

Parallel Computing & OpenMP – Parallel Regions

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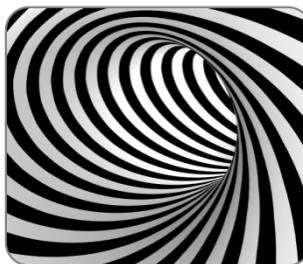
“Foundation of HPC” course



DATA SCIENCE &
SCIENTIFIC COMPUTING
2021-2022 @ Università di Trieste



OpenMP Outline



Parallel
Regions

Parallel
Loops

Advanced
Parallelism



NUMA
AWARENESS



Parallel Regions



```
#pragma omp parallel _some_clauses_here_
single-line-here
```

```
#pragma omp parallel _some_clauses_here_
{ ... }
```

```
#pragma omp parallel for _some_clauses_here_
{ ... }
```

```
#pragma omp sections _some_clauses_here_
{ ... }
```

```
#pragma omp task _some_clauses_here_
{ ... }
```

A parallel region can be as short as a single line

There are no limits on the size of the code included within {}.

The specific construct about `for` loops

A more general work-sharing construct

This allows task-based parallelism

The region starts at the opening { brace and ends at the closing } one.

An implicit synchronization barrier is present at the end of the region (*).

However, for efficiency reasons, it may be that the threads do not die at the end of the region but remain alive until the father process dies

(*) we'll see the details about synchronization later on



| The threads factory

When you create a parallel region, through an OpenMP directive, a pool of threads is created.

Each one receives an ID, ranging from 0 to $n-1$, where n is the number of threads.

Their stack and IP are separated from the others' ones and from the father-process' ones.

Can we check that?
What is the fate of the creating process (thread) ?

```
int      i;

register unsigned long long base_of_stack asm("rbp");
register unsigned long long top_of_stack asm("rsp");

printf( "\nmain thread (pid: %d, tid: %d) data:\n"
        "base of stack is: %p\n"
        "top of stack is : %p\n"
        "&i is           : %p\n"
        "    rbp - &i   : %td\n"
        "    &i - rsp    : %td\n"
        "\n\n",
        getpid(), syscall(SYS_gettid),
        (void*)base_of_stack,
        (void*)top_of_stack,
        &i,
        (void*)base_of_stack - (void*)&i,
        (void*)&i - (void*)top_of_stack );

#pragma omp parallel private(i)
{
    int me = omp_get_thread_num();
    unsigned long long my_stackbase;
    __asm__("mov %%rbp,%0" : "=mr" (my_stackbase));

    printf( "\tthread (tid: %ld) nr %d:\n"
            "\t\tmy base of stack is %p ( %td from main\'s stack )\n",
            "\t\tmy i address is %p\n"
            "\t\t\ttd from my stackbase and %td from main\'s\n",
            syscall(SYS_gettid), me,
            (void*)my_stackbase, (void*)base_of_stack - (void*)my_stackbase,
            &i, (void*)&i - (void*)my_stackbase,
            (void*)&i - (void*)base_of_stack);
}
```





| The threads factory

serial section

parallel region

```
main thread (pid: 26291, tid: 26291) data:  
base of stack is: 0x7ffe6f51efc0  
top of stack is : 0x7ffe6f51ef60  
&i is           : 0x7ffe6f51ef7c  
    rbp - &i     : 68  
    &i - rsp     : 28  
  
thread (tid: 26291) nr 0:  
    my base of stack is 0x7ffe6f51eee0 ( 224 from main's stack )  
    my i address is 0x7ffe6f51eea0  
        -64 from my stackbase and -288 from main's  
thread (tid: 26293) nr 2:  
    my base of stack is 0x7f7342c82ba0 ( 597747680288 from main's stack )  
    my i address is 0x7f7342c82b60  
        -64 from my stackbase and -597747680352 from main's  
thread (tid: 26294) nr 3:  
    my base of stack is 0x7f7342880ba0 ( 597751882784 from main's stack )  
    my i address is 0x7f7342880b60  
        -64 from my stackbase and -597751882848 from main's  
thread (tid: 26292) nr 1:  
    my base of stack is 0x7f7343084b20 ( 597743477920 from main's stack )  
    my i address is 0x7f7343084ae0  
        -64 from my stackbase and -597743477984 from main's
```

pid and tid are the IDs of processes and threads for the OS.

Here you see that the master thread stops its activity and join the new pool of threads, while maintaining its own tid.

- ▶ Each thread has its own stack; thread 0 inherits its own stack, which has grown to host the data relative to OpenMP parallel region;
- ▶ the private `i`s are actually in the threads stacks, and so they are at different memory locations than the shared `i`.





| The threads factory



How large is the stack of each thread? The default value is system dependent.

That is the output of the `00_stack_and_scope` code example: it prints the BP and SP pointers of each thread. Then, we know how large is the threads' stack at a given moment (only an integer is defined, plus the system's thread structure), and how much memory is reserved for the stack to grow:

```
thread 0: bp @ 0x7ffc6edea2f0, tp @ 0x7ffc6edea280: 112 B
thread 1: bp @ 0x7fa371d18b20, tp @ 0x7fa371d18ab0: 112 B      --> -382202615648 B from top of thread 0
thread 2: bp @ 0x7fa371916ba0, tp @ 0x7fa371916b30: 112 B      --> -4202256 B from top of thread 1
thread 3: bp @ 0x7fa371514ba0, tp @ 0x7fa371514b30: 112 B      --> -4202384 B from top of thread 2
```

`you did not define OMP_STACKSIZE: try to set it and check what happens to the threads' stack pointers`

In this case, compiled with gcc, it seems that about 4MB are reserved for each thread. The same value holds with pgi, whereas clang seems to reserve 132MB (i.e., the addresses in the virtual space are spaced by 132MB).

