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Session 5: Parallel Programming with OpenMP

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Agenda

- 10:00 11:00 OpenMP fundamentals, parallel regions
- 11:00 11:30 Worksharing constructs
- 11:30 12:00 Break
- 12:00 12:15 Synchronization mechanisms in OpenMP
- 12:15 13:00 Practical: heat diffusion
- 13:00 14:00 Lunch
- 14:00 14:30 Tasking in OpenMP
- 14:30 15:30 Programming using a hybrid MPI/OpenMP approach
- 15:30 16:00 Break
- 16:00 17:00 Practical: heat diffusion

Part I

[OpenMP fundamentals, parallel regions](#page-3-0)

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Outline

- **[OpenMP Overview](#page-5-0)**
- **•** [The OpenMP model](#page-10-0)
- [Writing OpenMP programs](#page-14-0)
- **[Creating Threads](#page-19-0)**
- [Data-sharing attributes](#page-28-0) \bullet

Outline

- **[OpenMP Overview](#page-5-0)**
- [The OpenMP model](#page-10-0) \bullet
- [Writing OpenMP programs](#page-14-0) \bullet
- **[Creating Threads](#page-19-0)**
- [Data-sharing attributes](#page-28-0) \bullet

What is OpenMP?

- \bullet It's an API extension to the C, C₊₊ and Fortran languages to write parallel programs for shared memory machines
	- Current version is 3.0 (May 2008)
	- Supported by most compiler vendors
		- Intel,IBM,PGI,Sun,Cray,Fujitsu,HP,GCC,...
- Maintained by the Architecture Review Board (ARB), a consortium of industry and academia

http://www.openmp.org

A bit of history

Advantages of OpenMP

- Mature standard and implementations
	- Standardizes practice of the last 20 years
- Good performance and scalability
- Portable across architectures
- Incremental parallelization
- Maintains sequential version
- (mostly) High level language
	- Some people may say a medium level language :-)
- Supports both task and data parallelism
- Communication is implicit

 \Box

Disadvantages of OpenMP

- Communication is implicit
- Flat memory model
- **•** Incremental parallelization creates false sense of glory/failure
- No support for accelerators
- No error recovery capabilities
- Difficult to compose
- Lacks high-level algorithms and structures
- **o** Does not run on clusters

Outline

- **[OpenMP Overview](#page-5-0)**
- [The OpenMP model](#page-10-0) \bullet
- [Writing OpenMP programs](#page-14-0) \bullet
- **[Creating Threads](#page-19-0)**
- [Data-sharing attributes](#page-28-0) \bullet

OpenMP at a glance

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Execution model

Fork-join model

- OpenMP uses a fork-join model
	- The master thread spawns a team of threads that joins at the end of the parallel region
	- Threads in the same team can collaborate to do work

 \Box

Memory model

• OpenMP defines a relaxed memory model

- Threads can see different values for the same variable
- Memory consistency is only guaranteed at specific points
- Luckily, the default points are usually enough
- Variables can be shared or private to each thread

Outline

- **[OpenMP Overview](#page-5-0)**
- [The OpenMP model](#page-10-0) \bullet
- [Writing OpenMP programs](#page-14-0) \bullet
- **[Creating Threads](#page-19-0)**
- [Data-sharing attributes](#page-28-0) \bullet

OpenMP directives syntax

In Fortran

Through a specially formatted comment:

sentinel construct [clauses]

where sentinel is one of:

- **.** I SOMP or CSOMP or $*$ SOMP in fixed format
- **.** I SOMP in free format

In C/C_{++}

Through a compiler directive:

```
#pragma omp construct [clauses]
```
• OpenMP syntax is ignored if the compiler does not recognize OpenMP

OpenMP directives syntax

In Fortran

Through a specially formatted comment:

sentinel construct [clauses]

where sentinel is one of:

- **.** I SOMP or CSOMP or $*$ SOMP in fixed format
- **.** I SOMP in free format

$In \overline{C/C_{++}}$

Through a compiler directive:

#pragma omp construct [clauses]

• OpenMP syntax is ignored if the compiler does not recognize

We'll be using C/C++ syntax through this tutorial

Headers/Macros

$C/C++$ only

o omp.h contains the API prototypes and data types definitions

- The OPENMP is defined by OpenMP enabled compiler
	- Allows conditional compilation of OpenMP

