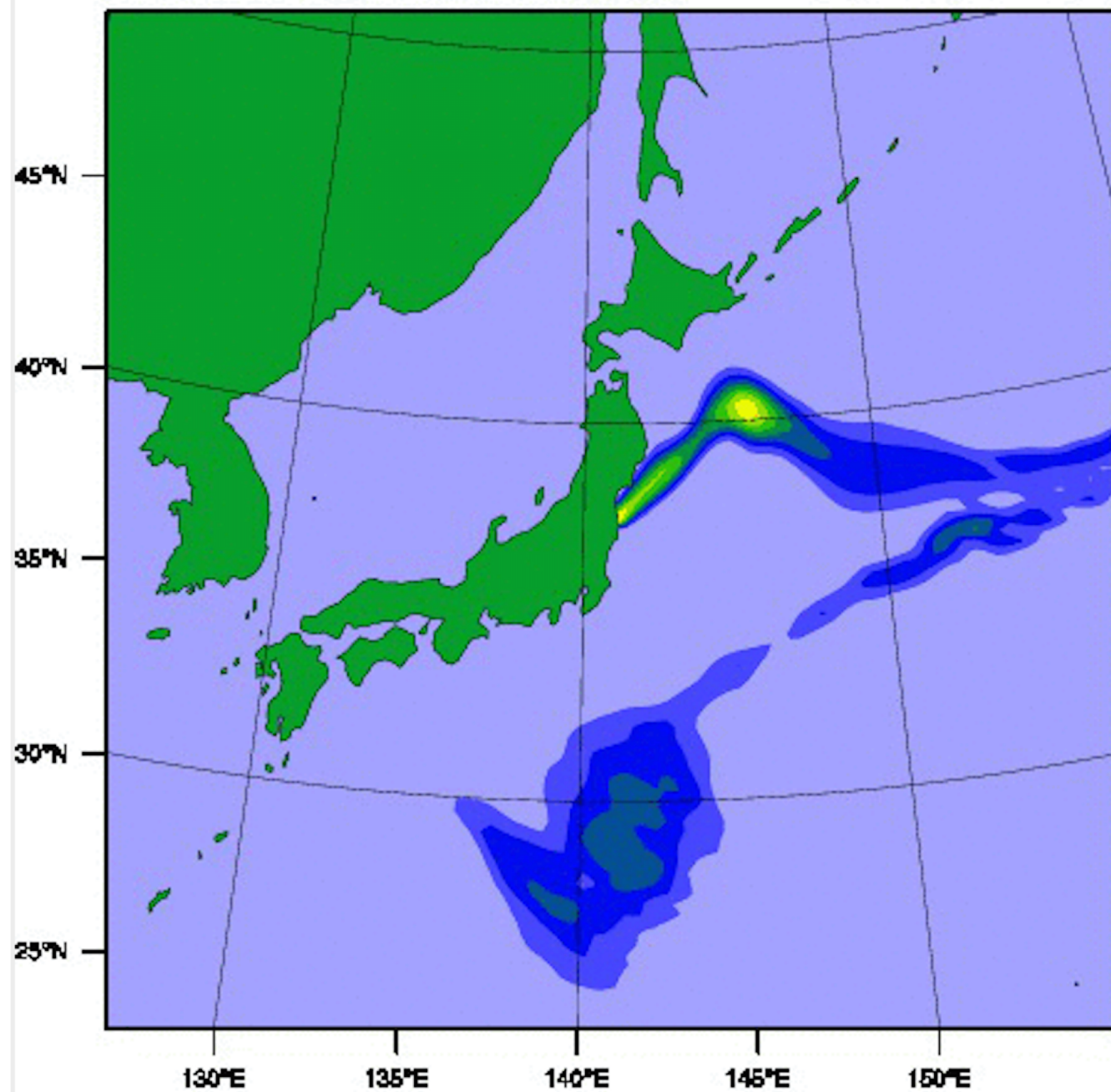
```
END IF
```

Tracer Example

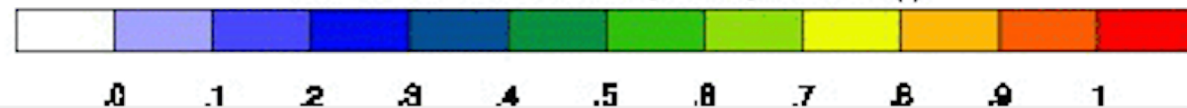
- Modify the test/em_real/namelist.input file
- Include the new settings for the tracer option required from the Registry file

```
&dynamics  
  tracer_opt = 3, 3, 3,
```

Fukushima 11-14 Mar 2011, 90-km, 100x100 (-)



Fukushima 11-14 Mar 2011, 30-km, 100x100 (-)



