

# Parallel Computing for Science & Engineering CS395T

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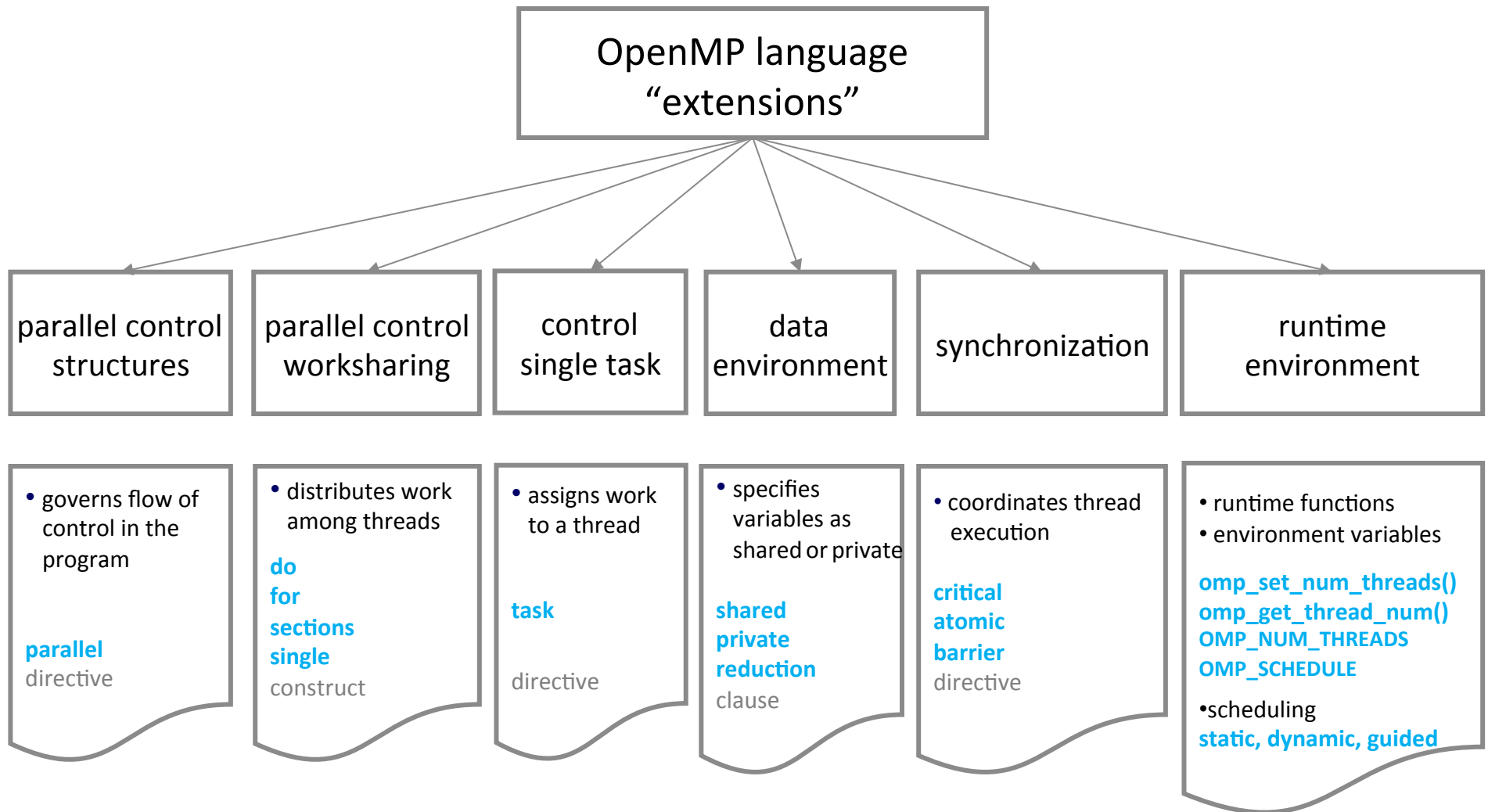
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# OpenMP Constructs



# OpenMP Worksharing -- Single

- SINGLE (or MASTER)
  - Block of code is executed only once by a single thread (or the master thread)
  - Implied barrier (ONLY single)
- Only one thread is executing the block at all!

