

# C-DAC Four Days Technology Workshop

*ON*

**Hybrid Computing – Coprocessors/Accelerators**  
**Power-Aware Computing – Performance of**  
**Applications Kernels**

**hyPACK-2013**  
**(Mode-1:Multi-Core)**

**Lecture Topic:**

**Multi-Core Processors : Shared Memory Prog:**  
**OpenMP Part-I**

*Venue : CMSD, UoHYD ; Date : October 15-18, 2013*

# Explicit Parallelism :Shared Memory Programming: The OpenMP Standard

## Lecture outline

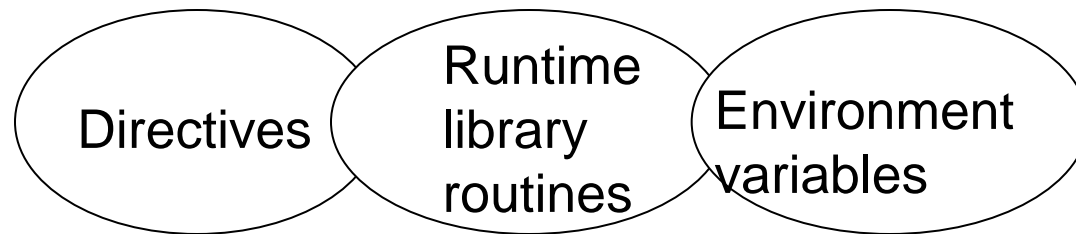
- ❖ Introduction to OpenMP
- ❖ OpenMP Programming Model
- ❖ OpenMP Constructs
  - Directives
  - Runtime Libraries
  - Environment variables

Source : Reference : [4], [6], [14],[17], [22], [28]

# OpenMP: Introduction

## OpenMP is an API for writing Multithreaded Applications

- ❖ It is a specification for



- ❖ Portable : Makes it easy to create multi-threaded programs in C,C++ and Fortran.
- ❖ Standardizes the SMP practice

Source : Reference : [4], [6], [14],[17], [22], [28]

# OpenMP: Introduction

## Why OpenMP ?

- ❖ Relatively easy to do parallelization for small parts of an application at a time.
- ❖ Impact on code quantity (e.g., amount of additional code required) and code quality (e.g., readability of parallel code)
- ❖ Feasibility of scaling an application to a large number of processes.
- ❖ Readability of the parallel code is high
- ❖ Availability of application development and debugging environment
- ❖ Standard and portable API

# Commonly Encountered Questions While Threading Application ?

- ❖ Where to thread ?
- ❖ How long would it take to thread?
- ❖ How much re-design / efforts is required?
- ❖ Is it worth threading the selected region ?
- ❖ What should the expected speedup be?
- ❖ Will the performance meet expectations?
- ❖ Will it scale if the more number of processor added?
- ❖ Which threading model is it?

# OpenMP: Introduction

## History

- ❖ First standard ANSI X3H5 in 1994
- ❖ OpenMP Standard SPECs started in 1997
- ❖ OpenMP Architecture review board
  - Compaq, HP, IBM, Sun Micro System, Intel Corp, Kuck & Associate Inc (KAI), SGI, US Dept. of Energy, ASCI program.

# OpenMP: Introduction

## Supporters

### ❖ Hardware vendors

➤ Intel, HP, SGI, IBM, SUN, Cray, AMD

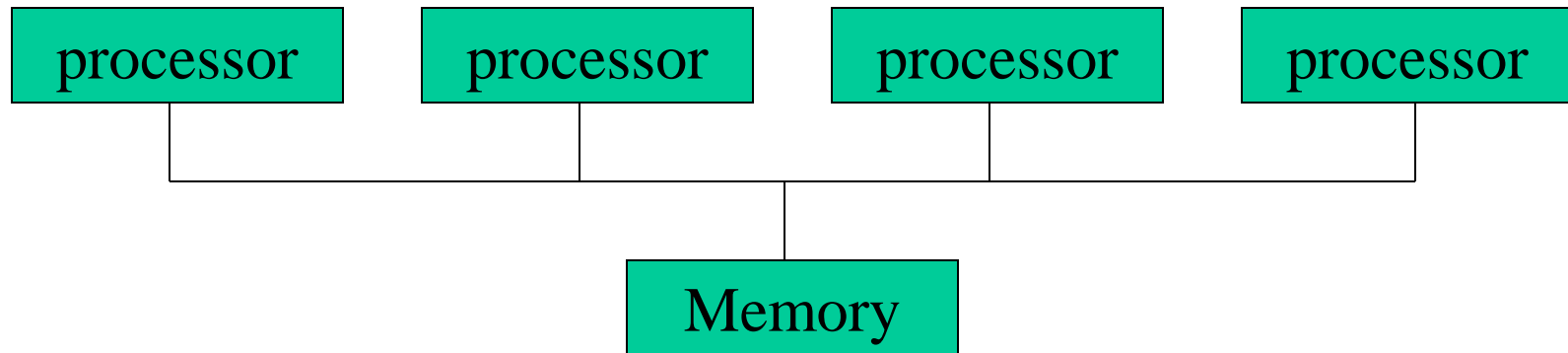
### ❖ Software tools vendors

➤ pathscale, Intel PGI, SGI, Sun, Absoft

# OpenMP: Programming Model

## Shared Memory Model

- ❖ Processes synchronize and communicate with each other through shared variables
- ❖ Supports *incremental parallelization*.

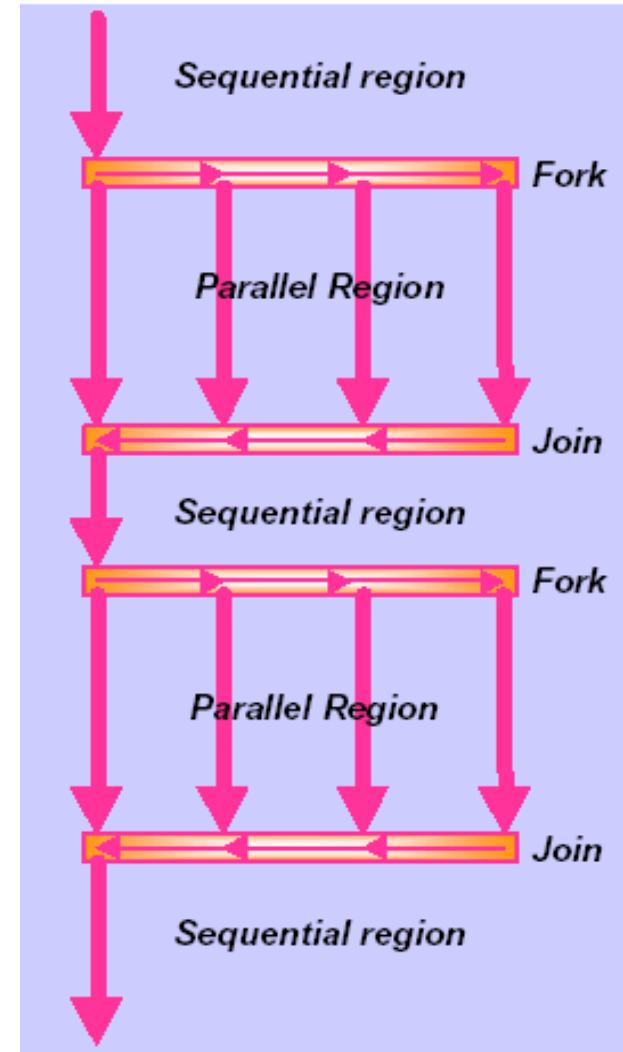




# OpenMP: Programming Model

## ❖ Fork – Join parallelism

- OpenMP uses fork and join model for parallel execution
- OpenMP programs begin with single process: **master thread**.
- FORK : Master thread creates a team of parallel threads
- JOIN: When the team threads complete the statements in parallel region, they synchronize and terminate leaving master thread.
- Parallelism is added incrementally



