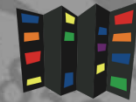


Introduction to OpenMP

Martin Čuma

*Center for High Performance
Computing University of Utah
m.cuma@utah.edu*

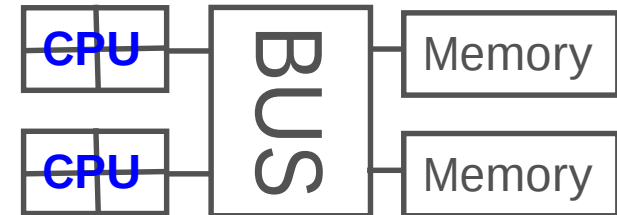


- Quick introduction.
- Parallel loops.
- Parallel loop directives.
- Parallel sections.
- Some more advanced directives.
- Summary.
- Survey

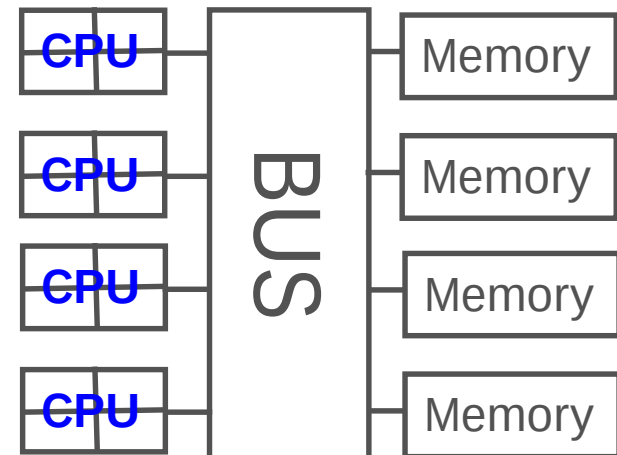
<https://www.surveymonkey.com/r/8HLJ3WK>

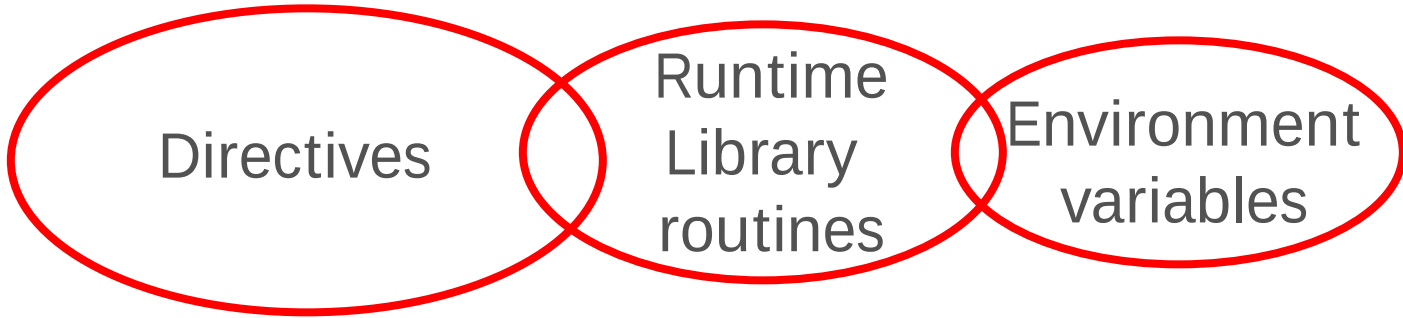
- All processors have access to local memory
- Simpler programming
- Concurrent memory access
- More specialized hardware
- CHPC :
Linux clusters 12 - 28 core nodes

Dual quad-core node



Many-core node (e.g. SGI)





