Parallelism (PAR) Short tutorial on OpenMP 5.0 tasking

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Part I

OpenMP Basics



Outline

- OpenMP overview
- OpenMP model
- Creating threads and accessing data
- Some API calls
- The single construct
- Thread synchronization
- Memory consistency



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What is OpenMP?

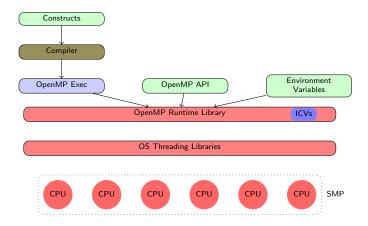
- ► It's an API extension to the C, C++ and Fortran languages to write parallel programs for shared memory machines
 - ► Current version is 5.0 (November 2018)
 - Supported by most compiler vendors and implementors
 - ► Intel, IBM, PGI, GCC, LLVM, ...
- Maintained by the Architecture Review Board (ARB), a consortium of industry and academia
- ► This mini-tutorial just covers part of the specification, for the complete reference please consult the documentation online

http://www.openmp.org



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OpenMP components





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OpenMP components

Constructs

These form the major elements of OpenMP programming

- Create threads and tasks
- Share the work amongst threads and accelerators (not covered in this mini-tutorial)
- Synchronize threads and memory, and wait for termination of tasks

Library routines

To control and query the parallel execution environment (internal control variables - ICVs)

Environment variables

The execution environment can also be set before the program execution is started



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OpenMP directives syntax

Through a compiler directive:

 OpenMP syntax is ignored if the compiler does not have the appropriate compilation flag activated

Structured block

Most directives apply to:

- A block of one or more statements
- One entry point, one exit point (no branching in or out allowed)



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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



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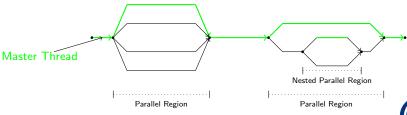


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Execution model

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





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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
 - ▶ If not ... there is a mechanism to guarantee it! (described at the end of Part I)
- Variables can be shared or private to each thread



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Directive

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

where some of the clauses are:

- ▶ num_threads(expression)
- ▶ if(expression)
- ► shared(var-list)← Coming shortly!
- ▶ private(var-list) ←
- ► firstprivate(var-list)←
- ► reduction(operator:var-list) ← We'll see it later



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Specifying the number of threads

- ► The nthreads-var ICV is used to determine the number of threads to be used for encountered parallel regions
 - It is a list of positive integer values, its first element specifying the number of processors for the next nesting level
 - When a parallel construct is encountered, and the generating task's nthreads-var list contains multiple elements, the generated task(s) inherit the value of nthreads-var as the list obtained by deletion of the first element
 - ► If the generating task's nthreads-var list contains a single element, the generated task(s) inherit that list as the value of nthreads-var



Specifying the number of threads

- ► The nthreads-var list can be defined from the execution command line by setting the OMP_NUM_THREADS environment variable
 - ► Example: setenv OMP_NUM_THREADS 4,3,2
- ► After that, the threads-var list can be modified through:
 - the omp_set_num_threads API, which sets the value of the first element of the current list
 - ▶ the num_threads clause, which causes the implementation to ignore the ICV and use the value of the clause for that region.



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The if clause

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



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

How many threads are used in each parallel region above?



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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
- By default, variables are implicitly shared



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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



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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



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```
int x=1;
#pragma omp parallel XXXXXX num_threads(2)
{
    x++;
    printf("%d\n",x);
}
printf("%d\n",x);
What does appear on the screen if XXXXXXX is shared(x),
```



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private(x) or firstprivate(x)?

Example: computation of PI

```
static long num_steps = 100000;
double step:
void main ()
   int i:
   double x, pi, sum = 0.0;
   step = 1.0/(double) num_steps;
   for (i=1; i \le num\_steps; i++){
      x = (i - 0.5) * step;
      sum = sum + 4.0/(1.0+x*x);
   pi = step * sum;
```



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Example: computation of PI (not equivalent to sequential!)