Fortran only

• The omp lib module contains the subroutine and function definitions

Structured Block

Definition

Most directives apply to a structured block:

- **Block of one or more statements**
- One entry point, one exit point
	- No branching in or out allowed
- Terminating the program is allowed

Outline

- **[OpenMP Overview](#page-5-0)**
- [The OpenMP model](#page-10-0) \bullet
- [Writing OpenMP programs](#page-14-0) \bullet
- [Creating Threads](#page-19-0) \bullet
- [Data-sharing attributes](#page-28-0) \bullet

The parallel construct

Directive

#pragma omp parallel [clau se s] structured block

where clauses can be:

- **num_threads(expression)**
- **if(expression)**

The parallel construct

Specifying the number of threads

- The number of threads is controlled by an internal control variable (**ICV**) called **nthreads-var**.
- When a parallel construct is found a parallel region with a maximum of **nthreads-var** is created
	- Parallel constructs can be nested creating nested parallelism
- **The nthreads-var** can be modified through
	- the **omp_set_num_threads** API called
	- **the OMP_NUM_THREADS** environment variable
- Additionally, the **num_threads** clause causes the implementation to ignore the ICV and use the value of the clause for that region.

The parallel construct

Avoiding parallel regions

- Sometimes we only want to run in parallel under certain conditions
	- E.g., enough input data, not running already in parallel, ...
- The **if** clause allows to specify an *expression*. When evaluates to false the **parallel** construct will only use 1 thread
	- Note that still creates a new team and data environment

Example

```
void main () {
  #pragma omp parallel
     . . .
  omp_set_num_threads ( 2 ) ;
  #pragma omp parallel
     . . .
  #pragma omp parallel num_threads( random ()%4+1) if( 0 )
     . . .
}
```


Example

Example

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Example

API calls

Other useful routines

int **omp_get_num_threads**() Returns the number of threads in the current team

int **omp** get thread num() Returns the id of the thread in the current team

int **omp** get num procs() Returns the number of processors in the machine

int omp qet max threads() Returns the maximum number of threads that will be used in the next parallel region double **omp get wtime**() Returns the number of seconds since an arbitrary point in the past

Outline

- **[OpenMP Overview](#page-5-0)**
- [The OpenMP model](#page-10-0) \bullet
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- **[Creating Threads](#page-19-0)**
- [Data-sharing attributes](#page-28-0) \bullet

Data environment

A number of clauses are related to building the data environment that the construct will use when executing.

- **shared**
- **private**
- **firstprivate**
- **default**
- **threadprivate**

 \Box

Shared

When a variable is marked as **shared**, the variable inside the construct is the same as the one outside the construct.

- In a parallel construct this means all threads see the same variable
	- but not necessarily the same value
- Usually need some kind of synchronization to update them correctly
	- OpenMP has consistency points at synchronizations

Data-sharing attributes

Example

```
int x=1;
#pragma omp parallel shared( x ) num_threads( 2 )
{
   X++;print(f("d\nu", x);}
print(f("d\nu", x);
```


Example

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Private

When a variable is marked as **private**, the variable inside the construct is a new variable of the same type with an undefined value.

- **•** In a parallel construct this means all threads have a different variable
- Can be accessed without any kind of synchronization

Data-sharing attributes

Example

```
int x=1;
#pragma omp parallel private( x ) num_threads( 2 )
{
   X++;print(f("d\nu", x);}
print(f("d\nu", x);
```


Example


```
int x=1;
#pragma omp parallel private( x ) num_threads( 2 )
{
   X + +:
    print(f("d\nu", x);
}
print f("d\nu", x); \longleftarrow Prints 1
```


Firstprivate

When a variable is marked as **firstprivate**, the variable inside the construct is a new variable of the same type but it is initialized to the original variable value.

- **•** In a parallel construct this means all threads have a different variable with the same initial value
- Can be accessed without any kind of synchronization

Data-sharing attributes

```
int x=1;
#pragma omp parallel firstprivate( x ) num_threads( 2 )
{
   X++;print(f("d\nu", x);}
print(f("d\nu", x);
```


Example

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What is the default?

