

# Shared Memory machines and OpenMP programming

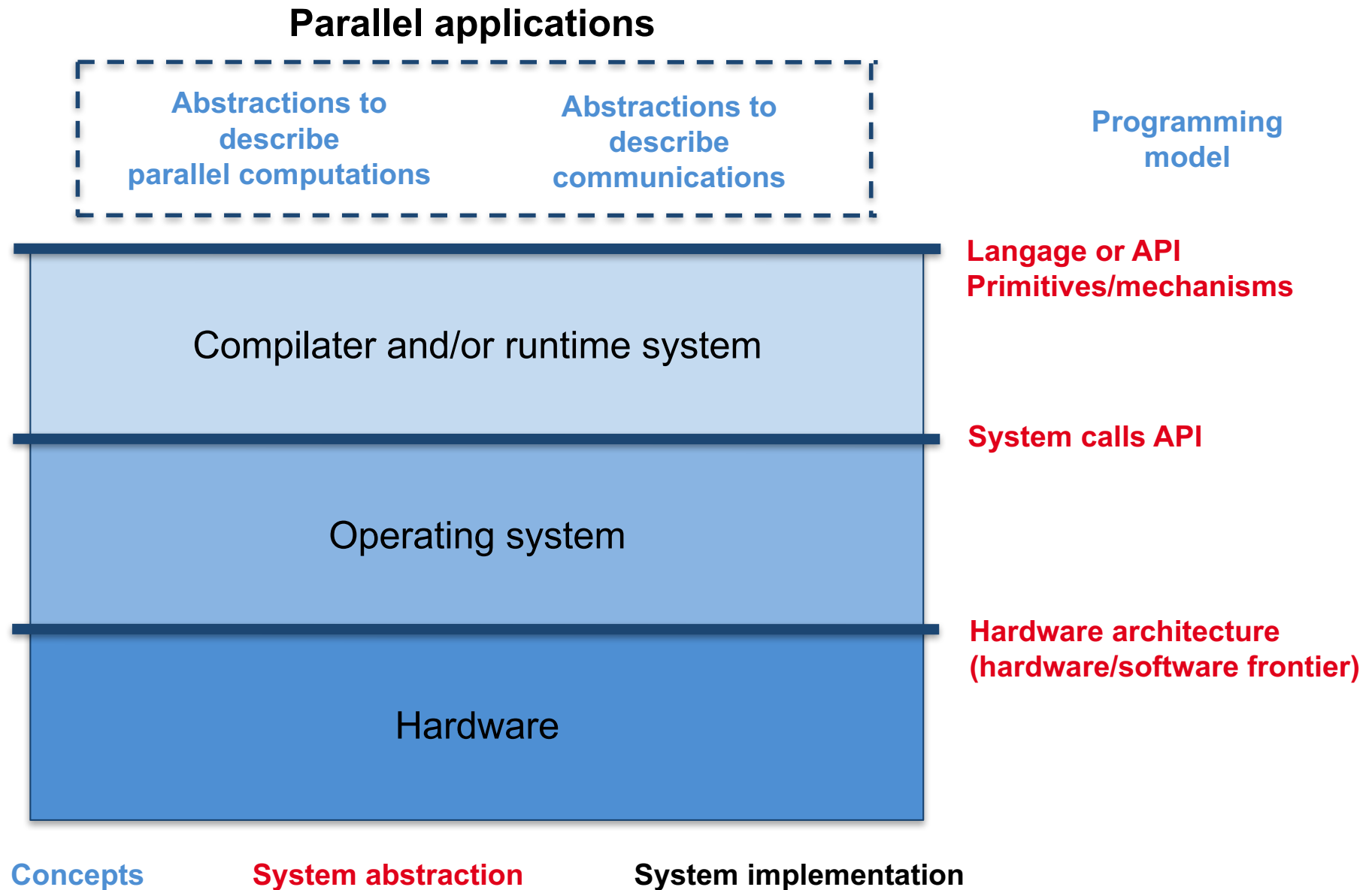
Frédéric Desprez

INRIA

# Some references

- **OpenMP web site**
  - <http://www.openmp.org>
  - <http://www.openmp.org/specifications/>
- **OpenMP lecture**, François Broquedis (Corse), CERMACS School 2016
  - <http://smai.emath.fr/cemracs/cemracs16/programme.php>
- **OpenMP lecture**, Françoise Roch (Grenoble)
- **IDRIS lecture and lab work**
  - <http://www.idris.fr/formations/openmp/>
- **Using OpenMP , Portable Shared Memory Model**, Barbara Chapman
- **Parallel Programming in C with MPI and OpenMP**, M.J. Quinn
- **Programming Models for Parallel Computing**, P. Balaji
- **Parallel Programming – For Multicore and Cluster System**, T. Rauber, G. Rünger

