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# Outline

- Introduction
- Computing Overview
- WRF Software Overview

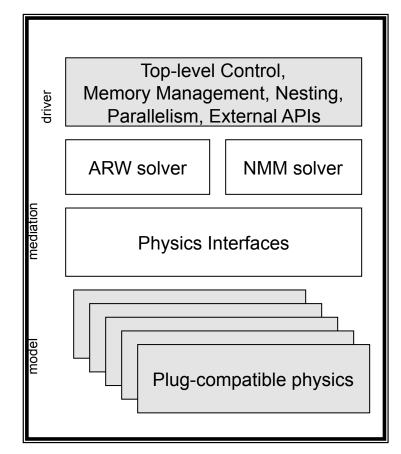
## Introduction – WRF Software Characteristics

- Developed from scratch beginning around 1998, primarily Fortran and C
- Requirements emphasize flexibility over a range of platforms, applications, users, performance
- WRF develops rapidly. First released Dec 2000; current release WRF v3.3.1 (Sep 2011); next release WRF v3.4 (April 2012)
- Supported by flexible efficient architecture and implementation called the WRF Software Framework

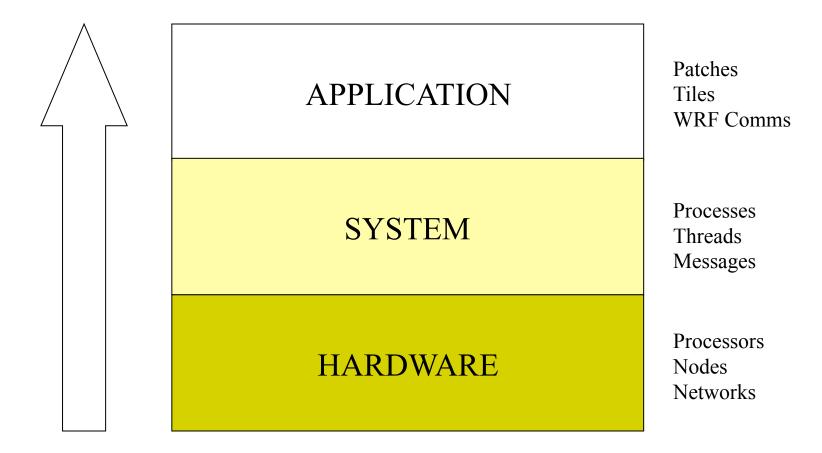
# Introduction - WRF Software Framework Overview

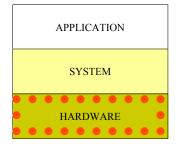
- Implementation of WRF Architecture
  - Hierarchical organization
  - Multiple dynamical cores
  - Plug compatible physics
  - Abstract interfaces (APIs) to external packages
  - Performance-portable
- Designed from beginning to be adaptable to today's computing environment for NWP

http://box.mmm.ucar.edu/wrf/WG2/bench/



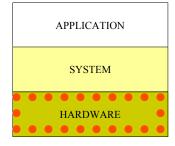
# **Computing Overview**





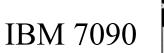
# Hardware: The Computer

- The 'N' in NWP
- Components
  - Processor
    - A program counter
    - Arithmetic unit(s)
    - Some scratch space (registers)
    - Circuitry to store/retrieve from memory device
    - Cache
  - Memory
  - Secondary storage
  - Peripherals
- The implementation has been continually refined, but the basic idea hasn't changed much