# OpenMP: Programming Model

## ❖ Threads based parallelization

- Open MP is based on the existence of multiple threads in the shared memory programming paradigm

## ❖ Explicit parallelization

- It is an explicit programming model, and offers full control over parallelization to the programmer

## ❖ Compiler directive based

- All of OpenMP parallelization is supported through the use of compiler directives

## ❖ Nested parallelism support

- The API support placement of parallel construct inside other parallel construct

Source : Reference : [4], [6], [14],[17], [22], [28]

# OpenMP: Programming Model

## ❖ Dynamic threads

- The API provides dynamic altering of number of threads (Depends on the implementation)

## How do threads interact?

### ❖ OpenMP is shared memory model.

- Threads communicate by sharing variables

### ❖ Unintended sharing of data can lead to race conditions:

- Race condition : when the program's outcome changes as the threads are scheduled differently
- To control race conditions: Use synchronization to protect data conflicts

# OpenMP : Fortran Directives Format

## Format

sentinel	directive-name	[clause ...]
<p>All Fortran OpenMP directives must begin with a sentinel. The accepted sentinels depend upon the type of Fortran source. Possible sentinels are:</p> <p><b>!\$OMP</b> <b>C\$OMP</b> <b>*\$OMP</b></p>	<p>A valid OpenMP directive. Must appear after the sentinel and before any clauses.</p>	<p>Optional. Clauses can be in any order, and repeated as necessary unless otherwise restricted.</p>

## Example

```
!$OMP PARALLEL SHARED(ALPHA) PRIVATE(BETA)
```

Source : Reference : [4], [6], [14],[17], [22], [28]

# OpenMP : C/C++ Directives Format

## Format

sentinel	directive-name	[clause ...]	newline
Required for all OpenMP C/C++ directives  #pragma omp	A valid OpenMP directive. Must appear after the <b>pragma</b> and before any clauses.	Optional. Clauses can be in any order, and repeated as necessary unless otherwise restricted.	Required. Proceeds the structured block which is enclosed by this directive.

## Example

```
#pragma omp parallel shared(alpha), private(beta)
```

# OpenMP : C/C++ Directives Format

## Format : Compiler Directive

### ❖ C/C++

```
sentinel directive-name clause
```

Ex.

```
#pragma omp parallel shared(alpha), private(beta)
```

### ❖ Fortran

```
sentinel directive-name clause
```

Ex.

```
!$OMP directive [clause, ...]  
!C$OMP directive [clause, ...]  
!*$OMP directive [clause, ...]
```

# OpenMP : C/C++ General Code Structure

```
# include <omp.h>
main() {
int var1, var2, var3;
```

*serial code*

.....

Beginning of parallel region, fork a team of threads.

*Specify variable scoping*

```
#pragma omp parallel private (var1, var2), shared(var3)
```

```
{
```

*Parallel region executed by all threads*

.....

.....

*All threads join master thread and disband*

```
}
```

*Resume Serial code*

```
}
```

# OpenMP : C/C++ General Code Structure

```
# include <omp.h>
```

```
main() {
```

```
int var1, var2, var3; Create thread here for this parallel region
```

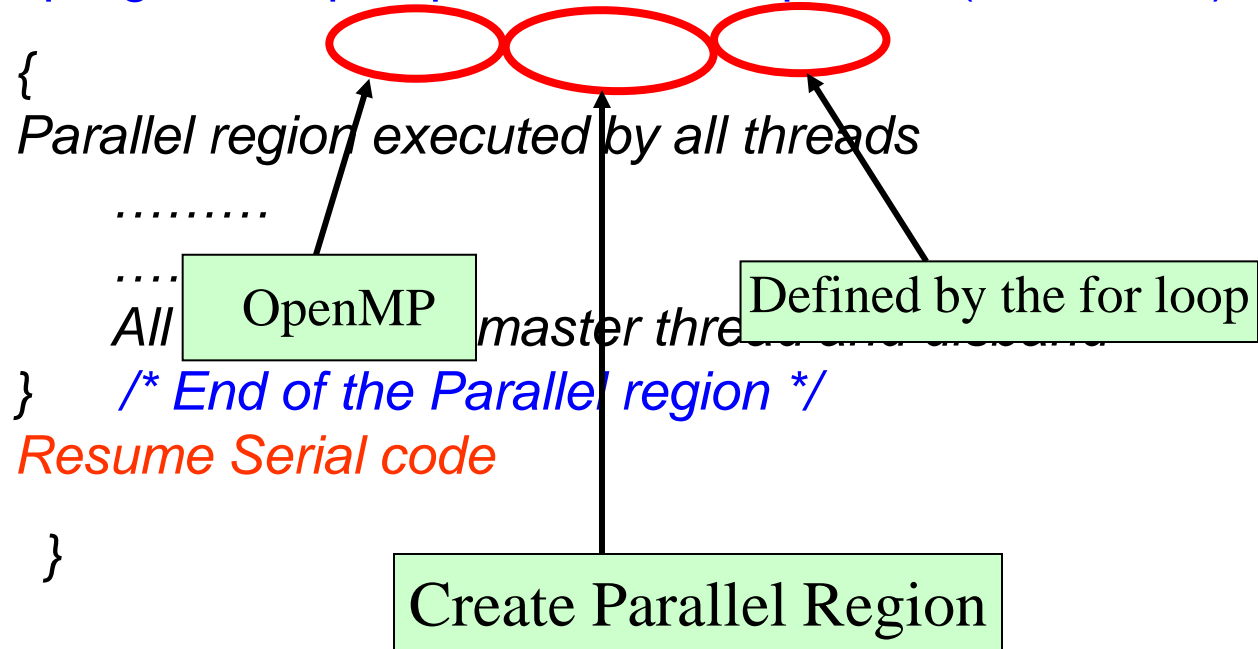
*serial*

*code*

.....

Beginning of parallel region, fork a team of threads.

```
#pragma omp parallel for private (var1, var2)
```





# OpenMP : FORTRAN General Code Structure

```
PROGRAM HELLO  
INTEGER VAR1, VAR2, VAR3
```

*Serial code*

.....

.....

*Beginning of parallel region, fork a team of threads.*

Specify variable scoping

```
!$OMP PARALLEL PRIVATE (VAR1, VAR2), SHARED(VAR3)
```

*Parallel region executed by all threads*

.....

.....

*All threads join master thread and disband*

```
!$OMP END PARALLEL
```

.....

*Resume Serial code*

```
END
```

# OpenMP : How is OpenMP typically used? (C/C++)

- ❖ OpenMP is usually used to parallelize loops:
  - Find your most time consuming loops.
  - Split them up between threads.

**Split-up this loop between multiple threads**

```
void main()
{
    double Res[1000];

    for(int i=0;i<1000;i++) {
        do_huge_comp(Res[i]);
    }
}
```

**Sequential Program**

```
#include "omp.h"
void main()
{
    double Res[1000];

    #pragma omp parallel for
    for(int i=0;i<1000;i++) {
        do_huge_comp(Res[i]);
    }
}
```

**Parallel Program**

# OpenMP : How is OpenMP typically used? (Fortran)

- ❖ OpenMP is usually used to parallelize loops:
  - Find your most time consuming loops.
  - Split them up between threads.