However, you can control the size of the threads' stack using the environmental variable `OMP_STACKSIZE`:

```
export OMP_STACKSIZE=N
```

Where N is a number, followed by a letter: 'B', 'K', 'M', 'G' mean bytes, kilobytes, megabytes and gigabytes respectively (if no letter is specified, the number is interpreted as kilobytes).



| The threads factory

```
export OMP_STACKSIZE=32K
```

```
thread 0: bp @ 0x7ffd1a2bbc70, tp @ 0x7ffd1a2bbbb0: 160 B
thread 1: bp @ 0x7f18b4be4af0, tp @ 0x7f18b4be4a50: 160 B      --> -980954214624 B from top of thread 0
thread 2: bp @ 0x7f18b4bdab70, tp @ 0x7f18b4bdaad0: 160 B      --> -40672 B from top of thread 1
thread 3: bp @ 0x7f18b4bd0b70, tp @ 0x7f18b4bd0ad0: 160 B      --> -40800 B from top of thread 2
```

you defined OMP_STACKSIZE as 32K: try to change that value and check what happens to the threads' stack pointers

```
export OMP_STACKSIZE=1M
```

```
thread 0: bp @ 0x7ffc6f8b97f0, tp @ 0x7ffc6f8b9750: 160 B
thread 1: bp @ 0x7fd5bf930af0, tp @ 0x7fd5bf930a50: 160 B      --> -166161058912 B from top of thread 0
thread 2: bp @ 0x7fd5bec96b70, tp @ 0x7fd5bec96ad0: 160 B      --> -13213408 B from top of thread 1
thread 3: bp @ 0x7fd5beb94b70, tp @ 0x7fd5beb94ad0: 160 B      --> -1056608 B from top of thread 2
```

you defined OMP_STACKSIZE as 1M: try to change that value and check what happens to the threads' stack pointers



| The threads factory

How many threads can be created by a single process?
Also this limit is system-dependent.

On Linux, it depends on the amount of total physical memory: basically, the maximum number of threads is the amount of physical memory divided by the memory amount needed to describe and run a thread, times a factor that accounts for the fact that you don't want all the memory allocated only to make the threads alive.
Within this limit, you can change the behaviour of your OpenMP program by using the env variable OMP_THREAD_LIMIT.

