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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



Outline

- OpenMP Overview
- The OpenMP model
- Writing OpenMP programs
- Creating Threads
- Data-sharing attributes



Outline

- OpenMP Overview
- The OpenMP model
- Writing OpenMP programs
- Creating Threads
- Data-sharing attributes



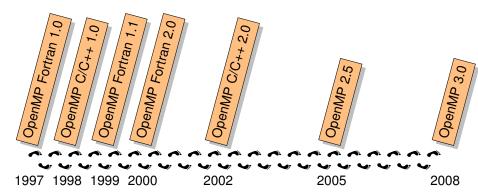
What is OpenMP?

- 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



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
- Does not run on clusters

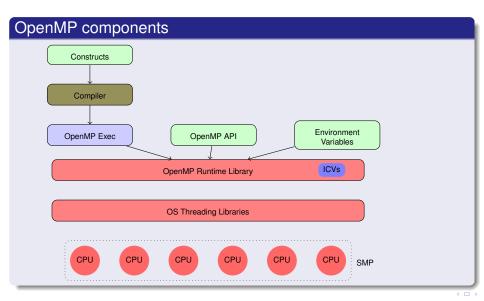


Outline

- OpenMP Overview
- The OpenMP model
- Writing OpenMP programs
- Creating Threads
- Data-sharing attributes



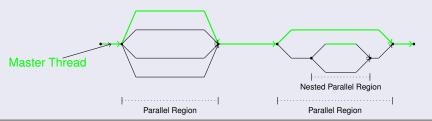
OpenMP at a glance



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





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
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OpenMP directives syntax

In Fortran

Through a specially formatted comment:

```
sentinel construct [clauses]
```

where sentinel is one of:

- !\$OMP or C\$OMP or *\$OMP in fixed format
- ! \$OMP 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:

- !\$OMP or C\$OMP or ★\$OMP in fixed format
- ! \$OMP in free format

In 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

- 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
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The parallel construct

```
Directive
         #pragma omp parallel [clauses]
                                     structured block
where clauses can be:
              num threads(expression)
              • if(expression)

    shared(var-list)←

                                                                                                                                                                                                                                                                                                  Coming shortly!
              • private(var-list)←
              firstprivate(var-list)

    default(none|shared| private  firstprivate  f

    reduction(var-list)
    ← We'll see it later

    copyin(var-list)← Not today

                                                                                                                                                                                                                                                                                                                                                                  Only in Fortran
```

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 <u>num_threads</u> 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
  void main () {
    #pragma omp parallel
    omp_set_num_threads(2);
    #pragma omp parallel
                A team of two threads here.
    #pragma omp parallel num_threads(random()%4+1) if(0)
```



```
void main () {
    #pragma omp parallel
    ...
    omp_set_num_threads(2);
    #pragma omp parallel
    ...
    #pragma omp parallel num threads(random()%4+1) if(0)
    ...
    A team of 1 thread here.
}
```



API calls

Other useful routines

int omp_get_num_threads()

int omp_get_thread_num()

int omp_get_num_procs()

int omp_get_max_threads()

double omp_get_wtime()

Returns the number of threads in the current team

Returns the id of the thread in the current team

Returns the number of processors in the machine

Returns the maximum number of threads that will be used in the next parallel region

Returns the number of seconds since an arbitrary point in the past



Outline

- OpenMP Overview
- The OpenMP model
- Writing OpenMP programs
- Creating Threads
- Data-sharing attributes



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
- lastprivate
- copvin←
- Out of our scope today

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



Example



```
int x=1;
#pragma omp parallel shared(x) num_threads(2)
{
    x++;
    printf("%d\n",x);
}
printf("%d\n",x);
Prints 2 or 3
```



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



Example

```
int x=1;
#pragma omp parallel private(x) num_threads(2)
   X++;
   printf("%d\n",x);
printf("%d\n",x);
```







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