# Introduction



# Introduction

**Programming model:** how to write (and describe) a parallel program

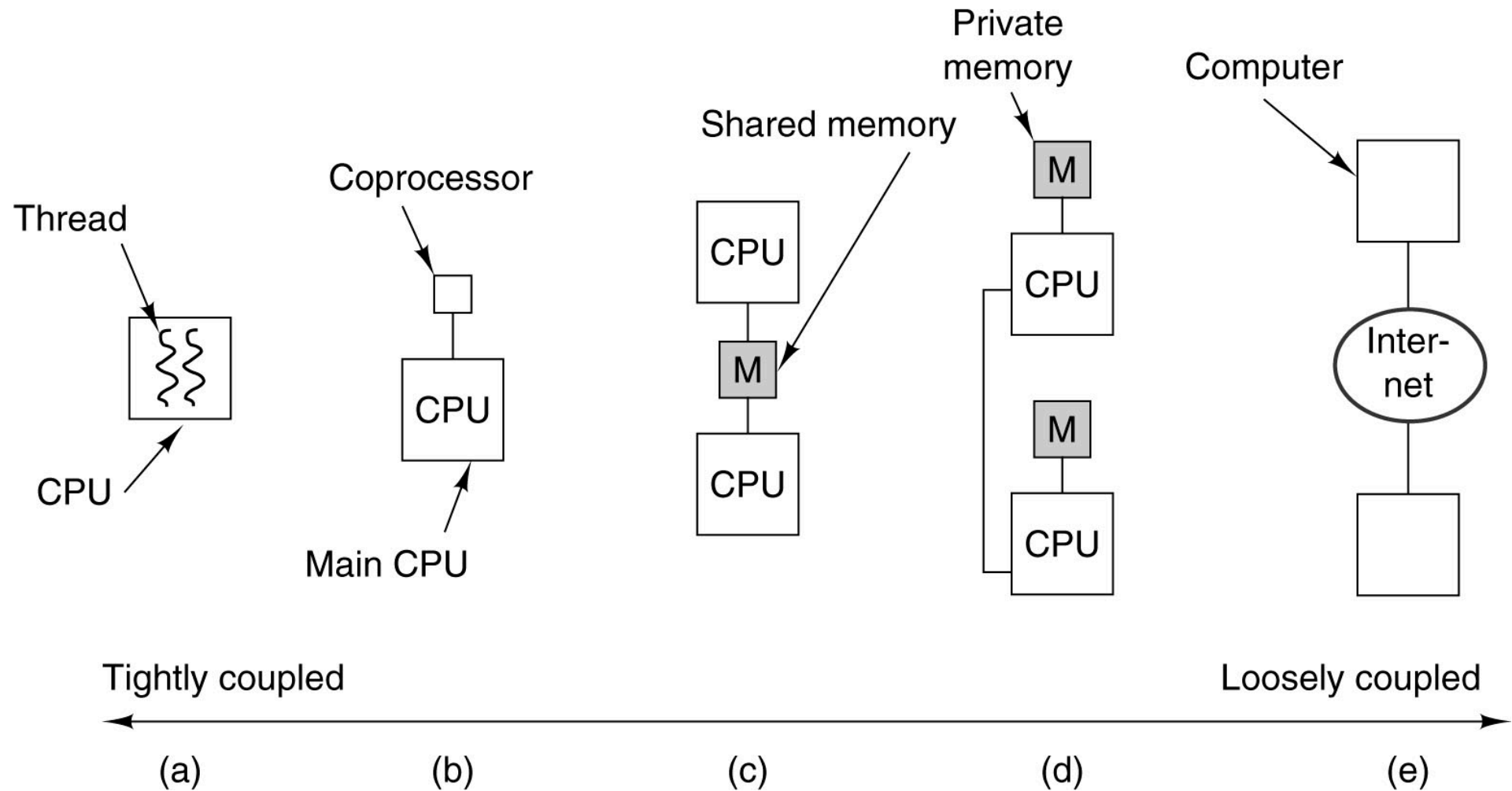
## **We will learn MPI (Message Passing Interface)**

- The programmer manages everything (data distribution, computation distribution, processors synchronization, data exchanges)
- **Avantages**
  - Greater control from the programmer
  - Performances (if the code is well written!)
- **Drawbacks**
  - Parallelisme assembly code
  - Performance portability
  - Less transparency

## **Other solution**

- Give more work to the compiler and the runtime system!

# Parallel architectures



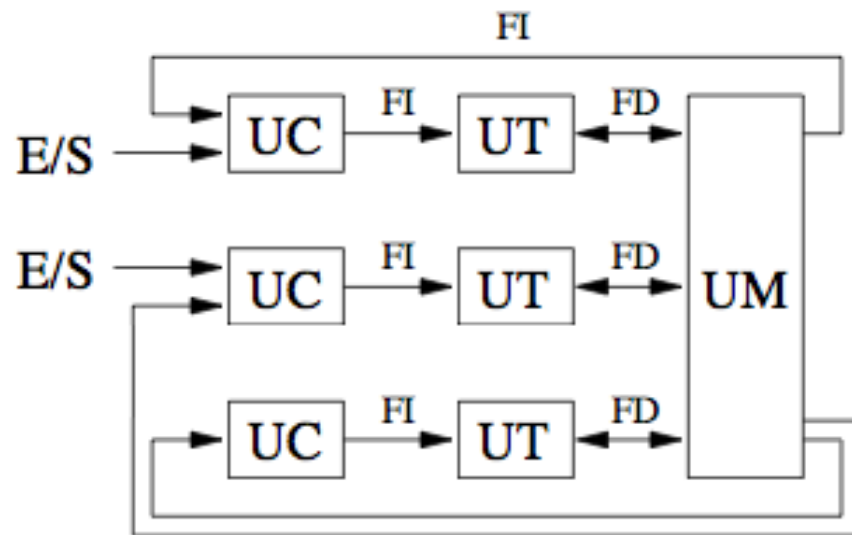
# MIMD: Multiple Instructions stream, multiple data stream

Multi-Processor Machines

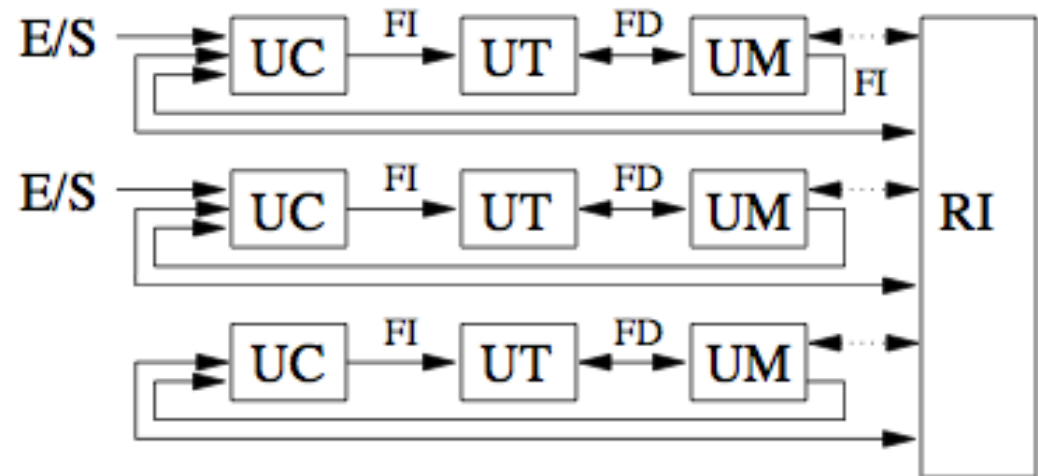
Each processor runs its own code asynchronously and independently

**Two sub-classes**

**Shared memory**



**Distributed memory**



A mix between SIMD and MIMD: SPMD (**Single Program, Multiple Data**)