**Split-up this loop between multiple threads**

program example  
double precision Res(1000)  
  
do l=1,1000  
  call huge\_comp(Res(l))  
end do  
end

**Sequential Program**

program example  
double precision Res(1000)  
  
**C\$OMP PARALLEL DO**  
do l=1,1000  
  call huge\_comp(Res(l))  
end do  
end

**Parallel Program**

# OpenMP : Constructs

## Main categories of OpenMP's constructs:

### ❖ Directives

- Parallel Regions
- Work-sharing
- Data Environment
- Synchronization

### ❖ Runtime library functions

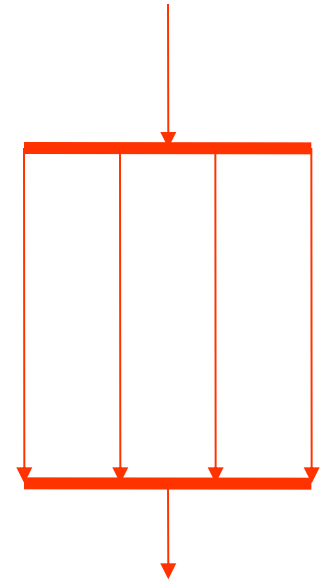
- Execution Environment Functions
- Lock functions
- Timing routines

### ❖ Environment variables

# OpenMP : 'PARALLEL' Region Construct

**A Parallel Region is a block of code executed by all threads simultaneously**

- The master thread always has thread ID 0
- Thread adjustment (if enabled) is only done before entering a parallel region
- Parallel regions can be nested, but support for this is implementation dependent



All thread perform identical task

# OpenMP : PARALLEL Region Construct

## Example

Each thread redundantly executes the code within the structured block

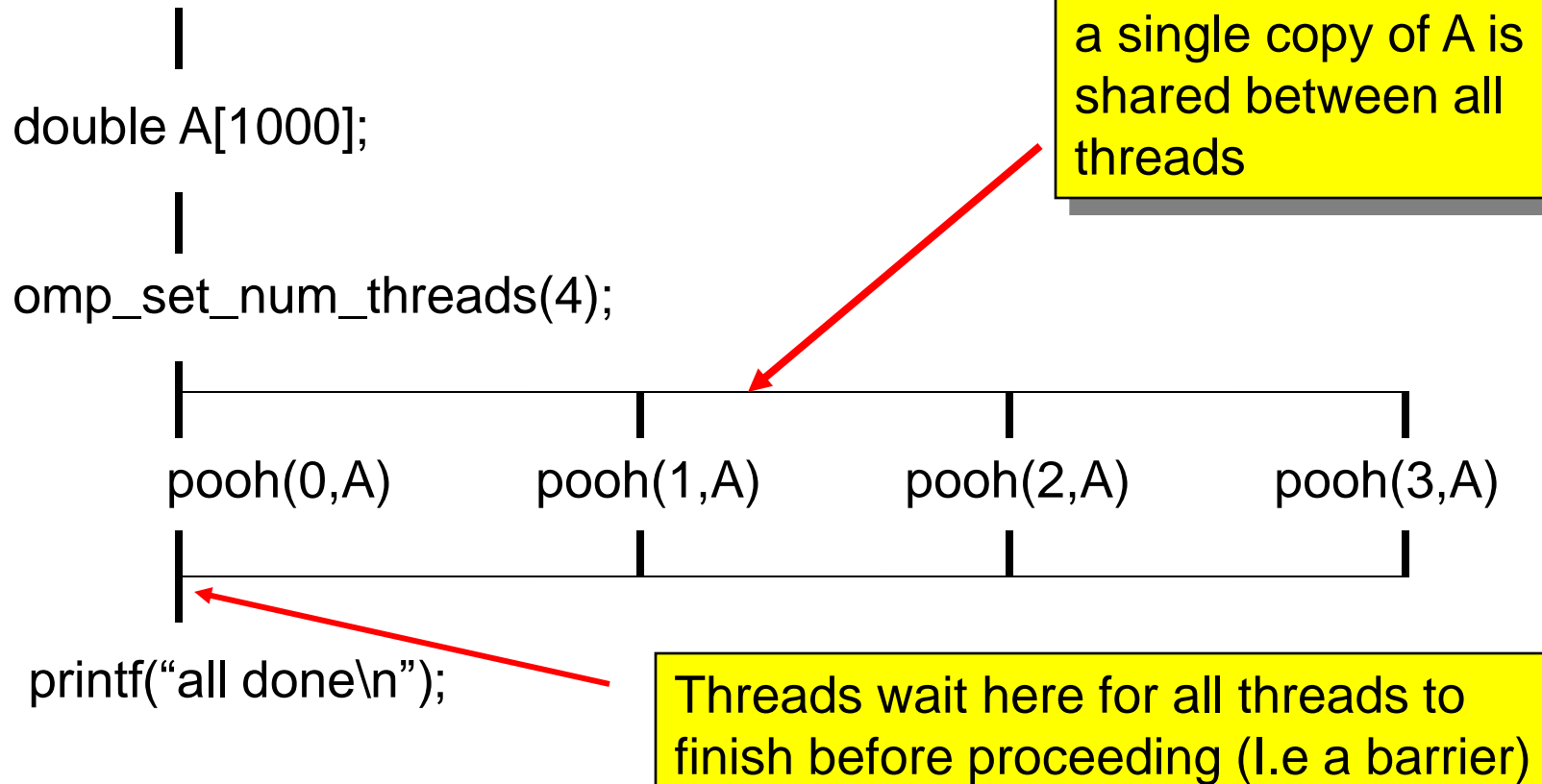
```
double A[1000];
omp_set_num_threads(4);
#pragma omp parallel
{
    int ID = omp_thread_num();
    pooh(ID,A);
}
Printf("all done\n");
```

Runtime function to request a certain number of threads

Runtime function returning a thread ID

# OpenMP : PARALLEL Region Construct

Each thread executes the same code redundantly.



# OpenMP : Work-sharing Construct

**It distributes the execution of the associated statement among the members of the team that encounter it**

- ❖ Work sharing construct do not launch new threads
- ❖ There is no barrier upon entry to work-sharing construct.
- ❖ There is an implied barrier at the end of a worksharing construct

## Restrictions

- ❖ Must be enclosed in the parallel region for parallel execution
- ❖ Must be encountered by all the members of the team or none of them



# OpenMP : Work-sharing Construct

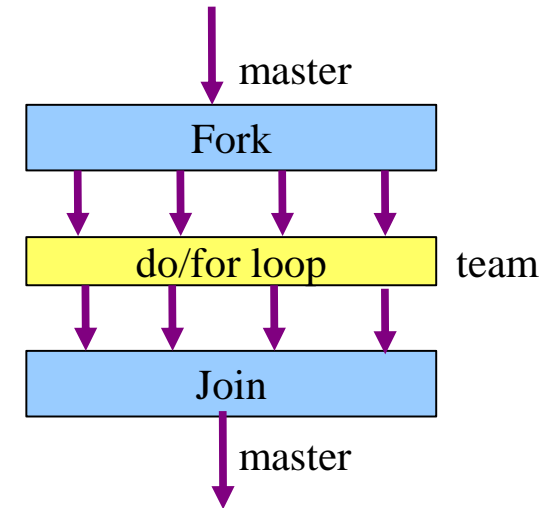
**OpenMP defines the following work-sharing constructs.**

- ❖ **for** directive
- ❖ **sections** directive
- ❖ **single** directive

# OpenMP : Work-sharing construct 'for'

- ❖ OpenMP is usually used to parallelize loops:
  - Find your most time consuming loops.
  - Split them up between threads.

