Introduction to OpenMP

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CS594: Scientific Computing for Engineers

February 29, 2012



Acknowledgment

These slides are adapted from the Lawrence Livermore
 OpenMP Tutorial by Blaise Barney at

https://computing.llnl.gov/tutorials/openMP/



What is OpenMP?

- An Application Program Interface (API) that may be used to explicitly direct multi-threaded, shared memory parallelism
- Comprises three primary API components
 - Compiler Directives
 - Runtime Library Routines
 - Environment Variables
- Portable
 - The API is specified for C/C++ and Fortran
 - Has been implemented for most major platforms including Unix/ Linux platforms and Windows NT



What is OpenMP? (cont.)

- Standardized
 - Jointly defined and endorsed by a group of major computer hardware and software vendors
 - Expected to become an ANSI standard later???
- What does OpenMP stand for?
 - Short version: **Open Multi-Processing**
 - Long version: Open specifications for Multi-Processing via collaborative work between interested parties from the hardware and software industry, government and academia.



OpenMP is NOT

- Meant for distributed memory parallel systems (by itself)
- Necessarily implemented identically by all vendors
- Guaranteed to make the most efficient use of shared memory
- Required to check for data dependencies, data conflicts, race conditions, or deadlocks
- Required to check for code sequences that cause a program to be classified as non-conforming
- Meant to cover compiler-generated automatic parallelization and directives to the compiler to assist such parallelization
- Designed to guarantee that input or output to the same file is synchronous when executed in parallel (The programmer is responsible for synchronizing input and output).



References

- OpenMP website: <u>openmp.org</u> API specifications, FAQ, presentations, discussions, media releases, calendar, membership application and more...
- Wikipedia: en.wikipedia.org/wiki/OpenMP
- Barbara Chapman, Gabriele Jost, and Ruud van der Pas: Using OpenMP. The MIT Press, 2008.
- Compiler documentation
 - IBM: <u>www-4.ibm.com/software/ad/fortran</u>
 - Cray: <u>http://docs.cray.com/</u> (Cray Fortran Reference Manual)
 - Intel: www.intel.com/software/products/compilers/
 - PGI: <u>www.pgroup.com</u>
 - PathScale: <u>www.pathscale.com</u>
 - GNU: <u>gnu.org</u>



History of OpenMP

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- In the early 90's, vendors of shared-memory machines supplied similar, directive-based, Fortran programming extensions.
 - The user would augment a serial Fortran program with directives specifying which loops were to be parallelized.
 - The compiler would be responsible for automatically parallelizing such loops across the SMP processors.
- Implementations were all functionally similar, but were divergent.
- First attempt at a standard was the draft for ANSI X3H5 in 1994. It was never adopted, largely due to waning interest as distributed memory machines became popular.



History of OpenMP (cont.)

- The OpenMP standard specification started in the spring of 1997, taking over where ANSI X3H5 had left off, as newer shared memory machine architectures started to become prevalent.
- Led by the OpenMP Architecture Review Board (ARB). Original ARB members included: (Disclaimer: all partner names derived from the <u>OpenMP web site</u>)
 - Compaq / Digital
 - Hewlett-Packard Company
 - Intel Corporation
 - International Business Machines (IBM)
 - Kuck & Associates, Inc. (KAI)
 - Silicon Graphics, Inc.
 - Sun Microsystems, Inc.
 - U.S. Department of Energy ASCI program

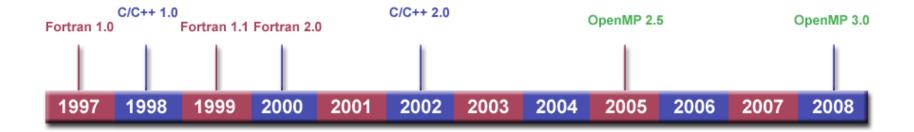


Other Contributors

- Endorsing application developers
 - ADINA R&D, Inc.
 - ANSYS, Inc.
 - Dash Associates
 - Fluent, Inc.
 - ILOG CPLEX Division
 - Livermore Software Technology Corporation (LSTC)
 - MECALOG SARL
 - Oxford Molecular Group PLC
 - The Numerical Algorithms Group Ltd.(NAG)
- Endorsing software vendors
 - Absoft Corporation
 - Edinburgh Portable Compilers
 - GENIAS Software GmBH
 - Myrias Computer Technologies, Inc.
 - The Portland Group, Inc. (PGI)



OpenMP Release History





Goals of OpenMP

- Standardization
 - Provide a standard among a variety of shared memory architectures/ platforms
- Lean and mean
 - Establish a simple and limited set of directives for programming shared memory machines
 - Significant parallelism can be implemented by using just 3 or 4 directives.
- Ease of Use
 - Provide capability to incrementally parallelize a serial program, unlike message-passing libraries which typically require an all or nothing approach
 - Provide the capability to implement both coarse-grain and fine-grain parallelism
- Portability
 - Supports Fortran (77, 90, and 95), C, and C++
- Public forum for API and membership

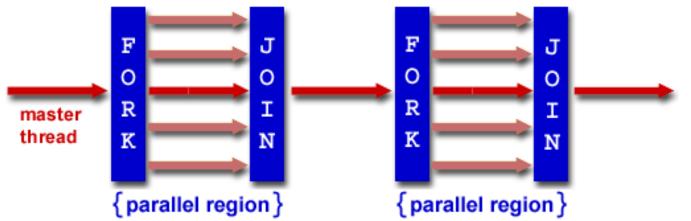


OpenMP Programming Model

- Shared memory, thread-based parallelism
 - OpenMP is based upon the existence of multiple threads in the shared memory programming paradigm.
 - A shared memory process consists of multiple threads.
- Explicit Parallelism
 - OpenMP is an explicit (not automatic) programming model, offering the programmer full control over parallelization.
- OpenMP uses the fork-join model of parallel execution.
- Compiler directive based
 - Most OpenMP parallelism is specified through the use of compiler directives which are imbedded in C/C++ or Fortran source code.
- Nested parallelism support
 - The API provides for the placement of parallel constructs inside of other parallel constructs.
 - Implementations may or may not support this feature.