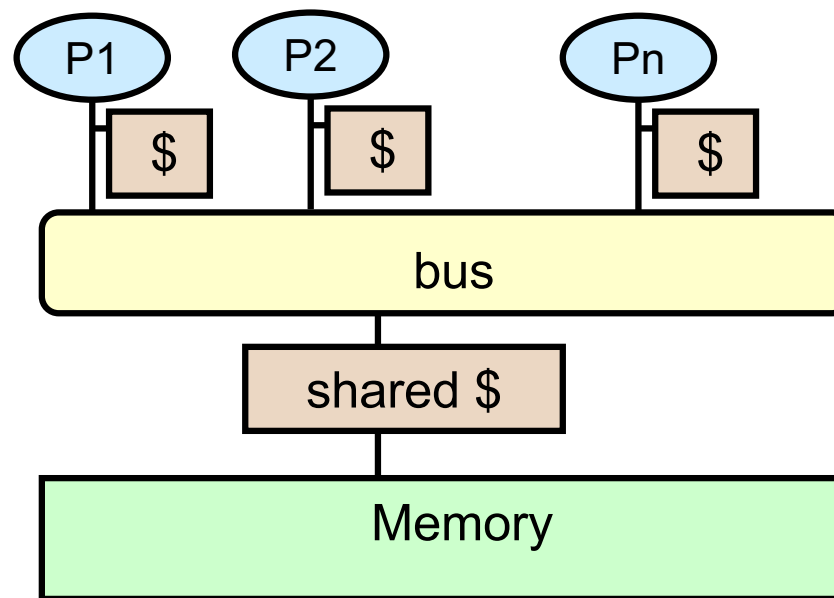
# Shared memory machine model

Processors are connected to a large shared memory

- Also known as *Symmetric Multiprocessors* (SMPs)
- SGI, Sun, HP, Intel, SMPs IBM
- Multicore processors (except that caches are shared)

Scalability issues for large numbers of processors

- Usually  $\leq 32$  processors
- Uniform memory access (*Uniform Memory Access*, UMA)
- Lower cost for caches compared to the main memory



Note: \$ = cache

# HPC architecture are getting more and more hierarchical

- **Parallelism is everywhere !**

- At the architecture level

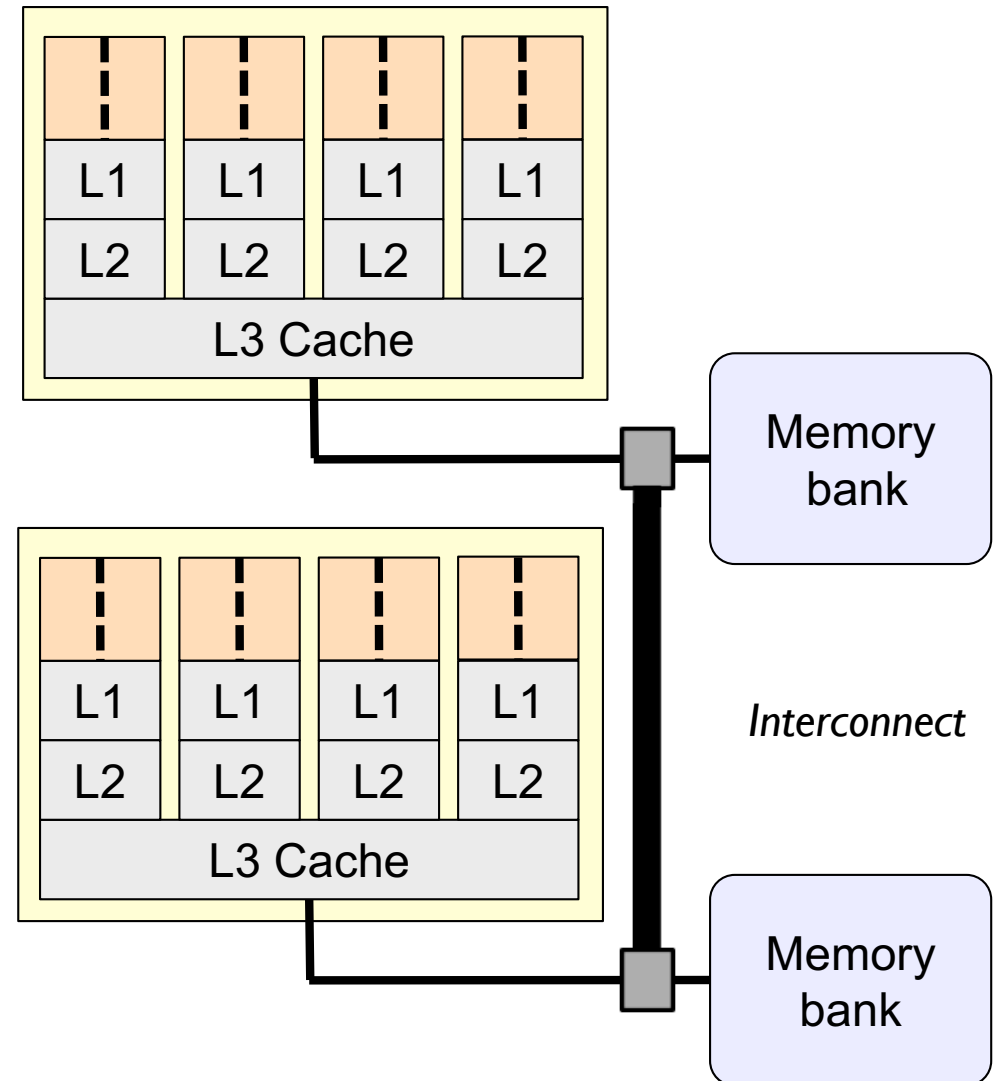
- SMP
    - NUMA

- At the processor level

- Multicore chips

- **Current (solid) trend: back to the cc-NUMA era**

- AMD Hypertransport or Intel QuickPath to connect multicore chips together in a NUMA fashion





# How to program these parallel machines?

- **The « good old » thread library**
  - A way to achieve the best performance for a particular instance of a problem (architecture, application, data set)
  - Not portable, most of the time...

## The « user-friendly » (...) parallel programming environments

- MPI
  - Standard for distributed programming
- OpenMP
  - De-facto standard for shared-memory programming
- and all these great programming languages I won't talk about today
  - Cilk+, TBB, Charm++, UPC, X10, Chapel, OpenCL, CUDA, OpenACC, ...