**Split-up this loop between multiple threads**



```
void main()
{
    double Res[1000];

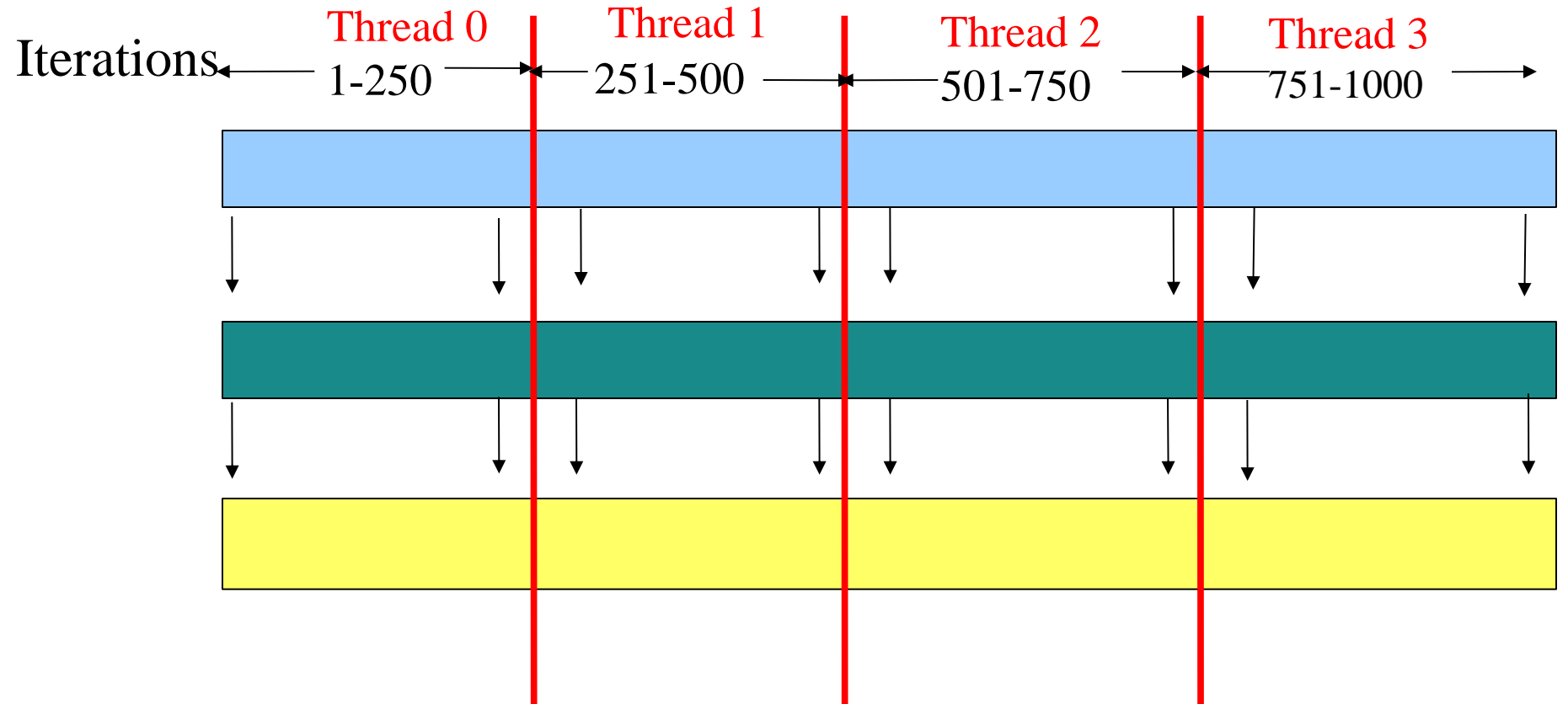
    for(int i=0;i<1000;i++) {
        do_huge_comp(Res[i]);
    }
}
```

**Sequential Program**

```
#include "omp.h"
void main()
{
    double Res[1000];
    #pragma omp parallel for
    for(int i=0;i<1000;i++) {
        do_huge_comp(Res[i]);
    }
}
```

**Parallel Program**

# OpenMP : Work-sharing construct 'for'



# OpenMP : Work-sharing Construct

## for directive

for directive identifies the iterative work-sharing construct.

```
#pragma omp for [clause[[,]clause]...] new-line  
for-loop
```

Clause is one of the following:

*private(variable list)*

*firstprivate (variable list)*

*lastprivate (variable list)*

*reduction (variable list)*

*ordered , nowait*

# OpenMP : Work-sharing Construct

## for directive

The “for” Work-Sharing construct splits up loop iterations among the threads in a team

```
#pragma omp parallel
```

```
#pragma omp for
```

```
    for (I=0;I<N;I++) {  
        NEAT_STUFF(I);  
    }
```

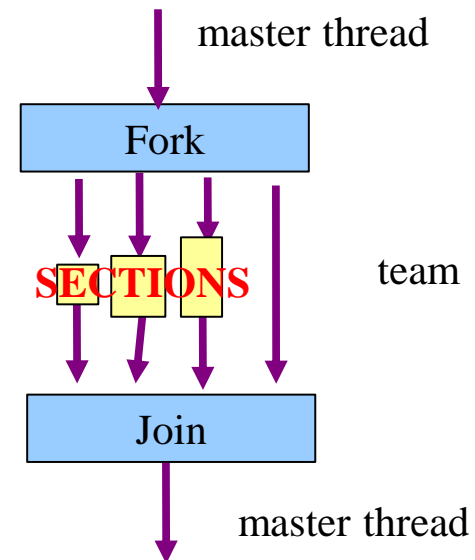
By default, there is a barrier at the end of the “omp for”.

# OpenMP : Work-sharing Construct

## sections directive

**sections directive gives different structured blocks to each thread.**

```
#pragma omp parallel
#pragma omp sections {
#pragma omp section
    x_calculation(); // thread 1 work
#pragma omp section
    y_calculation(); // thread 2 work
.....
}
```



# OpenMP : Work-sharing Construct

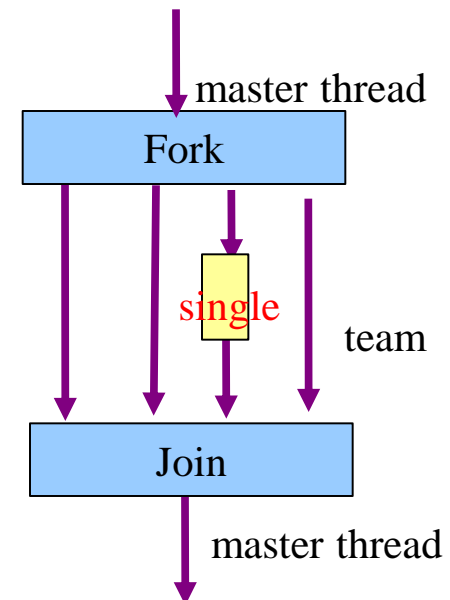
## single directive

This identifies that the associated structured block is to be executed by only one thread in the team (It can be any thread including master thread).

```
#pragma omp single [clause[[, clause] ...] new-line
```

```
structured-block
```

```
Example : !$omp single  
          call read_array(in, len)  
          !$omp end single
```



# OpenMP : Data Environment

## Default Storage attributes

### ❖ Shared Memory programming model:

- Most variables are shared by default

### ❖ Global variables and SHARED among threads

- Fortran : COMMON blocks, SAVE variables, MODULE variables.
- C: File scope variables, static

### ❖ But not everything is shared...

- Stack variables in sub-programs called from parallel regions are PRIVATE
- Automatic variables within a statement block are PRIVATE.