You can check the OS' limit with

```
cat /proc/sys/kernel/threads-max
```

or by calling the omp library function

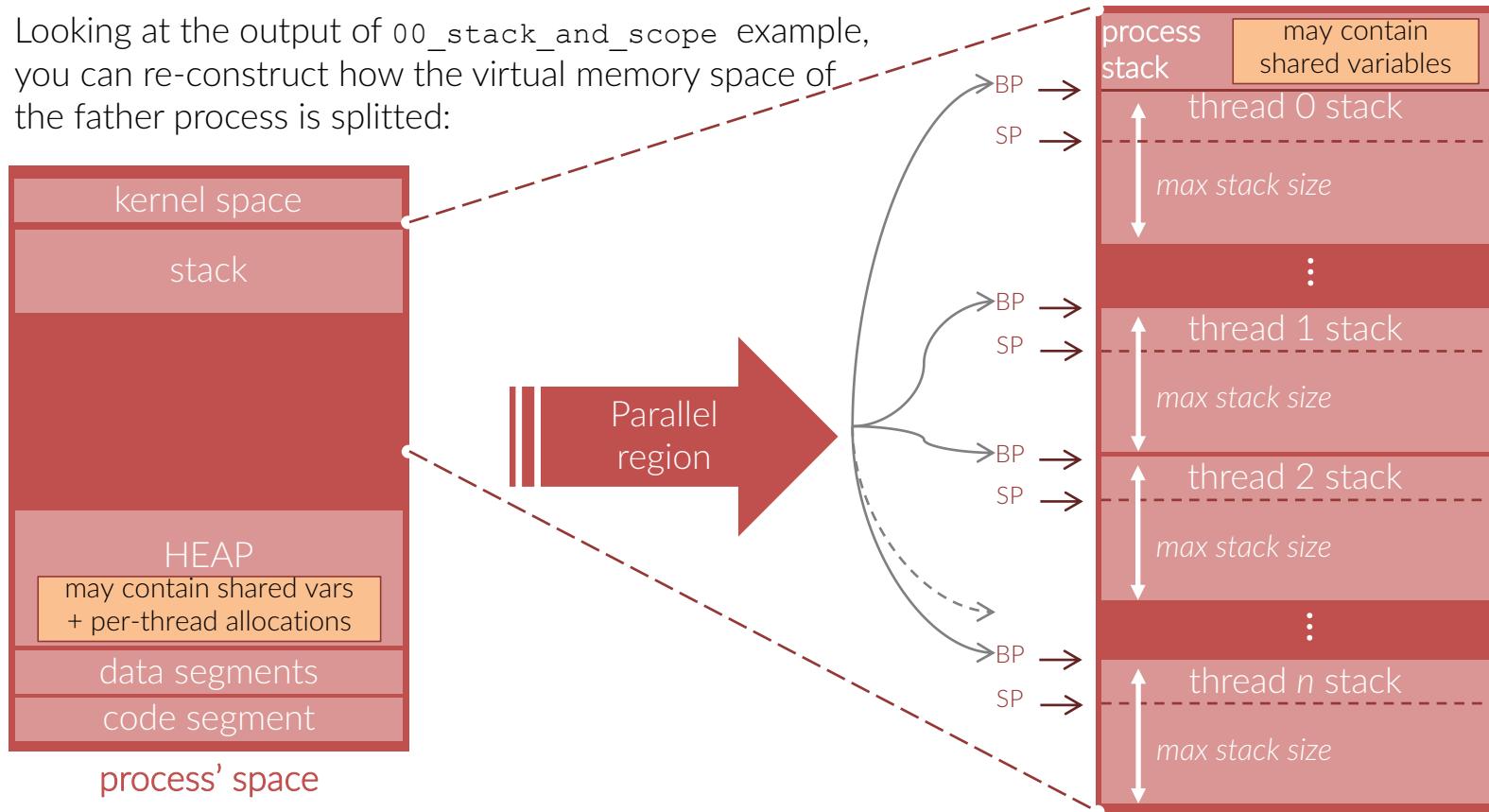
```
int omp_get_max_threads()
```



| The threads factory



Looking at the output of `00_stack_and_scope` example, you can re-construct how the virtual memory space of the father process is splitted:





| The threads factory

Looking at the output of `00_stack_and_scope` example,
you can re-construct how the virtual memory space of

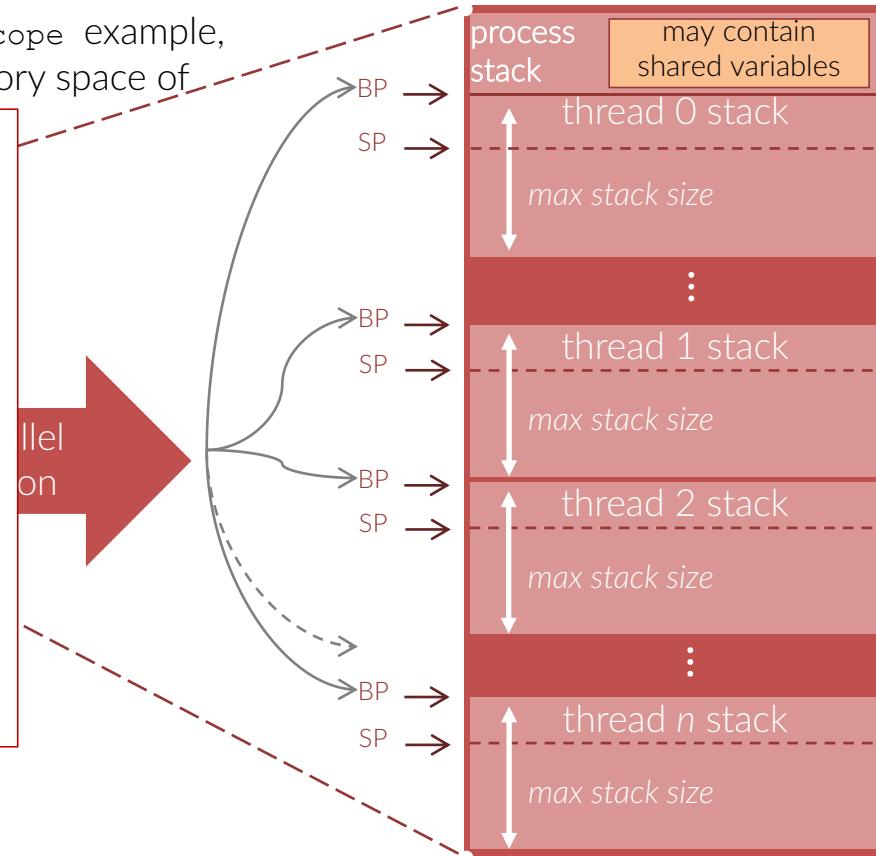
Be aware:

When the threads are created, the space needed for their stack is also allocated. If that space is quite large, you may run out of memory.

Conversely, if the stack is not large enough, you may incur into `seg fault` error (due to lack of memory in the stack)

code segment

process' space

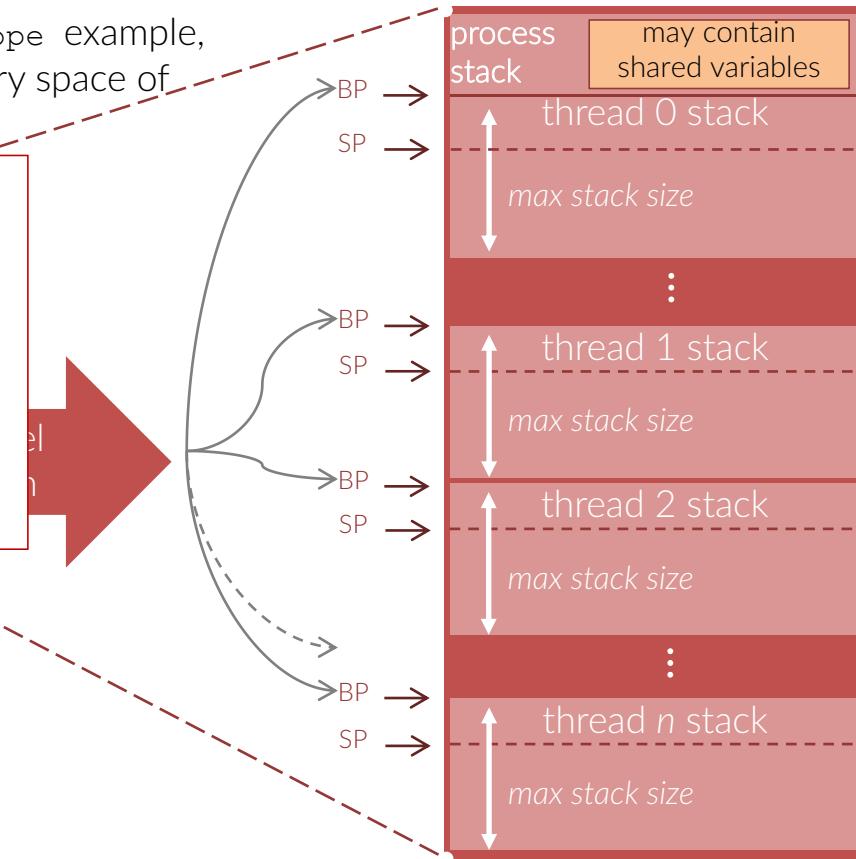
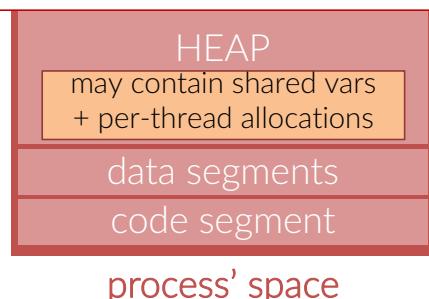