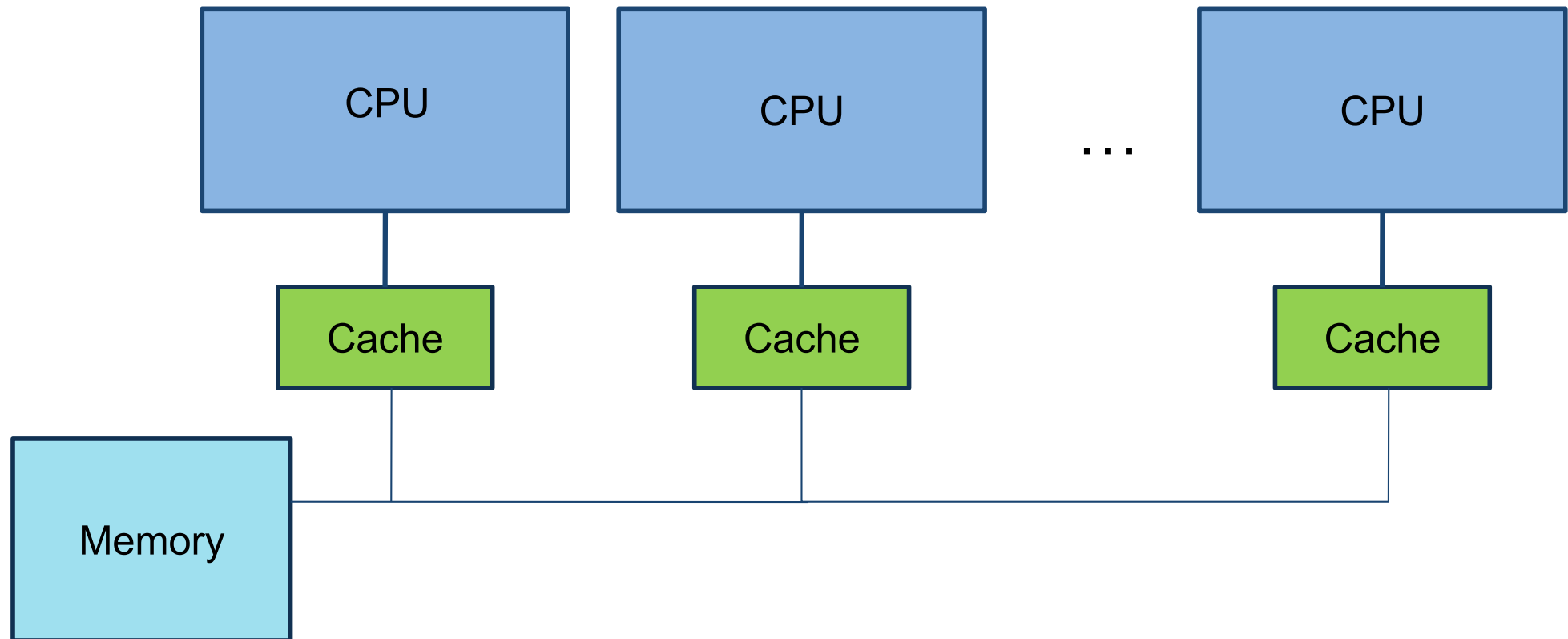
# Multi-task programming model on shared memory architecture

- Several tasks are executed in parallel
- Memory is shared (physically or virtually)
- Communication between tasks is done by reads and writes in the shared memory.
- Eg. The general-purpose multi-core processors share a common memory
  - Tasks can be assigned to distinct cores

# Multi-task programming model on shared memory architecture

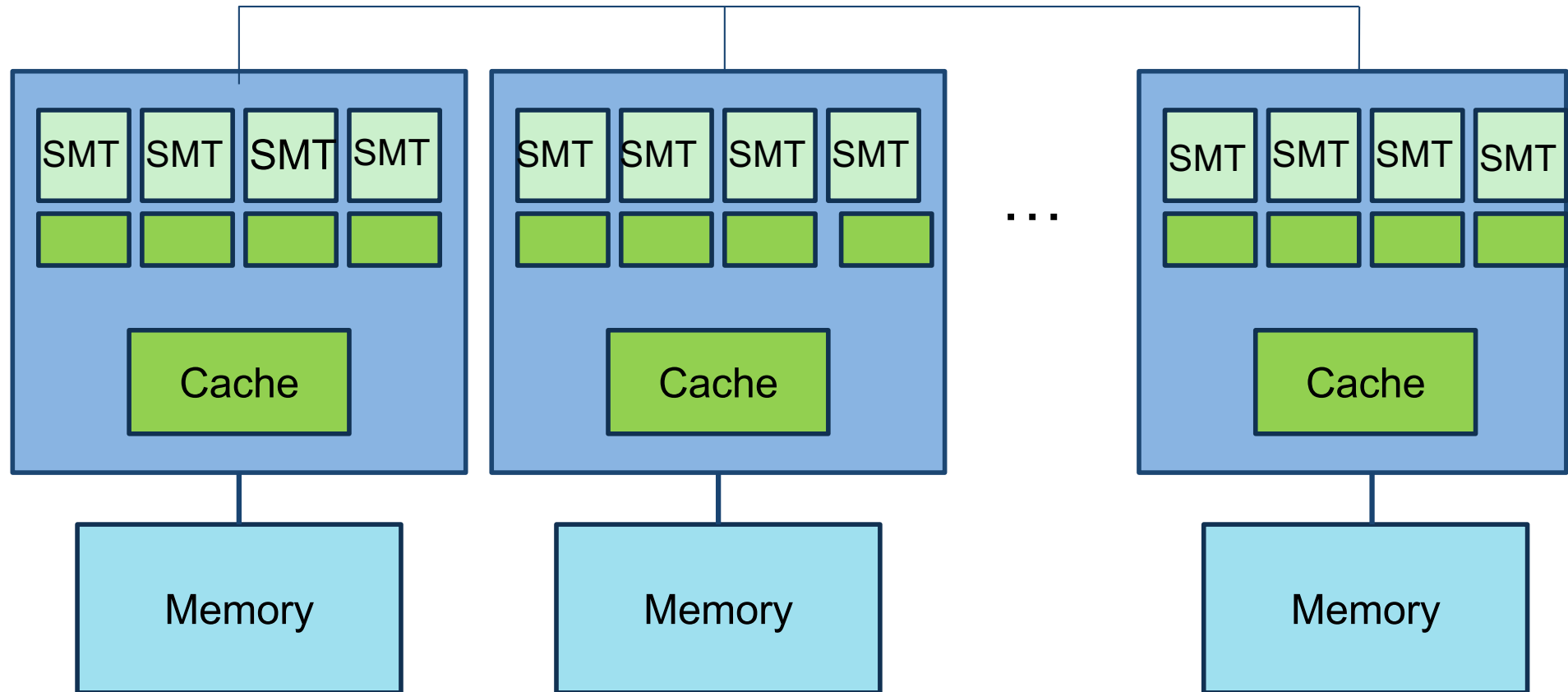
- The Pthreads library: POSIX thread library, adopted by most operating systems
  - The writing of a code requires a considerable number of lines specifically dedicated to threads
  - Example: parallelizing a loop involves
    - Declare thread structures,
    - create threads,
    - compute loop boundaries,
    - assign them to threads, ...
- OpenMP: a simpler alternative for the programmer