# OpenMP : Data Environment

## Example : Storage attributes

```
program sort
```

```
common /input/ A(10)
```

```
integer index(10)
```

```
call input
```

```
C$OMP PARALLEL
```

```
call work(index)
```

```
C$OMP END PARALLEL
```

```
print *, index(1)
```

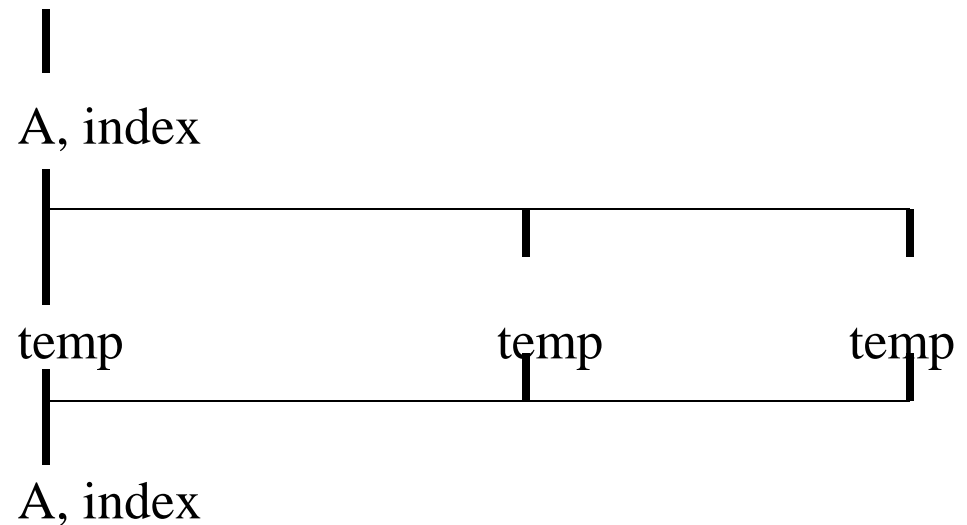
A, index are shared by all threads

temp is local to each thread

```
subroutine work
```

```
common /input/ A(10)
```

```
real temp(10)
```



# OpenMP : Data Environment

## Changing the storage attributes

One can selectively change storage attributes constructs using the following clauses\*

- SHARED** declares variables to be shared among all threads in the team
- PRIVATE** declares variables to be private to each thread.
- FIRSTPRIVATE** performs initialization of private variables
- LASTPRIVATE** performs finalization of private variables
- REDUCTION** performs a reduction on the variables subject to given operator.

The default status can be modified with:

**DEFAULT (PRIVATE | SHARED)**

# OpenMP : Data Environment

## private clause

Creates a local copy of variable for each thread.

- The value is **un-initialized**
- Private copy is **not** storage associated with the original

program sample

```
IS = 0
```

```
!$OMP PARALLEL DO PRIVATE(IS)
```

```
DO J=1,1000
```

```
IS = IS + J
```

```
ENDDO
```

```
print *, IS
```

IS was not initialized

IS is undefined at this point

# OpenMP : Data Environment

- ❖ First private initializes each thread's copy of a private variable to the value of the master copy.
- ❖ Last private writes back to the master's copy the value of private copy that executed the Sequentially last iteration.

program closer

```
IS = 0
```

```
C$OMP PARALLEL DO  
FIRSTPRIVATE(IS) LASTPRIVATE(IS)
```

```
DO J=1,1000
```

```
IS = IS + J
```

```
ENDDO
```

```
print *, IS
```

**Each thread gets its own IS with an initial value of 0**

**IS is defined as its value at the last iteration (i.e. for J=1000)**

## Example : Firstprivate & lastprivate

```
int x;
x = 0;
#pragma omp parallel for firstprivate(x)
for (i = 0; i < 10000; i++) {
    x = x + i;
}
printf( "x is %d\n", x );
```

Initialise x to zero

Each thread gets its own is with an initial value of 0

Print out value of x

**Oops! The value x is undefined!**  
Need lastprivate(x) to copy value back out to master

# OpenMP : Data Environment

**Example :**  
**Default Clause**

```
        itotal = 1000
#pragma omp parallel private(np, each)
{
        np = omp_get_num_threads()
        each = itotal/np
}
```

These two  
codes are  
equivalent

```
        itotal = 1000
#pragma omp parallel Default(private) shared(itotal)
{
        np = omp_get_num_threads()
        each = itotal/np
}
```

## Default clause

This clause is used for changing the default status of the variables.

### ❖ **default (private)**

- *Each* variable in *static* extent of the parallel region is made **private** as if specified in a private clause

### ❖ **default (shared)**

- *Each* variable in *static* extent of the parallel region is made **shared** as if specified in a shared clause

### ❖ **default (none)**

- *no* default for variables in static extent.

## OpenMP : Data Environment

**Example :**  
**Default Clause**

```
        itotal = 1000
C$OMP PARALLEL PRIVATE(np, each)
        np = omp_get_num_threads()
        each = itotal/np
C$OMP END PARALLEL
```

**These two  
codes are  
equivalent**

```
        itotal = 1000
C$OMP PARALLEL DEFAULT(PRIVATE)
        SHARED(itotal)
        np = omp_get_num_threads()
        each = itotal/np
C$OMP END PARALLEL
```



# OpenMP : Data Environment

## Reduction

- ❖ Another clause that effects the way variables are shared
  - `reduction(op:list)`
- ❖ The variables in “list” must be shared in the enclosing parallel region.
- ❖ Inside a parallel or a work sharing construct:
  - A local copy of each list variable is made and initialized depending on the “op” (e.g 0 for “+”)
  - Pair wise “op” is updated on the local value
  - Local copies are reduced into a single global copy at the end of the construct

# OpenMP Clauses

## Example : Reduction

```
#include <omp.h>
#define NUM_THREADS 2
void main() {
    int I;
    double ZZ, func(), res = 0.0;
    omp_set_num_threads(NUM_THREADS);
    #pragma omp parallel for reduction(+;res) private(ZZ)
        for(I=0;I<1000;I++)}
        ZZ=func(I);
        res = res + ZZ;
    }
}
```

