

Pemrograman OpenMP (2)

Kuliah#8 TSK617 Pengolahan Paralel - TA 2011/2012

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Tentang Kuliah #8 Pemrograman OpenMP

- ▶ Akan dibahas tentang **pemrograman paralel dengan OpenMP** menggunakan kompiler directive
 - ▶ Arsitektur memori: shared (SMP, symmetric multi-processor)
 - ▶ Model programming: thread
- ▶ Pokok Bahasan: (kuliah #8 akan membahas item yang ditebalkan)
 1. Pengantar OpenMP
 2. Membuat Thread
 3. Sinkronisasi dengan *critical*, *atomic*
 4. **Loop dan Worksharing**
 5. **Sinkronisasi dengan *barrier*, *single*, *master*, *ordered***
 6. **Environment Data**
 7. Menjadwalkan *for* dan *section*
 8. Model memori
 9. OpenMP 3.0

- ▶ Setelah mempelajari bab ini, mahasiswa akan mampu:
 1. [C2] Mahasiswa memahami konsep pemrograman paralel menggunakan OpenMP
 2. [C3] Mahasiswa akan mampu membuat program paralel dari program serial menggunakan compiler-directive dan pustaka-pustaka OpenMP
 3. [C5] Mahasiswa akan mampu memprogram suatu aplikasi komputasi matrik menggunakan OpenMP serta menghitung faktor speedupnya
- ▶ Link
 - ▶ Website: <http://didik.blog.undip.ac.id/2012/02/25/kuliah-tsk-617-pengolahan-paralel-2011/>
 - ▶ Email: didik@undip.ac.id

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Acknowledgment

- ▶ Materi dan gambar didapat dari:
 - ▶ Tim Mattson, Larry Meadows (2008): “A Hands-on Introduction to OpenMP“
 - ▶ Website: <http://openmp.org/wp/resources/>

Bahasan

Pemrograman
OpenMP (2)

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Loop Paralel

- SPMD vs Worksharing

- Loop Construct

- Working with Loop

- Reduction

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- Single Worksharing Construct

- Synchronization: Ordered

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SPMD vs Worksharing

- ▶ A parallel construct by itself creates an SPMD or “Single Program Multiple Data” program
 - ▶ each thread **redundantly executes the same code**
- ▶ How do you split up pathways through the code between threads within a team?
 - ▶ This is called **worksharing**
 - ▶ Loop construct
 - ▶ Sections/section constructs
 - ▶ Single construct
 - ▶ Task construct

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Loop Construct

- ▶ The loop worksharing construct **splits up loop iterations** among the threads in a team
- ▶ Loop construct name:
 - ▶ C/C++: **for**
 - ▶ Fortran: **do**

```
#pragma omp parallel
{

    #pragma omp for
    for (i=0;i<N;i++){
        NEAT_STUFF(i);
    }

}
```

- ▶ The variable *i* is made “**private**” to each thread **by default**. You could do this **explicitly** with a “private(*i*)” clause

```
#pragma omp parallel private(i)
{

    #pragma omp for
    for (i=0;i<N;i++){
        NEAT_STUFF(i);
    }

}
```

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Construct Worksharing Loop: Contoh

Sequential code

```
for(i=0;i<N;i++) { a[i] = a[i] + b[i]; }
```

OpenMP parallel region

```
#pragma omp parallel  
{  
  
    int id, i, Nthrds, istart, iend;  
    id = omp_get_thread_num();  
    Nthrds = omp_get_num_threads();  
    istart = id * N / Nthrds;  
    iend = (id+1)*N/Nthrds;  
    if (id == Nthrds-1) iend = N;  
    for(i=istart;i<iend;i++) {  
        a[i] = a[i] + b[i];  
    }  
  
}
```

OpenMP parallel region
and a worksharing for
construct

```
#pragma omp parallel  
#pragma omp for  
    for(i=0;i<N;i++) {  
        a[i] = a[i] + b[i];  
    }
```

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Kombinasi Construct Paralel/Worksharing

- ▶ OpenMP shortcut: Put the “parallel” and the worksharing directive **on the same line**

```
double res[MAX]; int i;
#pragma omp parallel
{
    #pragma omp for
    for(i=0;i<MAX;i++) {
        res[i] = huge();
    }
}
```

```
double res[MAX]; int i;
#pragma omp parallel for
    for(i=0;i<MAX;i++) {
        res[i] = huge();
    }
```