# Multi-task programming on UMA architectures



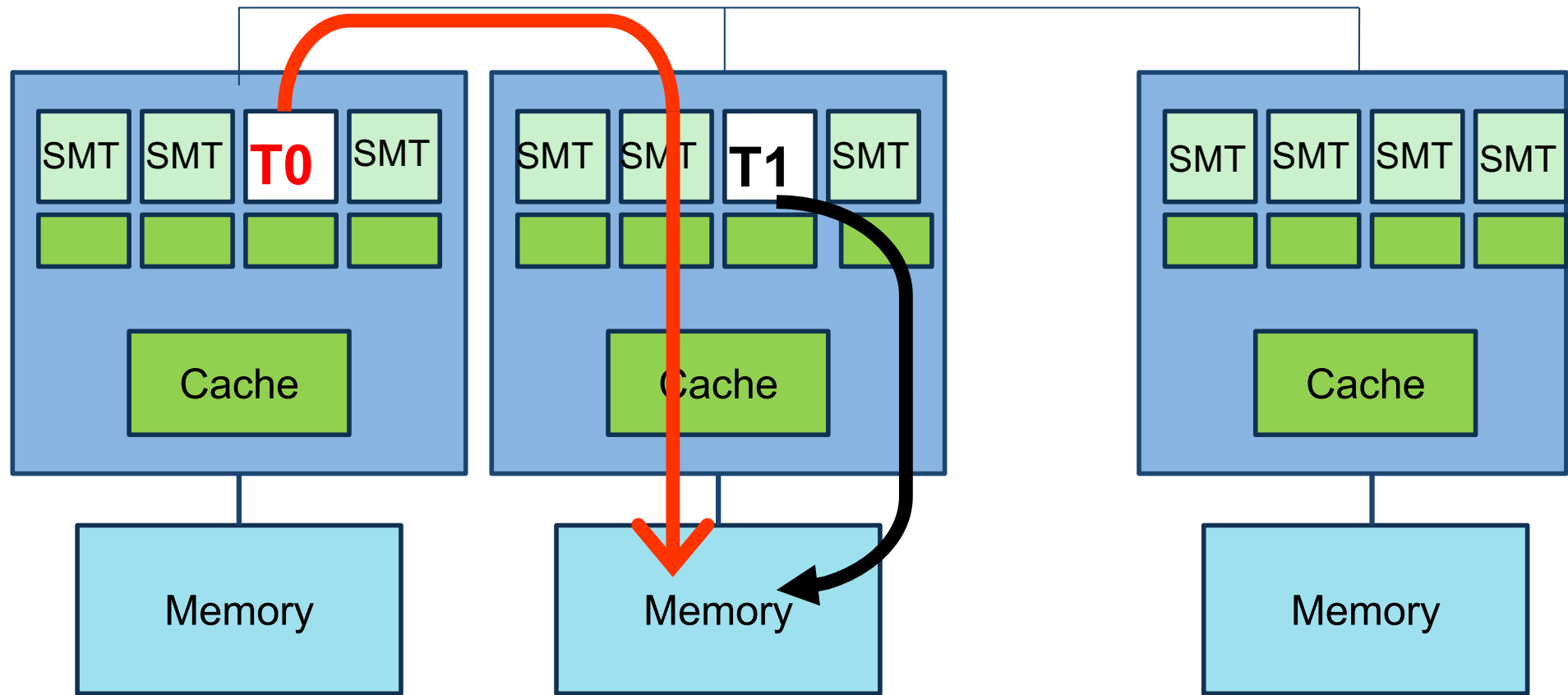
- Memory is shared
  - Uniform Memory Access Architectures (UMA)
  - An inherent problem: memory contentions

# Multi-task programming on UMA multicore architectures



- Memory is directly attached to multicore chips
  - Non-Uniform Memory Access architectures (NUMA)

# Multi-task programming on UMA multicore architectures



**Distant access**

**Local access**

# OpenMP

- A de-facto standard API to write shared memory parallel applications in C, C++ and Fortran
- Consists of compiler directives, runtime routines and environment variables
- Specification maintained by the OpenMP Architecture Review Board (<http://www.openmp.org>)
- Current version of the specification: 4.5 (November 2015)



# Advantages of OpenMP

- A **mature** standard
  - Speeding-up your applications since 1998
- **Portable**
  - Supported by many compilers, ported on many architectures
- Allows **incremental parallelization**
- Imposes low to **no overhead on the sequential execution** of the program
  - Just tell your compiler to ignore the OpenMP pragmas and you get back to your sequential program
- Supported by a wide and active community
  - The specifications have been moving fast since revision 3.0 (2008) to support:
    - new kinds of parallelism (tasking)
    - new kinds of architectures (accelerators)