```
static long num_steps = 100000;
double step:
#include <omp.h>
#define NUM_THREADS 2
void main ()
   int i. id:
   double \times, pi, sum=0.0;
   step = 1.0/(double) num_steps;
   omp_set_num_threads(NUM_THREADS);
   #pragma omp parallel private(x, i)
   for (i=1; i \le num\_steps; i++){
      x = (i - 0.5) * step;
      sum = sum + 4.0/(1.0+x*x);
   pi = sum * step:
```



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Some useful routines

int omp_get_num_threads()

int omp_get_thread_num()

void omp_set_num_threads()

int omp_get_max_threads()

double omp_get_wtime()

Returns the number of threads in the current team. 1 if outside a parallel region

Returns the id of the thread in the current team. id between 0 and omp_get_num_threads()-1

Sets the number of threads to be used in parallel regions at the next nesting level

Returns the number of threads that could be used in parallel regions at the next nesting level

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

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Example: computation of PI (data race!)

```
static long num_steps = 100000;
double step:
#include <omp.h>
#define NUM_THREADS 2
void main ()
   int i. id:
   double \times, pi, sum=0.0;
   step = 1.0/(double) num_steps;
   omp_set_num_threads(NUM_THREADS);
   #pragma omp parallel private(x, i, id)
      id = omp_get_thread_num();
      for (i=id+1; i \le num\_steps; i=i+NUM\_THREADS) {
         x = (i - 0.5)*step;
         sum = sum + 4.0/(1.0+x*x);
   pi = sum * step:
```





Example: computation of PI (measuring execution time)

```
static long num_steps = 100000;
double step:
#include <omp.h>
#define NUM_THREADS 2
void main ()
   int i. id:
   double \times, pi, sum=0.0;
   double TimeStart, TimeEnd;
   TimeStart = omp_get_wtime();
   step = 1.0/(double) num_steps;
   omp_set_num_threads(NUM_THREADS);
   #pragma omp parallel private(x, i, id)
   pi = sum * step;
   TimeEnd = omp_get_wtime();
   printf("Wall_clock_time_=_%.20f\n", TimeEnd-TimeStart);
```



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Only one thread in the team doing the work

The single construct

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

- where clauses can be:
 - private
 - ► firstprivate
 - ▶ nowait
- Only one thread of the team executes the structured block
- ► There is an implicit barrier at the end



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The nowait clause

When single has a nowait clause then the implicit barrier at the end of the single is removed.

► This allows to overlap the execution of non-dependent work inside single with what continues after it.

```
#pragma omp single nowait
for ( i = 0; i < n ; i++ )
    v[i] = 0;
#pragma omp single

for ( i = 0; i < n ; i++ )
    a[i] = 0;</pre>
First and second single regions are independent so we can overlap them
```



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Example: computation of PI with single

```
void main ()
   #pragma omp parallel private(x, i) firstprivate(sum)
      #pragma omp single nowait
         for (i=1; i \le num\_steps; i++) {
            x = (i - 0.5) * step;
            sum = sum + 4.0/(1.0+x*x);
         pi = sum * step;
      if (sum == 0.0)
         printf("Life_is_good_when_there_is_nothing_to_do_...
```

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Why synchronization?

OpenMP is a shared memory model

- ► Threads communicate by sharing variables
- Unintended sharing of data causes race conditions (i.e. the execution outcome may change as threads are scheduled differently)
- Threads need to synchronize to impose some ordering in their sequence of actions

Some OpenMP synchronization mechanisms:

- barrier
- critical
- ► atomic
- ► Use of locks through API



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



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Barrier

Example



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Exclusive access: 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
- Multiple mutual exclusion regions by providing them with a name
 - Only those with the same name synchronize



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Example: computation of PI

```
void main ()
   int i. id:
   double \times. pi. sum = 0.0:
   step = 1.0/(double) num_steps;
   omp_set_num_threads(NUM_THREADS);
   #pragma omp parallel private(x, i, id)
      id = omp_get_thread_num();
      for (i=id+1; i \le num\_steps; i=i+NUM\_THREADS) {
         x = (i - 0.5) * step;
         #pragma omp critical
         sum = sum + 4.0/(1.0+x*x);
   pi = sum * step;
```



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Critical construct

```
int x=1,y=0;
#pragma omp parallel num_threads(4)
{
#pragma omp critical (x)
    x++;
#pragma omp critical (y)
    y++;
Different names: One thread can
update x while another updates y
```



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Exclusive access: atomic construct