# Hardware has not changed much...

### A computer in 1960





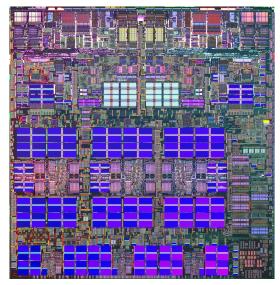
6-way superscalar

36-bit floating point precision

 $\sim \! 144 \text{ Kbytes}$ 

~50,000 flop/s 48hr 12km WRF CONUS in 600 years

#### A computer in 2008

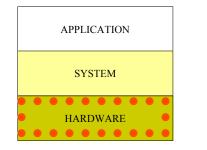


Dual core, 4.7 GHz chip 64-bit floating point precision 1.9 MB L2, 36 MB L3

Upto 16 GB per processor

~5,000,000,000 flop/s 48 12km WRF CONUS in 52 Hours

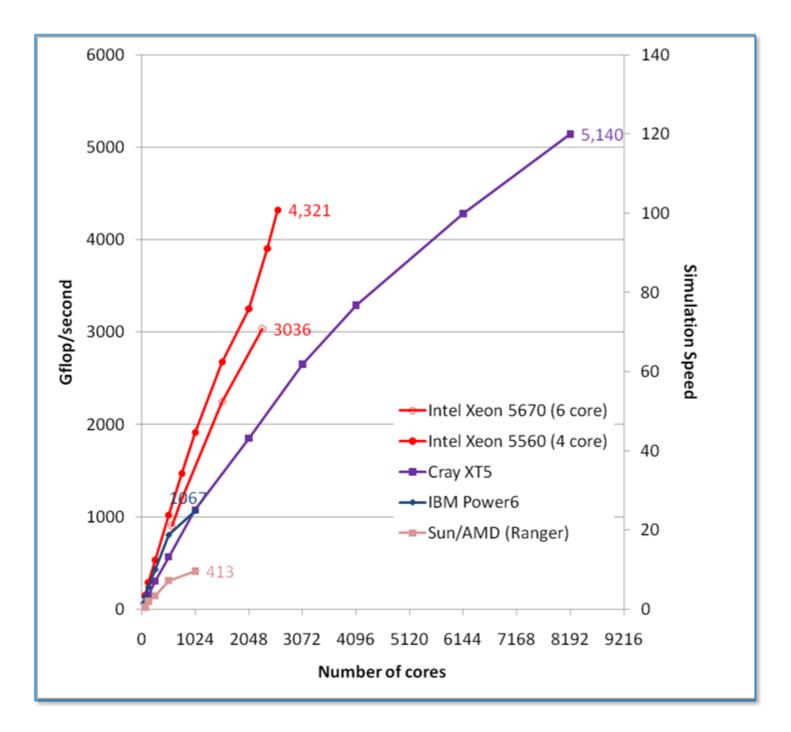
IBM P6



...how we use it has

- Fundamentally, processors haven't changed much since 1960
- Quantitatively, they haven't improved nearly enough
  - 100,000x increase in peak speed
  - 100,000x increase in memory size
- We make up the difference with <u>parallelism</u>
  - Ganging multiple processors together to achieve  $10^{11-12}$  flop/second
  - Aggregate available memories of  $10^{11-12}$  bytes

~1,000,000,000,000 flop/s ~250 procs 48-h,12-km WRF CONUS in under 15 minutes



## Examples

• If the machine consists of 4 nodes, each with 4 processors, how many different ways can you run a job to use all 16 processors?

	1 MPI	1 MPI
<ul> <li>4 MPI processes, each with 4 threads</li> <li>setenv OMP_NUM_THREADS 4</li> <li>mpirun -np 4 wrf.exe</li> </ul>	4 threads	4 threads
<ul> <li>8 MPI processes, each with 2 threads</li> </ul>	1 MPI	1 MPI
setenv OMP_NUM_THREADS 2		
mpirun -np 8 wrf.exe	4 threads	4 threads
<ul> <li>16 MPI processes, each with 1 thread</li> </ul>		

mpirun -np 16 wrf.exe

setenv OMP\_NUM\_THREADS 1

### Examples

• If the machine consists of 4 nodes, each with 4 processors, how many different ways can you run a job to use all 16 processors?

	2 MPI	2 MPI
<ul> <li>4 MPI processes, each with 4 threads</li> </ul>	2 threads	2 threads
setenv OMP_NUM_THREADS 4	2 threads	2 threads
mpirun -np 4 wrf.exe		
<ul> <li>8 MPI processes, each with 2 threads</li> </ul>	2 MPI	2 MPI
setenv OMP_NUM_THREADS 2 mpirun -np 8 wrf.exe	2 threads	2 threads
<ul> <li>16 MPI processes, each with 1 thread</li> </ul>	2 threads	2 threads

setenv OMP\_NUM\_THREADS 1
mpirun -np 16 wrf.exe

### Examples

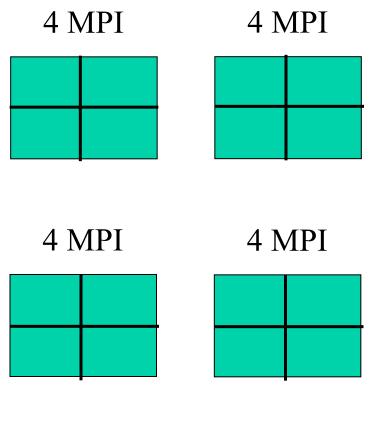
- If the machine consists of 4 nodes, each with 4 processors, how many different ways can you run a job to use all 16 processors?
  - 4 MPI processes, each with 4 threads

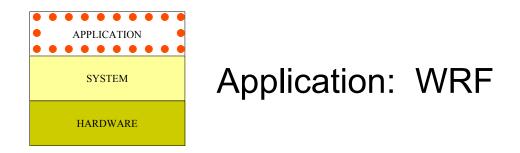
setenv OMP\_NUM\_THREADS 4
mpirun -np 4 wrf.exe

- 8 MPI processes, each with 2 threads

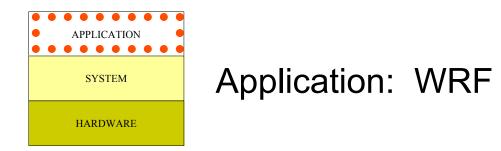
setenv OMP\_NUM\_THREADS 2
mpirun -np 8 wrf.exe

16 MPI processes, each with 1 thread
 setenv OMP\_NUM\_THREADS 1
 mpirun -np 16 wrf.exe

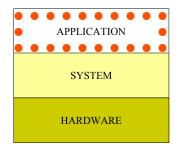




- WRF can be run serially or as a parallel job
- WRF uses *domain decomposition* to divide total amount of work over parallel processes