```
int x=1;
#pragma omp parallel firstprivate(x) num_threads(2)
{
     x++;
     printf("%d\n",x);
}
printf("%d\n",x);
Prints 1
```



What is the default?

- Static/global storage is shared
- Heap-allocated storage is shared
- Stack-allocated storage inside the construct is private
- 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



```
int x,y;
#pragma omp parallel private(y)
{
    x =
    y =
    #pragma omp parallel private(x)
    {
        x =
        y =
        y =
    }
}
```



```
Example
```

```
int x,y;
#pragma omp parallel private(y)
               X is shared
    #pragma om y is private ivate(X)
        X =
```



```
int x,y;
#pragma omp parallel private(y)
    Х
    #pragma omp parallel private(X)
                   X is private
                   y is shared
```



Threadprivate storage

The threadprivate construct

#pragma omp threadprivate(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



Threaprivate storage

```
char* foo ()
{
    static char buffer[BUF_SIZE];
    #pragma omp threadprivate(buffer)
```



return buffer;

Threaprivate storage

```
char* foo ()
{
    static char buffer[BUF_SIZE];
    #pragma omp threadprivate(buffer)

    return buffer;
}
Creates one static
copy of buffer per
thread
```



Threaprivate storage

```
char* foo ()
{
    static char buffer[BUF_SIZE];
    #pragma omp threadprivate(buffer)

    return buffer;
}
Now foo can be called by multiple threads at the same time
```



Part II

Worksharing constructs



Outline

The worksharing concept

Loop worksharing



Outline

The worksharing concept

Loop worksharing



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
- loop worksharing
- Section← We'll see them later
- workshare

Restriction: worksharings cannot be nested

Outline

The worksharing concept

Loop worksharing



Loop parallelism

The for construct

```
#pragma omp for [clauses]
for( init-expr ; test-expr ; inc-expr )
```

where clauses can be:

- private
- firstprivate
- lastprivate(variable-list)
- reduction(operator:variable-list)
- schedule(schedule-kind)
- nowait
- ollapse(n)
- 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
```

```
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;
}</pre>
```



```
void foo (int *m, int N, int M)
{
  int i;
  #pragma omp parallel for(privation for ( i = 0; i < N; i++ )
      for ( j = 0; j < M; j++ )
            m[i][j] = 0;
}</pre>
New created threads cooperate to execute all the iterations of the loop
```





```
void foo (int *m, int N, int M)
  int i;
  #pragma omp parallel for private(j)
  for (i = 0; i < N: i++)
     for ( j <del><= 0;</del> Must be explicitly privatized
```



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







Removing dependences

```
 \begin{array}{lll} x &=& 0;\\ & \text{for } ( & i &=& 0; & i &< n; & i++ &) \\ & & & v[i] &=& x;\\ & & & x &+=& dx; & \\ & & & & \\ \end{array}  Each iteration x depends on the previous one. Can't be parallelized
```



Removing dependences

```
x = 0;
for ( i = 0; i < n; i++ )

\begin{cases}
x = i * dx; & \text{But } x \text{ can be rewritten in terms of } i. \\
v[i] = x;
\end{cases}
Now it can be parallelized
```



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

- 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: +,-,*,|,||,&,&&,^
- 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.



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

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

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

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



```
#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;
```



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;</pre>
```

On a side note, you would be better by fusing the loops in this case



```
#pragma omp for nowait dent!. No guarantees that the pre-
for ( i = 0; i < n ; i++ )
    v[i] = 0;

#pragma omp for

for ( i = 0; i < n ; i++ )
    a[i] = v[i]*v[i];
```



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(static,2) nowait
for ( i = 0; i < n ; i++ )
   v[i] = 0;
#pragma omp for schedule(static,2)
for ( i = 0; i < n ; i++ )
   a[i] = v[i]*v[i];</pre>
```



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

```
#pragma omp for collapse(2)

for ( i = 0; i < N; i++ ) 

for ( j = 0; j < M; j++ ) 

foo (i,j);
```





Coffee time! :-)



Part III

Basic Synchronizations



Outline

Thread barriers

Exclusive access



Why synchronization?

