

Parallel Programming in OpenMP

Outline

- ❑ OpenMP Overview
- ❑ OpenMP
 - ❑ Basics
 - ❑ Syntax: Directives
 - ❑ Syntax: Clauses
 - ❑ Example

OpenMP – Overview

What is OpenMP?

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OpenMP – Overview

- ❑ OpenMP: stands for
 - Open Multi Processing
- ❑ parallel programming model for shared memory multiprocessors
- ❑ 'de-facto' standard, not an industry *standard*
- ❑ not a new language, but
 - ❑ compiler directives
 - ❑ support function library

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OpenMP – Overview

- ❑ OpenMP development is community driven
- ❑ Architecture Review Board (ARB):
 - ❑ hardware and software vendors
 - ❑ government and academia
- ❑ Official OpenMP website:
 - ❑ <https://www.openmp.org/>

OpenMP – Overview

- ❑ standard versions:
 - ❑ C/C++: version 2.0 (March 2002)
 - ❑ Fortran: version 2.0 (November 2000)
 - ❑ version 2.5 (Fortran and C/C++ / May 2005)
 - ❑ version 3.0 (May 2008)
 - ❑ version 3.1 (July 2011)
 - ❑ ...

OpenMP – Overview

- ❑ 4.0 (Jul 2013)
 - ❑ exception handling
- ❑ 4.5 – update (Nov 2015)
 - ❑ support for accelerators (GPUs, Xeon Phi)
- ❑ 5.0 (Nov 2018)
 - ❑ is implemented in most compilers now
- ❑ 5.1 – update (Nov 2020)
- ❑ The OpenMP standard specifications:
 - ❑ <http://www.openmp.org/specifications/>

OpenMP – Overview

OpenMP Literature:

- ❑ “Using OpenMP – The Next Step“ by R. van der Pas, E. Stotzer and C. Terboven, MIT Press (2017) – available via findit.dtu.dk
- ❑ “The OpenMP Common Core” by T.G. Mattson, Y. He, and A.E. Koniges, MIT Press (2019)
- ❑ “Using OpenMP”, B. Chapman, G. Jost, R. van der Pas, MIT Press (2007)

OpenMP Basics

Basic elements of OpenMP

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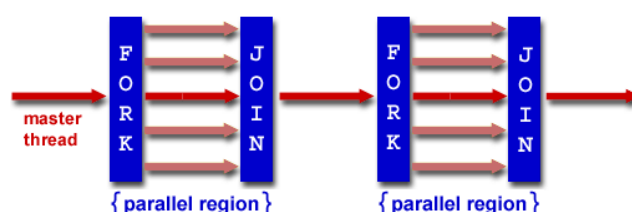
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OpenMP Basics

OpenMP uses the “Fork-Join Model”:

- ❑ All programs begin as a single process: the **master thread**.
- ❑ FORK: the master thread creates a team of parallel threads (**parallel region**).
- ❑ JOIN: synchronization and termination of the **worker threads**.



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OpenMP Basics

OpenMP is mostly based on compiler directives:

- ❑ C/C++:

```
#pragma omp directive [clause]
    {...code block...}
```

- ❑ Fortran:

```
!$OMP directive [clause]
    ...code block...
!$OMP end directive
```

OpenMP Basics

The OpenMP API has also

- ❑ a set of support library functions:

`omp_... ()`

e.g. `omp_get_thread_num()`

- ❑ control via environment variables:

`OMP_...`

e.g. `OMP_NUM_THREADS`

OpenMP Basics

First OpenMP version of “Hello world”:

```
#include <stdio.h>

int main(int argc, char *argv[]) {
    #pragma omp parallel
    {
        printf("Hello world!\n");
    } /* end parallel */
    return(0);
}
```

OpenMP Basics

Second version of “Hello world”:

```
#include <stdio.h>
#ifdef _OPENMP
#include <omp.h>
#endif

int main(int argc, char *argv[]) {
    int t_id = 0;
    #pragma omp parallel private(t_id)
    {
        #ifdef _OPENMP
        t_id = omp_get_thread_num();
        #endif
        printf("Hello world from %d!\n", t_id);
    } /* end parallel */
    return(0);
}
```