# OpenMP model characteristics

## Advantages

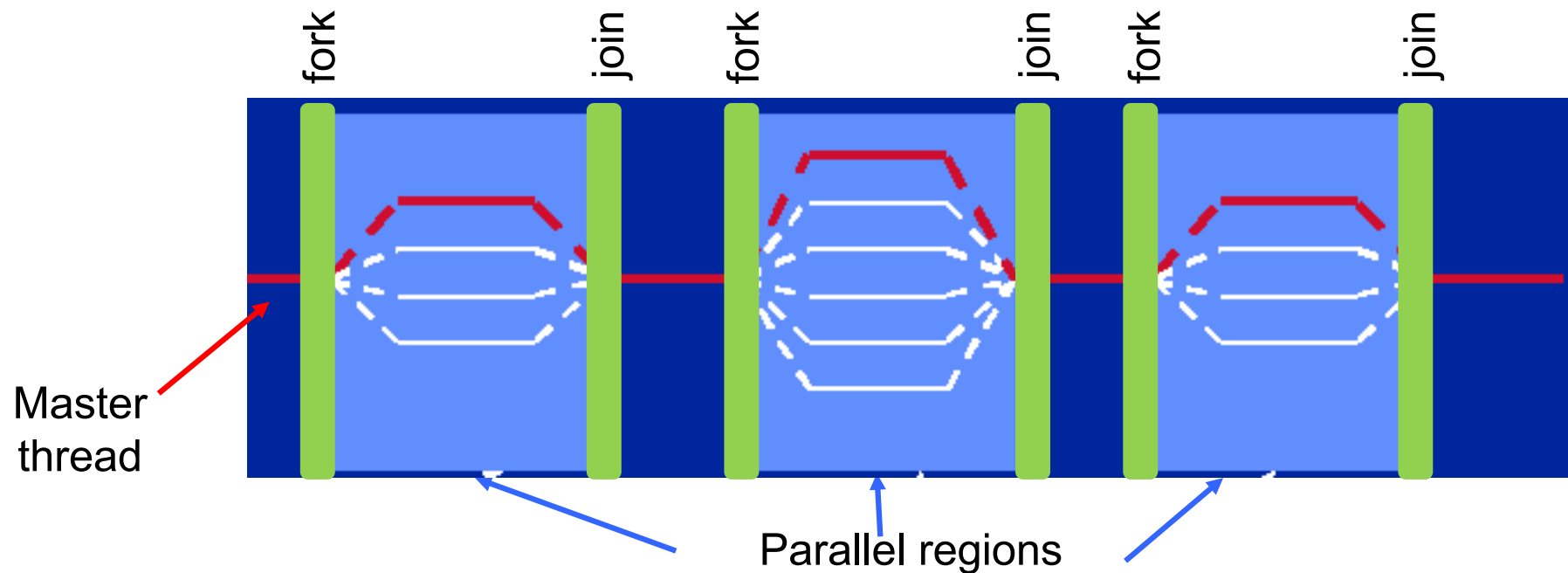
- Transparent and portable thread management
- Easy programming

## Drawbacks

- Data locality problem
- Shared but non-hierarchical memory
- Efficiency not guaranteed (impact of the material organization of the machine)
- Limited scalability, moderate parallelism

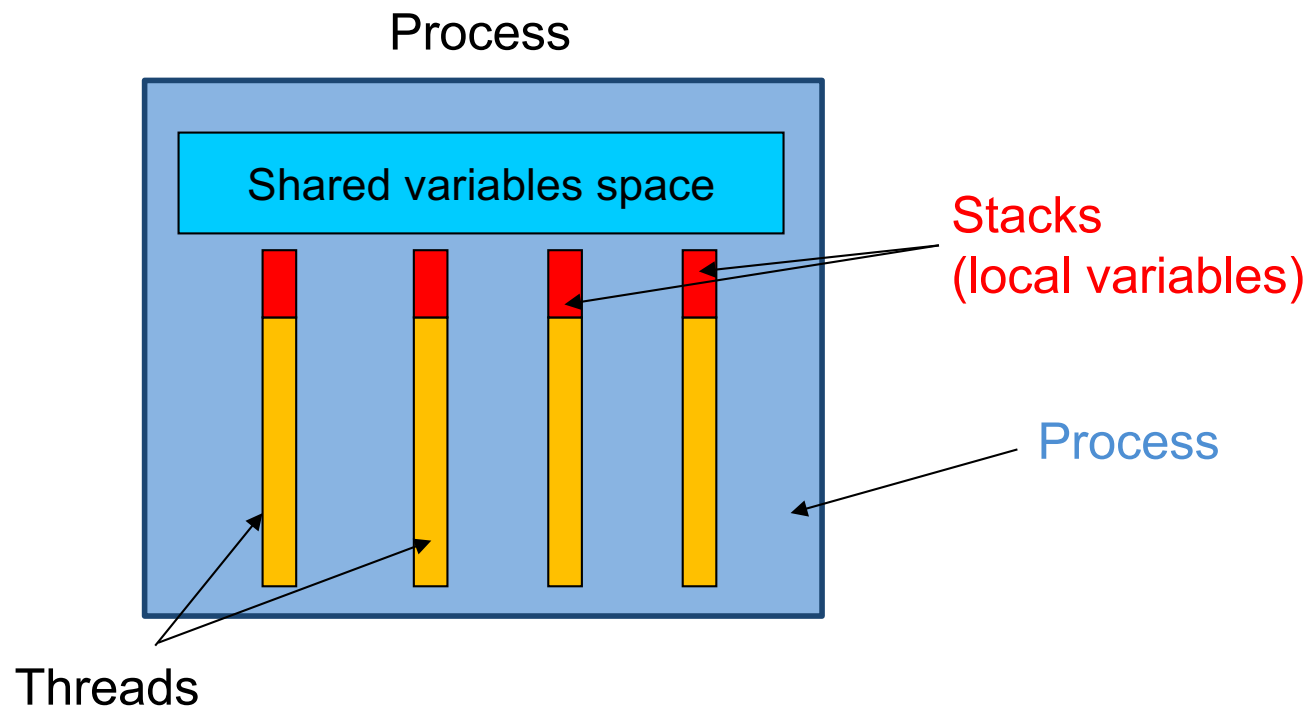
# Introduction: execution model

- An OpenMP program is executed by a **unique process** (on one or many cores)
- **Fork-Join Parallelism**
  - Master thread spawns a team of threads as needed
  - Parallelism is **added incrementally**: that is, the sequential program evolves into a parallel program
    - Entering a parallel region will **create** some threads (*fork*)
    - Leaving a parallel region will **terminate** them (*join*)
    - Any statement executed outside parallel regions are executed **sequentially**