- Static/global storage is **shared**
- **•** Heap-allocated storage is **shared**
- Stack-allocated storage inside the construct is **private**
- **o** Others
	- **If there is a default clause, what the clause says**
		- **none** means that the compiler will issue an error if the attribute is not explicitly set by the programmer
	- Otherwise, depends on the construct
		- For the **parallel** region the default is **shared**


```
i n t x , y ;
#pragma omp parallel private( y )
{
    x =V =#pragma omp parallel private( x )
     {
         x =V =}
}
```



```
i n t x , y ;
#pragma omp parallel private( y )
{
    x =y =#pragma omp parallel private( x )
     {
         x = xy =}
}
                     x is private
                     y is shared
```


Threadprivate storage

The threadprivate construct

```
#pragma omp thread private (var−list)
```
- Can be applied to:
	- Global variables
	- **•** Static variables
	- **Class-static members**
- Allows to create a per-thread copy of "global" variables.
- **threadprivate** storage persist across **parallel** regions if the number of threads is the same

 \Box

Threaprivate storage

```
char∗ foo ( )
{
  static char buffer [BUF SIZE];
  #pragma omp threadprivate (buffer)
  . . .
  return buffer;
}
```


Threaprivate storage

Threaprivate storage

Part II

[Worksharing constructs](#page-49-0)

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[The worksharing concept](#page-51-0)

• [Loop worksharing](#page-53-0)

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Outline

[The worksharing concept](#page-51-0)

• [Loop worksharing](#page-53-0)

Worksharings

Worksharing constructs divide the execution of a code region among the threads of a team

- Threads cooperate to do some work
- Better way to split work than using thread-ids
- Lower overhead than using **tasks**
	- But, less flexible

In OpenMP, there are four worksharing constructs:

- **•** single
- **o** loop worksharing

section⇔ (We'll see them later o workshare

Restriction: worksharings cannot be nested

 $(1 - 1)$

Outline

• [The worksharing concept](#page-51-0)

• [Loop worksharing](#page-53-0)

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Loop parallelism

The for construct

```
#pragma omp for [ clau se s ]
   for ( init -expr ; test -expr ; inc -expr )
```
where clauses can be:

- **•** private
- **•** firstprivate
- **lastprivate(variable-list)**
- **reduction(operator:variable-list)**
- **schedule(schedule-kind)**
- **nowait**
- **collapse(***n***)**
- o ordered We'll see it later

How it works?

The iterations of the loop(s) associated to the construct are divided among the threads of the team.

- Loop iterations must be independent
- Loops must follow a form that allows to compute the number of iterations
- Valid data types for inductions variables are: integer types, pointers and random access iterators (in C++)
	- The induction variable(s) are automatically privatized
- The default data-sharing attribute is **shared**

It can be merged with the **parallel** construct:

```
#pragma omp parallel for
```


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Example

```
void foo (int ∗m, int N, int M)
{
  int i;
  #pragma omp parallel for private( j )
  for (i = 0; i < N; i++)for ( j = 0; j < M; j_{++} )
        m[i][j] = 0;}
```


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Example

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```
void foo (int ∗m, int N, int M)
{
  int i;
  #pragma omp parallel for private( j )
  for ( i = 0; i < N; i + 1for ( j \leftarrow 0; -\lfloor \frac{M}{M} \rfloor has be explicitly privatized
          m[i][j] = b;
}
```


```
void foo ( std :: vector <int > &v )
{
  #pragma omp parallel for
  for ( std:: vector <int >:: iterator it = v. begin();
         it < v.end() ;
        it ++ )* it = 0;
}
```


[Loop worksharing](#page-63-0)

Removing dependences

Example

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[Loop worksharing](#page-64-0)

Removing dependences

The lastprivate clause

When a variable is declared **lastprivate**, a private copy is generated for each thread. Then the value of the variable in the last iteration of the loop is copied back to the original variable.

A variable can be both **firstprivate** and **lastprivate**

The reduction clause

A very common pattern is where all threads accumulate some values into a single variable

- \bullet E.g., $n + =$ v[i], our pi program, ...
- Using **critical** or **atomic** is not good enough
	- Besides being error prone and cumbersome
- Instead we can use the **reduction** clause for basic types.
	- Valid operators are: $+, -, *, |, |, |, 8, 8, 8, \rangle$
	- The compiler creates a **private** copy that is properly initialized
	- At the end of the region, the compiler ensures that the **shared** variable is properly (and safely) updated.

We can also specify **reduction** variables in the **parallel** construct.

 \Box

The reduction clause