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
- taskwait
- ordered← We'll see them later
- locks



Outline

Thread barriers

Exclusive access



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

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



Barrier

```
#pragma omp parallel

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

Forces all foo occurrences too happen before all bar occurrences
}
```



Barrier

```
#pragma omp parallel
{
    foo();
#pragma omp barrier
    bar():
}
Implicit barrier at the end of the parallel region
```



Outline

Thread barriers

Exclusive access



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







```
Example
int x=1;
#pragma omp parallel num_threads(2)
#pragma omp critical
    x++;← Only one thread at a time here
                        Prints 3!
printf("%d\n",x);\leftarrow
```







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





```
Example
int x=1;
#pragma omp parallel num_threads(2)
#pragma omp atomic
    x++;\leftarrow Only one thread at a time updates x here
printf("%d\n",x);
```







```
Example
int x=1;
#pragma omp parallel num_threads(2)
#pragma omp critical
                              Different threads can update x at
    X + + ; \leftarrow
                              the same time!
#pragma omp atomic
    X++:←
printf("%d\n",x);
```



```
Example
int x=1;
#pragma omp parallel num_threads(2)
#pragma omp critical
    X++;
#pragma omp atomic
    X++;
                         Prints 3,4 or 5 :()
printf("%d\n",x);\leftarrow
```



Part IV

Practical: heat diffusion



Outline

Heat diffusion



Outline

Heat diffusion



Before you start

Enter the OpenMP directory to do the following exercises.



Description of the Heat Diffusion app Hands-on

Parallel loops

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

- Annotate the jacobi, redblack, and gauss functions with OpenMP
- 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



Outline

- OpenMP tasks
- Task synchronization
- The single construct
- Task clauses
- Common tasking problems



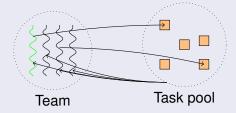
Outline

- OpenMP tasks
- Task synchronization
- The single construct
- Task clauses
- Common tasking problems



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, ...



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
 - Initialized at creation time
 - internal control variables (ICVs)
- Threads of the team cooperate to execute them



Creating tasks

The task construct

```
#pragma omp task [clauses]
    structured block
```

Where clauses can be:

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



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

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



In practice...

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

In practice...

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

In practice...

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

In practice...

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

In practice...

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

In practice...

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

In practice...

Example

```
int a:
void foo() {
  int b,c;
  #pragma omp parallel shared(b)
  #pragma omp parallel private(b)
        int d:
        #pragma omp task
            int e:
                 shared
                firstprivate
            C = shared
            d = firstprivate
            e = private
111
```

Tip: default (none) is your friend if you do not see it clearly

process(e); ← e is firstprivate

List traversal



Outline

- OpenMP tasks
- Task synchronization
- The single construct
- Task clauses
- Common tasking problems



Task synchronization

There are two main constructs to synchronize tasks:

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



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

```
void traverse list ( List I )
  Element e:
  for (e = 1 \rightarrow first; e; e = e \rightarrow next)
     #pragma omp task
        process(e);
  #pragma omp taskwait
          All tasks guaranteed to be completed here
```



Taskwait

```
void traverse_list ( List I )
{
    Element e;
    for ( e = I -> first; e ; e = e -> next )
        #pragma omp task
        process(e);
    #pragma omp taskwait
}
Now we need some threads
to execute the tasks

#pragma omp taskwait
}
```



Completing the picture

```
List |
#pragma omp parallel
    traverse_list(|);
```



Completing the picture

Example List | #pragma omp parallel traverse_list(|); This will generate multiple traversals



Completing the picture

Example List I #pragma omp parallel We need a way to have a single traverse_list(I);← thread execute traverse list