**!\$OMP single**

```
glob_count = glob_count + 1  
print *, glob_count
```

**!\$OMP end single**

**#pragma single**

```
{  
    glob_count++;  
    printf("%d\n", glob_count);
```

```
}
```

# Mutual exclusion – atomic and critical directives

When each thread must execute a section/statement of code serially (only one thread at a time can execute it) the region must be marked with CRITICAL/ATOMIC directives.

Use the ATOMIC directive if executing only one operation.

All threads are executing the block one after another

```
!$omp parallel shared(sum,x,y)
...
!$omp critical
    call update(x)
    call update(y)
    sum=sum+1
!$omp end critical
...
!$omp end parallel
```

```
!$omp parallel
...
!$omp atomic
    sum=sum+1
...
!$omp end parallel
```

Syntax:

\$pragma omp critical [*name*]

!\$omp critical [*name*]

!\$omp end critical [*name*]

# Mutual exclusion – atomic and critical directives

When each thread must execute a section/statement of code serially (only one thread at a time can execute it) the region must be marked with CRITICAL/ATOMIC directives.

Use the ATOMIC directive if executing only one operation.

```
#pragma parallel shared(sum,x,y) {  
    ...  
    #pragma omp critical{  
        update(x) ;  
        update(y) ;  
        sum=sum+1 ;  
    }  
    ...  
}
```

```
#pragma omp parallel{  
    ...  
    #pragma omp atomic  
        sum=sum+1 ;  
    ...  
}
```

Syntax:

\$pragma omp critical [*name*]

!\$omp critical [*name*]

!\$omp end critical [*name*]

# OpenMP Synchronization -- NOWAIT

Where a work-sharing region ends, a barrier is implied - all threads must reach the barrier before any can proceed. By using the NOWAIT clause at the end of each loop inside the parallel region, an unnecessary synchronization of threads can be avoided.

```
!$omp parallel
...

!$omp do
  do i=1,n
    call work(i)
  enddo
!$omp end do nowait

!$omp do schedule(dynamic,m)
  do i=1,n
    x(i)=y(i)+z(i)
  enddo

!$omp end parallel
```

```
#pragma omp parallel
{...

#pragma omp for nowait
  for(i=0; i<n; i++)
    work(i);

#pragma omp schedule(dynamic,m)
  for(i=0; i<n; i++)
    x(i)=y(i)+z(i);

}
```

# Runtime Library API

## Functions

## Operation

<code>omp_get_max_threads</code>	from <code>OMP_NUM_THREADS</code>
<code>omp_get_num_threads()</code>	Number of Threads in team, N.
<code>omp_get_thread_num()</code>	Thread ID. {0 -> N-1}
<code>omp_get_num_procs()</code>	Number of machine hardware threads
<code>omp_in_parallel()</code>	True if in parallel region & multiple thread executing
<code>omp_set_num_threads(#)</code>	Changes Number of Threads for parallel region.

# API Example: Pseudo code

Assume: 'export OMP\_NUM\_THREADS=3'

```
print omp_num_procs()
print omp_max_threads()

omp parallel
  print omp_num_threads()
  print omp_get_thread_num()
omp end parallel

omp_set_num_threads(6)
print omp_num_procs()
print omp_max_threads()

omp parallel num_threads(2)
  print omp_num_threads()
  print omp_get_thread_num()
omp end parallel
```

16  
3

3 printed 3 times  
0, 1, or 2 each printed once

?  
?

?, printed how often?  
?



# API example

Or can use OPENMP pragma (Fortran: compile with -fpp)

```
#include <omp.h>
```

```
#pragma omp parallel private(i,n)
{
  i = omp_get_thread_num( );
  n = omp_get_num_threads( );
}
```

```
use omp_lib
```

```
!$omp parallel private(i,n)
```

```
    i = omp_get_thread_num( )
    n = omp_get_num_threads( )
```

```
!$omp end parallel
```

**What happens if you don't compile with `-qopenmp`?**

# OpenMP Conditional Compilation

Fortran triggers: !\$

```
i=0; n=1
!$omp parallel private(i,n)
!$  i = omp_get_thread_num( )
!$  n = omp_get_num_threads( )
      call sub(i,n)
!$omp end parallel
```