| The threads factory

Looking at the output of `00_stack_and_scope` example, you can re-construct how the virtual memory space of the father process is splitted:

In general, this advocates for the same wise variables placement either on the stack or on the heap that is valid for single-thread application





Parallel Regions

```
int nthreads      = 1;
int my_thread_id = 0;

#ifndef _OPENMP
#pragma omp parallel
{
    my_thread_id = omp_get_thread_num();
    #pragma master
    nthreads     = omp_get_num_threads();
}
#endif
```

returns the ID of the calling thread



returns the number of threads active
in that parallel region

By default rules, variables declared outside a parallel region are *shared*, whereas those declared inside a region are *private*.
Then both `nthreads` and `my_thread_id` are shared.

As a consequence, in this example:

- all threads are writing in `my_thread_id`, in undefined order
- only the master thread is writing in `nthreads`

The value of `my_thread_id` is unpredictable, because it depends on the run-time order by which the threads access it and by each a thread accesses it again to write it.





By default the number of threads spawned in a parallel regions is the number of cores available.
However, you can vary that number in several way

- `OMP_NUM_THREADS` environmental variable
- `#pragma omp parallel num_threads(n)`
- `omp_set_num_threads(n);`
`#pragma omp parallel`





Controlling the number of threads



By default the number of threads spawned in a parallel regions is the number of cores available.
However, you can vary that number in several way

- `OMP_NUM_THREADS` environmental variable
- `#pragma omp parallel num_threads(n)`
- `omp_set_num_threads(n);`
`#pragma omp parallel`

That works if `OMP_DYNAMIC` variable is set to `TRUE`, otherwise the number of threads spawned is strictly equal to `OMP_NUM_THREADS`.
This setting can be changed through `omp_set_dynamic(true)`





| The ordering of the threads



```
#pragma omp parallel
{
    int my_thread_id = omp_get_thread_num();
    printf( "greetings from thread num %d\n",
            my_thread_id );
}
```

The order by which the greetings appear in the output is not deterministic.

In general, unless either you explicitly requires so – and that is possible only for some constructs, or you directly code it, there is no given order, since the threads are independent entities.

How would you build-up an enforcement of the order in the parallel region shown here on the left ?



| The ordering of the threads



```
int order = 0;  
  
#pragma omp parallel  
{  
    int myid = omp_get_thread_num();  
  
    #pragma omp critical  
    if ( order == myid ) {  
        printf( "greetings from thread num %d\n",  
                my_thread_id );  
        order++; }  
}
```



parallel_regions/
04_order_of_threads_wrong.c

The CRITICAL directive defines a region which is executed by all threads but by a single thread at a time. Which starts to sound like an “ordering”.

However, although the threads queue up at the begin of the region, no particular order is imposed to that queue. As a consequence, the result is unpredictable, aside the fact that thread 0 will print the message and increase order.

For instance, if thread 2 is queued immediately after, it will execute the `if` evaluation which will fail (`order` would be equal to 1) and the `thread 2` will not print the message at all.



| The ordering of the threads



```
int order = 0;

#pragma omp parallel
{
    int myid = omp_get_thread_num();
    int done = 0

    while (!done) {
        #pragma omp critical
        if ( order == myid ) {
            printf( "greetings from thread num %d\n",
                    my_thread_id );
            order++; done=1; } }
}
```



parallel_regions/
05_order_of_threads.c

Foundations of HPC

In this second implementation, the `while` cycle enforces the threads that fails the `if` condition to keep trying until the correct order is ensured.

Once a thread has executed the `critical` region, it then exits the `while` and join the implicit synchronization barrier at the end of the parallel region.

A note: without the `critical` directive, this code could work as well on most platform. However, it is still unreliable since it is not guaranteed that each thread complete the region *before* the successive threads enter in it. For instance, the compiler could have been reshuffled the `order++` instruction placing it before the print and so the thread `n+1` could enter the region *before* of the print of the thread `n` and its message could appear as first.



| The ordering of the threads

```
int nthreads = 1;

#pragma omp parallel
{
    int myid = omp_get_thread_num();
    #pragma omp master
    int nthreads = omp_get_num_threads();
    #pragma omp barrier

    #pragma omp for ordered
    for ( int i = 0; i < nthreads; i++ )
        #pragma omp ordered
        printf( "greetings from thread num %d\n",
                my_thread_id );
}
```

In this third implementation, we use the clause `ordered` which force a work-sharing loop to be executed in the order of the loop iteration. `ordered` may also be a stand-alone directive that specifies cross-iteration dependences.