Directives

Runtime
Library
routines

Environment
variables

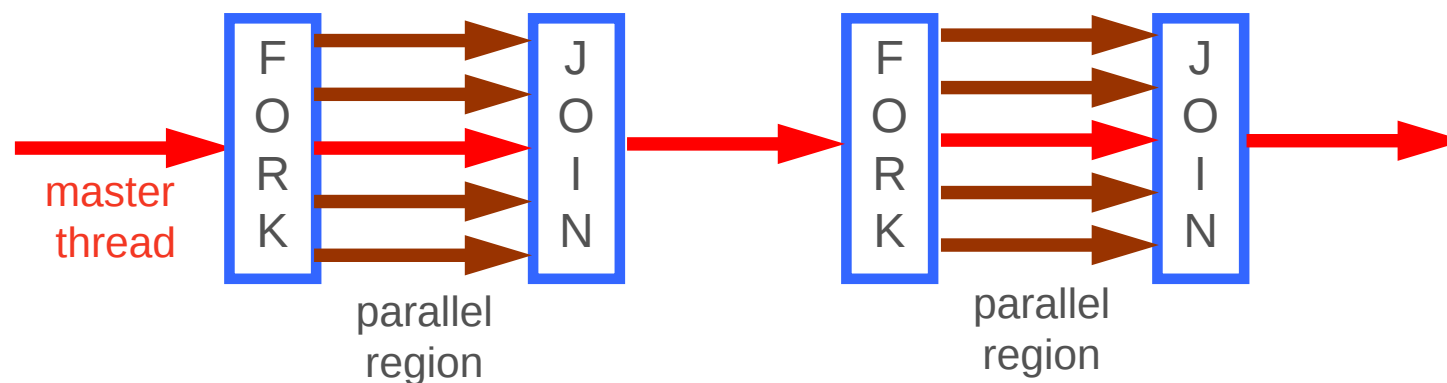
- Compiler directives to parallelize
 - Fortran – source code comments

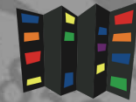
```
!$omp parallel/!$omp end parallel
```
 - C/C++ - #pragmas

```
#pragma omp parallel
```
- Small set of subroutines, environment variables

```
!$  iam = omp_get_num_threads()
```

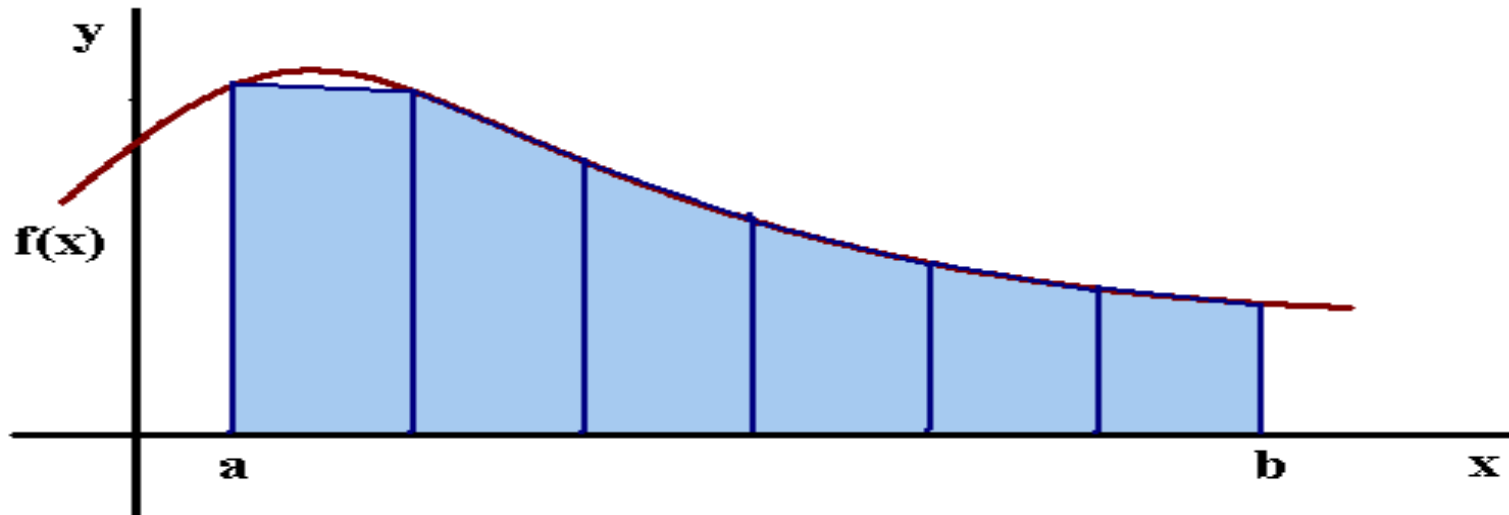
- Shared memory, thread based parallelism
- Explicit parallelism
- Nested parallelism support
- Fork-join model





$$\int_a^b f(x) \approx \sum_{i=1}^n \frac{1}{2} h [f(x_{i-1}) + f(x_i)] =$$

$$\frac{1}{2} h [f(x_0) + f(x_n)] + \sum_{i=1}^{n-1} h [f(x_i)]$$



```
program trapezoid
```

```
integer n, i
```

```
double precision a, b, h, x, integ, f
```

1.