Fork-Join Model



- All OpenMP programs begin as a single process: the master thread. The master thread executes sequentially until the first parallel region construct is encountered.
- FORK: the master thread then creates a *team* of parallel threads
- The statements in the program that are enclosed by the parallel region construct are then executed in parallel among the team threads.
- JOIN: When the team threads complete the statements in the parallel region construct, they synchronize and terminate, leaving only the master thread



I/O

- OpenMP specifies nothing about parallel I/O. This is particularly important if multiple threads attempt to write/ read from the same file.
- If every thread conducts I/O to a different file, the issues are not as significant.
- It is entirely up to the programmer to ensure that I/O is conducted correctly within the context of a multi-threaded program.



Memory Model

- OpenMP provides a "relaxed-consistency" and "temporary" view of thread memory (in their words). In other words, threads can "cache" their data and are not required to maintain exact consistency with real memory all of the time.
- When it is critical that all threads view a shared variable identically, the programmer is responsible for insuring that the variable is FLUSHed by all threads as needed.
- More on this later...



OpenMP Code Structure - Fortran

PROGRAM HELLO INTEGER VAR1, VAR2, VAR3 Serial code

Beginning of parallel section. Fork a team of threads. Specify variable scoping \$OMP PARALLEL PRIVATE(VAR1, VAR2) SHARED(VAR3)

Parallel section executed by all threads

All threads join master thread and disband \$OMP END PARALLEL

Resume serial code

END

OpenMP Code Structure – C/C++

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

.

.

Beginning of parallel section. Fork a team of threads. Specify variable scoping #pragma omp parallel private(var1, var2) shared(var3)



Fortran Directives Format

<sentinel> <directive-name> [clause...]

- All Fortran OpenMP directives must begin with a sentinel. The accepted sentinels depend on the type of Fortran source (see next two slides).
- A valid OpenMP directive must appear after the sentinel and before any clauses.
- Optional clauses can be in any order and repeated as necessary unless otherwise restricted.
- Example:

!\$OMP PARALLEL DEFAULT(SHARED) PRIVATE(BETA,PI)



Fortran Directives Format (2)

- Fixed Form Source
 - **!\$OMP C\$OMP *\$OMP** are accepted sentinels and must start in column 1
 - All Fortran fixed form rules for line length, white space, continuation and comment columns apply for the entire directive line
 - Initial directive lines must have a space/zero in column 6.
 - Continuation lines must have a non-space/zero in column 6.
- Free Form Source
 - **!\$OMP** is the only accepted sentinel. Can appear in any column, but must be preceded by white space only.
 - All Fortran free form rules for line length, white space, continuation and comment columns apply for the entire directive line
 - Initial directive lines must have a space after the sentinel.
 - Continuation lines must have an ampersand as the last non-blank character in a line. The following line must begin with a sentinel and then the continuation directives.



Fortran Directives Format (3)

- General Rules
 - Comments cannot appear on the same line as a directive
 - Only one directive-name may be specified per directive
 - Fortran compilers that are OpenMP enabled generally include a command line option that instructs the compiler to activate and interpret all OpenMP directives.
 - Several Fortran OpenMP directives come in pairs and have the form shown below. The "end" directive is optional but advised for readability.



C/C++ Directives Format

#pragma omp <directive-name> [clause,...] \n

• Newline at the end required and precedes the structured block that follows the directive

Example:

#pragma omp parallel default(shared) private(beta,pi)



C/C++ Directive Format (2)

- General Rules:
 - Case sensitive
 - Directives follow conventions of the C/C++ standards for compiler directives.
 - Only one directive-name may be specified per directive.
 - Each directive applies to at most one succeeding statement, which must be a structured block.
 - Long directive lines can be "continued" on succeeding lines by escaping the newline character with a backslash ("\") at the end of a directive line.



Directive Scoping

- Static (Lexical) Extent
 - The code textually enclosed between the beginning and the end of a structured block following a directive.
 - The static extent of a directives does not span multiple routines
 or code files
- Orphaned Directive
 - An OpenMP directive that appears independently from another enclosing directive is said to be an *orphaned* directive. It exists outside of another directive's static (lexical) extent.
 - Will span routines and possibly code files
- Dynamic Extent
 - The dynamic extent of a directive includes both its static (lexical) extent and the extents of its orphaned directives.



Directive Scoping Example

PROGRAM TEST

\$0MP PARALLEL

\$0MP DO

DO I=...

CALL SUB1

ENDDO

CALL SUB2

\$OMP END PARALLEL

. . .

STATIC EXTENT

The DO directive occurs within an enclosing parallel region.

SUBROUTINE SUB1

!\$OMP CRITICAL

!\$OMP END CRITICAL END

SUBROUTINE SUB2

!\$OMP SECTIONS

!\$OMP END SECTIONS

END

ORPHANED DIRECTIVES

The CRITICAL and SECTIONS directives occur outside an enclosing parallel region.