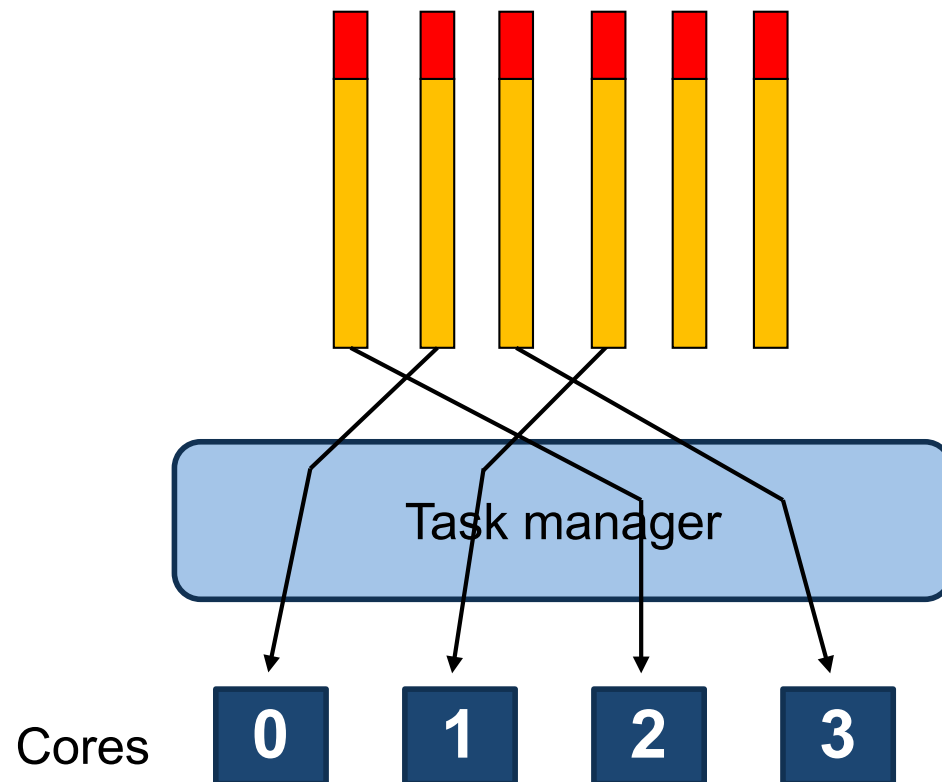
# Introduction: threads

- Threads access the same resources as the main process
- They have a stack (stack, stack pointer and clean instructions pointer)

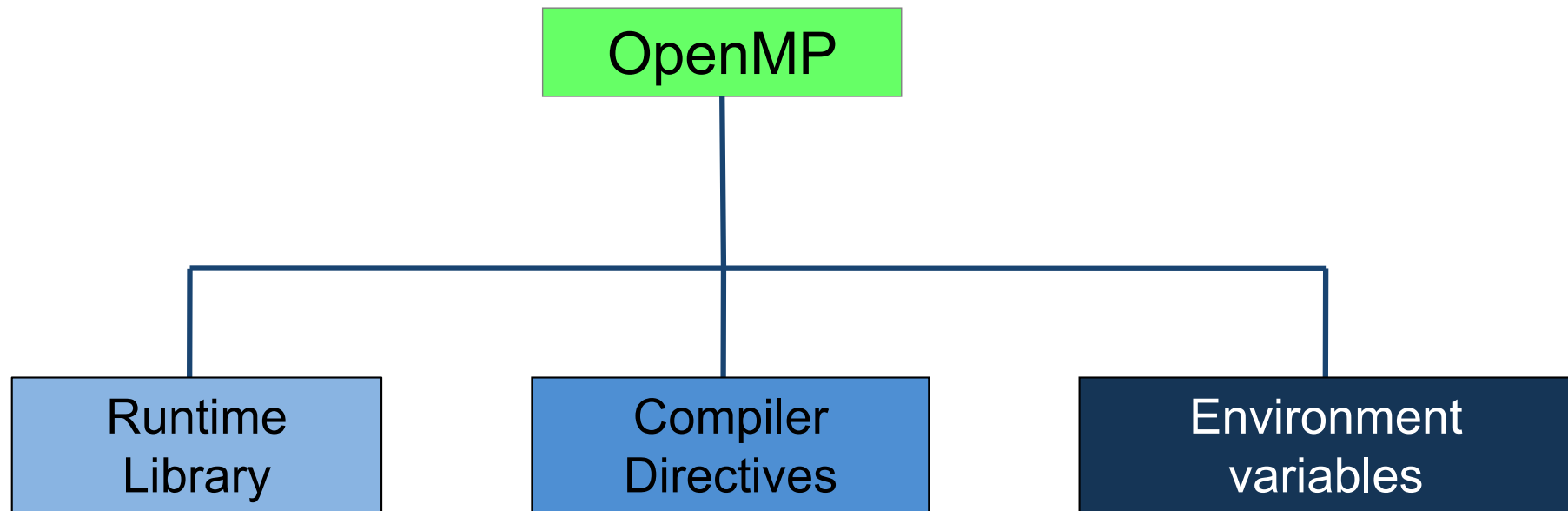


# Introduction: execution of an OpenMP program on a multicore

The task management system of the operating system assigns the tasks on the cores



# OpenMP structure: software architecture



# OpenMP structure: directives/pragmas formats

- directive [clause[clause]..]

Fortran

```
!$OMP PARALLEL PRIVATE(a,b) &  
!$OMP FIRSTPRIVATE(c,d,e)  
...  
!$OMP END PARALLEL
```

C/C++

```
#pragma omp parallel private(a,b)  
                        firstprivate(c,d,e)  
{  
    ...  
}
```

- The line is interpreted if openmp option to the compiler call otherwise comment  
→ portability

# OpenMP structure: prototyping

We have

- A Fortran 95 module OMP\_LIB
- An C/C++ input file omp.h

that define the prototypes of all the functions of the OpenMP library

## Fortran

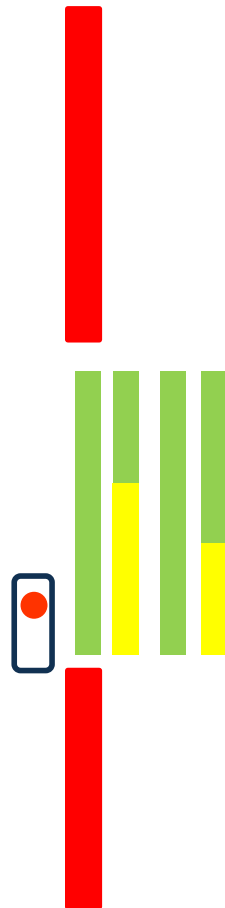
```
Program example
!$ USE OMP_LIB
!$OMP PARALLEL PRIVATE(a,b) &
...
tmp= OMP_GET_THREAD_NUM()
!$OMP END PARALLEL
```

## C/C++

```
#include <omp.h>
```



# OpenMP structure: construction of a parallel region



fortran

```
PROGRAM example
```

```
!$ USE OMP_LIB
```

```
Integer :: a, b, c
```

```
! Sequential code sequential executed by  
the master
```

```
!$OMP PARALLELPRIVATE(a,b) &  
!$OMP SHARED(c)
```

```
.  
! Parallel zone executed by all the  
! threads
```

```
!$OMP END PARALLEL
```

```
! Sequential code
```

```
END PROGRAM example
```

C/C++

```
#include <omp.h>
```

```
main () {
```

```
Int a,b,c;
```

```
/* Sequential code sequential executed by  
the master */
```

```
#pragma omp parallel private(a,b) \  
shared(c)
```

```
{  
    /* Parallel zone executed by all the  
    threads */  
}
```

```
/* Sequential code */
```

```
}
```

# Hello world !

```
void main()  
{  
    int ID = 0;  
    printf(" hello(%d) ", ID);  
    printf(" world(%d) \n", ID);  
}
```

# OpenMP's Hello world !

```
#include "omp.h"
void main()
{
    #pragma omp parallel
    {
        int ID = omp_get_thread_num();
        printf(" hello(%d) ", ID);
        printf(" world(%d) \n", ID);
    }
}
```

OpenMP include file

Parallel region with default number of threads

Runtime library function to return a thread ID.

