Parallel Programming in

OpenMP

Outline

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



What is OpenMP?



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



- OpenMP development is community driven
- Architecture Review Board (ARB):
 - hardware and software vendors
 - government and academia
- Founded back in 1997
 - unification of different vendor "standards"
- Official OpenMP website:
 - https://www.openmp.org/



- 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)



- □ 4.0 (Jul 2013)
 - exception handling
- □ 4.5 update (Nov 2015)
 - first support for accelerators ("GPU offloading")
- □ 5.0 (Nov 2018)
 - is implemented in most compilers now
- □ 5.2 update (Nov 2021)
- The OpenMP standard specifications:
 - http://www.openmp.org/specifications/



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)

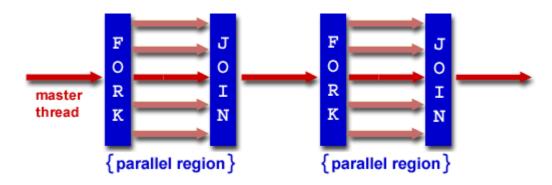


Basic elements of OpenMP



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.





OpenMP is mostly based on compiler directives:

□ C/C++:

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

Fortran:

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



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



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);
}
```



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);
```



```
$ ./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
 - Dynamic thread adjustment
- Dynamic thread adjustment

- Data scoping
- Nested parallelism

Nested parallelism

- no. of threads
- Orphaning

API for timers & locking

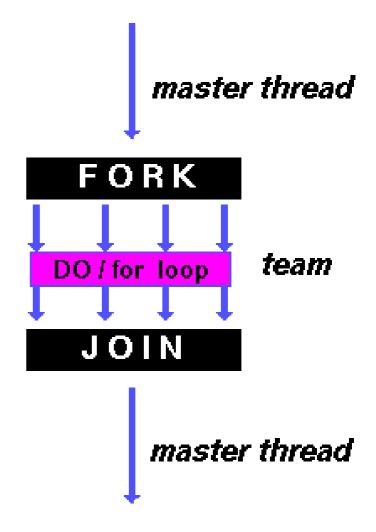


Work-sharing



Work-sharing constructs – 1:

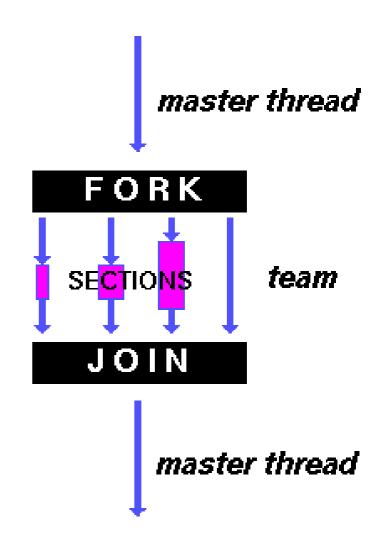
- do/for
- loop parallelism
- most common





Work-sharing constructs – 2:

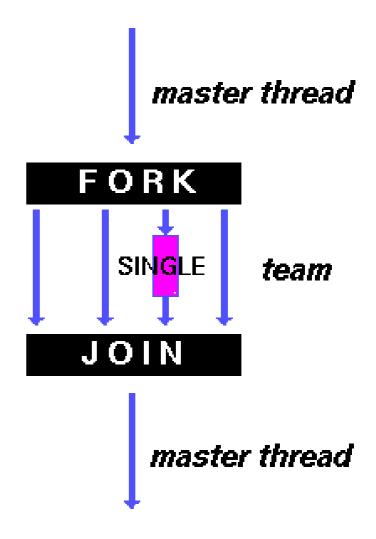
- sections
- functional parallelism
- typically independent calculations





Work-sharing constructs – 3:

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





Important rules for work-sharing constructs:

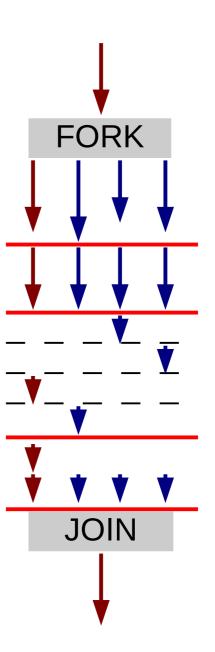
- must be enclosed in a parallel region
- must be encountered by all team members
- must be encountered in the same order



Synchronization



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





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



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!



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();
}
```



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)
                                 for has to follow
  #pragma omp for
                                 the pragma – no
  for (i=0; i < N; i++)
                                 {... }!
    c[i] = a[i] + b[i];
```



} /* end of parallel region */

Work-sharing – Loop parallelism:

Another version: combined "parallel for"



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



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 */
```



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 */
```



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];
```



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



The collapse(n) clause

a way to parallelize loop nests

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



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 */</pre>
```



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 */
```



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 */
```



Synchronization – Ordered:

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



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					



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)



doacross loops:

loops with dependencies

```
# of dependences
#pragma omp parallel
                                      dependence
    #pragma omp for ordered(1)
                                      vector
    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
```



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



Synchronization – Barrier:

synchronizes all threads in a team

```
#pragma omp parallel
{
    ...
    #pragma omp barrier
    ...
}    /* end of parallel section */
```



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 programming in C/C++
Part II: Clauses



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 Data scope attribute clauses:

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



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!



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!



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



Two examples

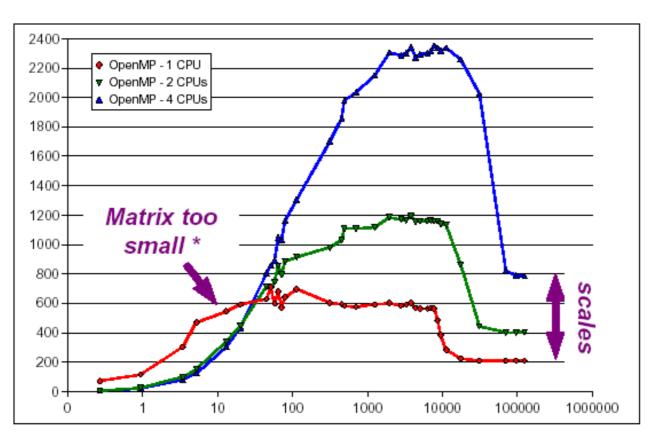


Matrix times vector

```
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;
```







Memory Footprint (KByte)

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

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



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
                                   sequential code!
   sum += f(x);
```

Race condition!



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); function evalution in parallel
#pragma omp critical
   sum += fx;
```



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
```



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)



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 Excercises – I

 \blacksquare Write an OpenMP code to calculate π , using

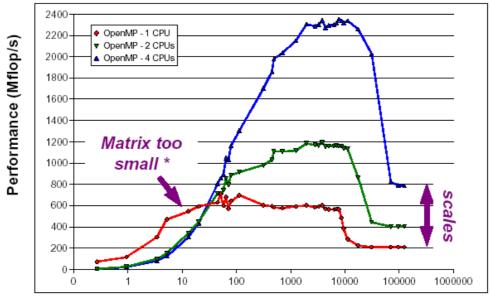
$$\pi = \int_{0}^{1} \frac{4}{(1+x^{2})} dx \approx \frac{1}{N} \sum_{i=1}^{N} \frac{4}{1+(\frac{i-0.5}{N})^{2}}$$

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



OpenMP Excercises – II

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



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