DYNAMIC EXTENT

The CRITICAL and SECTIONS directives occur within the dynamic extent of the DO and PARALLEL directives.

PARALLEL Region Construct

- Block of code that will be executed by multiple threads
- Fundamental OpenMP parallel construct
- When a thread reaches a PARALLEL directive, it creates a team of threads and becomes the master of the team. The master is a member of that team and has thread number 0 within that team.
- Starting from the beginning of this parallel region, the code is duplicated and all threads will execute that code.
- There is an implied barrier at the end of a parallel section. Only the master thread continues execution past this point.
- If any thread terminates within a parallel region, all threads in the team will terminate, and the work done up until that point is undefined.



PARALLEL Region Syntax - Fortran

block

!\$OMP END PARALLEL



PARALLEL Region Syntax – C/C++

#pragma omp parallel [clause ...] newline if (scalar_expression) private (list) shared (list) default (shared | none) firstprivate (list) reduction (operator: list) copyin (list) num threads (integer-expression)

structured_block



PARALLEL Region – Number of Threads

- The number of threads in a parallel region is determined by the following factors, in order of precedence:
 - 1. Evaluation of the **IF** clause
 - 2. Setting of the **NUM_THREADS** clause
 - 3. Use of the **omp_set_num_threads()** library function
 - 4. Setting of the **OMP_NUM_THREADS** environment variable
 - 5. Implementation default usually the number of cores on a node.
- Threads are numbered from 0 (master thread) to N-1



Nested Parallel Regions

- Use the omp_get_nested() library function to determine if nested parallel regions are enabled.
- The two methods available for enabling nested parallel regions (if supported) are:
 - The **omp_set_nested()** library routine
 - Setting of the **OMP_NESTED** environment variable to TRUE
- If not supported, a parallel region nested within another parallel region results in the creation of a new team, consisting of one thread, by default.



PARALLEL Region Restrictions

- A parallel region must be a structured block that does not span multiple routines or code files
- It is illegal to branch into or out of a parallel region
- Only a single IF clause is permitted
- Only a single NUM_THREADS clause is permitted



PARALLEL Region Example - Fortran

PROGRAM HELLO

INTEGER NTHREADS, TID, OMP_GET_NUM_THREADS,

+ OMP_GET_THREAD_NUM

C Fork a team of threads with each thread having a private TID variable

!\$OMP PARALLEL PRIVATE(TID)

C Obtain and print thread id

TID = OMP_GET_THREAD_NUM()

PRINT *, 'Hello World from thread = ', TID

C Only master thread does this

IF (TID .EQ. 0) THEN NTHREADS = OMP_GET_NUM_THREADS() PRINT *, 'Number of threads = ', NTHREADS END IF C All threads join master thread and disband !\$OMP END PARALLEL

END



PARALLEL Region Example - C

```
#include <omp.h>
main () {
int nthreads, tid;
/* Fork a team of threads with each thread having a private tid variable */
#pragma omp parallel private(tid)
   /* Obtain and print thread id */
   tid = omp get thread num();
   printf("Hello World from thread = %d\n", tid);
   /* Only master thread does this */
   if (tid == 0) {
     nthreads = omp get num threads();
     printf("Number of threads = %d\n", nthreads);
   /* All threads join master thread and terminate */
}
```



Work-sharing Constructs

- A work-sharing construct divides the execution of the enclosed code region among the members of the team that encounter it.
- Work-sharing constructs do not launch new threads
- There is no implied barrier upon entry to a work-sharing construct, however there is an implied barrier at the end of a work sharing construct.
- A work-sharing construct must be enclosed dynamically within a parallel region in order for the directive to execute in parallel.
- Work-sharing constructs must be encountered by all members of a team or none at all
- Successive work-sharing constructs must be encountered in the same order by all members of a team.



Types of Work-sharing Constructs

• DO / for

- Shares iterations of a loop across the team
- Represents a type of "data parallelism"

SECTIONS

- Breaks work into separate, discrete sections
- Each section is executed by a thread.
- Can be used to implement a type of "functional parallelism"

• SINGLE

• Serializes a section of code



Fortran DO Directive Syntax

!\$OMP DO [clause ...]
SCHEDULE (type [,chunk])
ORDERED
PRIVATE (list)
FIRSTPRIVATE (list)
LASTPRIVATE (list)
SHARED (list)
REDUCTION (operator | intrinsic : list)
COLLAPSE (n)

do_loop

!\$OMP END DO [NOWAIT]



C/C++ for Directive Syntax

#pragma omp for [clause ...] newline schedule (type [,chunk]) ordered private (list) firstprivate (list) lastprivate (list) shared (list) reduction (operator: list) collapse (n) nowait

for loop



SCHEDULE Clause

• Describes how iterations of the loop are divided among the threads in the team. The default schedule is implementation dependent.

STATIC

Loop iterations are divided into pieces of size *chunk* and then statically assigned to threads. If chunk is not specified, the iterations are evenly (if possible) divided contiguously among the threads.

DYNAMIC

Loop iterations are divided into pieces of size *chunk*, and dynamically scheduled among the threads; when a thread finishes one chunk, it is dynamically assigned another. The default chunk size is 1.

GUIDED

For a chunk size of 1, the size of each chunk is proportional to the number of unassigned iterations divided by the number of threads, decreasing to 1. For a chunk size with value k (greater than 1), the size of each chunk is determined in the same way with the restriction that the chunks do not contain fewer than k iterations (except for the last chunk to be assigned, which may have fewer than k iterations). The default chunk size is 1.