OpenMP Basics

```
$ ./hello2
Hello world from 0!
```

```
$ OMP_NUM_THREADS=4 ./hello2
Hello world from 0!
Hello world from 3!
Hello world from 1!
Hello world from 2!
```

- ❑ Note: The order of execution will be different from run to run!
- ❑ The default no. of threads depends on the OpenMP implementation

OpenMP Components

Directives

Environment variables

Runtime

- | | | |
|--------------------|-----------------------------|-----------------------------|
| ❑ Parallel regions | ❑ no. of threads | ❑ no. of threads |
| ❑ Worksharing | ❑ Scheduling | ❑ Scheduling |
| ❑ Synchronization | ❑ Dynamic thread adjustment | ❑ Dynamic thread adjustment |
| ❑ Data scoping | ❑ Nested parallelism | ❑ Nested parallelism |
| ❑ no. of threads | | ❑ API for timers & locking |
| ❑ Orphaning | | |

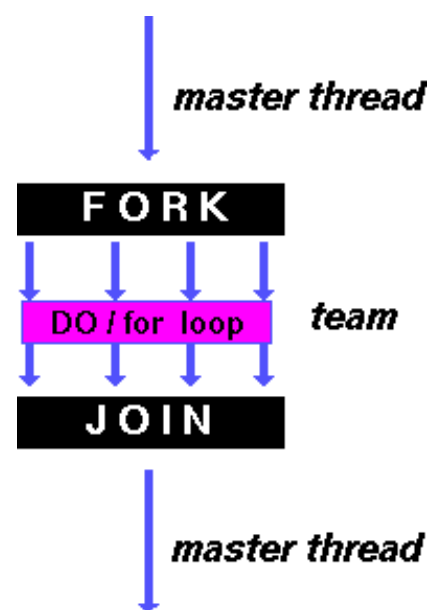
OpenMP Basics

Work-sharing

OpenMP Basics

Work-sharing constructs – 1:

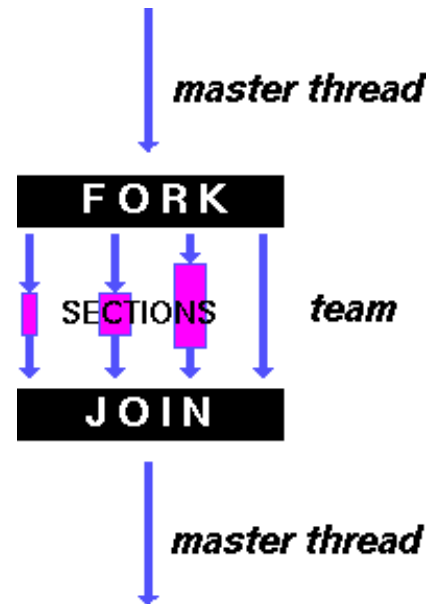
- ❑ **do/for**
- ❑ loop parallelism
- ❑ most common



OpenMP Basics

Work-sharing constructs – 2:

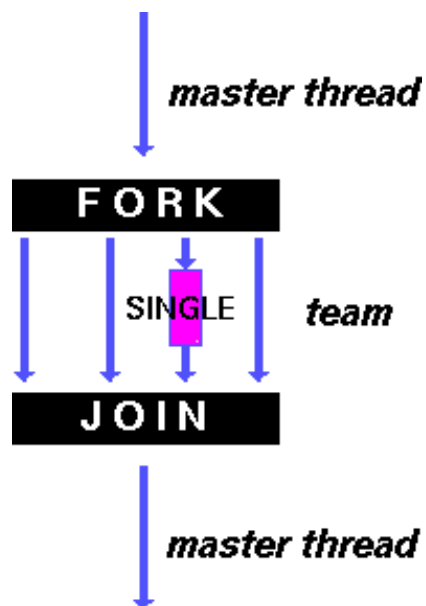
- sections
- functional parallelism
- typically independent calculations



OpenMP Basics

Work-sharing constructs – 3:

- single
- work assigned to one thread only
- typically I/O



OpenMP Basics

Important rules for work-sharing constructs:

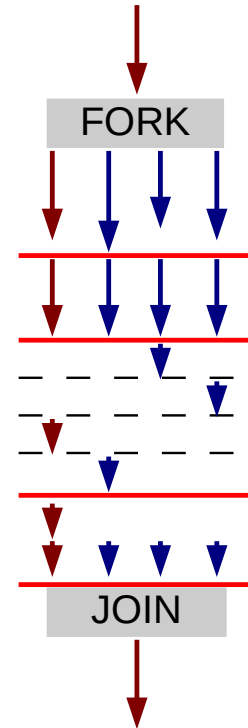
- ❑ must be enclosed in a parallel region
- ❑ must be encountered by all team members
- ❑ must be encountered in the same order

OpenMP Basics

Synchronization

OpenMP Basics

- ❑ most synchronization in OpenMP is implicit, but sometimes explicit synchronization is needed:
- ❑ barriers
- ❑ critical regions
- ❑ master only
- ❑ explicit locking



OpenMP Syntax

OpenMP programming in C/C++ Part I: Directives

OpenMP Syntax

OpenMP directives – general form:

```
#pragma omp directive [[clause] \  
  [clause] ...]  
{  
  <statements>  
} /* end of omp directive */
```

- ❑ **Note:** There is **no** “omp end” pragma!
- ❑ Best practice: add a comment at the end of the structured block!

OpenMP Syntax

Parallel region:

- ❑ Starts a team of parallel threads
- ❑ executes code in parallel
- ❑ synchronize/terminate threads

```
main() {  
  A();  
  #pragma omp parallel  
  {  
    B();    // all threads do B()!  
  } /* end omp parallel */  
  C();  
}
```

OpenMP Syntax

Work-sharing – Loop parallelism:

- ❑ OpenMP implements parallel do/for-loops only!

```
int i;
float a[N], b[N], c[N];

for (i=0; i < N; i++)
    a[i] = b[i] = i * 1.0;

#pragma omp parallel shared(a,b,c) private(i)
{
    #pragma omp for
    for (i=0; i < N; i++)
        c[i] = a[i] + b[i];
} /* end of parallel region */
```

*for has to follow
the pragma – no
{... }!*

OpenMP Syntax

Work-sharing – Loop parallelism:

- ❑ Another version: combined “parallel for”

```
int i;
float a[N], b[N], c[N];

for (i=0; i < N; i++)
    a[i] = b[i] = i * 1.0;

#pragma omp parallel for shared(a,b,c) \
    private(i)
for (i=0; i < N; i++)
    c[i] = a[i] + b[i];
```

OpenMP Syntax

Work-sharing – Fortran 95 array syntax

- ❑ Fortran 95 allows to address parts of or whole arrays – and the compiler will translate this into loops.
- ❑ A special Fortran directive:

```
double precision, dimension() :: A, B
double precision, dimension(N,M) :: C

!$ OMP WORKSHARE
...
A(1:M) = A(1:M) + B(1:M)
C = 0.00
...
!$ OMP END WORKSHARE
```

OpenMP Syntax

Work-sharing – Functional parallelism:

- ❑ Parallel sections:

```
#pragma omp parallel shared(a,b,c) private(i)
{
    #pragma omp sections
    {
        #pragma omp section
        for (i=0; i < N/2; i++)
            c[i] = a[i] + b[i];

        #pragma omp section
        for (i=N/2; i < N; i++)
            c[i] = a[i] + b[i];
    } /* end of sections */
} /* end of parallel region */
```

OpenMP Syntax

Work-sharing – Single thread execution:

- ❑ Work done by one thread only

```
#pragma omp parallel shared(a,b,c) private(i)
{
    #pragma omp single
    { read_array(a); read_array(b); }

    #pragma omp for
    for (i=0; i < N; i++)
        c[i] = a[i] + b[i];

    #pragma omp single
    { write_array(c); }
} /* end of parallel section */
```

OpenMP Syntax

Work-sharing – conditional parallelism:

- ❑ the `if(...)` clause

```
int i;
float a[N], b[N], c[N];

for (i=0; i < N; i++)
    a[i] = b[i] = i * 1.0;

#pragma omp parallel if (N > 10000) \
    shared(a,b,c) private(i)
{
    #pragma omp for
    for (i=0; i < N; i++)
        c[i] = a[i] + b[i];
}
```


OpenMP Syntax

The `num_threads(...)` clause:

```
#pragma omp parallel ... num_threads(int_expr)
{
    ...
}
```

- ❑ only one `num_threads` clause per parallel directive
- ❑ `int_expr` is evaluated before the parallel region is entered

OpenMP syntax

The `collapse(n)` clause

- ❑ a way to parallelize loop nests

```
subroutine sub()
!$omp do collapse(2) private(i,j,k)
  do k = kl, ku, ks
    do j = jl, ju, js
      do i = il, iu, is
        call bar(a,i,j,k)
      enddo
    enddo
  enddo
!$omp end do
end subroutine
```

collapse the two outer loops over k and j

OpenMP Syntax

Synchronization – Critical region:

- specifies a region of code that must be executed by only **one** thread at a time!
- can be named

```
#pragma omp parallel private(loc_sum)
{
    ...
    #pragma omp for
    for(int i = 0; i < n; i++)
        loc_sum += x[i];

    #pragma omp critical (cr_sum)
    sum += loc_sum;
} /* end of parallel section */
```

OpenMP Syntax

Synchronization – Atomic construct:

- specifies a single operation(!) that must be executed by only **one** thread at a time!
- restricted syntax (see OpenMP standard)

```
int x = 0;

#pragma omp parallel shared(x)
{
    ...
    #pragma omp atomic
    x = x + 1;
    ...
} /* end of parallel section */
```

OpenMP Syntax

Synchronization – Master region:

- ❑ specifies a region of code that is executed by the master thread only!
- ❑ ignored by others – no implicit barriers!

```
#pragma omp parallel
{
    ...
    #pragma omp master
    {
        printf("Hello\n");
    }
    ...
} /* end of parallel section */
```

OpenMP Syntax

Synchronization – Ordered:

- ❑ executes code block in the sequential order
- ❑ everything outside can run in parallel
- ❑ only within a parallel do/for loop

```
#pragma omp parallel for \
    private(i,myf) ordered
for(i = 0; i < n; i++) {
    myf = i+1;
    #pragma omp ordered
    {
        sum = sum + myf;
    }
} /* end of parallel for */
```

```
#pragma omp parallel for \
    private(i,myf)
for(i = 0; i < n; i++) {
    myf = i+1;
    #pragma omp critical
    {
        sum = sum + myf;
    }
} /*end of parallel for*/
```

OpenMP Syntax

Output:

non-ordered	vs.	ordered
sum[1] = 1		sum[1] = 1
sum[6] = 7		sum[2] = 3
sum[2] = 9		sum[3] = 6
sum[7] = 16		sum[4] = 10
sum[3] = 19		sum[5] = 15
sum[8] = 27		sum[6] = 21
sum[4] = 31		sum[7] = 28
sum[9] = 40		sum[8] = 36
sum[10] = 50		sum[9] = 45
sum[5] = 55		sum[10] = 55
Result: 55		Result: 55

OpenMP Syntax

Ordered with dependences (OpenMP 4.5)

- ❑ new clause `depend(...)` enhances ordered by adding dependences
 - ❑ `#pragma omp ordered depend(sink: vec)`
 - ❑ start the ordered region with the dependences given in `vec`
 - ❑ `#pragma omp ordered depend(source)`
 - ❑ ends the ordered execution defined by `sink` above
- ❑ this new feature allows to parallelize e.g. loops with dependences (see next slide)

OpenMP Syntax

doacross loops:

- ❑ loops with dependencies

```
#pragma omp parallel
{
    #pragma omp for ordered(1)
    for (int i=1; i<n; i++) {

        #pragma omp ordered depend(sink:i-1)
        a[i] = a[i-1] + b[i];
        #pragma omp ordered depend(source)

        c[i] = 2*a[i];
    } // End of for-loop
} // End of parallel region
```

of dependences
↓
dependence vector
↓

OpenMP Syntax

ordered depend(...) and doacross loops:

- ❑ provides logic in the OpenMP runtime, that else would have to be done manually
- ❑ helps to “resolve” static dependences and allows part of the code to run in parallel
- ❑ useful, if the parallel part of the loop(s) is more computational intensive than the ordered part
- ❑ hint: might need a different loop scheduling

OpenMP Syntax

Synchronization – Barrier:

- synchronizes all threads in a team

```
#pragma omp parallel
{
    ...

    #pragma omp barrier

    ...
} /* end of parallel section */
```

OpenMP Syntax

Synchronization – Implied barriers:

- exit from parallel region
- exit from `omp for/omp do/omp workshare`
- exit from sections
- exit from single

No *implied* barrier on the master construct,
neither on entry nor on exit!

OpenMP Syntax

OpenMP programming in C/C++ Part II: Clauses



OpenMP Syntax

Data scoping clauses:

- ❑ Understanding and the use of data scoping is really essential.
- ❑ Most problems/errors are due to wrong data scoping.
- ❑ Most variables are shared by default (shared memory programming model).
- ❑ Private variables: loop indices, stack of subroutines.



OpenMP Syntax

OpenMP Data scope attribute clauses:

- ❑ private
- ❑ shared
- ❑ default
- ❑ reduction
- ❑ firstprivate
- ❑ lastprivate
- ❑ copyin

OpenMP Syntax

The “private” clause:

- ❑ declares variables private to each thread:
`#pragma omp directive private (list)`
- ❑ a new variable is declared once for each thread
- ❑ all references are replaced with references to the newly declared variable
- ❑ variables declared private are uninitialized for each thread!

OpenMP Syntax

The “shared” clause:

- ❑ declares variables to be shared among all threads:

```
#pragma omp directive shared (list)
```

- ❑ a shared variable exists in only one memory location and all threads have read/write access to that address
- ❑ proper access to the variable is left to the programmer – that's YOU!

OpenMP Syntax

The “default” clause:

- ❑ allows the programmer to specify the default scope for all variables:

```
#pragma omp dir default(shared|none)
```

- ❑ C/C++ knows only those two types
- ❑ only one default clause per parallel region
- ❑ Best practice: use default(none) and scope all your variables explicitly

OpenMP Example

Two examples

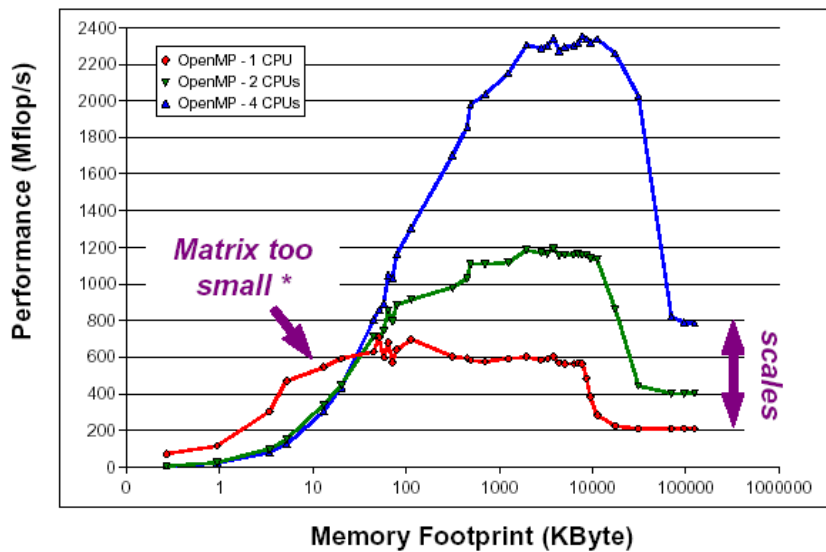
OpenMP Example

Matrix times vector

```
void
mxv(int m, int n, double *a, double *b, double *c)
{
    int i, j;
    double sum;

    #pragma omp parallel for default(none) \
        shared(m,n,a,b,c) private(i,j,sum)
    for (i=0; i<m; i++) {
        sum = 0.0;
        for (j=0; j<n; j++)
            sum += b[i*n+j] * c[j];
        a[i] = sum;
    }
}
```