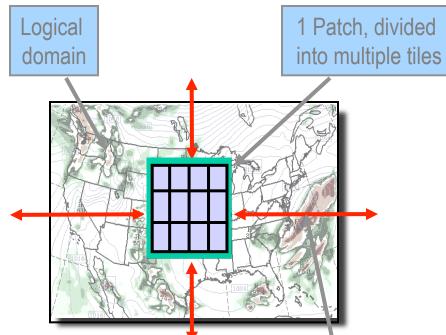


- The decomposition of the application over processes has two levels:
  - The *domain* is first broken up into rectangular pieces that are assigned to MPI (distributed memory) processes. These pieces are called *patches*
  - The *patches* may be further subdivided into smaller rectangular pieces that are called *tiles*, and these are assigned to *shared-memory threads* within the process.



### Parallelism in WRF: Multi-level Decomposition

- Single version of code for efficient execution on:
  - Distributed-memory
  - Shared-memory (SMP)
  - Clusters of SMPs
  - Vector and microprocessors



#### Model domains are decomposed for parallelism on two-levels

*Patch:* section of model domain allocated to a distributed memory node, this is the scope of a mediation layer solver or physics driver.

Inter-processor communication

*Tile:* section of a patch allocated to a shared-memory processor within a node; this is also the scope of a model layer subroutine.

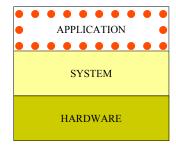
Distributed memory parallelism is over patches; shared memory parallelism is over tiles within patches

When Needed?	Communication is required between patches when a horizontal index is incremented or decremented on the right-hand-side of an assignment.
Why?	On a patch boundary, the index may refer to a value that is on a different patch.
	Following is an example code fragment that requires communication between patches
Signs in code	Note the tell-tale +1 and -1 expressions in indices for <b>rr</b> , <b>H1</b> , and <b>H2</b> arrays on right-hand side of assignment.
	These are <i>horizontal data dependencies</i> because the indexed operands may lie in the patch of a neighboring processor. That neighbor's updates to that element of the array won't be seen on this processor.

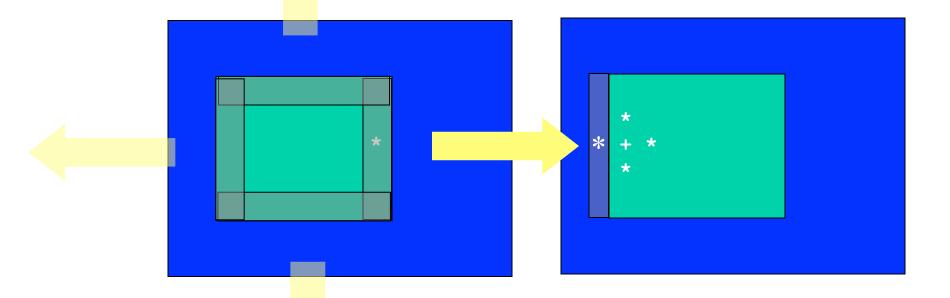
(module diffusion.F )

SUBROUTINE horizontal diffusion s (tendency, rr, var, . . .

```
DO j = jts,jte
DO k = kts, ktf
DO i = its, ite
   mrdx=msft(i,j)*rdx
   mrdy=msft(i,j)*rdy
   tendency(i,k,j) = tendency(i,k,j) -
                                                                  &
         (mrdx*0.5*((rr(i+1,k,j)+rr(i,k,j))*H1(i+1,k,j)-
                                                                  £
                    (rr(i-1,k,j)+rr(i,k,j)) * H1(i,k,j)) +
                                                                  £
         mrdy*0.5*((rr(i,k,j+1)+rr(i,k,j))*H2(i,k,j+1)-
                                                                  &
                    (rr(i,k,j-1)+rr(i,k,j)) * H2(i,k,j)) -
                                                                  £
         msft(i,j) * (Hlavg(i,k+1,j) - Hlavg(i,k,j) +
                                                                  &
                     H2avg(i,k+1,j)-H2avg(i,k,j)
                                                                  3
                               )/dzetaw(k)
                                                                  &
         )
ENDDO
ENDDO
ENDDO
```

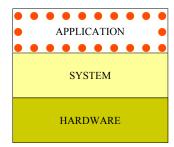


Halo updates

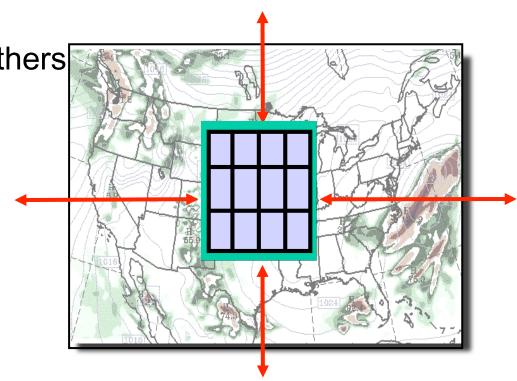


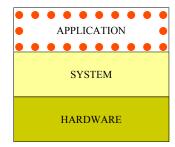
memory on one processor

memory on neighboring processor

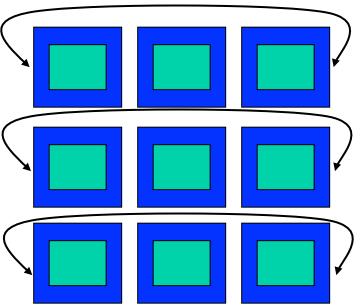


- Halo updates
- Periodic boundary updates
- Parallel transposes
- Nesting scatters/gathers

