

Topics in Computing for Astronomy:

Parallel Programming with OpenMP (I)

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Preliminaries

- ▶ The Mad SDK (contains gfortran, an OpenMP-capable Fortran 90/95/2003/2008 compiler)

<http://www.astro.wisc.edu/~townsend/static.php?ref=madsdk>

- ▶ This talk & accompanying code examples:

<http://www.astro.wisc.edu/~townsend/static.php?ref=talks>



What is OpenMP?

- ▶ An Application Programming Interface (API) for writing multithreaded programs
- ▶ Consists of a set of compiler directives, library routines and environment variables
- ▶ Supports Fortran, C and C++
- ▶ Developed by vendors (AMD, IBM, Intel, ...)
- ▶ Mainly targeted at Symmetric Multiprocessing systems



What is a Symmetric Multiprocessing (SMP)?

- ▶ A multiprocessor system with
 - ▶ centralized, shared memory
 - ▶ a single operating system
 - ▶ two or more homogeneous processors
- ▶ Modern desktops/laptops are SMP systems:
 - ▶ Multiple cores in each physical processor
 - ▶ Multiple physical processors



What is Multithreading?

- ▶ The ability of a program to have multiple, independent *threads* executing different sequences of instructions
- ▶ The threads can access the resources (e.g., memory) of the parent process
- ▶ The threads run concurrently on a multiprocessor systems



OpenMP Core Syntax

- ▶ Constructs use compiler directives specified using special Fortran comments:

Fortran 77

```
C$omp parallel do
```

Fortran 90+

```
!$omp parallel
```

- ▶ Most constructs apply to a “structured block” with one point of entry and one point of exit
- ▶ Additional library functions and subroutines accessed through the `omp_lib` module



Example: Hello World

```
program hello_world
```

```
hello_world.f90
```

```
use omp_lib  
implicit none
```

```
!$omp parallel num_threads(4)  
print *, 'Hello from thread', omp_get_thread_num()  
!$omp end parallel
```

```
end program hello_world
```

```
% gfortran -fopenmp -o hello_world hello_world.f90
```

```
% ./hello_world
```

```
Hello from thread      0  
Hello from thread      2  
Hello from thread      3  
Hello from thread      1
```



Example: Hello World

```
program hello.
```

Make
OpenMP routines
available

```
hello_world.f90
```

```
use omp_lib  
implicit none
```

```
!$omp parallel num_threads(4)  
print *, 'Hello from thread', omp_get_thread_num()  
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```
end program hello_world
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Start
a parallel region with
4 threads

```
!$omp parallel num_threads(4)  
print *, 'Hello from thread', omp_get_thread_num()  
!$omp end parallel
```

```
end program hello_world
```

```
% gfortran -fopenmp -o hello_world hello_world.f90
```

```
% ./hello_world
```

```
Hello from thread      0  
Hello from thread      2  
Hello from thread      3  
Hello from thread      1
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Example: Hello World

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program hello.
```

Make
OpenMP routines
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use omp_lib  
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Start
a parallel region with
4 threads

Get the ID of
the thread

```
!$omp parallel num_threads(4)  
print *, 'Hello from thread', omp_get_thread_num()  
!$omp end parallel
```

```
end program hello_world
```

```
% gfortran -fopenmp -o hello_world hello_world.f90
```

```
% ./hello_world
```

```
Hello from thread      0  
Hello from thread      2  
Hello from thread      3  
Hello from thread      1
```



Example: Hello World

```
program hello.
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Make
OpenMP routines
available

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hello_world.f90
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implicit none
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Start
a parallel region with
4 threads

Get the ID of
the thread

```
!$omp parallel num_threads(4)  
print *, 'Hello from thread', omp_get_thread_num()  
!$omp end parallel
```

```
end program hell
```

Tell gfortran to
recognize !\$omp directives

```
% gfortran -fopenmp -o hello_world hello_world.f90
```

```
% ./hello_world
```

```
Hello from thread      0  
Hello from thread      2  
Hello from thread      3  
Hello from thread      1
```



Example: Hello World

```
program hello.
```

Make
OpenMP routines
available

```
hello_world.f90
```

```
use omp_lib  
implicit none
```

Start
a parallel region with
4 threads

Get the ID of
the thread

```
!$omp parallel num_threads(4)  
print *, 'Hello from thread', omp_get_thread_num()  
!$omp end parallel
```

```
end program hell
```

Tell gfortran to
recognize !\$omp directives

```
% gfortran -fopenmp -o hello_world hello_world.f90  
% ./hello_world
```

