# Multi-Core Programming with OpenMP

### Clemens Grelck

University of Amsterdam Informatics Institute Computer Systems Architecture Group

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# Outline of the Day

- Multi-core computer architecture (brief)
- ▶ OpenMP the Basics
- ▶ Lunch Break
- ▶ OpenMP the Advanced Story



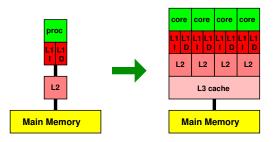
### What we are talking about:

- Bread and butter multi-core processors
- ► Intel XEON family
- AMD Opteron family
- maybe Oracle T3/T4 (Niagara)



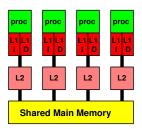
#### **Characteristics:**

- Consequence of Murphy's Law: abundance of transistors
- Put multiple fully-fledged previous generation processors on a single die.
- ► Add generous caches
- Hardware cache coherence



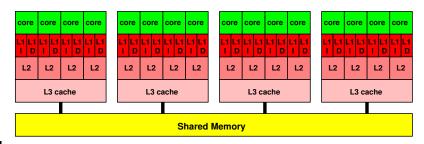
### Symmetric multiprocessing:

- ▶ Idea dates back to 1980s
- Assemble multiple processors on one board, 2-way or 4way
- Processor-local caches
- Shared main memory
- Hardware cache coherence



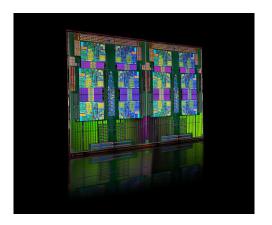
### Multi-core symmetric multiprocessors:

- Combining symmetric multiprocessing with multi-core processors
- Unprecedented parallelism on single board
  - ▶ 48 AMD Opteron cores
  - 24 hyperthreaded Intel Xeon cores
- Hardware cache coherence



# 12-core AMD Opteron

### One chip, two dies:





# A Note on Hyperthreading

#### Idea:

- Duplicate architectural state: (control) registers
- ▶ Do not duplicate functional units
- Let one core appear as two logical cores to the OS
- Make more efficient use of resources by computing two streams of instructions simultaneously
- ▶ If one stalls the other takes over

# A Note on Hyperthreading

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- Duplicate architectural state: (control) registers
- Do not duplicate functional units
- Let one core appear as two logical cores to the OS
- Make more efficient use of resources by computing two streams of instructions simultaneously
- ▶ If one stalls the other takes over

### Disadvantage:

- Requires the right instruction mix
- ► Together with symmetric multithreading OS must have specific support for hyperthreading instead of just seeing two cores
- ► Cache thrashing



# How to program such systems





# Design Rationale of OpenMP

#### Ideal:

- Automatic parallelisation of sequential code.
- ► No additional parallelisation effort for development, debugging, maintenance, etc.



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#### **Problem:**

- Data dependences are difficult to assess.
- Compilers must be conservative in their assumptions.

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#### **Problem:**

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- Compilers must be conservative in their assumptions.

### Way out:

- Take or write ordinary sequential program.
- ► Add annotations/pragmas/compiler directives that guide parallelisation.
- ▶ Let the compiler generate the corresponding code.



# Short History of OpenMP

#### **Towards OpenMP:**

- Mid 1980s: First commercial SMPs appear.
  - ANSI standard X3H5 (not adopted)
- ► Late 1980s to early 1990s:
  - Distributed memory architectures prevail.
- ► Early 1990s until today:
  - ► Shared memory architectures significantly gain popularity.
  - Different vendor-specific compiler extensions.
- **▶** 1996/97:
  - ► Development of a new industry standard: OpenMP
  - Actively supported by all major hardware vendors: Intel, HP, SGI, IBM, SUN, Compaq, ...
  - OpenMP Architectural Review Board: http://www.openmp.org/



# Short History of OpenMP

### **OpenMP Standardisation:**

- ▶ 1996 : Fortran Version 1.0
- ▶ 1998 : C/C++ Version 1.0
- ▶ 1999 : Fortran Version 1.1
- ▶ 2000 : Fortran Version 2.0
- ▶ 2002 : C/C++ Version 2.0
- ▶ 2005 : Joint Version 2.5
- ▶ 2008 : Joint Version 3.0