RUNTIME

The scheduling decision is deferred until runtime by the environment variable OMP_SCHEDULE. It is illegal to specify a chunk size for this clause.

AUTO

The scheduling decision is delegated to the compiler and/or runtime system.



Other DO/for Clauses

- **NO WAIT / nowait**: If specified, then threads do not synchronize at the end of the parallel loop.
- **ORDERED**: Specifies that the iterations of the loop must be executed as they would be in a serial program.
- COLLAPSE: Specifies how many loops in a nested loop should be collapsed into one large iteration space and divided according to the schedule clause. The sequential execution of the iterations in all associated loops determines the order of the iterations in the collapsed iteration space.
- Other clauses are described in detail later.



Fortran DO Directive Example

```
PROGRAM VEC ADD DO
      INTEGER N, CHUNKSIZE, CHUNK, I
      PARAMETER (N=1000)
      PARAMETER (CHUNKSIZE=100)
      REAL A(N), B(N), C(N)
      Some initializations
DOI = 1, N
         A(I) = I * 1.0
         B(I) = A(I)
      ENDDO
      CHUNK = CHUNKSIZE
!$OMP PARALLEL SHARED(A,B,C,CHUNK) PRIVATE(I)
!$OMP DO SCHEDULE(DYNAMIC,CHUNK)
      DOI = 1, N
         C(I) = A(I) + B(I)
      ENDDO
!$OMP END DO NOWAIT
!$OMP END PARALLEL
      END
```



C for Directive Example

#include <omp.h>
#define CHUNKSIZE 100
#define N 1000

```
main () {
int i, chunk;
float a[N], b[N], c[N];
/* Some initializations */
for (i=0; i < N; i++)
    a[i] = b[i] = i * 1.0;
chunk = CHUNKSIZE;
#pragma omp parallel shared(a,b,c,chunk) private(i)
    {
    #pragma omp for schedule(dynamic,chunk) nowait
    for (i=0; i < N; i++)
        c[i] = a[i] + b[i];
    } /* end of parallel section */
}</pre>
```



SECTIONS Directive – Fortran Syntax

!\$OMP SECTIONS [clause ...]

PRIVATE (list) FIRSTPRIVATE (list) LASTPRIVATE (list) REDUCTION (operator | intrinsic : list)

!\$OMP SECTION

block

!\$OMP SECTION

block

!\$OMP END SECTIONS [NOWAIT]



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SECTIONS Directive – C/C++ Syntax

#pragma omp sections [clause ...] newline
 private (list)
 firstprivate (list)
 lastprivate (list)
 reduction (operator: list)
 nowait

#pragma omp section *newline*

structured_block

#pragma omp section newline

structured_block

}



SECTIONS Directive Example - Fortran

```
PROGRAM VEC_ADD_SECTIONS
     INTEGER N. I
     PARAMETER (N=1000)
     REAL A(N), B(N), C(N), D(N)
Some initializations
     DOI = 1, N
        A(I) = I * 1.5
         B(I) = I + 22.35
      ENDDO
!$OMP PARALLEL SHARED(A,B,C,D), PRIVATE(I)
!$OMP SECTIONS
!$OMP SECTION
      DO I = 1, N
         C(I) = A(I) + B(I)
      ENDDO
!$OMP SECTION
      DO I = 1, N
         D(I) = A(I) * B(I)
      ENDDO
!$OMP END SECTIONS NOWAIT
!$OMP END PARALLEL
      END
```



SECTIONS Directive Example - C

```
#include <omp.h>
#define N 1000
main () {
int i:
float a[N], b[N], c[N], d[N];
/* Some initializations */
for (i=0; i < N; i++) {
 a[i] = i * 1.5;
 b[i] = i + 22.35;
#pragma omp parallel shared(a,b,c,d) private(i)
 #pragma omp sections nowait
    #pragma omp section
    for (i=0; i < N; i++)
      c[i] = a[i] + b[i];
    #pragma omp section
    for (i=0; i < N; i++)
      d[i] = a[i] * b[i];
    } /* end of sections */
  } /* end of parallel section */
}
```



Questions for Thought

- What happens if the number of threads and the number of SECTIONs are different? What if there are more threads than SECTIONs? fewer threads than SECTIONs?
- Which thread executes which SECTION?



Synchronization Constructs

 Motivation: Consider a simple example where two threads on two different processors are both trying to increment a variable x at the same time (assume x is initially 0).

THREAD 1:	THREAD 2:	
increment(x)	increment(x)	
{	{	
x = x + 1;	x = x + 1;	
}	}	
THREAD 1:	THREAD 2:	
10 LOAD A, (x address)	10 LOAD A, (x address)	
20 ADD A, 1	20 ADD A, 1	
30 STORE A, (x address)	30 STORE A, (x address)	

- The incrementation of x must be synchronized between the two threads to insure that the correct result is produced.
- OpenMP provides a variety of synchronization sonstructs that control how the execution of each thread proceeds relative to other team threads.

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CRITICAL Directive

• The CRITICAL directive specifies a region of code that must be executed by only one thread at a time.

Fortran	!\$OMP CRITICAL <i>[name]</i> <i>block</i> !\$OMP END CRITICAL
C/C++	#pragma omp critical <i>[name] newline</i> <i>structured_block</i>

- If a thread is currently executing inside a CRITICAL region and another thread reaches that CRITICAL region and attempts to execute it, it will block until the first thread exits that CRITICAL region.
- The optional name enables multiple different CRITICAL regions to exist.
 - Names act as global identifiers. Different CRITICAL regions with the same name are treated as the same region.
 - All unnamed CRITICAL regions are treated as the same region.