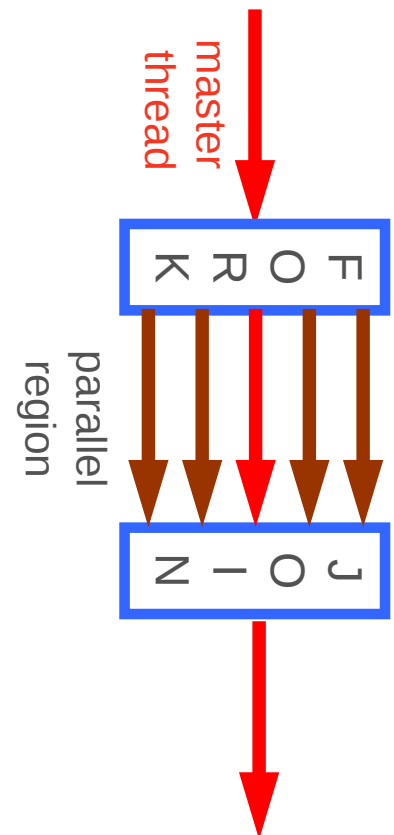
```
print*, "Input integ. interval, no. of trap:"  
read(*,*)a, b, n  
h = (b-a)/n  
integ = 0.
```

2.

```
!$omp parallel do reduction(+:integ) private(x)  
do i=1,n-1  
x = a+i*h  
integ = integ + f(x)  
enddo
```

3.

```
integ = integ + (f(a)+f(b))/2.  
integ = integ*h  
print*, "Total integral = ", integ  
end
```



```
em001:>%module load pgi
em001:>%pgf77 -mp=numa trap.f -o trap
em001:>%setenv OMP_NUM_THREADS 12
em001:>%trap
Input integ. interval, no. of trap:
0 10 100
Total integral =      333.350000000000001
```



- Fortran

```
!$omp parallel do [clause [, clause]]  
[!$omp end parallel do]
```

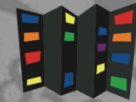
- C/C++

```
#pragma omp parallel for [clause [clause]]
```

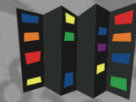
- Loops must have precisely determined *trip count*
 - no do-while loops
 - no change to loop indices, bounds inside loop (C)
 - no jumps out of the loop (Fortran – exit, goto; C – break, goto)
 - cycle (Fortran), continue (C) are allowed
 - stop (Fortran), exit (C) are allowed



- Control execution of parallel loop
 - `scope (shared, private)`
sharing of variables among the threads
 - `if`
whether to run in parallel or in serial
 - `schedule`
distribution of work across the threads
 - `collapse(n)`
combine nested loops into a single loop for better parallelism
 - `ordered`
perform loop in certain order
 - `copyin`
initialize private variables in the loop



- `private` – each thread creates a private instance
- not initialized upon entry to parallel region
undefined upon exit from parallel region
- default for loop indices, variables declared inside parallel loop
- `shared` – all threads share one copy
- update modifies data for all other threads
- default everything else
- Changing default behavior
- `default (shared | private | none)`



- `firstprivate/lastprivate` clause
 - initialization of a private variable
`!$omp parallel do firstprivate(x)`
 - finalization of a private variable
`!$omp parallel do lastprivate(x)`
- `reduction` clause
 - performs global operation on a variable
`!$omp parallel do reduction (+ : sum)`