# OpenMP: Synchronization

❖ OpenMP has the following constructs support synchronization

➤ Atomic

➤ Critical section

➤ Barrier

➤ Flush

➤ Ordered

# OpenMP: Synchronization constructs

## ❖ Some of the OpenMP synchronization constructs

- Single

- Master

- Atomic

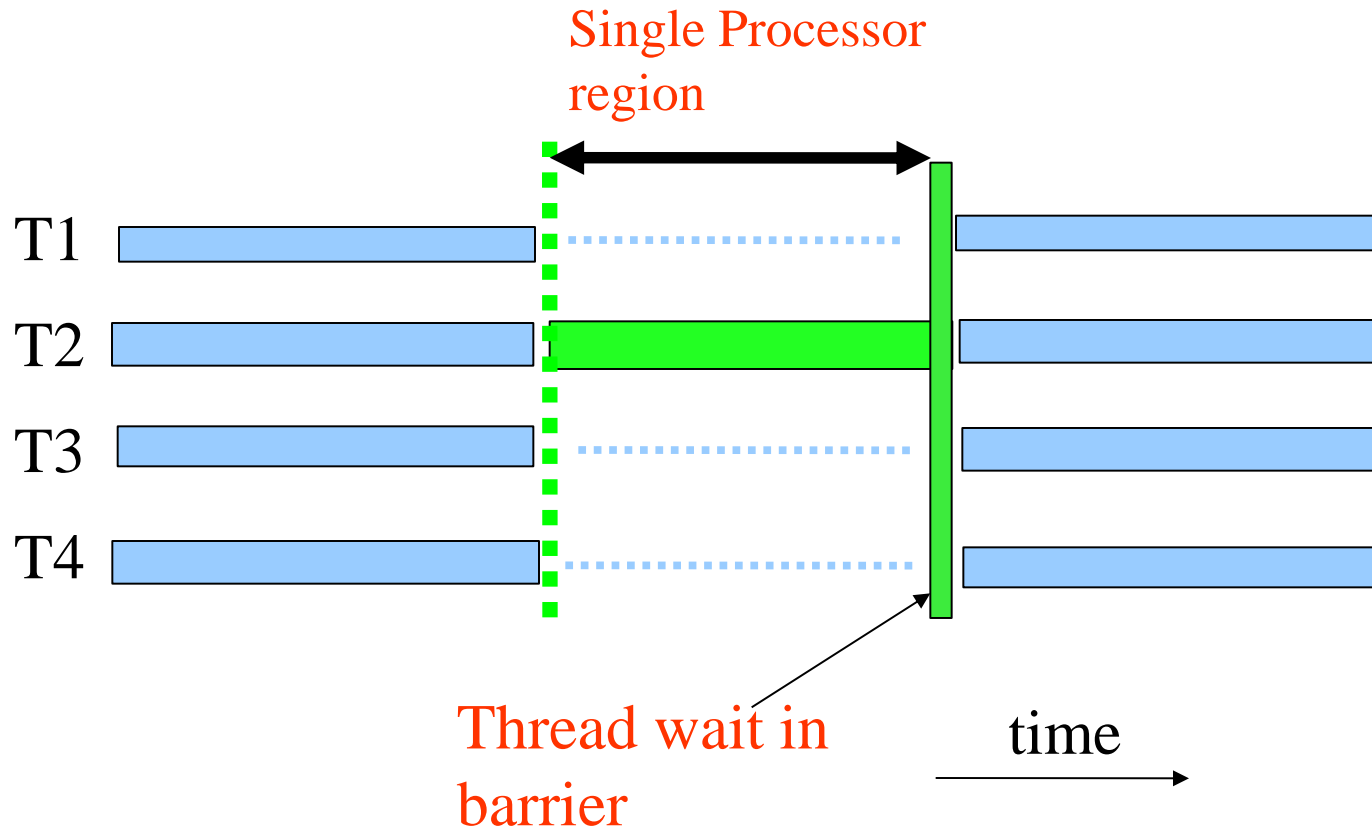
- Critical section

- Barrier

- Ordered

# SINGLE and MASTER Construct

Usually, there is a barrier at the end of the region



# SINGLE and MASTER Construct

Only one thread in a team executes the code enclosed

```
#pragma omp single [clause[[,] clause] ...]  
{  
    <code-block>  
}
```

Only the master thread executes the code block

```
#pragma omp master  
{  
    <code-block>  
}
```

There is no implied  
barrier on the  
entry & exit

# Synchronization : Critical Section and Atomic

**Only one thread at a time can enter a critical section**

If sum is a shared variable, this loop can not run in parallel

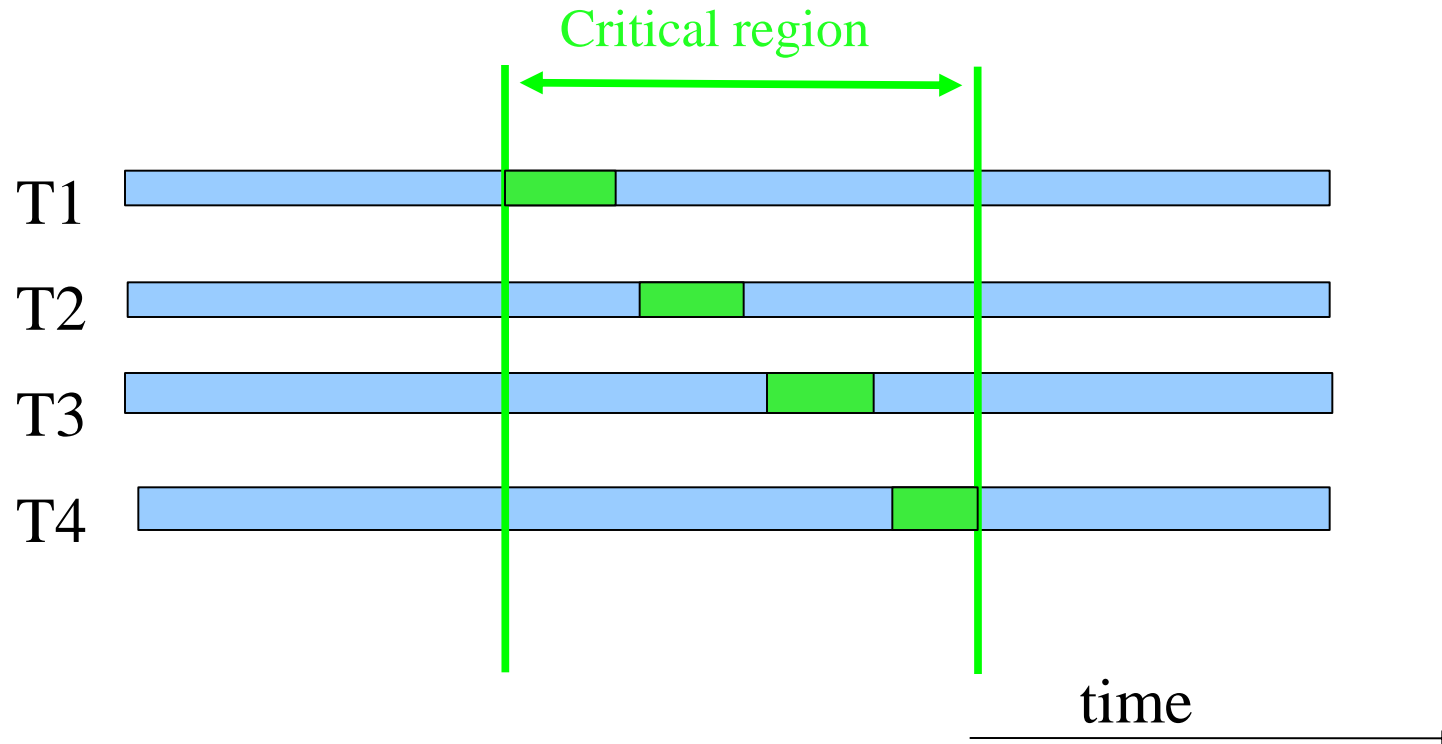
```
for (i=0; i < N; i++){  
    ....  
    sum += a[i];  
    .... }
```

We can use a critical region for this:

```
for (i=0; i < N; i++){  
    one at a time can proceed  
    .....  
    sum += a[i];  
    .....  
    next in line, please  
}
```

# Synchronization : Critical Section and Atomic

- Useful to avoid a race condition, or to perform I/O
- Be aware that your parallel computation may be serialized





# Synchronization : Critical Section and Atomic

All threads execute the code, but only one at a time:

```
#pragma omp critical [(name)]  
{<code-block>}
```

There is no implied  
barrier on entry or exit

```
#pragma omp atomic  
<statement>
```

This is a lightweight,  
special form of a  
critical section

# Synchronization : Critical Section and Atomic

Only one thread at a time can enter a critical section

```
#pragma omp parallel for  
for (i=0; i < N; i++){  
  #pragma omp critical  
    sum = sum + a[i];  
}
```

one thread at a time



# Synchronization construct: Barrier

Suppose we run each of these two loops in parallel over  $i$ :

```
for (i=0; i < N; i++)  
  a[i] = b[i] + c[i];
```

```
for (i=0; i < N; i++)  
  d[i] = a[i] + b[i];
```

**This may give us a wrong answer ?**

# Synchronization construct : Barrier

We need to have updated all of a[ ] first, before using a[ ]

```
for (i=0; i < N; i++)  
a[i] = b[i] + c[i];
```