CRITICAL Directive Example - Fortran

PROGRAM CRITICAL INTEGER X X = 0ISOMP PARALLEL SHARED(X) SOMP CRITICAL X = X + 1ISOMP END CRITICAL ISOMP END PARALLEL END



CRITICAL Directive Example - C

```
#include <omp.h>
main()
{
    int x;
    x = 0;
#pragma omp parallel shared(x)
    {
        #pragma omp critical
        x = x + 1;
        } /* end of parallel region */
}
```



Other Synchronization Constructs

- MASTER directive
 - Specifies a region that is to be executed only by the master thread of the team
 - All other threads on the team skip this section of code. Fortran: !\$OMP MASTER

block

!\$OMP END MASTER

C/C++: #pragma omp master newline

structured block

- BARRIER directive
 - Synchronizes all threads in the team
 - When a BARRIER directive is reached, a thread will wait at that point until all other threads have reached that barrier. All threads then resume executing in parallel the code that follows the barrier.

Fortran: **!**\$OMP BARRIER

C/C++: #pragma omp barrier newline

- ATOMIC directive
 - specifies that a specific memory location must be updated atomically, rather than letting multiple threads attempt to write to it
 - Appies only to a single, immediately following statement Fortran: !\$OMP ATOMIC

C/C++: #pragma omp atomic newline



FLUSH Directive

- Identifies a synchronization point at which the implementation must provide a consistent view of memory
- Thread-visible variables are written back to memory at this point.
- Necessary to instruct the compiler that a variable must be written to/read from the memory system, i.e. that a variable cannot be kept in a local CPU register
 - Keeping a variable in a register in a loop is very common when producing efficient machine language code for a loop.



FLUSH Directive (2)

Fortran: **!**\$OMP FLUSH (*list*)

C/C++: #pragma omp flush (list) newline

- The optional list contains a list of named variables that will be flushed in order to avoid flushing all variables. For pointers in the list, the pointer itself is flushed, not the object to which it points.
- Implementations must ensure any prior modifications to thread-visible variables are visible to all threads after this point; i.e., compilers must restore values from registers to memory, hardware might need to flush write buffers, etc.
- The FLUSH directive is implied for the directives shown in the table below. The directive is not implied if a NOWAIT clause is present.



FLUSH Directive (3)

 The FLUSH directive is implied for the directives shown in the table below. The directive is not implied if a NOWAIT clause is present.

Fortran	C/C++
BARRIER	barrier
END PARALLEL	parallel – upon entry and exit
CRITICAL and END CRITICAL	critical – upon entry and exit
END DO	ordered – upon entry and exit
END SECTIONS	for – upon exit
END SINGLE	sections – upon exit
ORDERED and END ORDERED	single – upon exit



ORDERED Directive

- Specifies that iterations of the enclosed loop will be executed in the same order as if they were executed on a serial processor
- Threads will need to wait before executing their chunk of iterations if previous iterations haven't completed yet.
- Used within a DO / for loop with an ORDERED clause
- The ORDERED directive provides a way to "fine tune" where ordering is to be applied within a loop. Otherwise, it is not required.
- An ORDERED directive can only appear in the dynamic extent of the following directives:
 - DO or PARALLEL DO (Fortran)
 - for or parallel for (C/C++)
- Only one thread is allowed in an ordered section at any time
- It is illegal to branch into or out of an ORDERED block.
- An iteration of a loop must not execute the same ORDERED directive more than once, and it must not execute more than one ORDERED directive.
- A loop that contains an ORDERED directive must be a loop with an ORDERED clause.



ORDERED Directive Syntax

Fortran	<pre>!\$OMP DO ORDERED [clauses] [loop region] !\$OMP ORDERED [block] !\$OMP END ORDERED [end of loop region] !\$OMP END DO</pre>
C/C++	<pre>#pragma omp for ordered [clauses] (loop region) #pragma omp ordered newline structured_block (endo of loop region)</pre>



ORDERED Example - C

```
#include <stdio.h>
#include <omp.h>
static float a[1000], b[1000], c[1000];
                                                        }
void test(int first, int last)
{
  int i;
  #pragma omp for schedule(static) ordered
  for (i = first; i <= last; ++i) {
     // Do something here.
     if (i % 2)
    {
        #pragma omp ordered
        printf("test() iteration %d\n", i);
     }
                                                          }
                                                        }
}
```

```
void test2(int iter)
```

```
#pragma omp ordered
printf("test2() iteration %d\n", iter);
}
int main()
```

```
int i;
#pragma omp parallel
{
   test(1, 8);
   #pragma omp for ordered
   for (i = 0 ; i < 5 ; i++)
      test2(i);</pre>
```



Data Scope Attribute Clauses

- Also called data sharing attribute clauses
- Because OpenMP is based upon the shared memory programming model, most variables are shared by default.
- Global variables include:
 - Fortran: COMMON blocks, SAVE variables, MODULE variables
 - C: File scope variables, static
- Private variables include:
 - Loop index variables
 - Stack variables in subroutines called from parallel regions
 - Fortran: Automatic variables within a statement block
- Clauses used to explicitly define how variables should be scoped include:
 - PRIVATE
 - FIRSTPRIVATE
 - LASTPRIVATE
 - SHARED
 - DEFAULT
 - REDUCTION
 - COPYIN