- Anti-dependence

race between statement S_1 writing and S_2 reading

- removal: **privatization**, multiple do loops

- Output dependence

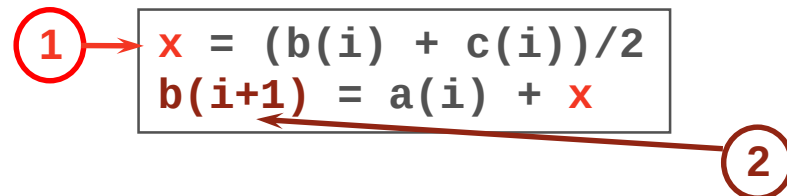
values from the last iteration used outside the loop

- removal: `lastprivate` clause

- Flow dependence

data at one iteration depend on data from another iteration

- removal: reduction, rearrangement, often impossible



- Serial trapezoidal rule

```
integ = 0.
do i=1,n-1
  x = a+i*h
  integ = integ + f(x)
enddo
```

- Parallel solution

```
integ = 0.
```

```
!$omp parallel do private(x) reduction (+:integ)
```

```
do i=1,n-1
  x = a+i*h
  integ = integ + f(x)
enddo
```

Thread 1

$x = a + i * h$

$integ = integ + f(x)$

Thread 2

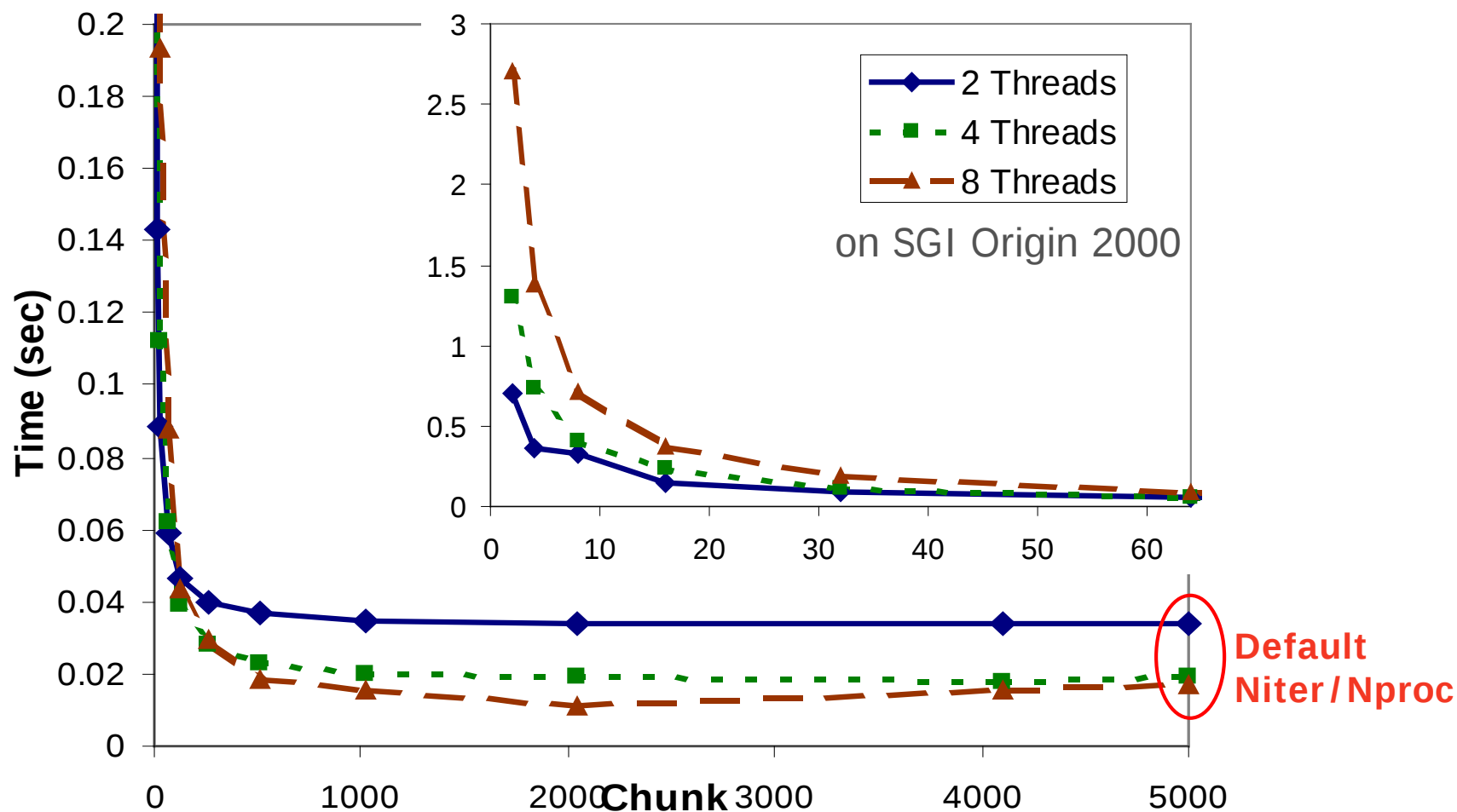
$x = a + i * h$

$integ = integ + f(x)$

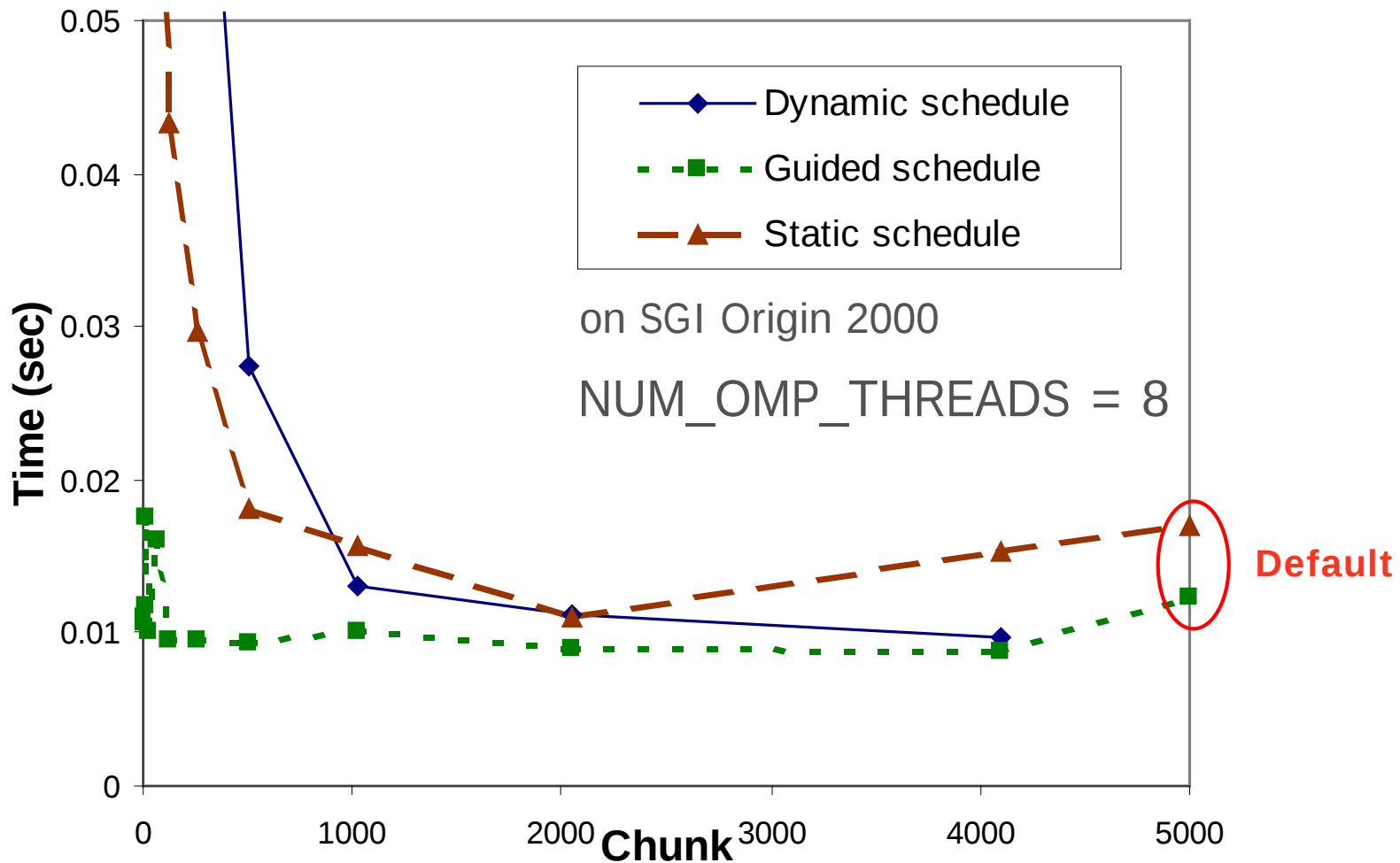
- Threads distribute work
- Need to collect work at the end
 - sum up total
 - find minimum or maximum
- Reduction clause – global operation on a variable
 - `!$omp parallel do reduction(+:var)`
 - `#pragma omp parallel for reduction(+:var)`
- Allowed operations - commutative
 - +, *, max, min, logical