# OpenMP at a Glance

### OpenMP as a programming interface:

- Compiler directives
- ► Library functions
- ► Environment variables

### C/C++ version:

```
#pragma omp name [clause]*
structured block
```

#### Fortran version:

```
!$ OMP name [ clause [, clause]*]
code block
!$ OMP END name
```

# Hello World with OpenMP

```
#include "omp.h"
#include <stdio.h>
int main()
  printf( "Starting execution with %d threads:\n",
          omp_get_num_threads());
  #pragma omp parallel
    printf( "Hello world says thread %d of %d.\n",
            omp_get_thread_num(),
            omp_get_num_threads());
  }
  printf( "Execution of %d threads terminated.\n",
          omp_get_num_threads());
  return(0);
```



# Hello World with OpenMP

### **Compilation:**

```
gcc -fopenmp hello_world.c
```

### **Output using 4 threads:**

```
Starting execution with 1 threads: Hello world says thread 2 of 4. Hello world says thread 3 of 4. Hello world says thread 1 of 4. Hello world says thread 0 of 4. Execution of 1 threads terminated.
```



# Hello World with OpenMP

### Using 4 threads:

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Starting execution with 1 threads: Hello world says thread 2 of 4. Hello world says thread 3 of 4. Hello world says thread 1 of 4. Hello world says thread 0 of 4. Execution of 1 threads terminated.
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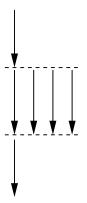
#### Who determines number of threads?

- Environment variable: export OMP\_NUM\_THREADS=4
- ► Library function: void omp\_set\_num\_threads( int)



# OpenMP Execution Model

### Classical Fork/Join:



Master thread executes serial code.

Master thread encounters parallel directive.

Master and slave threads concurrently execute parallel block.

Implicit barrier, wait for all threads to finish.

Master thread resumes serial execution.

# Simple Loop Parallelisation

### **Example: element-wise vector product:**

```
void elem_prod( double *c, double *a, double *b, int len)
{
  int i;

  #pragma omp parallel for

  for (i=0; i<len; i++)
  {
    c[i] = a[i] * b[i];
  }
}</pre>
```

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### Prerequisite:

▶ No data dependence between any two iterations.



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```

### Prerequisite:

- ▶ No data dependence between any two iterations.
- ► Caution: YOU claim this property !!



# Directive #pragma omp parallel for

### What the compiler directive does for you:

- It starts additional worker threads depending on OMP\_NUM\_THREADS.
- ▶ It divides the iteration space among all threads.
- It lets all threads execute loop restricted to their mutually disjoint subsets.
- It synchronizes all threads at an implicit barrier.
- It terminates worker threads.

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- It terminates worker threads.

#### Restrictions:

- ▶ The directive must directly precede for-loop.
- ▶ The for-loop must match a constrained pattern.
- ▶ The trip-count of the for-loop must be known in advance.



### Shared and Private Variables

### Example:

```
#pragma omp parallel for
for (i=0; i<len; i++)
{
   res[i] = a[i] * b[i];
}</pre>
```

- ► Shared variable: one instance for all threads
- ▶ Private variable: one instance for each thread

### Shared and Private Variables

### Example:

```
#pragma omp parallel for
for (i=0; i<len; i++)
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Who decides that res, a, b, and len are shared variables, whereas i is private ??

### Shared and Private Variables

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Who decides that res, a, b, and len are shared variables, whereas i is private ??

#### **Default rules:**

- All variables are shared.
- Only loop variables of parallel loops are private.



#### Mandelbrot set:

```
double x, y;
int i, j, max = 200;
int depth[M,N];
...
for (i=0; i<M; i++) {
   for (j=0; j<N; j++) {
      x = (double) i / (double) M;
      y = (double) j / (double) N;
      depth[i,j] = mandelval( x, y, max);
} }</pre>
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### Properties to check:

No data dependencies between loop iterations ?

#### Mandelbrot set:

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- ▶ No data dependencies between loop iterations? YES!
- ► Trip-count known in advance?

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- ▶ No data dependencies between loop iterations ? YES!
- Trip-count known in advance? YES!
- ► Function mandelval without side-effects?



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- Trip-count known in advance? YES!
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- ▶ Only loop variable i needs to be private?