Data Scope Attribute Clauses (2)

- Used in conjunction with several directives (PARALLEL, DO/for, and SECTIONS) to control the scoping of enclosed variables.
- Provide the ability to control the data environment during execution of parallel constructs
 - Define how and which data variables in the serial section of the program are transferred to the parallel sections of the program (and back)
 - Define which variables will be visible to all threads in the parallel sections and which variables will be privately allocated to all threads
- Effective only within their lexical/static extent



PRIVATE and SHARED Clauses

- PRIVATE Clause
 - · Declares variables in its list to be private to each thread
 - A new object of the same type is declared once for each thread in the team.
 - All references to the original object are replaced with references to the new object.
 - Variables declared PRIVATE should be assumed to be uninitialized for each thread.
- SHARED Clause
 - Declares variables in its list to be shared among all threads in the team
 - A shared variable exists in only one memory location and all threads can read or write to that address
 - It is the programmer's responsibility to ensure that multiple threads properly access SHARED variables (such as via CRITICAL sections)



FIRSTPRIVATE and LASTPRIVATE Clauses

- FIRSTPRIVATE Clause
 - Combines the behavior of the PRIVATE clause with automatic initialization of the variables in its list
 - Listed variables are initialized according to the value of their original objects prior to entry into the parallel or work-sharing construct.
- LASTPRIVATE Clause
 - Combines the behavior of the PRIVATE clause with a copy from the last loop iteration or section to the original variable object
 - The value copied back into the original variable object is obtained from the last (sequentially) iteration or section of the enclosing construct.
 - For example, the team member that executes the final iteration for a DO section, or the team member that executes the last SECTION of a SECTIONS context, performs the copy with its own values.



PRIVATE Variables Example

```
main()
ł
   int A = 10;
   int B, C;
   int n = 20;
   #pragma omp parallel
       #pragma omp for private(i) firstprivate(A) lastprivate(B)
       for (int i=0; i<n; i++)
       {
           B = A + i; /* A undefined unless declared firstprivate */
           ....
       C = B; /* B undefined unless declared lastprivate */
   } /* end of parallel region */
}
                                            ICL
```

DEFAULT Clause

- Allows the user to specify a default scope for all variables in the lexical extent of any parallel region
- Specific variables can be exempted from the default using the PRIVATE, SHARED, FIRSTPRIVATE, LASTPRIVATE, and REDUCTION clauses.
- The C/C++ OpenMP specification does not include private or firstprivate as a possible default. However, actual implementations may provide this option.
- Using NONE as a default requires that the programmer explicitly scope all variables.

Fortran	DEFAULT (PRIVATE FIRSTPRIVATE SHARED NONE)
C/C++	default (shared none)



REDUCTION Clause

- Performs a reduction on the variables that appear in its list.
- A private copy for each list variable is created for each thread. At the end of the reduction, the reduction operator is applied to all private copies of the shared variable, and the final result is written to the global shared variable.
- Syntax:

Fortran	REDUCTION (operator intrinsic: list)
C/C++	reduction <i>(operator: list</i>)



REDUCTION Clause Example - Fortran

```
PROGRAM DOT PRODUCT
      INTEGER N, CHUNKSIZE, CHUNK, I
      PARAMETER (N=100)
      PARAMETER (CHUNKSIZE=10)
      REALA(N), B(N), RESULT
      Some initializations
DO | = 1. N
         A(I) = I * 1.0
         B(I) = I * 2.0
       ENDDO
       RESULT= 0.0
       CHUNK = CHUNKSIZE
!SOMP PARALLEL DO
!$OMP& DEFAULT(SHARED) PRIVATE(I)
!$OMP& SCHEDULE(STATIC,CHUNK)
!$OMP& REDUCTION(+:RESULT)
       DO I = 1, N
         RESULT = RESULT + (A(I) * B(I))
       ENDDO
!$OMP
       END PARALLEL DO NOWAIT
       PRINT *, 'Final Result= ', RESULT
       END
```



REDUCTION Clause Example - C

#include <omp.h>

```
main () {
int i, n, chunk;
float a[100], b[100], result;
/* Some initializations */
n = 100:
chunk = 10;
result = 0.0:
for (i=0; i < n; i++) {
  a[i] = i * 1.0;
  b[i] = i * 2.0;
}
#pragma omp parallel for default(shared) private(i) \
  schedule(static,chunk) reduction(+:result)
  for (i=0; i < n; i++)
    result = result + (a[i] * b[i]);
printf("Final result= %f\n",result);
}
```



THREADPRIVATE Directive

- Used to make global file scope variables (C/C++) or common blocks (Fortran) local and persistent to a thread through the execution of multiple parallel regions
- Must appear after the declaration of listed variables/common blocks
- Each thread then gets its own copy of the variable/common block, so data written by one thread is not visible to other threads.
- On first entry to a parallel region, data in THREADPRIVATE variables and common blocks should be assumed undefined, unless a COPYIN clause is specified in the PARALLEL directive
- THREADPRIVATE variables differ from PRIVATE variables because they are able to persist between different parallel sections of a code.
- Data in THREADPRIVATE objects is guaranteed to persist only if the dynamic threads mechanism is "turned off" and the number of threads in different parallel regions remains constant. The default setting of dynamic threads is undefined.