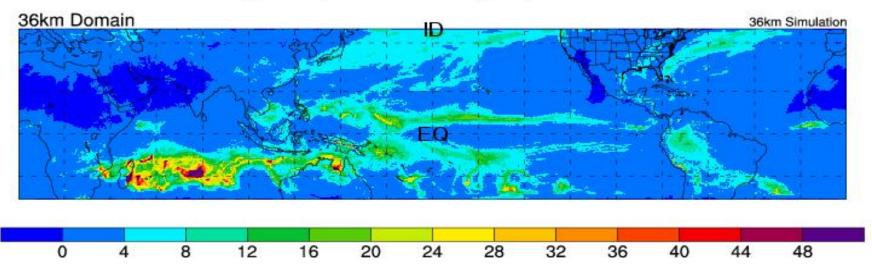


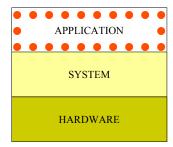


- Halo updates
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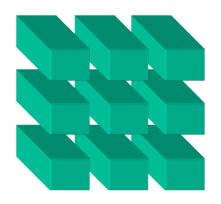


Average Daily Total rainfall (mm) - March 1997

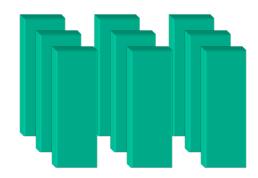




- Halo updates
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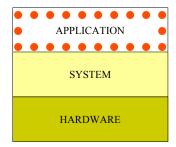
all y on patch