A note: without the `barrier` directive, it may happen that some threads (potentially all of them) arrive at the worksharing construct `for` with a wrong value of `nthreads`



parallel_regions/
05b_order_of_threads.c + 05c_order_of_threads.c



| Specializing execution in PR



```
#pragma omp parallel
{
    #pragma omp critical
    { code block
    }

    #pragma omp atomic
    assignment instruction

    #pragma omp single
    { code block
    }

    #pragma omp master
    { code block
    }
}
```

The `critical` directive ensures that only a thread at a time executes the block. All the threads will execute it, although in unspecified order.

Like `critical` but limited to the following single line.
The instruction can only be an assignment in the form
`x binop = expr`

Only one among the threads will enter the code block

Only the master thread will enter the code block



| Specializing execution in PR



```
#pragma omp parallel
{
    #pragma omp critical
    { code block
    }

    #pragma omp atomic
    assignment instruction

    #pragma omp single
    { code block
    }

    #pragma omp master
    { code block
    }
}
```

A barrier (in the form of queuing) is present at the entry point; no one at the end point, i.e. the threads exiting the region will continue the run.

Synchronization as in `critical` region.

Implicit synchronization at *the end*. Only one thread is inside the region, the others are waiting at the end of the region.

No explicit synchronization at all. Only the thread 0 is inside the region, the other ones can be everywhere else.



| Specializing execution in PR



TIPS

Rationale of

`#pragma omp atomic
assignment instruction`

When you deal with shared variables, ensuring that the workflow maintains the semantic correctness of your code is fundamental. For instance, the assignment

$$a += b$$

(where either `a`, `b` or both are shared) is meaningful only if the value of `a` (or `b`, or both) does not change in the middle of the instruction itself.

In fact, while a single assembly instruction is guaranteed not to be ever interrupted on the fly, we have to take into account that even a single C line translates in several asm instructions. So, the value of a shared variable could have been changed in the middle of that groups of asm instructions, breaking the correctness of the code (*).

To avoid that, it is necessary to “protect” the sensible regions.

An **atomic** assignment has, obviously, a much lower overhead than a **critical** region and must be preferred in the appropriate cases.

(*) std C and C++ also expose to the programmer a large bunch of atomic instructions



| Specializing execution in PR



TIPS

Rationale of

`#pragma omp atomic
assignment instruction`

clauses:
`read, write, update,`
`capture`

Special clauses for **atomic****read**`var = mem`causes an atomic read of the location designated by *mem* regardless of the native machine word size**write**`mem = var`causes an atomic write of the location designated by *mem* regardless of the native machine word size**update**

- `mem++, ++mem,`
- `mem--, --mem,`
- `mem binop= expr`

Causes an atomic update of the location designated by *mem* using the designated operator or intrinsic**capture**

- `val = mem++,`
- `val = ++mem,`
- `val = mem--,`
- `val = --mem,`
- `val = mem binop= expr`
- `{val = mem; mem binop= expr}`
- `{mem binop= expr; val = mem}`

Causes an atomic update of the location designated by *mem* using the designated operator or intrinsic while also capturing the original or final value of the location designated by *x* with respect to the atomic update.The original or final value of the location designated by *mem* is written in the location designated by *val*, depending on the form of the `atomic` construct, structured block, or statements, following the usual language semantics



| Specializing execution in PR



```
#pragma omp critical
{
    "A" code block
}
#pragma omp critical
{
    "B" code block
}
#pragma omp critical(my_loop)
{
    code block
}
```

Named critical regions

In OpenMP it exists a unique global `critical` section. Hence, when you define a `critical` section, it is logically considered to be part of the global one.

As a consequence *only one thread can be inside any of the unnamed `critical` sections*, which of course limits the performance when more than one region is present.

However, that can be cured by the *named regions*, defined as the last one in the code snippet here on the left.

In that example, only one thread can be either in region "A" or in region "B": so, if one thread is executing "A", another one is forced to delay entering "B" even if it would have been reading for it.

Instead, the named `critical` regions can be executed at the same time (by different threads, of course, and only by one thread at a time).



| Specializing execution in PR



```
#pragma omp critical
{
    ...
    some_assignment_to_x;
    ...
}

#pragma omp atomic
x op= value
```

critical and **atomic**

Consider that by design **atomic** protects memory regions, while **critical** protects code regions.

Hence they are not mutually exclusive: a thread may be executing the **critical** region while another may be executing the **atomic**.



| Specializing execution in PR



```
#pragma omp single  
{ ...code block... }
```

```
#pragma omp master  
{ ...code block... }
```

Is there any difference between using `single` or `master` ?

There obviously is if you're assigning some special workflow to `thread 0`.

If not, the entity of the difference depends on the implementation details.

In general, using `master` requires only a simple test on the thread id, while using `single` requires more synchronization (the `openmp` infrastructure has to keep track about what thread is in the region and so on).

In general, if you have some insight about the fact that the workload before that point may be systematically unbalanced, so that some thread will likely arrive first, using `single` is ok.
Otherwise, it may be better to use `master`.



| Work assignment

Inside a parallel region, the work can be assigned to each threads (actually, that's why PR are created..):

```
...
results[0] = heavy_work_0();
results[1] = heavy_work_1();
results[2] = heavy_work_2();
...