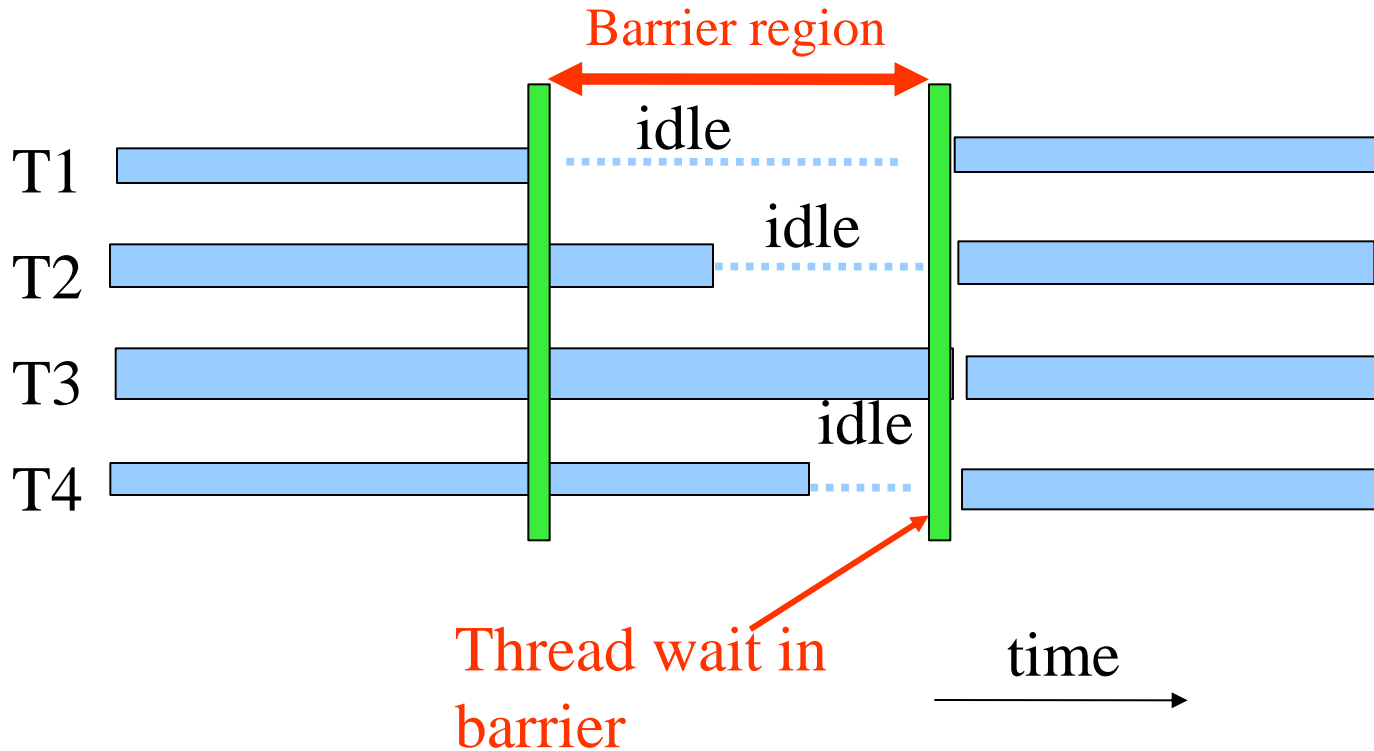
**wait !**

**barrier**

```
for (i=0; i < N; i++)  
d[i] = a[i] + b[i];
```

**All threads wait at the barrier point and only continue when all threads have reached the barrier point**

# Synchronization Construct :Barrier



Each thread waits until all others have reached this point:

```
#pragma omp barrier
```

# OpenMP: Synchronization

## Critical Section

- ❖ Only one thread at a time can enter a critical section

```
cur_max = MINUS_INFINITY
!$omp parallel do
  do I = 1 , n
    !omp critical
      if (a(I) .gt. Cur_max) then
        cur_max = a(I)
      endif
    !omp end critical
  endif
enddo
```

# OpenMP: Synchronization

```
#pragma omp parallel shared (A, B, C)
  private(d)
{
  id = omp_get_thread_num();
  A[id]=big_calc1(id);
#pragma omp barrier
#pragma omp for
  for(l=0;l<N;l++){ C[l]=big_calc3(l,A);}
#pragma omp for nowait
  for(l=0;l<N;l++){ B[l]=big_calc2(C,l);}
  A[id]=big_calc3(id);
}
```

**Barrier** : Each thread waits until all threads arrive

Implicit barrier at the end of a for work-sharing construct

No implicit barrier due to nowait

Implicit barrier at the end of a parallel region

# OpenMP: Synchronization

## Ordered

- ❖ The ordered construct enforces the sequential order for a block.

```
#pragma omp parallel private(tmp)
```

```
#pragma omp for ordered
```

```
    for (I=0;I<N;I++){
```

```
        tmp = NEAT_STUFF(I);
```

```
#pragma ordered
```

```
    res = consum(tmp);
```

```
}
```



## flush

- ❖ The FLUSH construct denotes a sequence point where a thread tries to create a consistent view of memory.
  - All memory operations (both reads and writes) defined prior to the sequence point must complete.
  - All memory operations (both reads and writes) defined after the sequence point must follow the flush.
  - Variables in registers or write buffers must be updated in memory.
- ❖ Arguments to flush specify which variables are flushed. No arguments specifies that all thread visible variables are flushed.

# OpenMP: Library Routines

## ❖ Lock routines

omp\_init\_lock(),      omp\_set\_lock(),  
omp\_destroy\_lock(),  
omp\_unset\_lock(),  
omp\_test\_lock()

## ❖ Runtime environment routines:

- Modify/Check the number of threads

omp\_set\_num\_threads(),  
omp\_get\_num\_threads(),  
omp\_get\_thread\_num(),  
omp\_get\_max\_threads()

## ❖ Runtime environment routines

➤ How many processors in the system?

–omp\_num\_procs()

➤ Turn on/off nesting and dynamic mode

omp\_set\_nested(),

omp\_get\_nested(),

omp\_set\_dynamic(),

omp\_get\_dynamic()

## ❖ Protect resources with locks

```
    omp_lock_t lck;  
    omp_init_lock(&lck);  
#pragma omp parallel private(tmp)  
{  
    id = omp_get_thread_num();  
    tmp = do_lots_of_work(id);  
    omp_set_lock(&lck);  
    printf("%d %d", id, tmp);  
    omp_unset_lock(&lck);  
}
```

## OpenMP: Library Routines

- ❖ **To fix the number of threads used in a program, first turn off dynamic mode and then set the number of threads.**

```
#include <omp.h>

Void main()
{
    omp_set_dynamic(0);
    omp_set_num_threads(4);
    #pragma omp parallel {
        Int id=omp_get_thread_num();
        do_lots_of-stuff(id); }
}
```

# OpenMP :Environment Variables

## ❖ Environment Variables

<b>Variables</b>	<b>Value</b>	<b>Description</b>
OMP_NUM_THREADS	4	Specify the no. of threads
OMP_DYNAMIC	TRUE or FALSE	Enable/disable dynamic adj of threads
OMP_NESTED	TRUE or FALSE	Enable/disable nested parallelism

# A more elaborate example

```
#pragma omp parallel if (n>limit) default(none) \  
    shared(n,a,b,c,x,y,z) private(f,i,scale)
```

```
{  
    f = 1.0;  
#pragma omp for  
    for (i=0; i<n; i++)  
        z[i] = x[i] + y[i];
```

```
#pragma omp for  
    for(i=0; i<n; i++)  
        a[i] = b[i] + c[i];
```

```
#pragma omp barrier
```

```
    ....  
    scale = sum(a,0,n) + sum(z,0,n) + f;  
    ....  
} /*-- End of parallel region --*/
```