66	
C/C++	#pragma omp threadprivate (list)
Fortran	!\$OMP THREADPRIVATE (/cb/,)

THREADPRIVATE Example - Fortran

PROGRAM THREADPRIV

INTEGER A, B, I, TID, OMP_GET_THREAD_NUM

REAL*4 X

COMMON /C1/ A

!\$OMP THREADPRIVATE(/C1/, X)

C Explicitly turn off dynamic threads

CALL OMP_SET_DYNAMIC(.FALSE.)

PRINT *, '1st Parallel Region:'

!\$OMP PARALLEL PRIVATE(B, TID)

TID = OMP_GET_THREAD_NUM()

- A = TID
- B = TID

```
X = 1.1 * TID + 1.0
```

PRINT *, 'Thread', TID, ': A, B, X=', A, B, X

!\$OMP END PARALLEL

PRINT *, 'Master thread doing serial work here'

67

PRINT *, '2nd Parallel Region: '

!\$OMP PARALLEL PRIVATE(TID)

TID = OMP_GET_THREAD_NUM()

PRINT *, 'Thread', TID,': A,B,X=',A,B,X

!\$OMP END PARALLEL

END



THREADPRIVATE Example - C

```
int a, b, i, tid;
float x;
#pragma omp threadprivate(a, x)
main () {
/* Explicitly turn off dynamic threads */
  omp set dynamic(0);
  printf("1st Parallel Region:\n");
  #pragma omp parallel private(b,tid)
  tid = omp get thread num();
  a = tid;
   b = tid:
  x = 1.1 * tid + 1.0;
   printf("Thread %d: a,b,x= %d %d %f\n",tid,a,b,x);
 } /* end of parallel region */
  printf("Master thread doing serial work here\n");
  printf("2nd Parallel Region:\n");
  #pragma omp parallel private(tid)
  {
   tid = omp get thread num();
   printf("Thread %d: a,b,x= %d %d %f\n",tid,a,b,x);
 } /* end of parallel section */
                              68
}
```

#include <omp.h>



COPYIN Clause

- Provides a means for assigning the same value to THREADPRIVATE variables for all threads in the team
- List contains the names of variables to copy. In Fortran, the list can contain both the names of common blocks and named variables.
- The master thread variable is used as the copy source. The team threads are initialized with its value upon entry into the parallel construct.

Fortran	COPYIN <i>(list)</i>
C/C++	copyin <i>(list)</i>



Run-time Library Routines

- The OpenMP standard defines an API for library calls that perform a variety of functions:
 - Query the number of threads/processors, set number of threads to use
 - General purpose locking routines (semaphores)
 - Portable wall clock timing routines
 - Set execution environment functions: nested parallelism, dynamic adjustment of threads.
 - For C/C++, it may be necessary to specify the include file "omp.h".
 - Note: Your implementation may or may not support nested parallelism and/or dynamic threads. If nested parallelism is supported, it is often only nominal, in that a nested parallel region may only have one thread.



OMP_SET_NUM_THREADS()

- Sets the number of threads that will be used in the next parallel region
- Must be a positive integer
- Can only be called from serial portion of the code
- Has precedence over the OMP_NUM_THREADS environment variable

Fortran	SUBROUTINE OMP_SET_NUM_THREADS(scalar_integer_expression)
C/C++	#include <omp.h> void omp_set_num_threads(int num_threads)</omp.h>



OMP_GET_NUM_THREADS() and OMP_GET_THREAD_NUM()

- OMP_GET_NUM_THREADS()
 - Returns the number of threads that are currently in the team executing the parallel region from which it is called

Fortran	INTEGER FUNCTION OMP_GET_NUM_THREADS()
C/C++	#include <omp.h></omp.h>
	int omp_get_num_threads(void)

• OMP_GET_THREAD_NUM()

- Returns the thread number of the thread, within the team, making this call. This number will be between 0 and OMP_GET_NUM_THREADS-1. The master thread of the team is thread 0.
- If called from a nested parallel region, or a serial region, this function will return 0.



OMP_GET_MAX_THREADS()

- Returns the maximum value that can be returned by a call to the OMP_GET_NUM_THREADS function
- Generally reflects the number of threads as set by the OMP_NUM_THREADS environment variable or the OMP_SET_NUM_THREADS() library routine.
- May be called from both serial and parallel regions of code



OMP_GET_THREAD_LIMIT() and OMP_GET_NUM_PROCS()

- OMP_GET_THREAD_LIMIT()
 - New with OpenMP 3.0
 - Returns the maximum number of OpenMP threads available to a program
- OMP_GET_NUM_PROCS()
 - Returns the number of processors that are available to the program



OMP_IN_PARALLEL()

- May be called to determine if the section of code currently executing is parallel or not
- For Fortran, this function returns .TRUE. if it is called from the dynamic extent of a region executing in parallel, and .FALSE. otherwise. For C/C++, it will return a nonzero integer if parallel, and zero otherwise.

Fortran	LOGICAL FUNCTION OMP_IN_PARALLEL()
C/C++	<pre>#include <omp.h> int omp_in_parallel(void)</omp.h></pre>



OMP_SET_DYNAMIC()

- Enables or disables dynamic adjustment (by the run time system) of the number of threads available for execution of parallel regions
- For Fortran, if called with .TRUE. then the number of threads available for subsequent parallel regions can be adjusted automatically by the run-time environment. If called with .FALSE., dynamic adjustment is disabled.
- For C/C++, if dynamic_threads evaluates to non-zero, then the mechanism is enabled, otherwise it is disabled.
- The OMP_SET_DYNAMIC subroutine has precedence over the OMP_DYNAMIC environment variable.
- The default setting is implementation dependent.
- Must be called from a serial section of the program

Fortran	SUBROUTINE OMP_SET_DYNAMIC(scalar_logical_expression)
C/C++	#include <omp.h></omp.h>
	<pre>void omp_set_dynamic(int dynamic_threads)</pre>



OMP_GET_DYNAMIC()

- Used to determine if dynamic thread adjustment is enabled or not
- For Fortran, this function returns .TRUE. if dynamic thread adjustment is enabled, and .FALSE. otherwise.
- For C/C++, non-zero will be returned if dynamic thread adjustment is enabled, and zero otherwise.