- ▶ Both codes are equivalent

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► Basic approach:

1. Find **compute intensive loops**
2. Make **the loop iterations independent** ..
 - So they can safely execute in any order without loop-carried dependencies
3. **Place** the appropriate OpenMP directive and test

```
int i, j, A[MAX];
j = 5;
for (i=0;i< MAX; i++) {
    j +=2;
    A[i] = big(j);
}
```

```
int i, A[MAX];
#pragma omp parallel for
for (i=0;i< MAX; i++) {
    int j = 7 + 2*i;
    A[i] = big(j);
}
```

- Remove loop dependency, i.e from variable j

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- ▶ How do we handle this case?

```
double ave=0.0, A[MAX]; int i;
for (i=0;i< MAX; i++) {
    ave + = A[i];
}
ave = ave/MAX;
```

- ▶ It is combining values into a single accumulation variable (**ave**)
 - ▶ There is a true dependence between loop iterations that **can't be trivially removed**
 - ▶ This is a very common situation
 - ▶ It is called a “**reduction**”.
- ▶ Support for reduction operations is included in **most** parallel programming environments

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Reduction

- ▶ OpenMP reduction clause:

```
reduction (op : list)
```

Op is an operator

- ▶ Inside a parallel or a work-sharing construct:
 - ▶ **A local copy** of each list variable is made and **initialized** depending on the “op” (e.g. 0 for “+”).
 - ▶ Compiler finds standard reduction expressions containing “**op**” and uses them to **update** the local copy.
 - ▶ Local copies are reduced into **a single** value and **combined** with the original global value.
- ▶ The variables in “list” must be shared in the enclosing parallel region.

```
double ave=0.0, A[MAX]; int i;  
#pragma omp parallel for reduction (+:ave)  
for (i=0;i< MAX; i++) {  
    ave + = A[i];  
}  
ave = ave/MAX;
```


Reduction operands/initial-values

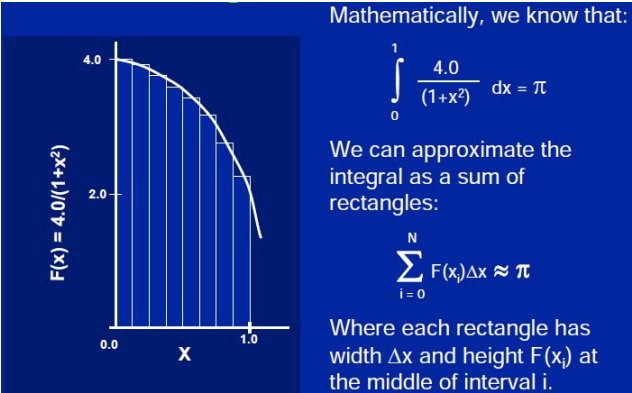
- ▶ Many different associative operands can be used with reduction
 - ▶ Initial values are the ones that make sense mathematically.

Operator	Initial Value
+	0
*	1
-	0

C/C++ Only	
Operator	Initial Value
&	~0
	0
^	0
&&	1
	0

Exercise

- ▶ Parallelize serial pi program with a loop construct
 - ▶ The goal is to **minimize the number changes** made to the serial program.



Parallelize Pi Program

```
/*Serial pi program*/  
#include <stdlib.h>  
#include <omp.h>  
define num_steps 1000000;  
double step;  
int main (int argc, char *argv[])  
{  
    int i, j;  
    double x, pi, sum = 0.0;  
    step = 1.0/num_steps;  
    for (i=0;i< num_steps; i++){  
        x = (i+0.5)*step;  
        sum += 4.0/(1.0+x*x);  
    }  
    pi = step * sum;  
    return EXIT_SUCCESS;  
}
```

```
/*Serial pi program*/  
#include <stdlib.h>  
#include <omp.h>  
#define num_steps 1000000;  
double step;  
int main (int argc, char *argv[])  
{  
    int i, j;  
    double x, pi, sum = 0.0;  
    step = 1.0/num_steps;  
    #pragma omp parallel for re-  
duction(+:sum)  
    for (i=0;i< num_steps; i++){  
        x = (i+0.5)*step;  
        sum += 4.0/(1.0+x*x);  
    }  
    pi = step * sum;  
    return EXIT_SUCCESS;  
}
```