Statement is executed by all threads

parallel loop (work will be distributed)

parallel loop (work will be distributed)

synchronization

Statement is executed by all threads

Parallel region

# Compilation & Execution of OpenMP programs

```
#include<stdio.h>
#include<omp.h>
main() {
#pragma omp parallel
{
    printf( "hello world from thread %d of
           %d\n", omp_get_thread_num(),
           omp_get_num_threads() );
}
}
```



# Compilation & Execution of OpenMP programs

## Compilation :

```
$ cc -o <objectFileName> <programName> <omp-compiler-flag>
```

Ex.

```
$gcc -o omp-hello-world omp-hello-world.c -fopenmp
```

## Setting the Number of Threads :

❖ Environment Variables :

```
$ export OMP_NUM_THREADS= <No. of threads>
```

❖ Environment variable can be overridden by the programmer :

```
omp_set_num_threads(int n)
```

# Sample Output

**From a Dual Socket Quad Core machine:**

hello world from thread 0 of 8

hello world from thread 2 of 8

hello world from thread 3 of 8

hello world from thread 7 of 8

hello world from thread 6 of 8

hello world from thread 1 of 8

hello world from thread 4 of 8

hello world from thread 5 of 8

# Commonly Encountered Questions While Threading Application ?

- ❖ What should the expected speedup be?
- ❖ Will the performance meet expectations?
- ❖ Will it scale if the more number of processor added?
- ❖ Which threading model is it?

# Commonly Encountered Questions While Threading Application ?

## Analysis

- ❖ Where to thread ?
  - thread the more time consuming section of code like loops
- ❖ How long would it take to thread?
  - Very minimum time just need to use some directives/library routine
- ❖ How much re-design / efforts is required?
  - Very less
- ❖ Is it worth threading the selected region ?
  - Appears to have minimal dependencies
  - Consuming over 90% of run time

## OpenMP :Conclusions

- ❖ Features and advantages of OpenMP is discussed.
- ❖ OpenMP programming models are covered.
- ❖ Parallelization using OpenMP is explained.
- ❖ Various OpenMP Constructs are discussed with examples.

Source : Reference : [4], [6], [14],[17], [22], [28]

## References

1. Andrews, Grogory R. **(2000)**, Foundations of Multithreaded, Parallel, and Distributed Programming, Boston, MA : Addison-Wesley
2. Butenhof, David R **(1997)**, Programming with POSIX Threads , Boston, MA : Addison Wesley Professional
3. Culler, David E., Jaswinder Pal Singh **(1999)**, Parallel Computer Architecture - A Hardware/Software Approach , San Francscico, CA : Morgan Kaufmann
4. Grama Ananth, Anshul Gupts, George Karypis and Vipin Kumar **(2003)**, Introduction to Parallel computing, Boston, MA : Addison-Wesley
5. Intel Corporation, **(2003)**, Intel Hyper-Threading Technology, Technical User's Guide, Santa Clara CA : Intel Corporation Available at : <http://www.intel.com>
6. Shameem Akhter, Jason Roberts **(April 2006)**, Multi-Core Programming - Increasing Performance through Software Multi-threading , Intel PRESS, Intel Corporation,
7. Bradford Nichols, Dick Buttlar and Jacqueline Proulx Farrell **(1996)**, Pthread Programming O'Reilly and Associates, Newton, MA 02164,
8. James Reinders, Intel Threading Building Blocks – **(2007)** , O'REILLY series
9. Laurence T Yang & Minyi Guo (Editors), **(2006)** *High Performance Computing - Paradigm and Infrastructure* Wiley Series on Parallel and Distributed computing, Albert Y. Zomaya, Series Editor
10. Intel Threading Methodology ; Principles and Practices Version 2.0 copy right **(March 2003)**, Intel Corporation

## References

11. William Gropp, Ewing Lusk, Rajeev Thakur **(1999)**, Using MPI-2, Advanced Features of the Message-Passing Interface, The MIT Press..
12. Pacheco S. Peter, **(1992)**, Parallel Programming with MPI, , University of Sanfrancisco, Morgan Kaufman Publishers, Inc., Sanfrancisco, California
13. Kai Hwang, Zhiwei Xu, **(1998)**, Scalable Parallel Computing (Technology Architecture Programming), McGraw Hill New York.
14. Michael J. Quinn **(2004)**, Parallel Programming in C with MPI and OpenMP McGraw-Hill International Editions, Computer Science Series, McGraw-Hill, Inc. Newyork
15. Andrews, Grogory R. **(2000)**, Foundations of Multithreaded, Parallel, and Distributed Progrmaming, Boston, MA : Addison-Wesley
16. SunSoft. Solaris multithreaded programming guide. SunSoft Press, Mountainview, CA, **(1996)**, Zomaya, editor. Parallel and Distributed Computing Handbook. McGraw-Hill,
17. Chandra, Rohit, Leonardo Dagum, Dave Kohr, Dror Maydan, Jeff McDonald, and Ramesh Menon, **(2001)**,Parallel Programming in OpenMP San Fracncisco Moraan Kaufmann
18. S.Kieriman, D.Shah, and B.Smaalders **(1995)**, Programming with Threads, SunSoft Press, Mountainview, CA. 1995
19. Mattson Tim, **(2002)**, Nuts and Bolts of multi-threaded Programming Santa Clara, CA : Intel Corporation, Available at : <http://www.intel.com>
20. I. Foster **(1995)**, Designing and Building Parallel Programs ; Concepts and tools for Parallel Software Engineering, Addison-Wesley (1995)
21. J.Dongarra, I.S. Duff, D. Sorensen, and H.V.Vorst **(1999)**, Numerical Linear Algebra for High Performance Computers (Software, Environments, Tools) SIAM, 1999

## References

22. OpenMP C and C++ Application Program Interface, Version 1.0". **(1998)**, OpenMP Architecture Review Board. October 1998
23. D. A. Lewine. *Posix Programmer's Guide: (1991)*, Writing Portable Unix Programs with the Posix. 1 Standard. O'Reilly & Associates, 1991
24. Emery D. Berger, Kathryn S McKinley, Robert D Blumofe, Paul R. Wilson, *Hoard : A Scalable Memory Allocator for Multi-threaded Applications* ; The Ninth International Conference on Architectural Support for Programming Languages and Operating Systems (ASPLOS-IX). Cambridge, MA, November **(2000)**. Web site URL : <http://www.hoard.org/>
25. Marc Snir, Steve Otto, Steyen Huss-Lederman, David Walker and Jack Dongarra, **(1998)** *MPI-The Complete Reference: Volume 1, The MPI Core, second edition* [MCMPI-07].
26. William Gropp, Steven Huss-Lederman, Andrew Lumsdaine, Ewing Lusk, Bill Nitzberg, William Saphir, and Marc Snir **(1998)** *MPI-The Complete Reference: Volume 2, The MPI-2 Extensions*
27. A. Zomaya, editor. *Parallel and Distributed Computing Handbook*. McGraw-Hill, **(1996)**
28. OpenMP C and C++ Application Program Interface, Version 2.5 **(May 2005)**", From the OpenMP web site, URL : <http://www.openmp.org/>
29. Stokes, Jon 2002 Introduction to Multithreading, Super-threading and Hyper threading *Ars Technica*, October **(2002)**
30. Andrews Gregory R. 2000, *Foundations of Multi-threaded, Parallel and Distributed Programming*, Boston MA : Addison – Wesley **(2000)**
31. Deborah T. Marr , Frank Binns, David L. Hill, Glenn Hinton, David A Koufaty, J . Alan Miller, Michael Upton, "Hyperthreading, Technology Architecture and Microarchitecture", Intel **(2000-01)**



**Thank You**  
*Any questions ?*