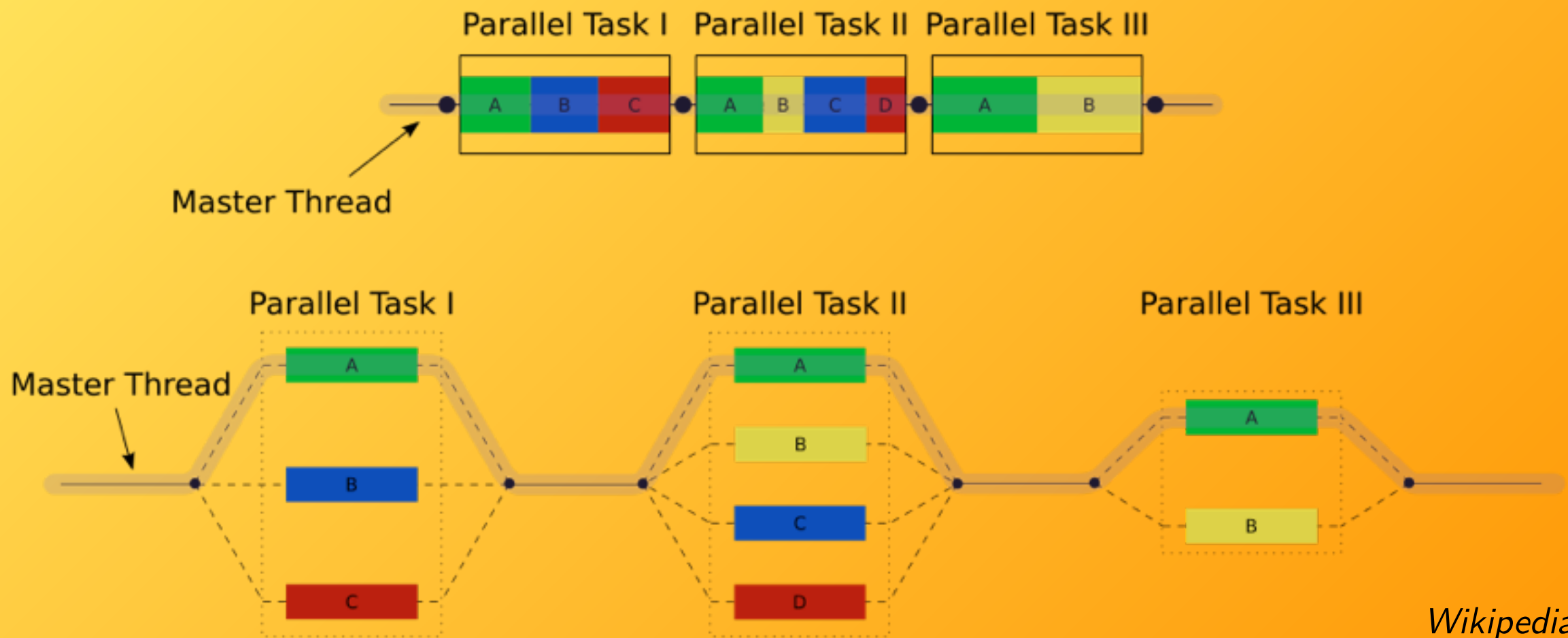
```
Hello from thread      0  
Hello from thread      2  
Hello from thread      3  
Hello from thread      1
```

Not sequential —
OpenMP does not guarantee
execution order!



The Fork-Join Thread Model

- ▶ At the start of a program, there is a single thread — the *master thread*
- ▶ When entering a parallel region, the master thread ‘forks’ into a team of threads
- ▶ When exiting a parallel region, the team joins back into a single master thread



Sharing Variables

```
program hello_world_shared
```

```
hello_world_shared.f90
```

```
use omp_lib  
implicit none
```

```
integer :: id
```

```
!$omp parallel num_threads(4)  
id = omp_get_thread_num() ←  
print *, 'Hello from thread', id  
!$omp end parallel
```

```
end program hello_world_shared
```

This overwrites
the same memory location!



Sharing Variables

```
program hello_world_shared
```

```
hello_world_shared.f90
```

```
use omp_lib  
implicit none
```

```
integer :: id
```

```
!$omp parallel num_threads(4)  
id = omp_get_thread_num()  
print *, 'Hello from thread', id  
!$omp end parallel
```

```
end program hello_world_shared
```

The id variable is
by default shared amongst
threads

This overwrites
the same memory location!