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Synchronization: Barrier

- **Barrier:** Each thread **waits** until all threads arrive

```
#pragma omp parallel shared (A, B, C) private(id)
{
    id=omp_get_thread_num();
    A[id] = big_calc1(id);
#pragma omp barrier
#pragma omp for
    for(i=0;i<N;i++){C[i]=big_calc3(i,A);}
        // implicit barrier at the end of worksharing construct
#pragma omp for nowait
    for(i=0;i<N;i++){ B[i]=big_calc2(C, i); }
        // no implicit barrier due to nowait
    A[id] = big_calc4(id);
} // implicit barrier at the end of a parallel region
```

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- ▶ The **master construct** denotes a structured block that is only executed by the **master** thread.
 - ▶ The other threads just skip it (**no synchronization** is implied).

```
#pragma omp parallel
{
    do_many_things();
#pragma omp master
    { exchange_boundaries(); }
#pragma omp barrier
    do_many_other_things();
}
```

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Single Worksharing Construct

- ▶ The single construct denotes a block of code that is executed by **only one thread** (not necessarily the master thread)
 - ▶ A barrier is implied at the end of the single block (can remove the barrier with a nowait clause)

```
#pragma omp parallel
{
    do_many_things();
#pragma omp single
    { exchange_boundaries(); }
    do_many_other_things();
}
```

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Synchronization: Ordered

- ▶ The ordered region executes in the **sequential order**

```
#pragma omp parallel private (tmp)
#pragma omp for ordered reduction(+:res)
for (i=0;i<N;i++){
    tmp = NEAT_STUFF(i);
#pragma ordered
    res += konsum(tmp);
}
```

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Synchronization: Lock Routines

▶ Simple Lock routines

- ▶ A simple lock is available if it is unset
*omp_init_lock(), omp_set_lock(), omp_unset_lock(),
omp_test_lock(), omp_destroy_lock()*

▶ Nested Locks

- ▶ A nested lock is available if **it is unset or if it is set** but owned by the thread executing the nested lock function
*omp_init_nest_lock(), omp_set_nest_lock(),
omp_unset_nest_lock(), omp_test_nest_lock(),
omp_destroy_nest_lock()*

▶ Note:

- ▶ A lock implies a memory fence (a “flush”) of all thread visible variables
- ▶ A thread always accesses the most recent copy of the lock, so you **don't need** to use a flush on the lock variable

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► Protect resources with locks

```
omp_lock_t lck;
omp_init_lock(&lck);
#pragma omp parallel private (tmp, id)
{
    id = omp_get_thread_num();
    tmp = do_lots_of_work(id);
    // wait here for your turn
    omp_set_lock(&lck);
    printf("%d %d", id, tmp);
    // release the lock so the next thread gets a turn
    omp_unset_lock(&lck);
}
// Free-up storage when done
omp_destroy_lock(&lck);
```

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▶ Runtime environment routines

- ▶ Modify/Check the number of threads
omp_set_num_threads(), *omp_get_num_threads()*,
omp_get_thread_num(), *omp_get_max_threads()*
- ▶ Are we in an active parallel region?
omp_in_parallel()
- ▶ Do you want the system to dynamically vary the number of threads from one parallel construct to another?
omp_set_dynamic(), *omp_get_dynamic()*;
- ▶ How many processors in the system?
omp_num_procs()

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Runtime Routine Usage

- ▶ To use a known, fixed number of threads in a program
 1. tell the system that you don't want dynamic adjustment of the number of threads
 2. set the number of threads
 3. save the number you got

```
#include <omp.h>
void main()
{
    int num_threads;
    // Disable dynamic adjustment of the number of threads
    omp_set_dynamic( 0 );
    // Request as many threads as you have processors
    omp_set_num_threads( omp_num_procs() );
    #pragma omp parallel
    {
        int id=omp_get_thread_num();
    #pragma omp single
        num_threads = omp_get_num_threads();
        do_lots_of_stuff(id);
    }
}
```

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Variabel Environment

- ▶ Set the default number of threads to use.

```
OMP_NUM_THREADS int_literal
```

- ▶ Control how “omp for schedule(RUNTIME)” loop iterations are scheduled

```
OMP_SCHEDULE “schedule[, chunk_size]”
```

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Data Environment: Atribut Storage Default

- ▶ Shared Memory programming model:
 - ▶ Most variables are shared by default
- ▶ Global variables are **SHARED** among threads
 - ▶ **File scope** variables, **static**
 - ▶ dynamically allocated memory (**ALLOCATE**, **malloc**, **new**)
- ▶ But not everything is shared...
 - ▶ Stack variables in functions called from parallel regions are **PRIVATE**
 - ▶ Automatic variables within a statement block are **PRIVATE**