```
int vector_sum (int n, int v[n])
{
   int i, sum = 0;
   #pragma omp parallel for reduction ( + : sum )
   {
      for ( i = 0; i < n; i_{++} )
         sum += v[i];
   }
   return sum;
}
```


The reduction clause

The schedule clause

The **schedule** clause determines which iterations are executed by each thread.

- If no **schedule** clause is present then is implementation defined There are several possible options as schedule:
	- **STATIC**
	- **STATIC,chunk**
	- **DYNAMIC[,chunk]**
	- **GUIDED[,chunk]**
	- **AUTO**
	- **RUNTIME**

The schedule clause

Static schedule

The iteration space is broken in chunks of approximately size *N*/*num* − *threads*. Then these chunks are assigned to the threads in a Round-Robin fashion.

Static,N schedule (Interleaved)

The iteration space is broken in chunks of size *N*. Then these chunks are assigned to the threads in a Round-Robin fashion.

Characteristics of static schedules

- **o** Low overhead
- Good locality (usually)
- Can have load imbalance problems

 $(1 - 1)$

The schedule clause

Dynamic,N schedule

Threads dynamically grab chunks of *N* iterations until all iterations have been executed. If no chunk is specified, $N = 1$.

Guided,N schedule

Variant of **dynamic**. The size of the chunks deceases as the threads grab iterations, but it is at least of size *N*. If no chunk is specified, $N = 1$.

Characteristics of dynamic schedules

- **•** Higher overhead
- Not very good locality (usually)
- Can solve imbalance problems

 $(1 - 1)$
The schedule clause

Auto schedule

In this case, the implementation is allowed to do whatever it wishes.

• Do not expect much of it as of now

Runtime schedule

The decision is delayed until the program is run through the **sched-nvar** ICV. It can be set with:

- **The OMP SCHEDULE** environment variable
- **•** The omp set schedule() API call

When a worksharing has a **nowait** clause then the implicit **barrier** at the end of the loop is removed.

• This allows to overlap the execution of non-dependent loops/tasks/worksharings

Example

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Example

```
#pragma omp for nowait
for ( i = 0; i < n ; i++ )
  v[i] = 0;#pragma omp for
for (i = 0; i < n; i++)a[i] = 0;
```
On a side note, you would be better by fusing the loops in this case

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Exception: static schedules

If the two (or more) loops have the same **static** schedule and all have the same number of iterations.

```
#pragma omp for schedule( s t a ti c , 2 ) nowait
for ( i = 0; i < n ; i_{++} )
   v[i] = 0;
#pragma omp for schedule( s t a ti c , 2 )
for (i = 0; i < n; i++)a[i] = v[i]*v[i];
```


The collapse clause

Allows to distribute work from a set of *n* nested loops.

- Loops must be perfectly nested
- The nest must traverse a rectangular iteration space

The collapse clause

Allows to distribute work from a set of *n* nested loops.

- Loops must be perfectly nested
- The nest must traverse a rectangular iteration space

Example

i and *j* loops are folded and iterations distributed among all threads. Both *i* and *j* are privatized

[Break](#page-80-0)

Coffee time! :-)

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Part III

[Basic Synchronizations](#page-81-0)

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• [Thread barriers](#page-84-0)

• [Exclusive access](#page-89-0)

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Mechanisms

Threads need to synchronize to impose some ordering in the sequence of actions of the threads. OpenMP provides different synchronization mechanisms:

- **barrier**
- **critical**
- **atomic**

Outline

• [Thread barriers](#page-84-0)

• [Exclusive access](#page-89-0)

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Thread Barrier

The barrier construct

#pragma omp barrier

- Threads cannot proceed past a barrier point until all threads reach the barrier AND all previously generated work is completed
- Some constructs have an implicit **barrier** at the end
	- E.g., the **parallel** construct

Barrier

Example

```
#pragma omp parallel
{
    foo();
#pragma omp barrier
    bar();
}
```


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Barrier

Example

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Barrier

Outline

• [Thread barriers](#page-84-0)

• [Exclusive access](#page-89-0)

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Exclusive access

The critical construct

#pragma omp critical [(name)] structured block

- Provides a region of mutual exclusion where only one thread can be working at any given time.
- By default all critical regions are the same, but you can provide them with names
	- Only those with the same name synchronize