- Parallelization costs CPU time
 - Nested loops
 - parallelize the outermost loop
 - `if` clause
 - parallelize only when it is worth it – above certain number of iterations:
- ```
!$omp parallel do if (n .ge. 800)
 do i = 1, n
 . . .
 enddo
```

- user-defined work distribution  
  `schedule (type[, chunk])`
- chunk – number of iterations contiguously  
  assigned to threads
- type
  - `static` – each thread gets a constant chunk
  - `dynamic` – work distribution to threads varies
  - `guided` – chunk size exponentially decreases
  - `runtime` – schedule decided at the run time

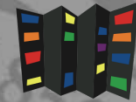


# Different schedule timings



# Example 2

## MPI-like parallelization



```
#include <stdio.h>
#include "omp.h"
#define min(a,b) ((a) < (b) ? (a) : (b))
```

```
int istart,iend;
```

1. 

```
#pragma omp threadprivate(istart,iend)
```

```
int main (int argc, char* argv[]){
int n,nthreads,iam,chunk; float a, b;
double h, integ, p_integ;
double f(double x);
double get_integ(double a, double h);
```

2. 

```
printf("Input integ. interval, no. of trap:\n");
scanf("%f %f %d",&a,&b,&n);
h = (b-a)/n;
integ = 0.;
```

```
3. #pragma omp parallel shared(integ)
 private(p_integ,nthreads,iam,chunk){
 nthreads = omp_get_num_threads();
 iam = omp_get_thread_num();
 chunk = (n + nthreads -1)/nthreads;
 istart = iam * chunk + 1;
 iend = min((iam+1)*chunk+1,n);

4. p_integ = get_integ(a,h);

5. #pragma omp critical
 integ += p_integ;
 }

6. integ += (f(a)+f(b))/2.;
 integ *= h;
 printf("Total integral = %f\n",integ);
 return 0;}
```

```
double get_integ(double a, double h)
{
 int i;
 double sum, x;

 sum = 0;
 for (i=istart; i<iend; i++)
 {
 x = a+i*h;
 sum += f(x);
 }
 return sum;
}
```

- Fortran

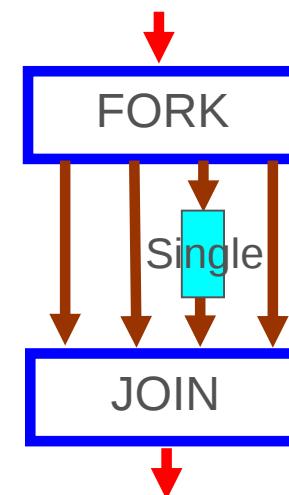
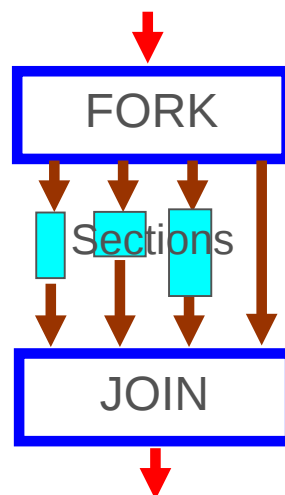
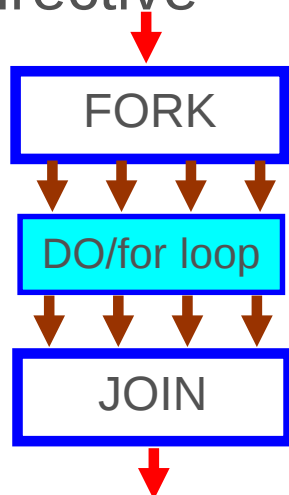
```
!$omp parallel ... !$omp end parallel
```