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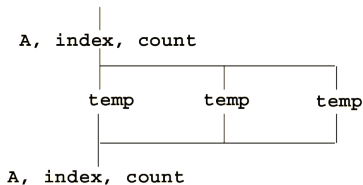
Data Sharing: Contoh

```
double A[10];
int main() {
    int index[10];
    #pragma omp parallel
        work(index);
    printf("%d\n", index[0]);
}
```

- ▶ Function work() is implemented on different file from function main()

```
extern double A[10];
void work(int *index) {
    double temp[10];
    static int count;
    ...
}
```

- ▶ A, index and count are shared by all threads
- ▶ temp is local to each thread



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Data Sharing: Mengubah Atribut Storage

- ▶ One can selectively change storage attributes for constructs using the following clauses:
 - ▶ SHARED
 - ▶ PRIVATE
 - ▶ FIRSTPRIVATE
- ▶ The final value of a private inside a parallel loop can be transmitted to the shared variable outside the loop with:
 - ▶ LASTPRIVATE
- ▶ The default attributes can be overridden with:
 - ▶ DEFAULT (PRIVATE | SHARED | NONE)
 - ▶ DEFAULT(PRIVATE) is Fortran only
- ▶ All data clauses apply to parallel constructs and worksharing constructs except “shared” which only applies to parallel constructs.

Data Sharing: Clause Private

- ▶ **private(var)** creates a new local copy of var for each thread
 - ▶ The value is uninitialized
 - ▶ In OpenMP 2.5 the value of the shared variable is undefined after the region

```
void wrong() {  
    int tmp = 0;  
    #pragma omp for private(tmp)  
    for (int j = 0; j < 1000; ++j)  
        tmp += j; // tmp was not initialized  
    printf("%d\n", tmp);  
    // tmp: 0 in 3.0, unspecified in 2.5  
}
```

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Data Sharing: Clause Private

When is the original variable valid?

- ▶ The original variable's value is unspecified in OpenMP 2.5.
- ▶ In OpenMP 3.0, if it is referenced outside of the construct
- ▶ Implementations may reference the original variable or a copy
 - ▶ A dangerous programming practice!

```
int tmp;
void danger() {
    tmp = 0;
    #pragma omp parallel private(tmp)
        work();
    printf(“%d\n”, tmp); // tmp has unspecified value
}
```

```
extern int tmp;
void work() {
    tmp = 5; // unspecified which copy of tmp
}
```

Data Sharing: Clause Firstprivate

- ▶ **Firstprivate** is a special case of private
 - ▶ **Initializes each private copy** with the corresponding value from the master thread

```
void useless() {  
    int tmp = 0;  
    #pragma omp for firstprivate(tmp)  
    // Each thread gets its own tmp with an initial value of 0  
    for (int j = 0; j < 1000; ++j)  
        tmp += j;  
    printf(“%d\n”, tmp);  
    // tmp: 0 in 3.0, unspecified in 2.5  
}
```

Loop Paralel

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Changing Storage
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Data Sharing: Clause Lastprivate

- ▶ Lastprivate passes the value of a private from the last iteration to a global variable

```
void closer() {  
    int tmp = 0;  
    #pragma omp parallel for firstprivate(tmp) \  
        lastprivate(tmp)  
    for (int j = 0; j < 1000; ++j)  
        // Each thread gets its own tmp with an initial value of 0  
        tmp += j;  
    printf(“%d\n”, tmp);  
    // tmp is defined as its value at the “last sequential” itera-  
    tion (i.e., for j=999)  
}
```

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Data Sharing: Testing Data Environment

- ▶ Consider this example of PRIVATE and FIRSTPRIVATE Variables A,B, and C = 1

```
#pragma omp parallel private(B) firstprivate(C)
```

- ▶ Are A,B,C local to each thread or shared inside the parallel region?
- ▶ What are their initial values inside and values after the parallel region?
- ▶ Inside this parallel region
 - ▶ A is shared by all threads; equals 1
 - ▶ B and C are local to each thread.
 - ▶ B's initial value is undefined
 - ▶ C's initial value equals 1
- ▶ Outside this parallel region
 - ▶ The values of B and C are unspecified in OpenMP 2.5, and in OpenMP 3.0 if referenced in the region but outside the construct

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Data Sharing: Clause Default

- ▶ The default storage attribute is **DEFAULT(SHARED)** (so no need to use it)
 - ▶ Exception: ***#pragma omp task***
 - ▶ Task default storage is **firstprivate**
- ▶ To change default: **use** DEFAULT(PRIVATE), Only on Fortran
 - ▶ each variable in the construct is made private as if specified in a private clause
 - ▶ mostly saves typing
- ▶ **DEFAULT(NONE)**: no default for variables in static extent
 - ▶ Must list storage attribute for each variable in static extent
 - ▶ Good programming practice!