```
#pragma omp atomic [update | read | write]
    expression
```

- Ensures that a specific storage location is accessed atomically, avoiding the possibility of multiple, simultaneous reading and writing threads
 - Atomic updates: x += 1, x = x foo(), x[index[i]]++
 - Atomic reads: value = *p
 - ► Atomic writes: *p = value
- Only protects the read/operation/write
- Usually more efficient than a critical construct
- Other clauses and forms for atomic are allowed in the specification



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First example: computation of PI

```
void main ()
   int i. id:
   double \times. pi. sum = 0.0:
   step = 1.0/(double) num_steps;
   omp_set_num_threads(NUM_THREADS);
   #pragma omp parallel private(x, i, id)
      id = omp_get_thread_num();
      for (i=id+1; i \le num\_steps; i=i+NUM\_THREADS) {
         x = (i - 0.5)*step;
         #pragma omp atomic
         sum = sum + 4.0/(1.0+x*x);
   pi = sum * step;
```





The reduction clause

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

```
reduction(operator:list)
```

- ► Valid operators are: $+,-,*,|,||,\&,\&\&,^n,min, max$
- ► The compiler creates a private copy of each variable in list that is properly initialized to the identity value
- ► At the end of the region, the compiler ensures that the shared variable is properly (and safely) updated with the partial values of each thread, using the specified operator



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First example: computation of PI

```
void main ()
   int i. id:
   double x, pi, sum;
   step = 1.0/(double) num_steps;
   omp_set_num_threads(NUM_THREADS);
   #pragma omp parallel private(x, i, id) reduction(+:sum)
      id = omp_get_thread_num();
      for (i=id+1; i \le num\_steps; i=i+NUM\_THREADS) {
         x = (i - 0.5) * step;
         sum = sum + 4.0/(1.0+x*x);
   pi = sum * step;
```



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Locks

OpenMP provides lock primitives for low-level synchronization

omp_init_lock Initialize the lock omp_set_lock Acquires the lock omp_unset_lock Releases the lock

omp_test_lock Tries to acquire the lock (won't block)

omp_destroy_lock Frees lock resources



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Locks

```
Example
```



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The flush construct

Relaxed consistency memory model

- ➤ A thread's temporary view of memory is not required to be consistent with memory at all times
- ► A value written to a variable can remain in the thread's temporary view until it is forced to memory at a later time
- ► Likewise, a read from a variable may retrieve the value from the thread's temporary view, unless it is forced to read from memory



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The flush construct

```
#pragma omp flush (list)
```

- ▶ It enforces consistency between the temporary view and memory for those variables in list
- Synchronization (implicit or explicit) constructs have an associated flush operation



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Task Parallelism in OpenMP



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Outline

- OpenMP tasks
- Task synchronization
- Taskloop construct
- Reductions in the tasking model



Outline

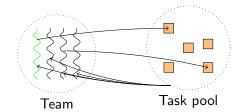
- OpenMP tasks
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Task parallelism model

- Tasks are work units whose execution may be deferred
 - they can also be executed immediately
- Threads of the team cooperate to execute them





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Creating tasks

Implicit and explicit tasks

- ► A parallel region creates tasks
 - One implicit task is created and assigned to each thread in the team of threads
- ▶ Each thread that encounters a task construct
 - Packages the code and data
 - Creates a new explicit task



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Creating (explicit) tasks

The task construct

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

Where some possible clauses are:

- shared
- ▶ private
- ► firstprivate
 - Values are captured at creation time
- ▶ if (expression)
- ► final(expression)
- ► mergeable



Example: list traversal



Example: list traversal

Completing the picture

We need threads to execute the tasks ...

... but not that many! This will generate multiple traversals



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Example: list traversal