Fortran	LOGICAL FUNCTION OMP_GET_DYNAMIC()
C/C++	#include <omp.h></omp.h>
	int omp_get_dynamic(void)



OMP_SET_NESTED()

- Used to enable or disable nested parallelism
- For Fortran, calling this function with .FALSE. will disable nested parallelism, and calling with .TRUE. will enable it.
- For C/C++, if nested evaluates to non-zero, nested parallelism is enabled; otherwise it is disabled.
- The default is for nested parallelism to be disabled.
- This call has precedence over the OMP_NESTED environment variable.

Fortran	SUBROUTINE OMP_SET_NESTED(scalar_logical_expression)
C/C++	<pre>#include <omp.h> void omp_set_nested(int nested)</omp.h></pre>



OMP_GET_NESTED()

- Used to determine if nested parallelism is enabled or not
- For Fortran, this function returns .TRUE. if nested parallelism is enabled, and .FALSE. otherwise.
- For C/C++, non-zero will be returned if nested parallelism is enabled, and zero otherwise.

Fortran	LOGICAL FUNCTION OMP_GET_NESTED
C/C++	Include <omp.h></omp.h>
	Int omp_get_nested(void)



Locking Routines

- OMP_INIT_LOCK()
 - Initializes a lock associated with the lock variable
 - The initial state is unlocked.
 - For Fortran, var must be an integer large enough to hold an address, such as INTEGER*8 on 64-bit systems.

Fortran	SUBROUTINE OMP_INIT_LOCK(var)
C/C++	#include <omp.h> void omp_init_lock(omp_lock_t *lock)</omp.h>

OMP_DESTROY_LOCK()

• Disassociates the given lock variable from any locks

Fortran	SUBROUTINE OMP_DESTROY_LOCK(var)
C/C++	#include <omp.h></omp.h>
	void omp_destroy_lock(omp_lock_t *lock)



Locking Routines (2)

- OMP_SET_LOCK()
 - Forces the executing thread to wait until the specified lock is available

Fortran	SUBROUTINE OMP_SET_LOCK(var)
C/C++	#include <omp.h></omp.h>
	void omp_set_lock(omp_lock_t *lock)

- OMP_UNSET_LOCK()
 - Releases the lock from the executing thread
- OMP_TEST_LOCK()
 - Attempts to set a lock, but does not block if the lock is unavailable



OMP_GET_WTIME()

- Provides a portable wall clock timing routine
- Returns a double-precision floating point value equal to the number of elapsed seconds since some point in the past
- Usually used in pairs with the value of the first call subtracted from the value of the second call to obtain the elapsed time for a block of code
- Designed to be per thread times, and therefore may not be globally consistent across all threads in a team

Fortran	DOUBLE PRECISION FUNCTION OMP_GET_WTIME()
C/C++	#include <omp.h></omp.h>
	double omp_get_wtime(void)



OpenMP Environment Variables

• OMP_SCHEDULE

- Applies only to DO, PARALLEL DO (Fortran) and for, parallel for (C/C++) directives which have their schedule clause set to RUNTIME. The value of this variable determines how iterations of the loop are scheduled on processors.
- For example:

setenv OMP_SCHEDULE "guided, 4" setenv OMP_SCHEDULE "dynamic"

• OMP_NUM_THREADS

- · Sets the number of threads to use during execution
- For example:

setenv OMP_NUM_THREADS 8



OpenMP Environment Variables (2)

• OMP_DYNAMIC

- Enables or disables dynamic adjustment of the number of threads available for execution of parallel regions
- Valid values are TRUE or FALSE.
- For example:

setenv OMP_DYNAMIC TRUE

- OMP_NESTED
 - Enables or disables nested parallelism
 - Valid values are TRUE or FALSE.
 - For example:

setenv OMP_NESTED TRUE



Thread Stack Size

- The OpenMP standard does not specify how much stack space a thread should have. Consequently, implementations will differ in the default thread stack size.
- Default thread stack size can be easy to exhaust. It can also be non-portable between compilers.
- Threads that exceed their stack allocation may or may not seg fault. An application may continue to run while data is being corrupted.
- Statically linked codes may be subject to further stack restrictions.
- A user's login shell may also restrict stack size.
- May need to increase thread stack size. How to do this is system-dependent.



Performance Expectations

- One might expect to get an N times speedup when running a program parallelized using OpenMP on an N processor/core platform. However, this is seldom the case due to the following reasons:
 - A large portion of the program may not be parallelized by OpenMP, which means that the theoretical upper limit of speedup is limited according to Amdahl's law.
 - N processors in a SMP may have N times the computation power, but the memory bandwidth usually does not scale up N times. Quite often, the original memory path is shared by multiple processors and performance degradation may be observed when they compete for the shared memory bandwidth.
 - Many other common problems affecting the final speedup in parallel computing also apply to OpenMP, such as load balancing and synchronization overhead.