```



```
#pragma omp parallel
{
    if( myid % 3 == 0)
        result = heavy_work_0( );
    else if ( myid % 3 == 1 )
        result = heavy_work_1( );
    else if ( myid % 3 == 2 )
        result = heavy_work_2( );

    if ( myid < 3 )
        results[myid] = result;
}
```

dynamic extent



parallel_region/
06_assign_work.c

run with a second
argument >0 to check
about it



parallel_region/
06_assign_work.c



| Conditional creation of PR

It is possible to spawn a parallel region only if some conditions are met:

- `#pragma omp parallel if(any valid C expression)`

```
#pragma omp parallel if( amount_of_work > high )
{
    if( myid % 3 == 0)
        result = heavy_work_0( );
    else if ( myid % 3 == 1 )
        result = heavy_work_1( );
    else if ( myid % 3 == 2 )
        result = heavy_work_2( );

    if ( myid < 3 )
        results[myid] = result;
}
```

If the condition is not fulfilled, the threads are not created and only the master thread will execute the code block.



parallel_region/
06_assign_work.c

if the first argument, that determines the amount of work, is ≤ 10 , the parallel region is **not** created



What is this useful for?

TIPS

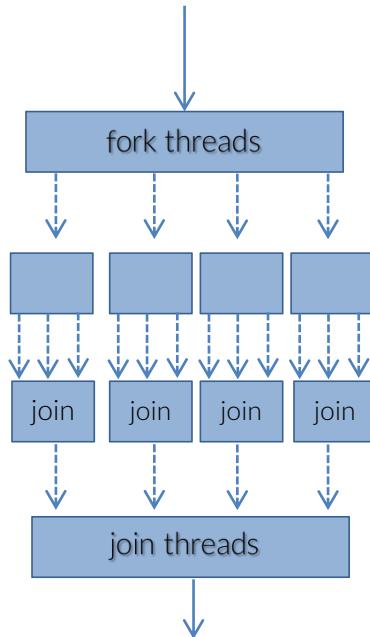
The overhead of the creation of a parallel region is of the order of $\sim 10 \mu\text{seconds}$ (that figure is dependent on the system, the compiler, the number of threads,...).

Then, you may want to create a parallel region (i.e. to spawn more threads) only if the serial execution of that code section is at least several times this overhead.



Nested Parallel Regions

NESTED PARALLELISM



Nested parallelism is explicitly_permitted in OpenMP. Whether it is *active* or not, depends on the value of some environmental variables:

`OMP_NESTED=<TRUE | FALSE>`

(* default value is FALSE)

`OMP_NUM_THREADS=N0, N1, ... >`

`OMP_MAX_ACTIVE_LEVELS=n`

Which you can change by calling the appropriate OMP_ functions

`omp_get_nested(), omp_set_nested(cond)`

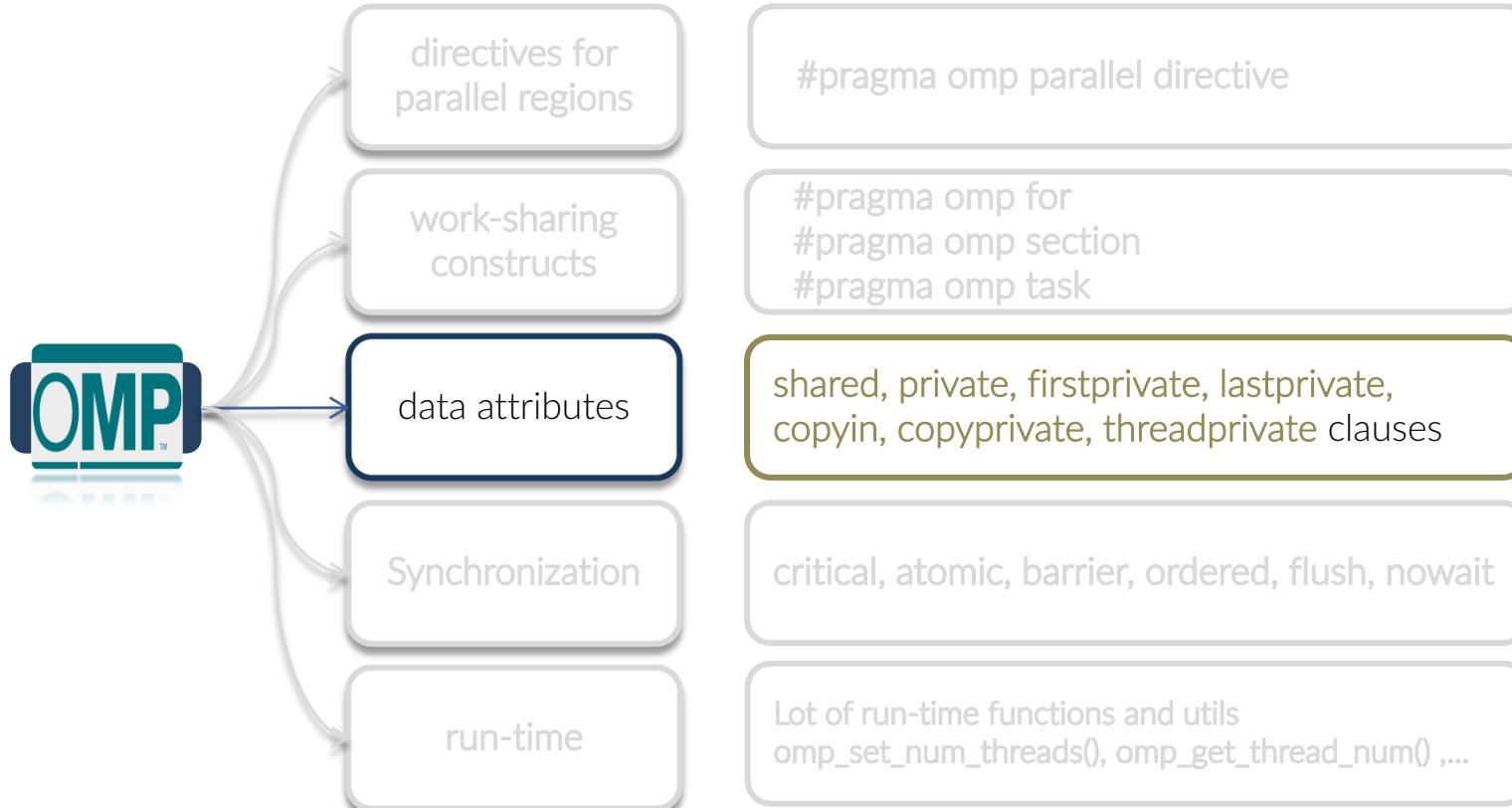
(*`OMP_NESTED` is deprecated in new OpenMP 5.0. You should use only `OMP_MAX_ACTIVE_LEVELS` and `OMP_NUM_THREADS`.