Not Sharing Variables

```
program hello_world_private
```

```
hello_world_private.f90
```

```
use omp_lib  
implicit none
```

```
integer :: id
```

```
!$omp parallel num_threads(4) private(id)  
id = omp_get_thread_num()  
print *, 'Hello from thread', id  
!$omp end parallel
```

```
end program hello_world_private
```



Not Sharing Variables

```
program hello_world_private
```

```
hello_world_private.f90
```

```
use omp_lib  
implicit none
```

```
integer :: id
```

```
!$omp parallel num_threads(4) private(id)  
id = omp_get_thread_num()  
print *, 'Hello from thread', id  
!$omp end parallel
```

```
end program hello_world_private
```

The id
variable is private; each
thread has its own copy



Read/Write Races

```
program count_threads
```

```
count_threads.f90
```

```
  use omp_lib  
  implicit none
```

```
  integer :: n
```

```
  n = 0
```

```
  !$omp parallel num_threads(4)
```

```
  n = n + 1
```

```
  !$omp end parallel
```

```
  print *, 'There were', n, 'threads'
```

```
end program count_threads
```



Read/Write Races

```
program count_threads
```

```
count_threads.f90
```

```
use omp_lib  
implicit none
```

```
integer :: n
```

```
n = 0
```

Bad: another
thread could update n
between read and write

```
!$omp parallel num_threads(4)
```

```
n = n + 1
```

```
!$omp end parallel
```

```
print *, 'There were', n, 'threads'
```

```
end program count_threads
```



Atomic Operations

```
program count_threads_atomic
```

```
count_threads_atomic.f90
```

```
use omp_lib  
implicit none
```

```
integer :: n
```

```
n = 0
```

```
!$omp parallel num_threads(4)
```

```
!$omp atomic
```

```
n = n + 1
```

```
!$omp end parallel
```

```
print *, 'There were', n, 'threads'
```

```
end program count_threads_atomic
```



Atomic Operations

```
program count_threads_atomic
```

```
count_threads_atomic.f90
```

```
use omp_lib  
implicit none
```

```
integer :: n
```

```
n = 0
```

atomic directive:
the read/write in the
following operation must be
executed as one

```
!$omp parallel threads(4)
```

```
!$omp atomic
```

```
n = n + 1
```

```
!$omp end parallel
```

```
print *, 'There were', n, 'threads'
```

```
end program count_threads_atomic
```



Synchronization Directives

- ▶ `!$omp atomic` is an example of a *synchronization directive*
- ▶ Other examples:
 - ▶ `!$omp critical` — only execute on one thread at a time
 - ▶ `!$omp master` — only execute on master thread
 - ▶ `!$omp barrier` — wait for all threads



Sharing the Load

```
program square_nums
```

```
square_nums.f90
```

```
use omp_lib  
implicit none
```

```
integer :: i, j(4)
```

```
!$omp parallel num_threads(4)  
i = omp_get_thread_num() + 1  
j(i) = i**2  
!$omp end parallel
```

```
print *, 'Square numbers:', j
```

```
end program square_nums
```



Sharing the Load

```
program square_nums
```

```
square_nums.f90
```

```
use omp_lib  
implicit none
```

```
integer :: i, j(4)
```

```
!$omp parallel num_threads(4)  
i = omp_get_thread_num() + 1  
j(i) = i**2  
!$omp end parallel
```

```
print *, 'Square numbers:', j
```

```
end program square_nums
```

Calculate array index
from thread id



Parallelizing DO loops

```
program square_nums_do
```

```
square_nums_do.f90
```

```
use omp_lib  
implicit none
```

```
integer :: i, j(13)
```

```
!$omp parallel num_threads(4)
```

```
!$omp do
```

```
do i = 1, SIZE(j)
```

```
    j(i) = i**2
```

```
enddo
```

```
!$omp end parallel
```

```
print *, 'Square numbers:', j
```

```
end program square_nums_do
```



Parallelizing DO loops

```
program square_nums_do
```

```
square_nums_do.f90
```

```
use omp_lib  
implicit none
```

Will work with
array of any size

```
integer :: i, j(13)
```

```
!$omp parallel num_threads(4)
```

```
!$omp do
```

```
do i = 1, SIZE(j)
```

```
    j(i) = i**2
```

```
enddo
```

```
!$omp end parallel
```

```
print *, 'Square numbers:', j
```

```
end program square_nums_do
```



Parallelizing DO loops

```
program square_nums_do
```

```
square_nums_do.f90
```

```
use omp_lib  
implicit none
```

```
integer :: i, j(13)
```