```
int x=1;
#pragma omp parallel num_threads( 2 )
{
#pragma omp critical
    X + +;
}
print(f("d\nu", x);
```


Example

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Example

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```
int x = 1, y = 0;
#pragma omp parallel num_threads( 4 )
{
#pragma omp critical ( x )
    x + +;
#pragma omp critical ( y )
    y++;}
```


Exclusive access

The atomic construct

#pragma omp atomic expression

• Provides an special mechanism of mutual exclusion to do read & update operations

• Only supports simple read & update expressions

• E.g., $x == 1$, $x = x - foo()$

- Only protects the read & update part
	- foo() not protected
- Usually much more efficient than a **critical** construct
- Not compatible with **critical**

```
int x=1;
#pragma omp parallel num_threads( 2 )
{
#pragma omp atomic
    X + +;
}
print(f("d\nu", x);
```


Example

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```
int x=1;
#pragma omp parallel num_threads( 2 )
{
#pragma omp critical
    x + +;
#pragma omp atomic
    X + +;
}
print(f("d\nu", x);
```


Example

 $(1 - 1)$

```
int x=1;
#pragma omp parallel num_threads( 2 )
{
#pragma omp critical
     x + +;
#pragma omp atomic
     X + +;
}
{\sf print(f(\texttt{"\%d}\setminus n", x);\longleftarrow \texttt{Prints 3,4 or 5 :()})}
```


Part IV

[Practical: heat diffusion](#page-103-0)

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• [Heat diffusion](#page-105-0)

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[Heat diffusion](#page-105-0)

Outline

• [Heat diffusion](#page-105-0)

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Before you start

Enter the OpenMP directory to do the following exercises.

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Description of the Heat Diffusion app Hands-on

Parallel loops

The file solver.c implements the computation of the Heat diffusion

- ¹ Annotate the jacobi, redblack, and gauss functions with OpenMP
- ² Execute the application with different numbers of processors, and compare the results

Bon appétit!*

*Disclaimer: actual food may differ from the image! :-)

Part V

[Task Parallelism in OpenMP](#page-109-0)

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Outline

- **•** [OpenMP tasks](#page-111-0)
- **•** [Task synchronization](#page-125-0)
- [The single construct](#page-133-0)
- **•** [Task clauses](#page-140-0)
- [Common tasking problems](#page-144-0)

Outline

- **•** [OpenMP tasks](#page-111-0)
- [Task synchronization](#page-125-0) \bullet
- [The single construct](#page-133-0) \bullet
- [Task clauses](#page-140-0) \bullet
- [Common tasking problems](#page-144-0)

Task parallelism in OpenMP

Task parallelism model

Parallelism is extracted from "several" pieces of code

- Allows to parallelize very unstructured parallelism
	- Unbounded loops, recursive functions, ...

 \Box

What is a task in OpenMP ?

- Tasks are work units whose execution may be deferred • they can also be executed immediately
- Tasks are composed of:
	- code to execute
	- a data environment
		- **o** Initialized at creation time
	- internal control variables (ICVs)
- Threads of the team cooperate to execute them

Creating tasks

The task construct

#pragma omp task [clau se s] structured block

Where clauses can be:

- **o** shared
- **o** private
- **•** firstprivate
	- Values are captured at creation time
- **o** default
- **if(expression)**
- **untied**

 \Box

When are task created?

Parallel regions create tasks

- One implicit task is created and assigned to each thread
	- So all task-concepts have sense inside the parallel region

Each thread that encounters a **task** construct

- Packages the code and data
- Creates a new explicit task

Default task data-sharing attributes

When there are no clauses ...

If no default clause

- o Implicit rules apply
	- e.g., global variables are shared
- **Otherwise...**
	- **firstprivate**
	- **shared** attribute is lexically inherited

Task default data-sharing attributes

In practice...

```
int a:
void foo () {
  i n t b , c ;
  #pragma omp parallel shared( b )
  #pragma omp parallel private( b )
  {
         int d;
         #pragma omp task
         {
             int e:
             a =h =c =d =e =} } }
```
Task default data-sharing attributes

In practice...

```
int a:
void foo () {
  i n t b , c ;
  #pragma omp parallel shared( b )
  #pragma omp parallel private( b )
  {
         int d;
         #pragma omp task
         {
             int e:
             a = shared
             h =c =d =e =} } }
```
Task default data-sharing attributes

In practice...