all z on patch

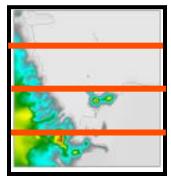


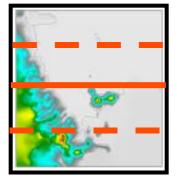
all x on patch

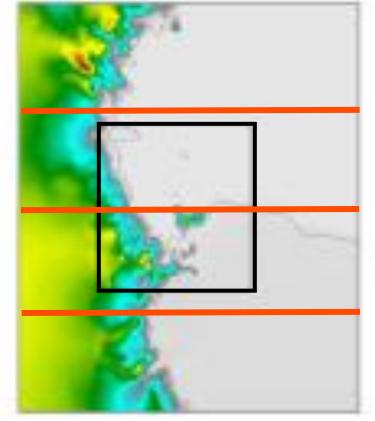


- Halo updates
- Periodic boundary updates
- Parallel transposes

Nesting scatters/gathers





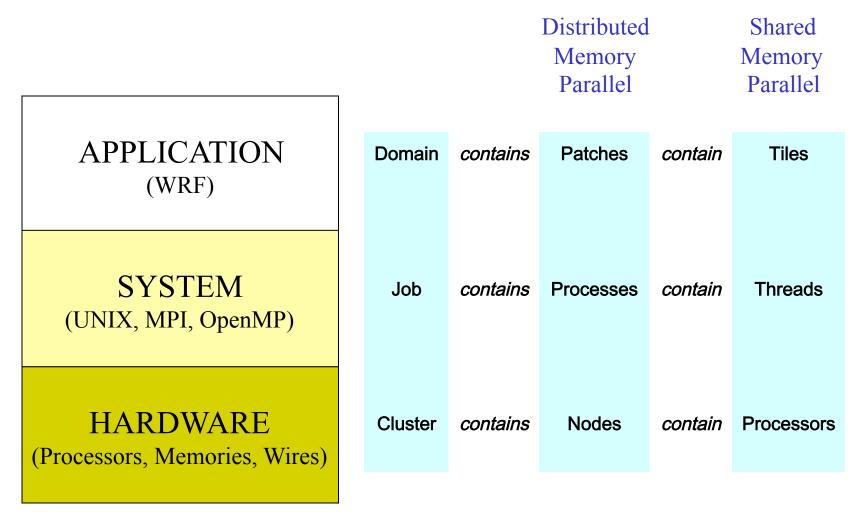


COARSE Ross Island 6.66 km

NEST:2.22 km

INTERMEDIATE: 6.66 km

### **Review – Computing Overview**

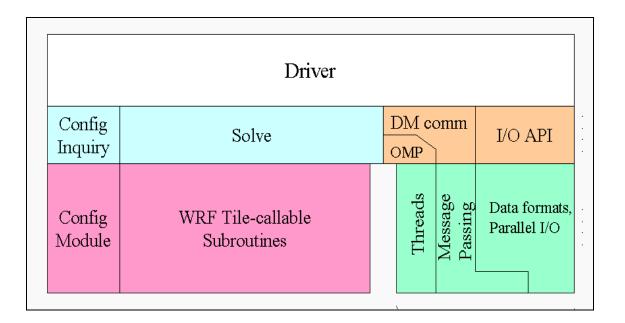


## Outline

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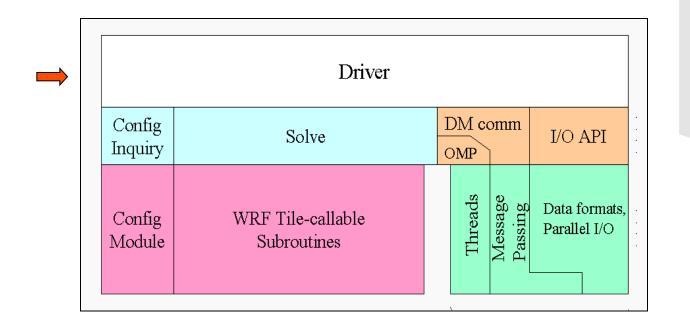
# WRF Software Overview

- Architecture
- Directory structure
- Model Layer Interface
- Data Structures
- I/O



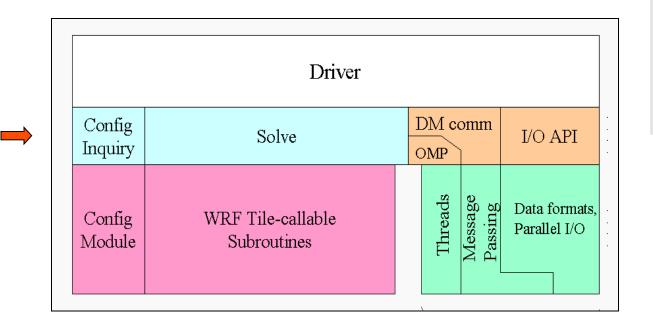
### Registry

- Hierarchical software architecture
  - Insulate scientists' code from parallelism and other architecture/ implementation-specific details
  - Well-defined interfaces between layers, and external packages for communications, I/O, and model coupling facilitates code reuse and exploiting of community infrastructure, e.g. ESMF.



Registry

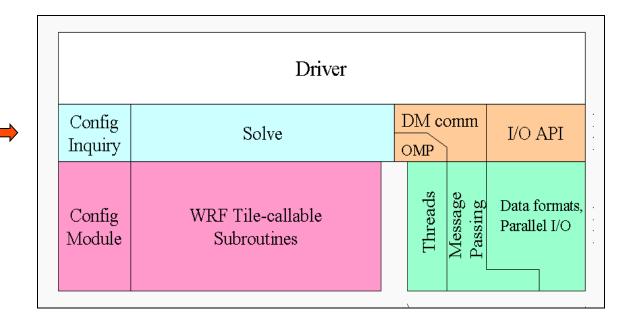
- Driver Layer
  - Domains: Allocates, stores, decomposes, represents abstractly as single data objects
  - Time loop: top level, algorithms for integration over nest hierarchy



- Mediation Layer
  - Solve routine, takes a domain object and advances it one time step

Registry

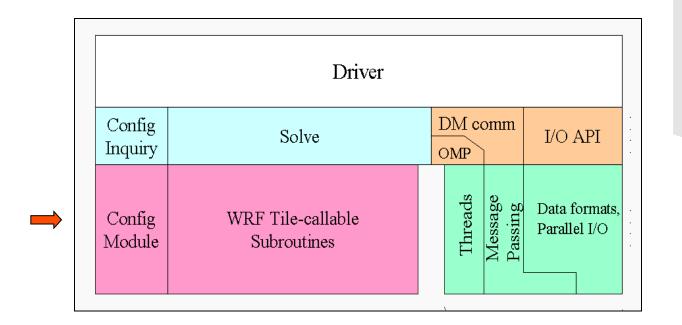
- Nest forcing, interpolation, and feedback routines



- Mediation Layer
  - The sequence of calls for doing a time-step for one domain is known in Solve routine

Registry

- Dereferences fields in calls to physics drivers and dynamics code
- Calls to message-passing are contained here as part of Solve routine

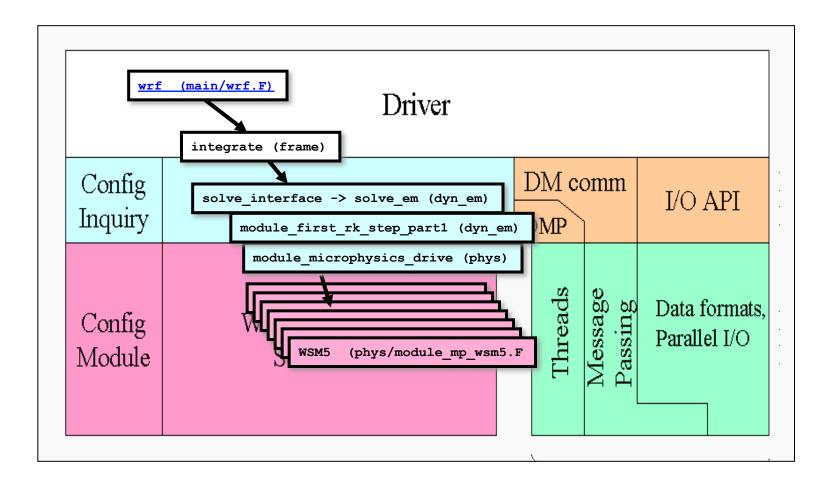


- Model Layer
  - Physics and Dynamics: contains the actual WRF model routines are written to perform some computation over an arbitrarily sized/ shaped, 3d, rectangular subdomain