Or can use OPENMP pragma (Fortran: compile with -fpp)

```
#include <omp.h>
i=0; n=1;
#pragma omp parallel private(i,n)
{
  #ifdef OPENMP
    i = omp_get_thread_num( ) ;
    n = omp_get_num_threads( );
  #endif
  sub(i,n);
}
```

```
use omp_lib
i=0; n=1
!$omp parallel private(i,n)

  #ifdef OPENMP
    i = omp_get_thread_num( )
    n = omp_get_num_threads( )
  #endif
  call sub(i,n)
!$omp end parallel
```

# Runtime Library API

## API Environment Variables

<b>OMP_NUM_THREADS</b>	Set to No. of Threads
<b>OMP_SCHEDULE</b>	Schedule for the run-time clause

### Example

```
export OMP_NUM_THREADS=4  
export OMP_SCHEDULE='static,4'  
./a.out
```

# References

- Some material identical to:  
[www.chpc.utah.edu/attachments/20110112.05/IntroOpenMP06.pdf](http://www.chpc.utah.edu/attachments/20110112.05/IntroOpenMP06.pdf)
- This one is a real tutorial and even has test modules:  
<http://www.citutor.org/login.php>
- The sites  
[www.llnl.gov/computing/tutorials/openMP/](http://www.llnl.gov/computing/tutorials/openMP/)  
[www.nersc.gov/assets/Uploads/XE62011OpenMP.pdf](http://www.nersc.gov/assets/Uploads/XE62011OpenMP.pdf)  
have good reference/tutorial pages for OpenMP.

# Examples

```
export OMP_NUM_THREADS=2  
./a.out
```

```
x and y : arrays with 2 elements  
x(1) = 2.; x(2) = 4.  
j = 0
```

```
!$omp parallel private(i)  
!$omp do  
do i=1 , 2  
    j = j + 1  
    y(i) = x(j)  
enddo  
!$omp end parallel
```

What values in x(1) and x(2)?

# Splitting dependences out of loop

## Loop fission

```
for(i=1; i<n; i++){  
    b[i] = a[i]*r[i]  
    a[i-1] = t[i-1]*s[i-1]  
}
```



```
for(i=1; i<n; i++){  
    b[i] = a[i]*r[i]  
}  
for(i=1; i<n; i++){  
    a[i-1] = t[i-1]*s[i-1]  
}
```

C/C++

```
do i=2,n  
    b(i) = a(i)*r(i)  
    a(i-1) = s(t-1)*t(t-1)  
enddo
```



```
do i=2,n  
    b(i) = a(i)*r(i)  
enddo  
do i=2,n  
    a(i-1) = s(t-1)*t(t-1)  
enddo
```

F90

# Examples

```
export OMP_NUM_THREADS=2  
./a.out
```

```
x : array with 2 elements  
x(1) = 2.; x(2) = 4.  
sum = 0
```

```
!$omp parallel  
!$omp do  
do i=1 , 2  
    sum = sum + x(i)  
enddo  
!$omp end parallel
```

# Examples

```
export OMP_NUM_THREADS=4  
./a.out
```

```
x : array with 8 elements  
'all elements set to 1'  
sum = 0
```

```
!$omp parallel reduction(+:sum)  
!$omp do  
do i=1 , 8  
    sum = sum + x(i)  
enddo  
!$omp end parallel
```



# Examples

```
export OMP_NUM_THREADS=4  
./a.out
```

```
x : array with 8 elements  
'all elements set to 1'  
sum = 0
```

```
!$omp parallel reduction(+:sum)  
!$omp do  
do i=1 , 8  
    sum = sum + x(i) * y(i) + fct(...)  
enddo  
!$omp end parallel
```

# Examples

```
export OMP_NUM_THREADS=4  
./a.out
```

```
x(4,4) : array with 4x4 elements
```

```
!$omp parallel  
!$omp do schedule(static,1)  
do i=1, 4  
  do j=1, i  
    x(i,j) = x(i,j) + 1.  
  enddo  
enddo nowait
```

```
!$omp do schedule(static,1)  
do i=1, 4  
  ...  
enddo
```