```
int a:
void foo () {
  i n t b , c ;
  #pragma omp parallel shared( b )
  #pragma omp parallel private( b )
  {
         int d;
         #pragma omp task
         {
             int e:
             a = shared
             b = firstprivate
             c =d =e =} } }
```
Task default data-sharing attributes

In practice...

```
int a:
void foo () {
  i n t b , c ;
  #pragma omp parallel shared( b )
  #pragma omp parallel private( b )
  {
         int d;
         #pragma omp task
         {
             int e:
             a = shared
             b = firstprivate
             c = shared
             d =e =} } }
```
Task default data-sharing attributes

In practice...

Example

```
int a:
void foo () {
  i n t b , c ;
  #pragma omp parallel shared( b )
  #pragma omp parallel private( b )
  {
         int d;
        #pragma omp task
         {
             int e:
             a = shared
             b = firstprivate
             c = shared
             d = firstprivate
             e =
```
} } }

Task default data-sharing attributes

In practice...

Example

```
int a:
void foo () {
  i n t b , c ;
  #pragma omp parallel shared( b )
  #pragma omp parallel private( b )
  {
         int d;
        #pragma omp task
         {
             int e:
             a = shared
             b = firstprivate
             c = shared
             d = firstprivate
             e = private
```
} } }

Task default data-sharing attributes

In practice...

Example

```
int a:
void foo ( ) {
  i n t b , c ;
  #pragma omp parallel shared( b )
  #pragma omp parallel private( b )
  {
         int d;
         #pragma omp task
         {
             int e:
             a = shared
             b = firstprivate
             c = shared
             d = firstprivate
             e = private
} } }
Tip: default (none) is your friend if you do not see it clearly
```


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List traversal

Example

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Outline

- **[OpenMP tasks](#page-111-0)**
- [Task synchronization](#page-125-0) \bullet
- [The single construct](#page-133-0)
- [Task clauses](#page-140-0) \bullet
- [Common tasking problems](#page-144-0)

[Task synchronization](#page-126-0)

Task synchronization

There are two main constructs to synchronize tasks:

- **barrier**
	- Remember: all previous work (including tasks) must be completed
- **taskwait**

[Task synchronization](#page-127-0)

Waiting for children

The taskwait construct

#pragma omp taskwait

Suspends the current task until all children tasks are completed **Just direct children, not descendants**

Taskwait

Example

}

#pragma omp taskwait

All tasks guaranteed to be completed here

Taskwait

Example

List |

```
#pragma omp parallel
   traverse\_list(1);
```


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List traversal Completing the picture

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Outline

- **[OpenMP tasks](#page-111-0)**
- [Task synchronization](#page-125-0) \bullet
- [The single construct](#page-133-0)
- [Task clauses](#page-140-0) \bullet
- [Common tasking problems](#page-144-0)

Giving work to just one thread

The single construct

- **#pragma omp single** [clau se s] structured block
	- o where clauses can be:
		- private
		- **•** firstprivate
		- nowait <-We'll see it later
		- copyprivate<
		Not today
	- o Only one thread of the team executes the structured block
	- There is an implicit **barrier** at the end


```
int main (int argc, char **argv)
{
    #pragma omp parallel
    {
       #pragma omp single
        {
           printf ("Hello_world!\n");
        }
    }
}
```


Example

List |

```
#pragma omp parallel
#pragma single
   traverse\_list(1);
```


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Example

List 1

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Example List 1

Xavier Martorell (BSC) [PATC Parallel Programming Workshop](#page-1-0) November 26-30, 2012 97 / 120

Outline

- **[OpenMP tasks](#page-111-0)**
- [Task synchronization](#page-125-0) \bullet
- [The single construct](#page-133-0) \bullet
- **[Task clauses](#page-140-0)**
- [Common tasking problems](#page-144-0)

Task scheduling

How it works?

Tasks are tied by default

- **Tied tasks are executed always by the same thread**
	- Not necessarily the creator
- **Tied tasks have scheduling restrictions**
	- Deterministic scheduling points (creation, synchronization, ...)
		- **•** Tasks can be suspended/resumed at these points
	- Another constraint to avoid deadlock problems
- Tied tasks may run into performance problems

The untied clause

A task that has been marked as **untied** has none of the previous scheduling restrictions:

- Can *potentially* switch to any thread
- Can *potentially* switch at any moment
- Bad mix with thread based features
	- thread-id, critical regions, threadprivate
- **•** Gives the runtime more flexibility to schedule tasks

The if clause

If the the expression of an if clause evaluates to false

- The encountering task is suspended
- The new task is executed immediately
	- with its own data environment
	- **•** different task with respect to synchronization
- The parent task resumes when the task finishes
- Allows implementations to optimize task creation
	- For very fine grain task you may need to do your own if

Outline

- **[OpenMP tasks](#page-111-0)**
- [Task synchronization](#page-125-0) \bullet
- [The single construct](#page-133-0) \bullet
- [Task clauses](#page-140-0) \blacksquare
- \bullet [Common tasking problems](#page-144-0)