```
Using single ...
    List I
    #pragma omp parallel
    #pragma omp single
    traverse_list(|);
    void traverse_list ( List I )
       Element e:
       for (e = I \rightarrow first; e; e = e \rightarrow next)
          #pragma omp task
          process(e);
```

One thread creates the tasks of the traversal. The rest (and this one once task generation is finished) cooperate to execute them





Default task data-sharing attributes

When no data clauses are specified, some rules apply:

- Global variables are shared
- ▶ Variables declared in the scope of a task are private
- ► The rest are firstprivate except when a shared attribute can be lexically inherited



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Task default data-sharing attributes

In practice...

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





The if clause: immediate task execution

- ▶ If 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
 - as a different task with respect to synchronization
 - ► The parent task resumes when the new task finishes
 - ► Allows implementations to optimize task creation



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The final clause: immediate task execution (nested)

- ▶ If the expression of a final clause evaluates to true
 - ► The generated task and all of its child tasks will be final
 - ► The execution of a final task is sequentially **included** in the generating task (executed immediately)
- When a mergeable clause is present on a task construct, and the generated task is an included task, the implementation may generate a merged task instead (i.e. no task and context creation for it).



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Final and mergeable tasks (data race!)

```
int fib(int n) {
  int i, j;

if (n<2)
    return n;
#pragma omp task shared(i) final(n <= THOLD) mergeable
i=fib(n-1);
#pragma omp task shared(j) final(n <= THOLD) mergeable
j=fib(n-2);
....
return i+j;
}</pre>
```



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Outline

- OpenMP tasks
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Task synchronization

There are two types of task barriers:

- taskwait
 - Suspends the current task waiting on the completion of child tasks of the current task. The taskwait construct is a stand-alone directive
- ► taskgroup
 - Suspends the current task at the end of structured block waiting on completion of child tasks of the current task and their descendent tasks



Taskwait

```
#pragma omp taskwait
```



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```
int fib(int n) {
  int i, j;
  if (n<2)
    return n:
  #pragma omp task shared(i) final(n <= THOLD) mergeable</pre>
  i=fib(n-1);
  #pragma omp task shared(i) final(n <= THOLD) mergeable</pre>
  j=fib(n-2);
  #pragma omp taskwait
  return i+i;
```



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```
#pragma omp taskgroup structured block
```



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Data sharing inside tasks

In addition one can use critical and atomic to synchronize the access to shared data inside the task

```
void process (Element e)
{
    ...
    #pragma omp atomic
    solutions_found++;
    ...
}
```



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Task dependences

Definition of dependences between sibling tasks (i.e. from the same father)

Task dependences are derived from the dependence type (in, out or inout) and its items in var_list. This list may include array sections



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Task dependences

- ➤ The in dependence-type: the generated task will be a dependent task of all previously generated sibling tasks that reference at least one of the list items in an out or inout dependence-type list
- ▶ The out and inout dependence-types: the generated task will be a dependent task of all previously generated sibling tasks that reference at least one of the list items in an in, out, or inout dependence-type list.



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Task dependences

Example: wave-front execution



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Creating (explicit) tasks from loop iterations

The taskloop construct

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

specifies that the iterations of one or more associated loops will be executed in parallel using OpenMP tasks. Implicit taskgroup synchronization associated with taskloop

Some clauses are used to specify data sharing attributes:

- ▶ shared(list)
- private(list)
- ► firstprivate(list)



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Creating (explicit) tasks from loop iterations

Other clauses to control task generation:

- grainsize(n)
- ▶ num_tasks(n)
- ► collapse(n)
- ▶ if (expression)
- ► final(expression)
- mergeable

Or to override the implicit taskgroup associated with the taskloop construct:

► nogroup



Taskloop example

```
Granularity: BS iterations per task
void vector_add(int *A, int *B, int *C, int n) {
   #pragma omp taskloop grainsize(BS)
   for (int i=0; i < n; i++)
       C[i] = A[i] + B[i];
void main() {
   #pragma omp parallel
   #pragma omp single
   ... vector_add(a, b, c, N); ...
or alternatively
#pragma omp taskloop num_tasks(n/BS)
```



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Outline

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Reduction clauses

Directives and clauses associated



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Task reductions examples (1)

Reductions with explicit tasks always occur in the environment of a taskgroup (explicit if not implicit), which delimits the scope of reduction operation:

```
#pragma omp parallel
#pragma omp single
{
    // implicit taskgroup in taskloop construct
    #pragma omp taskloop reduction(+: sum)
    for (i=0; i< SIZE; i++)
        sum += X[i];
}</pre>
```



Task reductions examples (2)

Reductions with explicit tasks always occur in the environment of a taskgroup (explicit if not implicit), which delimits the scope of reduction operation:





Parallelism (PAR) Short tutorial on OpenMP 5.0 tasking

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