Will work with
array of any size

```
!$omp parallel num_threads(4)
```

```
!$omp do
```

do directive
distributes loop iterations
amongst threads

```
do i = 1, SIZE(j)
```

```
  j(i) = i**2
```

```
enddo
```

```
!$omp end parallel
```

```
print *, 'Square numbers:', j
```

```
end program square_nums_do
```



A Shorthand: parallel do

```
program square_nums_pardo
```

```
square_nums_pardo.f90
```

```
use omp_lib  
implicit none
```

```
integer :: i, j(13)
```

```
!$omp parallel do num_threads(4)
```

```
do i = 1, SIZE(j)
```

```
    j(i) = i**2
```

```
enddo
```

```
!$omp end parallel do
```

```
print *, 'Square numbers:', j
```

```
end program square_nums_pardo
```



A Shorthand: parallel do

```
program square_nums_pardo
```

```
square_nums_pardo.f90
```

```
use omp_lib  
implicit none
```

```
integer :: i, j(13)
```

```
!$omp parallel do num_threads(4)
```

```
do i = 1, SIZE(j)
```

```
    j(i) = i**2
```

```
enddo
```

```
!$omp end parallel do
```

```
print *, 'Square numbers:', j
```

```
end program square_nums_pardo
```

parallel do
directive creates parallel region
AND distributes loop iterations
amongst threads



Work-Sharing Constructs

- ▶ `!$omp do` is an example of a work-sharing construct
- ▶ A work-sharing construct divides the execution of the enclosed region among the members of the thread team that encounters it
- ▶ Other examples:
 - ▶ `!$omp sections` — divide multiple enclosed blocks amongst threads of team
 - ▶ `!$omp workshare` — divide single enclosed block amongst threads of team
 - ▶ `!$omp single` — execute all work on a single thread



The workshare Construct

```
program square_nums_workshr
```

```
square_nums_workshr.f90
```

```
use omp_lib  
implicit none
```

```
integer :: i, j(13)
```

```
j = [(i,i=1,SIZE(j))]
```

```
!$omp parallel workshare num_threads(4)
```

```
j = j**2
```

```
!$omp end parallel workshare
```

```
print *, 'Square numbers:', j
```

```
end program square_nums_workshr
```



The workshare Construct

```
program square_nums_workshr
```

```
square_nums_workshr.f90
```

```
use omp_lib  
implicit none
```

```
integer :: i, j(13)
```

Initialize j using an
array assignment

```
j = [(i,i=1,SIZE(j))]
```

```
!$omp parallel workshare num_threads(4)
```

```
j = j**2
```

```
!$omp end parallel workshare
```

```
print *, 'Square numbers:', j
```

```
end program square_nums_workshr
```



The workshare Construct

```
program square_nums_workshr
```

```
square_nums_workshr.f90
```

```
use omp_lib  
implicit none
```

```
integer :: i, j(13)
```

Initialize j using an
array assignment

```
j = [(i,i=1,SIZE(j))]
```

These array
arithmetic/assignment operations
are distributed amongst the
team

```
!$omp parallel workshare num_threads(4)
```

```
j = j**2
```

```
!$omp end parallel workshare
```

```
print *, 'Square numbers:', j
```

```
end program square_nums_workshr
```



Restrictions on workshare Constructs

- ▶ Only the following operations permitted:
 - ▶ array assignments
 - ▶ scalar assignments
 - ▶ FORALL statements
 - ▶ FORALL constructs
 - ▶ WHERE statements
 - ▶ WHERE constructs
 - ▶ atomic constructs
 - ▶ critical constructs
 - ▶ parallel constructs



The sections Construct

```
program hello_sections
```

```
hello_sections.f90
```

```
use omp_lib  
implicit none
```

```
!$omp parallel sections num_threads(4)  
!$omp section  
print *, 'Hello from thread', omp_get_thread_num()  
!$omp section  
print *, 'Hola from thread', omp_get_thread_num()  
!$omp section  
print *, 'Aloha from thread', omp_get_thread_num()  
!$omp end parallel sections
```

```
end program hello_sections
```



The sections Construct

```
program hello_sections
```

```
hello_sections.f90
```

```
use omp_lib
```

```
implicit none
```

Each block delimited by !\$omp
section runs on a separate thread

```
!$omp parallel section private(i) num_threads(4)
```

```
!$omp section
```

```
print *, 'Hello from thread', omp_get_thread_num()
```

```
!$omp section
```

```
print *, 'Hola from thread', omp_get_thread_num()
```

```
!$omp section
```

```
print *, 'Aloha from thread', omp_get_thread_num()
```

```
!$omp end parallel sections
```

```
end program hello_sections
```