```
void search (int n, int j, bool *state)
{
    int i, res;
    if (n == i) {
      /* good solution, count it */
      s o lutions ++:
      re tu rn ;
    }
    /* try each possible solution */
    for (i = 0; i < n; i++)state[i] = i;
       if (ok(j+1, state))search(n, j+1, state);}
}
```


```
void search (int n, int j, bool *state)
{
    int i, res;
    if (n == i) {
      /* good solution, count it */
      s o lutions ++:
      re tu rn ;
     }
    /* try each possible solution */
    for (i = 0; i < n; i++)#pragma omp task
       state[i] = i;
       if ( ok(j+1, state)) {
          search(n, j+1, state);}
}
```


Example

```
void search (int n, int j, bool *state)
{
    int i, res;
    if (n == i)/* good solution, count it */
      s o lutions ++:
      re tu rn ;
    }
    /* try each possible solution */for (i = 0; i < n; i++)#pragma omp task
       state[i] = i;
       if ( ok(j+1, state)) {
         search(n, j+1, state);}
}
```
Data scoping

Because it's an orphaned task all variables are firstprivate

Example

```
void search (int n, int j, bool *state)
{
    int i, res;
    if (n == i)/* good solution, count it */
      s o lutions ++:
      re tu rn ;
    }
    /* try each possible solution */
    for (i = 0; i < n; i++)#pragma omp task
       state[i] = i;
       if ( ok(j+1, state)) {
          search(n, j+1, state);}
}
```
Data scoping

Because it's an orphaned task all variables are firstprivate

State is not captured

Just the pointer is captured not the pointed data

 $+$ $+$

Example

```
void search (int n, int j, bool *state)
{
    int i, res;
    if (n == i)/* good solution, count it */
      s o lutions ++:
      re tu rn ;
    }
    /* try each possible solution */
    for (i = 0; i < n; i++)#pragma omp task
       state[i] = i;
       if ( ok(j+1, state)) {
         search(n, j+1, state);}
}
```
Problem #1

Incorrectly capturing pointed data

Problem #1

Incorrectly capturing pointed data

Problem

firstprivate does not allow to capture data through pointers

Solutions

- **1** Capture it manually
- 2 Copy it to an array and capture the array with firstprivate


```
void search (int n, int j, bool *state)
{
    int i.res:
    if (n == i)/* good solution, count it */
      s o lutions ++;
      re tu rn ;
    /* try each possible solution */for (i = 0; i < n; i++)#pragma omp task
       bool ∗new state = alloca ( sizeof ( bool )∗n ) ;
       memcpy ( new_state, state, sizeof ( bool ) *n ) ;
       new state [i] = i;
       if (ok(i+1.new state)) {
          search (n, j+1, new state);
     }
}
```


Example

```
void search (int n, int j, bool *state)
{
    int i.res:
    if (n == i)/* good solution, count it */
      s o lutions ++;
      re tu rn ;
    /* try each possible solution */for (i = 0; i < n; i++)#pragma omp task
       bool ∗new state = alloca ( sizeof ( bool )∗n ) ;
       memcpy ( new_state, state, sizeof ( bool ) *n ) ;
        new state [i] = i;
        if ( ok(i+1), new state ) {
          search ( n, j+1, new state );
     }
}
```
Caution!

Will new state still be valid by the time memcpy is executed?

Example