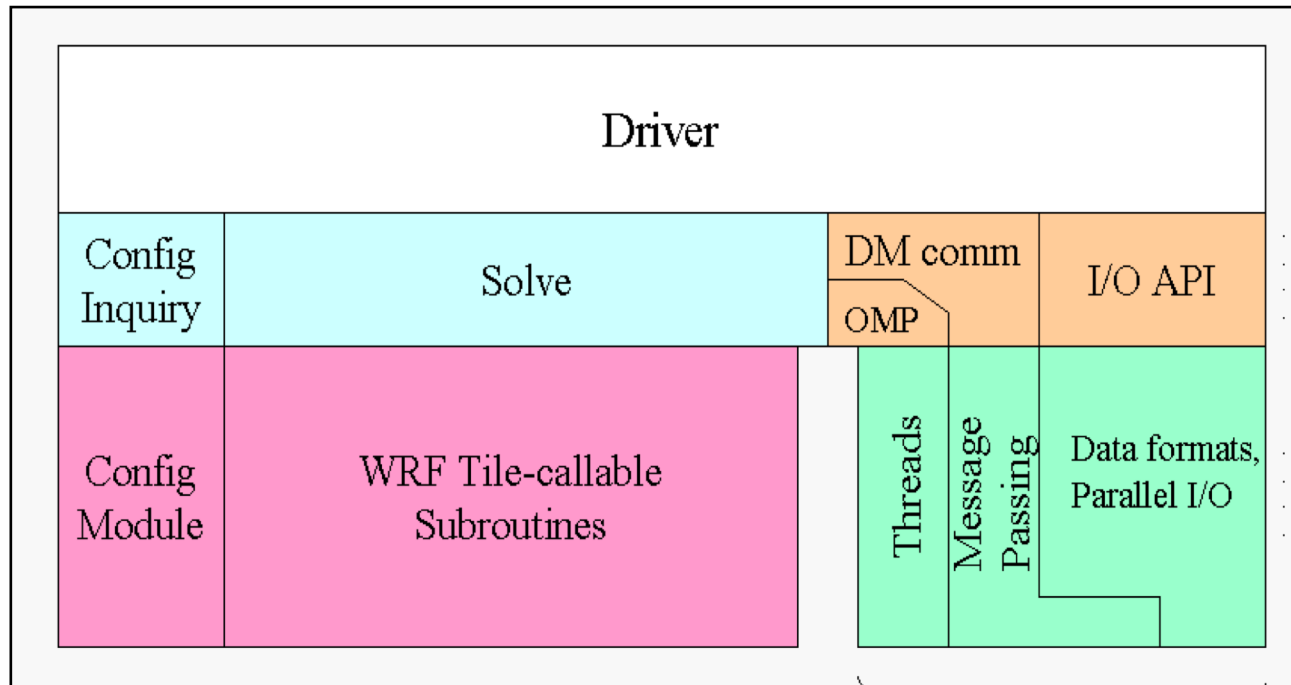
Review

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Outline for EXTRA Slides

- What is the WRF Registry
- Keyword syntax
- The BIG Three
 - state
 - rconfig
 - package

WRF Software Architecture



Text based file for real and WRF
 Active data dictionary
 Used with cpp to auto generate source
 Controls/defines
 Variables (I/O, comms, nesting)
 Communications
 namelist options

About 400k lines added to source
 Easy – 3x the size since initial release
 Compile-time option
 ./clean
 ./configure
 ./compile
 Registry.EM_COMMON (else lose changes)

Registry Keywords

- Currently implemented as a text file: **Registry/Registry.EM_COMMON**
- Types of entry:
 - *Dimspec*— Describes dimensions that are used to define arrays in the model
 - *State*— Describes state variables and arrays in the domain structure
 - *//*— Describes local variables and arrays in solve
 - *Typedef*— Describes derived types that are subtypes of the domain structure

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Registry Keywords

- Types of entry:
 - *Rconfig*— Describes a configuration (e.g. namelist) variable or array
 - *Package*— Describes attributes of a package (e.g. physics)
 - *Halo*— Describes halo update interprocessor communications
 - *Period*— Describes communications for periodic boundary updates
 - *Xpose*— Describes communications for parallel matrix transposes
 - *include*— Similar to a CPP #include file

Registry Keywords

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Registry State Entry

#	Type	Sym	Dims	Use	Tlev	Stag	IO	Dname	Descrip
state	real	tsk	ij	misc	1	-	i01rhusdf	"TSK"	"SKIN TEMP"

- Elements
 - *Entry*. The keyword “state”
 - *Type*. The type of the state variable or array (real, double, integer, logical, character, or derived)
 - *Sym*. The symbolic name of the variable or array
 - *Dims*. A string denoting the dimensionality of the array or a hyphen (-)
 - *Use*. A string denoting association with a solver or 4D scalar array, or a hyphen
 - *NumTlev*. An integer indicating the number of time levels (for arrays) or hyphen (for variables)