OpenMP Example



SunFire 6800
UltraSPARC III Cu @ 900 MHz
8 MB L2-cache

*) With the IF-clause in OpenMP this performance degradation can be avoided

OpenMP Example

Example: numerical integration of $f(x)$

```
int i, n;
double h, x, sum;

h = 1.0 / (double)n;
sum = 0.0;

#pragma omp parallel for default(none) \
    shared(n,h,sum) private(i,x)

for(i=1; i<=n; i++) {
    x = h * ((double)i + 0.5);
    #pragma omp critical
    sum += f(x);
}
```

sequential code!

Race condition!

OpenMP Example

Example: numerical integration of $f(x)$

Improvement 1

```
int i, n;
double h, x, fx, sum;

h = 1.0 / (double)n;
sum = 0.0;

#pragma omp parallel for default(none) \
    shared(n,h,sum) private(i,x,fx)
for(i=1; i<=n; i++) {
    x = h * ((double)i + 0.5);
    fx = f(x);
    #pragma omp critical
        sum += fx;
}
```

function evaluation in parallel

OpenMP Example

Example: numerical integration of $f(x)$

Improvement 2

```
int i, n; double h, x, t_sum, sum;

h = 1.0 / (double)n; sum = 0.0;
#pragma omp parallel default(none) \
    shared(n,h,sum) private(i,x,t_sum) {
    t_sum = 0.0;
    #pragma omp for
    for(i=1; i<=n; i++) {
        x = h * ((double)i + 0.5);
        t_sum += f(x);
    }
    #pragma omp critical
        sum += t_sum;
} // end omp parallel
```

OpenMP Syntax

The “reduction” clause:

- ❑ performs a reduction on the variables that appear on the list:

```
#pragma omp dir reduction(op: list)
```

- ❑ a private copy for each thread of all variables on the list is created
- ❑ at the end, the reduction operation is carried out and the result(s) written to the global variable(s)

OpenMP Example

Example: numerical integration of $f(x)$

- ❑ smart OpenMP solution

```
int i, n;
double h, x, sum;

h = 1.0 / (double)n;
sum = 0.0;

#pragma omp parallel for default(none) \
    shared(n,h) private(i,x) \
    reduction(+: sum)
for(i=1; i<=n; i++) {
    x = h * ((double)i + 0.5);
    sum += f(x);
}
```

OpenMP Exercises – I

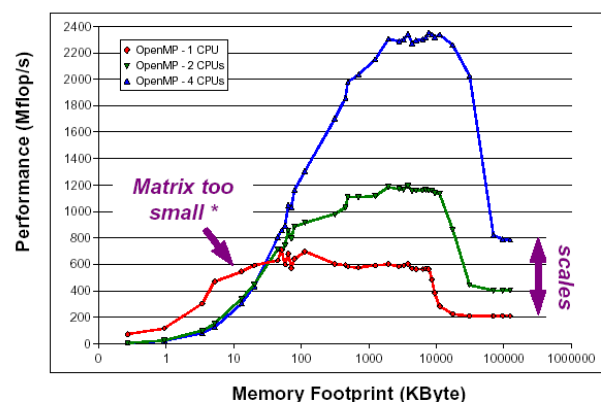
- Write an OpenMP code to calculate π , using

$$\pi = \int_0^1 \frac{4}{(1+x^2)} dx \approx \frac{1}{N} \sum_{i=1}^N \frac{4}{1 + \left(\frac{i-0.5}{N}\right)^2}$$

- implement the integrand as a function
- write your own reduction code
- use the OpenMP reduction clause
- compare the run-times

OpenMP Exercises – II

- Improve the matrix times vector example by adding an if-clause to the omp pragma – experiment with the threshold value!



SunFire 6800
UltraSPARC III Cu @ 900 MHz
8 MB L2-cache

*) With the IF-clause in OpenMP this performance degradation can be avoided