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Data Sharing: Task (OpenMP 3.0)

- ▶ The default for tasks is usually **firstprivate**, because the task may not be executed until later (and variables may have gone out of scope)
- ▶ Variables that are shared in all constructs starting from the innermost enclosing parallel construct are shared, because the barrier guarantees task completion

```
#pragma omp parallel shared(A) private(B)
{
    ...
    #pragma omp task
    {
        int C;
        compute(A, B, C);
    }
}
```

- ▶ A is **shared**, B is **firstprivate**, and C is **private**
- ▶ Task in detail will be presented in next lecture

Data Sharing: Threadprivate

- ▶ Makes global data private to a thread
 - ▶ In C: File scope and static variables, static class members
- ▶ Different from making them PRIVATE
 - ▶ with PRIVATE global variables are masked
 - ▶ THREADPRIVATE preserves global scope within each thread
- ▶ Threadprivate variables can be initialized using COPYIN or at time of definition (using language-defined initialization capabilities)

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Data Sharing: Contoh Threadprivate

- ▶ Use threadprivate to create a counter for each thread

```
int counter = 0;
#pragma omp threadprivate(counter)
int increment_counter()
{
    counter++;
    return (counter);
}
```

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Penyalinan Data: Copyin

- ▶ Initialize threadprivate data using a copyin clause
- ▶ Used with a single region to broadcast values of privates from one member of a team to the rest of the team

```
#include <omp.h>
void input_parameters (int, int); // fetch values of input parameters
void do_work(int, int);
void main()
{
    int Nsize, choice;
    #pragma omp parallel private (Nsize, choice)
    {
        #pragma omp single copyprivate (Nsize, choice)
            input_parameters (Nsize, choice);
    }
}
```

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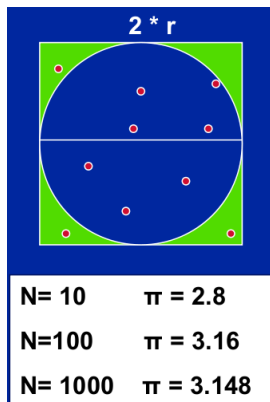
Changing Storage
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Penghitungan Monte Carlo

Menggunakan bilangan random untuk memecahkan problem

- ▶ Sample a problem domain to estimate areas, compute probabilities, find optimal values, etc.
- ▶ Example: Computing π with a digital dart board



- ▶ Throw darts at the circle/square
- ▶ Chance of falling in circle is proportional to ratio of areas:
 - ▶ $A_c = r^2 * \pi$
 - ▶ $A_s = (2 * r) * (2 * r) = 4 * r^2$
 - ▶ $P = A_c/A_s = \pi/4$
- ▶ Compute π by randomly choosing points, count the fraction that falls in the circle, compute π

Latihan: Penghitungan Monte Carlo

- ▶ File untuk program serial:
 - ▶ pi_mc.c: program pi dengan metode monte carlo
 - ▶ random.c: generator random sederhana
 - ▶ random.h: file header untuk generator random
- ▶ Buat program paralel menggunakan OpenMP!

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- ▶ Anda bebas:
 - ▶ untuk **Membagikan** — untuk menyalin, mendistribusikan, dan menyebarkan karya, dan
 - ▶ untuk **Remix** — untuk mengadaptasikan karya
- ▶ Di bawah persyaratan berikut:
 - ▶ **Atribusi** — Anda harus memberikan atribusi karya sesuai dengan cara-cara yang diminta oleh pembuat karya tersebut atau pihak yang mengeluarkan lisensi.
 - ▶ **Pembagian Serupa** — Jika Anda mengubah, menambah, atau membuat karya lain menggunakan karya ini, Anda hanya boleh menyebarkan karya tersebut hanya dengan lisensi yang sama, serupa, atau kompatibel.
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