- C/C++

```
#pragma omp parallel
```

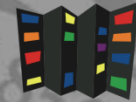
- SPMD parallelism – replicated execution
- must be a self-contained block of code – 1 entry, 1 exit
- implicit barrier at the end of parallel region
- can use the same clauses as in **parallel do/for**

- DO/for loop – distributes loop - do directive
- Sections – breaks work into separate, discrete sections - section directive
- Workshare – parallel execution of separate units of work - workshare directive
- Single/master – serialized section of code - single directive

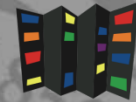




- Restrictions:
  - continuous block; no nesting
  - all threads must reach the same construct
  - constructs can be outside lexical scope of the parallel construct (e.g. subroutine)

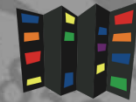


- Used to parallelize irregular, recursive algorithms
- All tasks run independent of each other in parallel, on up to `OMP_NUM_THREADS`
- Use `taskwait` to wait for all tasks to finish
- Each task has its own data space – use `mergeable` for shared variables to reduce storage needs
- Use `depend` to specify data dependencies
- Often started from `serial` section



- Calculate Fibonacci number using recursion

```
int fib(int n) {
 int i, j;
 if (n<2) return n;
 else {
 #pragma omp task shared(i)
 i=fib(n-1);
 #pragma omp task shared(j)
 j=fib(n-2);
 #pragma omp taskwait // wait for completion of child tasks
 return i+j;
 }
 #pragma omp parallel {
 fibn = fib(n); }
}
```



- global/common block variables are private only in lexical scope of the parallel region
  - possible solutions
    - pass private variables as function arguments
    - use `threadprivate` – identifies common block/global variable as private
    - `!$omp threadprivate (/cb/ [, /cb/] ...)`  
`#pragma omp threadprivate (list)`
    - use `copyin` clause to initialize the `threadprivate` variable
- e.g. `!$omp parallel copyin(istart,iend)`

- `critical` section
    - limit access to the part of the code to one thread at the time
- ```
!$omp critical [name]  
...  
!$omp end critical [name]
```
- `atomic` section
 - atomically updating single memory location
- ```
sum += x
```
- runtime library functions

- thread set/inquiry

`omp_set_num_threads(integer)`

`OMP_NUM_THREADS`

`integer omp_get_num_threads()`

`integer omp_get_max_threads()`

`integer omp_get_thread_num()`

- set/query dynamic thread adjustment

`omp_set_dynamic(logical)`

`OMP_DYNAMIC`

`logical omp_get_dynamic()`

- lock/unlock functions

`omp_init_lock()`

`omp_set_lock()`

`omp_unset_lock()`

`logical omp_test_lock()`

`omp_destroy_lock()`

- other

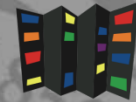
`integer omp_get_num_procs()`

`logical omp_in_parallel()`

`OMP_SCHEDULE`

- `barrier - !$omp barrier`
  - synchronizes all threads at that point
- `ordered - !$omp ordered`
  - imposes order across iterations of a parallel loop
- `master - !$omp master`
  - sets block of code to be executed only on the master thread
- `flush - !$omp flush`
  - synchronizes memory and cache on all threads

- nested parallel loops
- accelerator support (4.0)
- user defined reduction (4.0)
- thread affinity (4.0)



- parallel do/for loops
  - variable scope, reduction
  - parallel overhead, loop scheduling
- parallel regions
  - mutual exclusion
  - work sharing, tasking
  - synchronization

[http://www.chpc.utah.edu/short\\_courses/intro\\_openmp](http://www.chpc.utah.edu/short_courses/intro_openmp)

- Web

<http://www.openmp.org/>

<https://computing.llnl.gov/tutorials/openMP>

[http://openmp.org/mp-documents/OpenMP\\_Examples\\_4.0.1.pdf](http://openmp.org/mp-documents/OpenMP_Examples_4.0.1.pdf)

- Books

Chapman, Jost, van der Pas – Using OpenMP

Pacheco – Introduction to Parallel Computing

Survey <https://www.surveymonkey.com/r/8HLJ3WK>