`OMP_NESTED` is in fact redundant. However, at the moment of writing the compilers still support `OMP_NESTED`.





Managing the shared/private memory





| Managing the shared/private memory

data attributes

shared, private, firstprivate, lastprivate,
copyin, copyprivate, threadprivate clauses

These clauses, that you can use at the opening of parallel regions and constructs allow a fine tuning about how to manage variables that are shared, i.e. that are formally declared in a serial region of the code.



```
int      i, j, k;  
double *array;  
  
#pragma omp parallel  
{  
    // i,j,k and array  
    // are shared here  
    ...  
}
```

| Managing the shared/private memory

The basic rule is that everything that is defined in a serial region is inherited as *shared* in a parallel region that originates from the former serial region. *Global* variables are always shared (we'll see one exception).

In this case, `i, j, k` and `array` are all *shared* within the parallel region, meaning that all the threads inside that region can concurrently access them.

To be absolutely clear:

`&i, &j, &k` and `&array` are exactly the same for all the threads.



Managing the shared/private memory

private

SHARED

```
int      i, j, k;  
double *array;  
  
#pragma omp parallel private(i,k)  
{  
    // j and array  
    // are shared here  
    // i and k are unique to each  
    // thread, they live in each  
    // thread's stack.  
    // They are different than  
    // the original j and k  
    ...  
}
```

You can specify that a list of variables(*) that exists in the local stack at the moment of the creation of the parallel region are **private** to each thread inside the parallel region.

That means that the needed space will be reserved in the threads' stack to host variables of the same types. As such, those memory regions are *different* than the original ones, although within the parallel region those variables are referred in the source code with the same names.

(*) specified as a comma-separated list of names



private

TUTORIAL

```
int      i, j, k;
double *array;
array = (double*)malloc(...);

#pragma omp parallel private(array)
{
    // now array is unique to
    // each thread and it does NOT
    // point to the region pointed
    // by the original array pointer
    ...
}
```

In this case, array is declared as *private*. As before, it means that in the stack of every thread there will be 8 bytes dedicated to host a variable referred as `array` in the source code that however is different than the original array variable in the serial region.

Then if the code tries to use it to address the memory allocated in the serial region, there will be weird things going on.
(see the example

[OpenMP/parallel_regions/00_stack_and_scope.c](#)



| Managing the shared/private memory

firstprivate

shared

```
int      i, j, k;
double *array;
array = (double*)malloc(...);

#pragma omp parallel
firstprivate(array)
{
    // now array is unique to each thread,
    // BUT each copy is initialized to the
    // value that the original array has
    // at the entry of the parallel region.
    // As such, now array can be used to
    // access the previously allocated
    // memory.

    ...
}
```

If the clause **firstprivate()** is used instead, every variable listed is private in the same sense than before, but its value is not randomic but it is **initialized** at the value that the corresponding variable in the serial region has at the moment of entering in the parallel region.



parallel_regions/
09_clauses__firstprivate.c



Managing the shared/private memory

lastprivate

```
double last_result;  
  
#pragma omp parallel  
lastprivate(last_result)  
{  
    ...  
    #pragma omp for  
    for( int j = 0; j < some_value; j++ )  
        last_result = calculation( j, ... );  
    ...  
}  
  
other_calculations( last_result, ... );  
// at this point, last_result has the last  
// value from the last iteration in the  
// for loop in the parallel region
```

parallel_regions/
09_clauses_lastprivate.c

The clause **lastprivate()** pertains only to the **for** construct.

When used, every variable listed is private in the same sense than before, and its value is not initialized.