```
void search (int n, int j, bool *state)
{
    int i.res:
    if (n == i)/* good solution, count it */
      s o lutions ++;
      re tu rn ;
    /* try each possible solution */for (i = 0; i < n; i++)#pragma omp task
       bool ∗new state = alloca ( sizeof ( bool )∗n ) ;
       memcpy ( new_state, state, sizeof ( bool ) *n ) ;
       new state [j] = i;
        if ( ok(i+1), new state ) {
          search ( n, j+1, new state );
     }
}
```
Problem #2

Data can go out of scope!

Problem #2 Out-of-scope data

Problem

Stack-allocated parent data can become invalid before being used by child tasks

• Only if not captured with firstprivate

Solutions

- **1** Use firstprivate when possible
- 2 Allocate it in the heap
	- Not always easy (we also need to free it)
- ³ Put additional synchronizations
	- May reduce the available parallelism

 \Box

```
void search (int n, int j, bool *state)
    int i.res:
     if (n == i) {
       /* good solution, count it */
       s o lution s++:
       return:
     }
    /* try each possible solution */for (i = 0; i < n; i++)#pragma omp task
        bool * new state = alloca ( sizeof (bool) *n ) ;
        memcpy ( new state , state , sizeof ( bool )∗n ) ;
        new state \overline{[} \overline{]} = i;
        if ( ok ( j + 1, new state ) ) {
          search(n, j+1, new\_state);
     }
    #pragma omp taskwait
}
```


```
void search (int n, int j, bool *state)
    int i.res:
    if (n == i)/* good solution, c
      s o lutions ++re tu rn ;
    }
    /* try each possible solution */for (i = 0; i < n; i++)#pragma omp task
       bool * new state = alloca(<b>sizeof</b> (bool) *n);
       memcpy ( new state , state , sizeof ( bool )∗n ) ;
       new state [i] = i;
       if ( ok(i+1) new state ) {
          search(n, j+1, new\_state);
    }
    #pragma omp taskwait
}
                          Shared variable needs protected access
```


Example

```
void search (int n, int j, bool *state)
    int i.res:
    if (n == i) {
      /* good solution, count it */
      s o lution s++:
      re tu rn ;
     }
    /* try each possible solution */for (i = 0; i < n; i++)#pragma omp task
       bool * new state = alloca ( sizeof (bool) *n ) ;
       memcpy ( new state , state , sizeof ( bool )∗n ) ;
       new state [i] = i;
       if (ok(i+1.new state))search(n, j+1, new\_state);
     }
    #pragma omp taskwait
}
```
Solutions

- Use **critical**
- Use **atomic**
- Use **threadprivate**

Reductions for tasks


```
void search (int n, int j, bool *state)
    int i.res:
    if (n == i) {
      /* good solution, count it */
      mysolutions++;
      re tu rn ;
     }
    /* try each possible solution */for (i = 0; i < n; i++)#pragma omp task
     {
        bool * new state = alloca ( sizeof (bool) *n ) ;
       memcpy ( new state , state , sizeof ( bool ) ∗n ) ;
        new state [i] = i;
        if ( ok(i+1, new state) ) {
          search(n, j+1, new\_state);
        }
     }
    #pragma omp taskwait
```


Part VI

[Programming using a hybrid](#page-160-0) [MPI/OpenMP approach](#page-160-0)

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[MPI+OpenMP programming](#page-162-0)

Xavier Martorell (BSC) [PATC Parallel Programming Workshop](#page-1-0) November 26-30, 2012 111 / 120

[MPI+OpenMP programming](#page-162-0)

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Alternatives

MPI + computational kernels in OpenMP

Use OpenMP directives to exploit parallelism between communication phases

• OpenMP parallel will end before new communication calls

MPI inside OpenMP constructs

Call MPI from within for-loops, or tasks

• MPI needs to support multi-threaded mode

[MPI+OpenMP programming](#page-164-0)

Compiling MPI+OpenMP

MPI compiler driver gets the proper OpenMP option

- o mpicc -openmp
- mpicc -fopenmp

[Break](#page-165-0)

Coffee time! :-)

Part VII

[Practical: heat diffusion](#page-166-0)

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[MPI+OpenMP Heat diffusion](#page-168-0)

[MPI+OpenMP Heat diffusion](#page-168-0) \bullet

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Before you start

Enter the MPI+OpenMP directory to do the following exercises.

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Description of the Heat Diffusion app Hands-on

Parallel loops

The file solver.c implements the computation of the Heat diffusion.

- **1** Use MPI to distribute the work across nodes
- ² Annotate the jacobi, redblack, and gauss functions with OpenMP tasks
- ³ Execute the application with different numbers of nodes/processors, and compare the results