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}</pre>
```

- No data dependencies between loop iterations? YES!
- Trip-count known in advance? YES!
- Function mandelval without side-effects? YES!
- ➤ Only loop variable i needs to be private? NO !!!! Check x,y,j



#### Mandelbrot set:

```
double x, y;
int i, j, max = 200;
int depth[M,N];
...
#pragma omp parallel for private( x, y, j)
for (i=0; i<M; i++) {
   for (j=0; j<N; j++) {
      x = (double) i / (double) M;
      y = (double) j / (double) N;
      depth[i,j] = mandelval( x, y, max);
} }</pre>
```

#### Private clause:

- Directives may be refined by clauses.
- ▶ Private clause allows to tag any variable as private.
- Caution: private variables are **not** initialised outside parallel section!!



#### Mandelbrot set with additional counter:

```
int total = 0;
...
for (i=0; i<M; i++) {
  for (j=0; j<N; j++) {
    x = (double) i / (double) M;
    y = (double) j / (double) N;
    depth[i,j] = mandelval(x, y, max);
    total = total + depth[i,j];
} }</pre>
```

# Parallelisation of a Less, Less Simple Loop

#### Mandelbrot set with additional counter:

```
int total = 0;
...
for (i=0; i<M; i++) {
  for (j=0; j<N; j++) {
    x = (double) i / (double) M;
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    depth[i,j] = mandelval(x, y, max);
    total = total + depth[i,j];
} }</pre>
```

#### **Problems:**

- ▶ New variable total introduces data dependence.
- ▶ Data dependence could be ignored due to associativity.
- ▶ New variable total must be shared.
- ▶ Incrementation of total must avoid *race condition*.



# What is a race condition?

# Note: Adding two values is not atomic

#### Thread 1:

```
load reg1 <= &depth
load reg2 <= &total
add reg1, reg2 => reg1
store reg1 => &total
...
```



## What is a race condition?

#### Note: Adding two values is not atomic

#### Thread 1:

# Thread 2:

load reg1 <= &depth
load reg2 <= &total
add reg1, reg2 => reg1
store reg1 => &total
...

```
load reg1 <= &depth
load reg2 <= &total
add reg1, reg2 => reg1
store reg1 => &total
```

#### **Problem:**

- ► Threads operate with any interleaving.
- ► If different interleavings cause different results, we have a race condition.
- Program behaves non-deterministically.



## What is a race condition?

#### Note: Adding two values is not atomic

#### Thread 1:

```
load reg1 <= &depth
load reg2 <= &total
add reg1, reg2 => reg1
store reg1 => &total
```

#### Thread 2:

```
load reg1 <= &depth
load reg2 <= &total
add reg1, reg2 => reg1
store reg1 => &total
...
```

#### Solution:

. . .

- ► We must identify critical regions: sequences of instructions that must be executed without interleaving.
- ▶ We must establish mutual exclusion between critical regions.

# Parallelisation of a Less, Less Simple Loop

#### Mandelbrot set with additional counter:

```
int total = 0;
#pragma omp parallel for private( x, y, j)
for (i=0; i<M; i++) {
  for (j=0; j<N; j++) {
   x = (double) i / (double) M;
    v = (double) i / (double) N;
   depth[i,j] = mandelval( x, y, max);
   #pragma omp critical
       total = total + depth[i,j];
```



# Critical Regions

#### The critical directive:

- ▶ Directive must immediately precede new statement block.
- Statement block is executed without interleaving.
- ▶ Directive implements critical region.

## **Equivalence:**

# Critical Regions

#### The critical directive:

- ▶ Directive must immediately precede new statement block.
- Statement block is executed without interleaving.
- ▶ Directive implements critical region.

## **Equivalence:**

# Disadvantage:

- ▶ All critical regions in entire program are synchronised.
- ▶ Unnecessary overhead.



# Critical Regions

#### The named critical directive

- Critical regions may be associated with names.
- Critical regions with identical names are synchronised.
- Critical regions with different names are executed concurrently.

# **Equivalence:**

pthread\_mutex\_unlock( &name\_lock);

# **Reduction Operations**

## Specific solution: reduction clause

## **Properties:**

- ► Reduction clause only supports built-in reduction operations: +, \*, ^, &, |, &&, ||.
- ▶ User-defined reductions only supported via critical regions.
- ▶ Bit accuracy not guaranteed.



#### **Shared variables:**

- One instance shared between sequential and parallel execution.
- Value unaffected by transition.

#### Private variables:

- One instance during sequential execution.
- One instance per worker thread during parallel execution.
- No exchange of values.



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## New: Firstprivate variables:

- Like private variables, but ...
- ▶ Worker thread instances initialised with master thread value.

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## New: Firstprivate variables:

- Like private variables, but ...
- ▶ Worker thread instances initialised with master thread value.

# New: Lastprivate variables:

- ► Like private variables, but ...
- ► Master thread instance updated to value of worker thread instance that executes the last (in sequential order) iteration.



## Example:

#### Note:

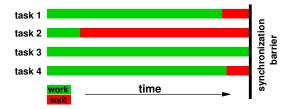
► The value of d depends on the number of iterations executed by the thread that is assigned iteration i=9 by the OpenMP loop scheduler.

#### **Definition:**

Loop scheduling determines which iterations are executed by which thread.

#### Aim:

Equal workload distribution



#### **Problem:**

Different situations require different techniques

#### The schedule clause:

```
#pragma omp parallel for schedule( <type > [, <chunk >])
for (...)
{
    ...
}
```

## **Properties:**

- Clause selects one out of a set of scheduling techniques.
- Optionally, a chunk size can be specified.
- ▶ Default chunk size depends on scheduling technique.



## Static scheduling:

```
#pragma omp parallel for schedule( static)
```

- ▶ Loop is subdivided into as many chunks as threads exist.
- ► Often called block scheduling.

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#### Static scheduling with chunk size:

```
#pragma omp parallel for schedule( static, <n>)
```

- ▶ Loop is subdivided into chunks of *n* iterations.
- ► Chunks are assigned to threads in a round-robin fashion.
- ► Also called block-cyclic scheduling.



# **Dynamic scheduling:**

#pragma omp parallel for schedule( dynamic, <n>)

- ▶ Loop is subdivided into chunks of *n* iterations.
- Chunks are dynamically assigned to threads on their demand.
- Also called self scheduling.
- Default chunk size: 1 iteration.

# **Properties:**

- Allows for dynamic load distribution and adjustment.
- Requires additional synchronization.
- Generates more overhead than static scheduling.



#### Dilemma of chunk size selection:

- Small chunk sizes mean good load balancing, but high synchronisation overhead.
- ► Large chunk sizes reduce synchronisation overhead, but result in poor load balancing.

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# Rationale of guided scheduling:

- In the beginning, large chunks keep synchronisation overhead small.
- When approaching the final barrier, small chunks balance workload.



# **Guided scheduling:**

#pragma omp parallel for schedule( guided, <n>)

- ► Chunks are dynamically assigned to threads on their demand.
- Initial chunk size is implementation dependent.
- ► Chunk size decreases exponentially with every assignment.
- ► Also called guided self scheduling.
- ▶ Minimum chunk size: *n* (default: 1)

# **Example:**

- ▶ Total number of iterations: 250
- ▶ Initial / minimal chunk size: 50 / 5
- ► Current chunk size: 80% of last chunk size: 50 40 32 26 21 17 14 12 10 8 6 5 5 4



## Conditional Parallelisation

#### **Problem:**

- Parallel execution of a loop incurs overhead:
  - creation of worker threads
  - scheduling
  - synchronisation barrier
- This overhead must be outweighed by sufficient workload.
- Workload depends on
  - loop body,
  - ▶ trip count.



## Conditional Parallelisation

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  - scheduling
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- Workload depends on
  - ▶ loop body,
  - trip count.

## **Example:**

```
if (len < 1000) {
  for (i=0; i<len; i++)
    res[i] = a[i] * b[i];
else {
  #pragma omp parallel for
  for (i=0; i<len; i++)
    res[i] = a[i] * b[i]:
```

## Conditional Parallelisation

# Introducing the if-clause:

```
if (len < 1000) {
   for (i=0; i<len; i++) {
      res[i] = a[i] * b[i];
   }
} else {
   #pragma omp parallel for
   for (i=0; i<len; i++) {
      res[i] = a[i] * b[i];
   }
}</pre>
```



```
#pragma omp parallel for if (len >= 1000)
for (i=0; i<len; i++) {
  res[i] = a[i] * b[i];
}</pre>
```

# **Eventually**

# Questions



Coffee

**Practicals** 

Lunch