Outline

- OpenMP tasks
- Task synchronization
- The single construct
- Task clauses
- Common tasking problems



Giving work to just one thread

The single construct

```
#pragma omp single [clauses]
    structured block
```

- where clauses can be:
 - private
 - firstprivate
 - nowait

 We'll see it later
 - copyprivate Not today
- Only one thread of the team executes the structured block
- There is an implicit barrier at the end



The single construct



The single construct



Completing the picture

```
List |
#pragma omp parallel
#pragma single
    traverse_list(|);
```



Completing the picture

```
Example
```

List I

```
#pragma omp parallel
#pragma single
    traverse list(|);←
```

One thread creates the tasks of the traversal



Completing the picture

```
Example

List |

#pragma omp parallel

#pragma single

traverse_list(|);

All threads cooperate to execute them
```



Outline

- OpenMP tasks
- Task synchronization
- The single construct
- Task clauses
- Common tasking problems



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
- Task synchronization
- The single construct
- Task clauses
- Common tasking problems



Example

```
void search (int n, int j, bool *state)
    int i, res;
    if (n == j) {
      /* good solution, count it */
      solutions++:
      return:
    /* try each possible solution */
    for (i = 0; i < n; i++)
       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 == j) {
      /* good solution, count it */
      solutions++:
      return:
    /* 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 == j) {
      /* good solution, count it */
      solutions++:
      return:
    /* 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 */
      solutions++;
      return:
    /* 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 == j) {
      /* good solution, count it */
      solutions++:
      return:
    /* 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

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



Example

```
void search (int n, int j, bool *state)
    int i, res;
    if (n == j) {
     /* good solution, count it */
      solutions++;
      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[j] = i;
       if (ok(j+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 */
      solutions++;
      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[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 == j) {
      /* good solution, count it */
      solutions++;
      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[i] = 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

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



Example

```
void search (int n, int j, bool *state)
    int i.res:
    if (n == i) {
     /* good solution, count it */
      solutions++:
      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 [i] = 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 == j) {
    /* good solution, c
    solutions++←
    Shared variable needs protected access
    return;
}

/* try each possible solution*/
for (i = 0; i < n; i++)
#pragma omp task</pre>
```

bool *new_state = alloca(sizeof(bool)*n);
memcpy(new state, state, sizeof(bool)*n);



#pragma omp taskwait

new_state[j] = i;
if (ok(j+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 */
      solutions++:
      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 [i] = i;
       if (ok(j+1,new state)) {
         search(n, j+1, new state);
    #pragma omp taskwait
```

Solutions

- Use critical
- Use atomic
- Use threadprivate



Reductions for tasks

```
Example
 int solutions = 0:
 int mysolutions=0:
                                               Use a separate counter for each thread
 #pragma omp threadprivate(mysolutions)←
 void start search ()
   #pragma omp parallel
     #pragma omp single
         bool initial state[n];
         search(n,0,initial state);
     #pragma omp atomic
                                            Accumulate them at the end
          solutions += mysolutions:
```

Example

```
void search (int n, int j, bool *state)
    int i.res:
    if (n == i) {
      /* good solution, count it */
      mysolutions++;
      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 [i] = i;
       if (ok(j+1,new state)) {
         search(n, j+1, new state);
    #pragma omp taskwait
```



Part VI

Programming using a hybrid MPI/OpenMP approach



Outline

MPI+OpenMP programming



Outline

MPI+OpenMP programming



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



Compiling MPI+OpenMP

MPI compiler driver gets the proper OpenMP option

- mpicc -openmp
- mpicc -fopenmp





Coffee time! :-)



Part VII

Practical: heat diffusion



Outline

MPI+OpenMP Heat diffusion



Outline

MPI+OpenMP Heat diffusion



Before you start

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



Description of the Heat Diffusion app Hands-on

Parallel loops

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

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