Registry State Entry

#	Type	Sym	Dims	Use	Tlev	Stag	IO	Dname	Descrip
state	real	tsk	ij	misc	1	-	i01r	hud	"TSK" "SKIN TEMP"

- Elements
 - *Stagger*: String indicating staggered dimensions of variable (X, Y, Z, or hyphen)
 - *IO*: String indicating whether and how the variable is subject to various I/O and Nesting
 - *DName*: Metadata name for the variable
 - *Units*: Metadata units of the variable
 - *Descrip*: Metadata description of the variable

State Entry: Defining a variable-set for an I/O stream

- Fields are added to a variable-set on an I/O stream in the Registry

#	Type	Sym	Dims	Use	Tlev	Stag	IO	Dname	Descrip
state	real	tsk	ij	misc	1	-	i01rhud	"TSK"	"SKIN TEMP"

- IO** is a string that specifies if the variable is to be available to initial, restart, or history I/O. The string may consist of '**h**' (subject to history output), '**i**' (initial dataset I/O), '**r**' (restart dataset I/O).
- The '**h**', '**r**', and '**i**' specifiers may appear in any order or combination.

State Entry: Defining a variable-set for an I/O stream

- Fields are added to a variable-set on an I/O stream in the Registry

#	Type	Sym	Dims	Use	Tlev	Stag	IO	Dname	Descrip
state	real	tsk	ij	misc	1	-	i01rhud	"TSK"	"SKIN TEMP"

- The '**h**' and '**i**' specifiers may be followed by an optional integer string consisting of '0', '1', ..., '9'
- Zero denotes that the variable is part of the principal input or history I/O stream.
- The characters '1' through '9' denote one of the auxiliary input or history I/O streams.
- Double digit streams require "{}" braces: **i01{19}{24}**

State Entry: Defining a variable-set for an I/O stream

The nesting info for each variable is co-located with the I/O

#	Type	Sym	Dims	Use	Tlev	Stag	IO	Dname	Descrip
state	real	tsk	ij	misc	1	-	i01rhud	"TSK"	"SKIN TEMP"

usdf refers to nesting options:

u = UP, d = DOWN, s = SMOOTH, f = FORCE

u – at end of each set of child time steps

d – at instantiation of child domain

f – at beginning of each set of child time steps

s – after each feedback

State Entry: Defining a variable-set for an I/O stream

Only variables involved with I/O,
communications, packages are required to
be state

Local variables inside of physics packages
are not controlled by the Registry

Rconfig Entry

#	Type	Sym	How set	Nentries	Default
<code>rconfig</code>	<code>integer</code>	<code>spec_bdy_width</code>	<code>namelist, bdy_control</code>	1	1

- This defines namelist entries
- Elements
 - *Entry*: the keyword “rconfig”
 - *Type*: the type of the namelist variable (integer, real, logical, string)
 - *Sym*: the name of the namelist variable or array
 - *How set*: indicates how the variable is set: e.g. namelist or derived, and if namelist, which block of the namelist it is set in

Rconfig Entry

#	Type	Sym	How set	Nentries	Default
rconfig	integer	spec_bdy_width	namelist, bdy_control	1	1

- This defines namelist entries
- Elements
 - *Nentries*: specifies the dimensionality of the namelist variable or array. If 1 (one) it is a variable and applies to all domains; otherwise specify max_domains (which is an integer parameter defined in module_driver_constants.F).
 - *Default*: the default value of the variable to be used if none is specified in the namelist; hyphen (-) for no default

Package Entry

- Elements
 - *Entry*: the keyword “**package**”,
 - *Package name*: the name of the package: e.g. “*kesslerscheme*”
 - *Associated rconfig choice*: the name of a rconfig variable and the value of that variable that chooses this package

```
# specification of microphysics options
package    passiveqv      mp_physics==0      -      moist:qv
package    kesslerscheme  mp_physics==1      -      moist:qv,qc,qr
package    linscheme      mp_physics==2      -
moist:qv,qc,qr,qi,qs,qg
package    ncepcloud3     mp_physics==3      -      moist:qv,qc,qr
package    ncepcloud5     mp_physics==4      -      moist:qv,qc,qr,qi,qs

# namelist entry that controls microphysics option
rconfig    integer      mp_physics    namelist,physics    max_domains    0
```

Package Entry

- Elements
 - *Package state vars*: unused at present; specify hyphen (-)
 - *Associated variables*: the names of 4D scalar arrays (*moist*, *chem*, *scalar*) and the fields within those arrays this package uses, and the state variables (*state:u_gc*, ...)

```
# specification of microphysics options
package    passiveqv      mp_physics==0      -      moist:qv
package    kesslerscheme  mp_physics==1      -      moist:qv,qc,qr
package    linscheme      mp_physics==2      -
moist:qv,qc,qr,qi,qs,qg
package    ncepcloud3      mp_physics==3      -      moist:qv,qc,qr
package    ncepcloud5      mp_physics==4      -      moist:qv,qc,qr,qi,qs

# namelist entry that controls microphysics option
rconfig integer  mp_physics  namelist,physics  max_domains  0
```