What is affected is the value that the original variable, the one that is declared in the master thread's stack, has after the parallel region. It will have the value that the private copy of it has in the last iteration of the for loop.

we'll see in detail how the for loop works and how the work is distributed within it



Managing the shared/private memory

threadprivate

```
int      myN;
double *array;
#pragma omp threadprivate(myN, array)

#pragma omp parallel
{
    // array does exist and it's private
    // to each thread
    array = ... allocate memory...;

}
// .. something serial here..
#pragma omp parallel
{
    // array does exist here and
    // the allocated memory is available
}
```



parallel_regions/
09_clauses_threadprivate.c

The clause **threadprivate** applies to global variables and has a global scope.

threadprivate variables are private variables that do exist all along the lifetime of the process.

i.e. they are private variables that do not die in between of two different parallel regions.

note: when using `threadprivate`, the dynamic thread creation is not allowed, i.e. the number of threads in each parallel region is constant.



```
int golden_values[3];
#pragma omp threadprivate(golden_values)

for( int i = 0; i < N; i++ )
{
    get_golden_values();
    #pragma omp parallel copyin(golden_values)
    {
        // each thread uses golden_values[]
    }
}
```



parallel_regions/
09_clauses_copyin.c, 09_clauses_copyin_clarify.c



Managing the shared/private memory

The clause **copyin()** applies to parallel and worksharing (e.g. `for`) constructs.

This clause basically provides a way to perform a **broadcast of threadprivate variables** from the master thread (i.e. the thread 0) to the corresponding `threadprivate` variables of other threads.

The copying happens at the creation of the region, before the associated structured block is executed.



```
#pragma omp parallel
{
    double seed[2];

#pragma omp single copyprivate(seed)
{
    // something happens here and the
    // thread that executes this block
    // initializes the seed[2] array
}

// at this point the values of the seed[2]
// array have been propagated to all the
// other threads
}
```

parallel_regions/
09_clauses_copyprivate.c

Managing the shared/private memory

The clause **copyprivate()** applies only to **single** construct.

This clause provides a mechanism to propagate the values of private variables, including threadprivate, from a thread to the others inside a parallel region.

The copying is ultimated before any threads leave the implicit barrier at then end of the construct.



You can configure the OpenMP behaviour through Internal Control Variables (ICV). Those are “conceptual” variables, meaning that you, as a programmer, are not interested in where and how those internal variables are implemented; you are only interested in how to access and use them.

These variables have different scopes:

- *Global scope* applies to the whole program, it is fixed once at the begin (there's only one, to my knowledge)
- *Device scope* OpenMP, since v4.0, can deploy threads on heterogeneous platforms (CPUs, GPUs, co-processors, ...); some ICV are naturally different on different devices
- *Task scope* most ICVs have a *local* scope, i.e. each OMP task^() holds its own values and can change them; when new tasks are created with either a `parallel` or `task` directives, they inherit the ICVs. If a task modifies the ICVs, this has effect only during its lifetime and can not be backward-propagated.

()a “task” is in general a target to be executed by a thread: for instance, a section of a for loop scheduled for a thread is, in this sense, a “task”.



The OpenMP environmental variables



ICV	Scope	Env. Variable	Accessibility from src	
<i>nthreads</i>	TASK	OMP_NUM_THREADS	get / set	Affect parallel region
<i>nested</i>	TASK	OMP_NESTED	get / set	
<i>max-active-levels</i>	TASK	OMP_MAX_ACTIVE_LEVELS	get / set	
<i>active-level</i>	TASK	-	get	
<i>dynamic</i>	TASK	OMP_DYNAMIC	get / set	
<i>stacksize</i>	DEVICE	OMP_STACKSIZE	-	Affect program execution
<i>threads-limit</i>	TASK	OMP_THREAD_LIMIT	get	
<i>waiting-policy</i>	DEVICE	OMP_WAIT_POLICY	-	
<i>binding</i>	TASK	OMP_PROC_BIND	set	
<i>placement</i>	TASK	OMP_PLACES	-	
<i>cancellation</i>	GLOBAL	OMP_CANCELLATION	get	
<i>default-device</i>	TASK	OMP_DEFAULT_DEVICE	get / set	
<i>run-schedule</i>	DEVICE	OMP_SCHEDULE	get / set	Affect auto-parallelization of loops
<i>default-schedule</i>	DEVICE	-	-	



Lab exercise / 1

I propose you to write a code here and now, in the next 20 minutes, using what you know on OpenMP up to know and drawing from the codes that you find in the examples that we have discussed.

Please, ask about any doubt you may have; let's make this an interactive lab.

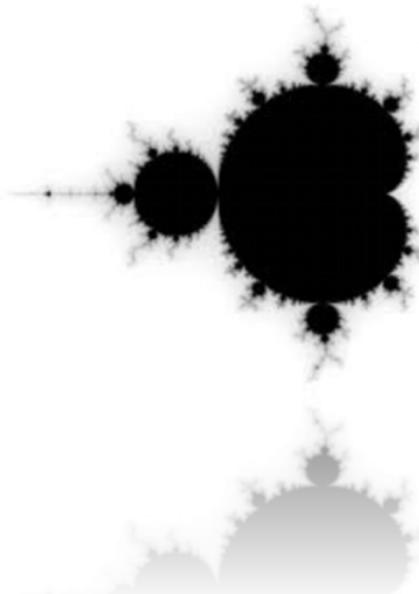


`parallel_regions/
10_lab_exercise.c, 10_lab_exercise.v2.c`

*let's check together the code
that is your starting point*



Lab exercise / 2



Let's aim to a more ambitious and, likely, funny, exercise.

Write a code that calculates the Mandelbrot set that, as you may know, is a fractal generated by the very simple equation

$$f_c(z) = z^2 + c$$

in the complex plane.

You find the details in

[OpenMP/parallel_regions/11_lab_exercise.md](#)

that's all, have fun

"So long
and thanks
forall the fish"