Registry

### Call Structure Superimposed on Architecture

module\_microphysics\_driver (phys)



# WRF Software Overview

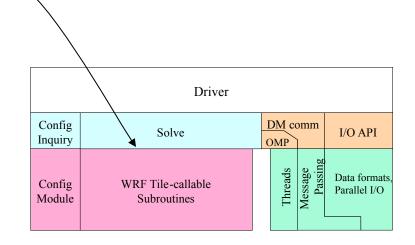
- Architecture
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#### **WRF Model Layer Interface – The Contract with Users**

All state arrays passed through argument list as simple (not derived) data types

Domain, memory, and run dimensions passed unambiguously in three dimensions

Model layer routines are called from mediation layer (physics drivers) in loops over tiles, which are multi-threaded



**WRF Model Layer Interface – The Contract with Users** 

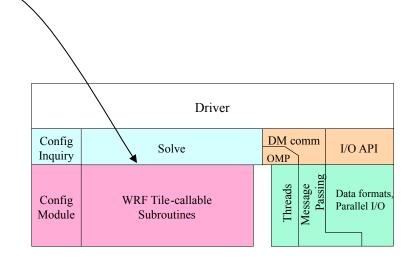
**Restrictions** on Model Layer subroutines:

No I/O, communication

No stops or aborts Use wrf\_error\_fatal

No common/module storage of decomposed data

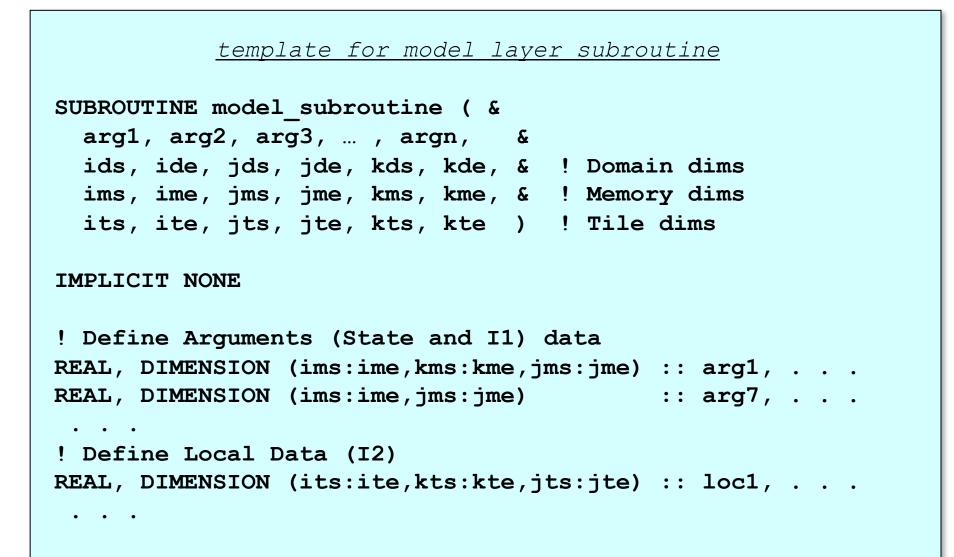
Spatial scope of a Model Layer call is one "tile"



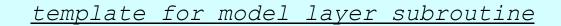
#### **WRF Model Layer Interface**

```
SUBROUTINE driver for some physics suite (
!$OMP DO PARALLEL
  DO ij = 1, numtiles
     its = i start(ij) ; ite = i end(ij)
     jts = j start(ij) ; jte = j end(ij)
     CALL model subroutine (arg1, arg2, . . .
           ids , ide , jds , jde , kds , kde ,
           ims , ime , jms , jme , kms , kme ,
           its , ite , jts , jte , kts , kte )
  END DO
END SUBROUTINE
```

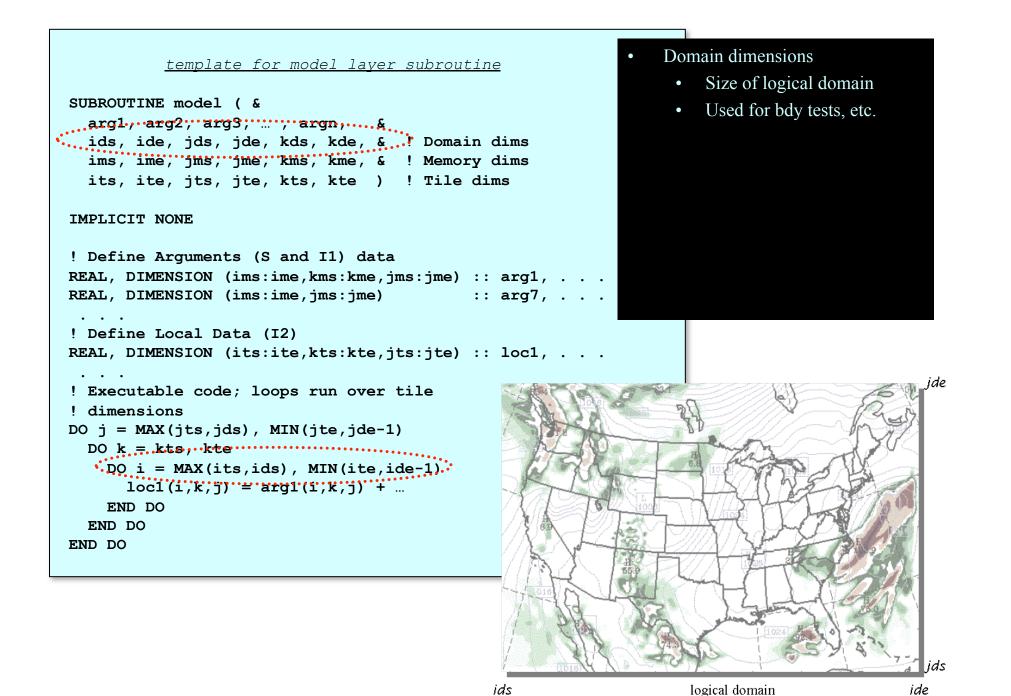
#### **WRF Model Layer Interface**

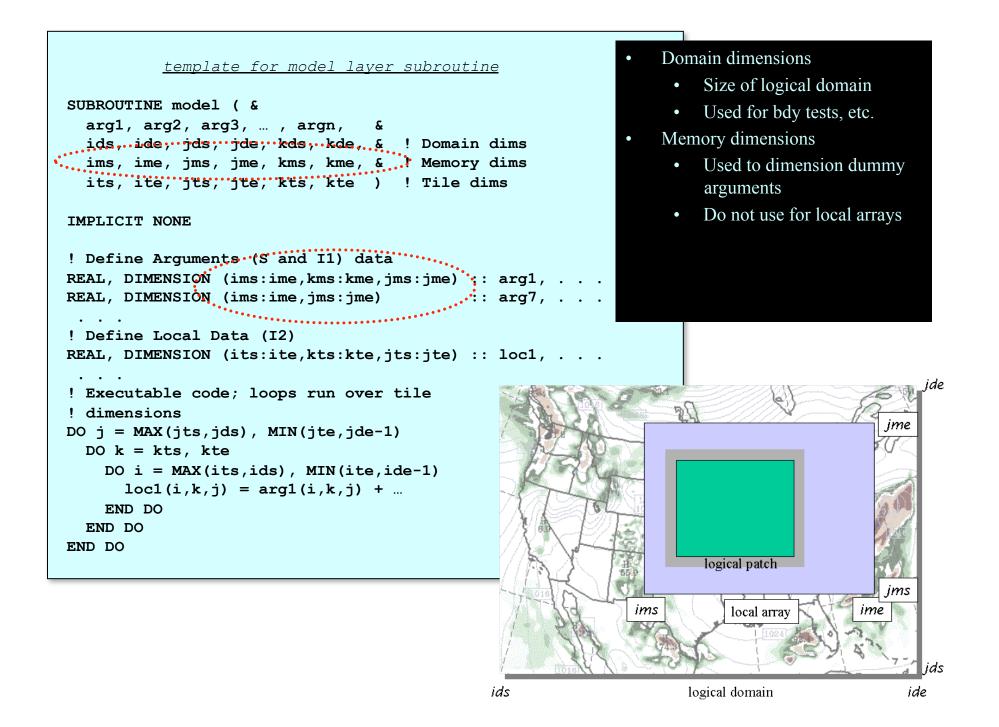


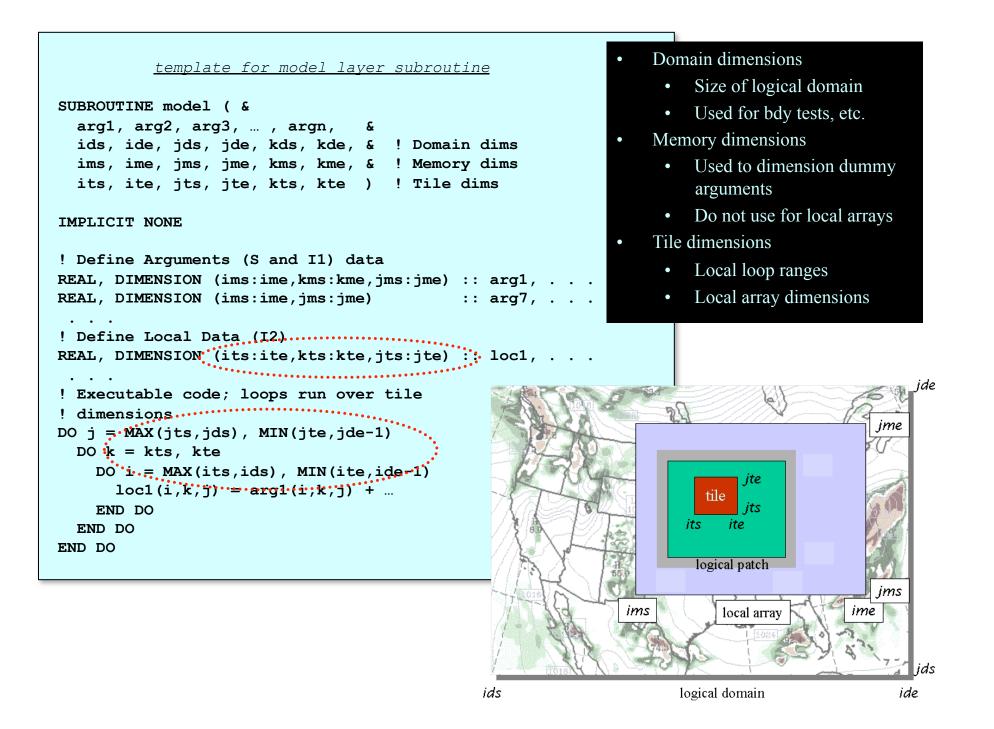
#### **WRF Model Layer Interface**

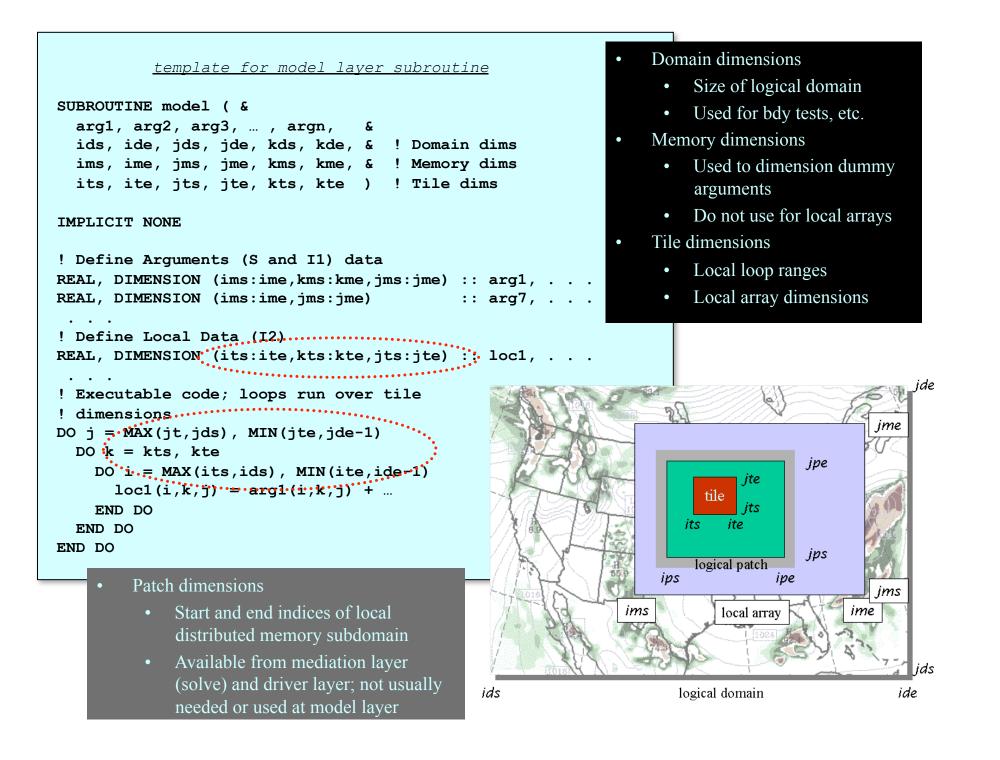


```
! Executable code; loops run over tile
! dimensions
DO j = jts, MIN(jte,jde-1)
DO k = kts, kte
DO i = its, MIN(ite,ide-1)
loc1(i,k,j) = arg1(i,k,j) + ...
END DO
END DO
END DO
```









# WRF Software Overview

- Architecture
- Directory structure
- Model Layer Interface
- Data Structures
- I/0

# WRF I/O

- Streams: pathways into and out of model
  - History + auxiliary output streams (10 and 11 are reserved for nudging)
  - Input + auxiliary input streams (10 and 11 are reserved for nudging)
  - Restart, boundary, and a special Var stream

# WRF I/O

- Attributes of streams
  - Variable set
    - The set of WRF state variables that comprise one read or write on a stream
    - Defined for a stream at compile time in Registry
  - Format
    - The format of the data outside the program (e.g. NetCDF), split
    - Specified for a stream at run time in the namelist

# WRF I/O

- Attributes of streams
  - Additional namelist-controlled attributes of streams
    - Dataset name
    - Time interval between I/O operations on stream
    - Starting, ending times for I/O (specified as intervals from start of run)

## **Outline - Review**

- Introduction
  - WRF started 1998, clean slate, Fortran + C
  - Targeted for research and operations
- WRF Software Overview
  - Hierarchical software layers
  - Patches (MPI) and Tiles (OpenMP)
  - Strict interfaces between layers
  - Contract with developers
  - I/O