End of the Parallel region

## Sample Output

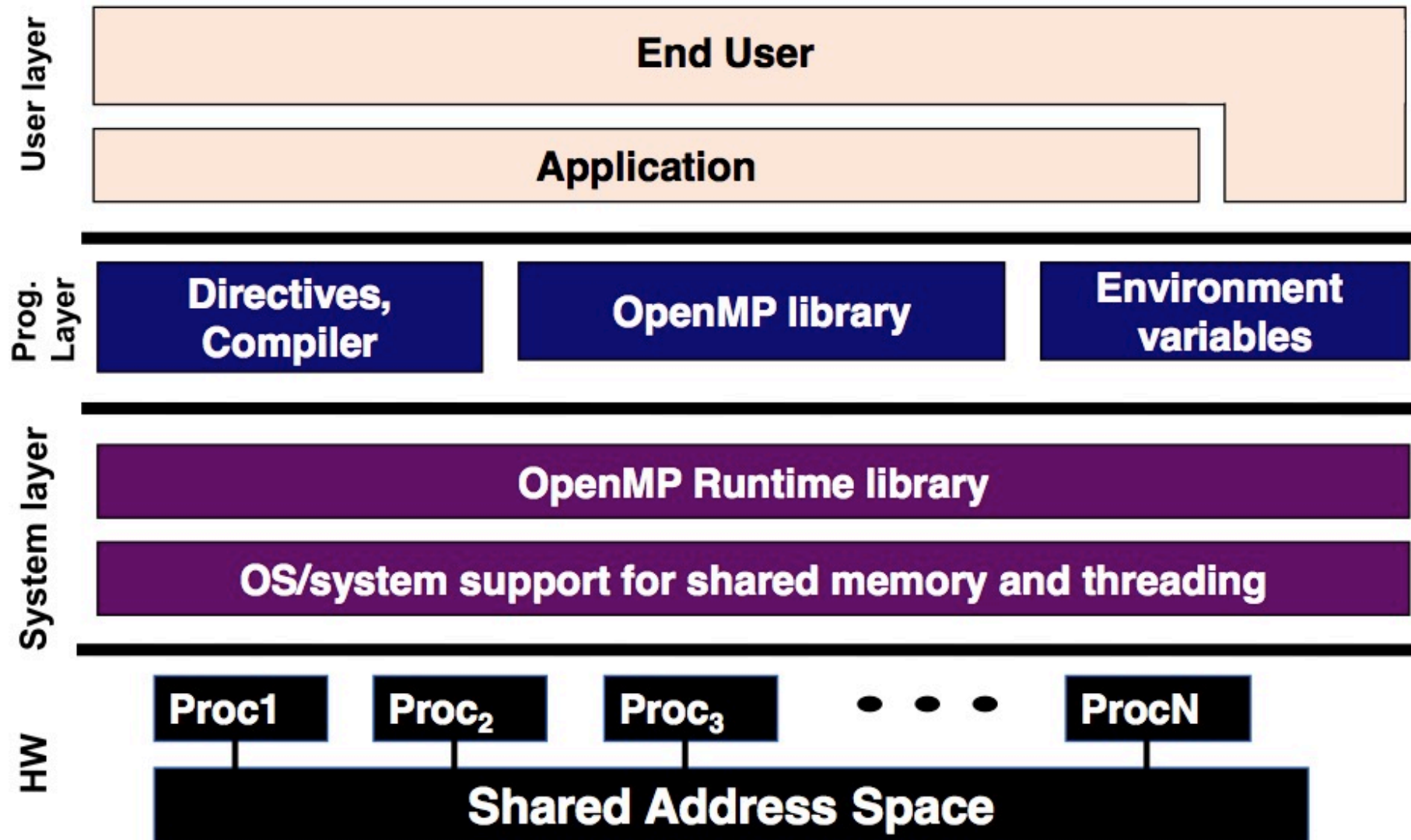
```
hello(1) hello(0) world(1)
world(0)
hello (3) hello(2) world(3)
world(2)
```

## Example: Hello world Solution Calling the OpenMP compiler

```
#include "omp.h"
void main()
{
#pragma omp parallel
{
    int ID = omp_get_thread_num();
    printf(" hello(%d) ", ID);
    printf(" world(%d) \n", ID);
}
}
```

<b>Linux and OS X</b>	<b>gcc -fopenmp</b>
<b>PGI Linux</b>	<b>pgcc -mp</b>
<b>Intel windows</b>	<b>icl /Qopenmp</b>
<b>Intel Linux and OS X</b>	<b>icpc -openmp</b>

# OpenMP Basic Defs: Solution Stack



# IF clause of the PARALLEL directive

Conditional creation of a parallel region **IF(logical\_expression)** clause

```
_____ fortran _____  
  
! Sequential code  
  
!$OMP PARALLEL IF(expr)  
  
! Parallel or sequential code depending of the expr value  
  
!$OMP END PARALLEL  
  
! Sequential code
```

The logical expression will be evaluated before entering the parallel region

# How do threads interact?

- OpenMP is a multi-threading, shared address model
  - Threads communicate by sharing variables
- Unintended sharing of data causes race conditions:
  - race condition: when the program's outcome changes as the threads are scheduled differently
- To control race conditions
  - Use synchronization to protect data conflicts
- Synchronization is expensive so
  - Change how data is accessed to minimize the need for synchronization

# OpenMP threads

## Number of threads definition

- Through an environment variable: OMP\_NUM\_THREADS
- Through the routine: OMP\_SET\_NUM\_THREADS()
- Through the clause NUM\_THREADS() of the PARALLEL directive

## Threads are numbered

- The number of threads is not necessary equal to the number of physical cores
- thread #0 is the master task
- OMP\_GET\_NUM\_THREADS(): number of threads
- OMP\_GET\_THREAD\_NUM(): thread number
- OMP\_GET\_MAX\_THREADS(): maximum number of threads



# OpenMP structure: compilation and execution

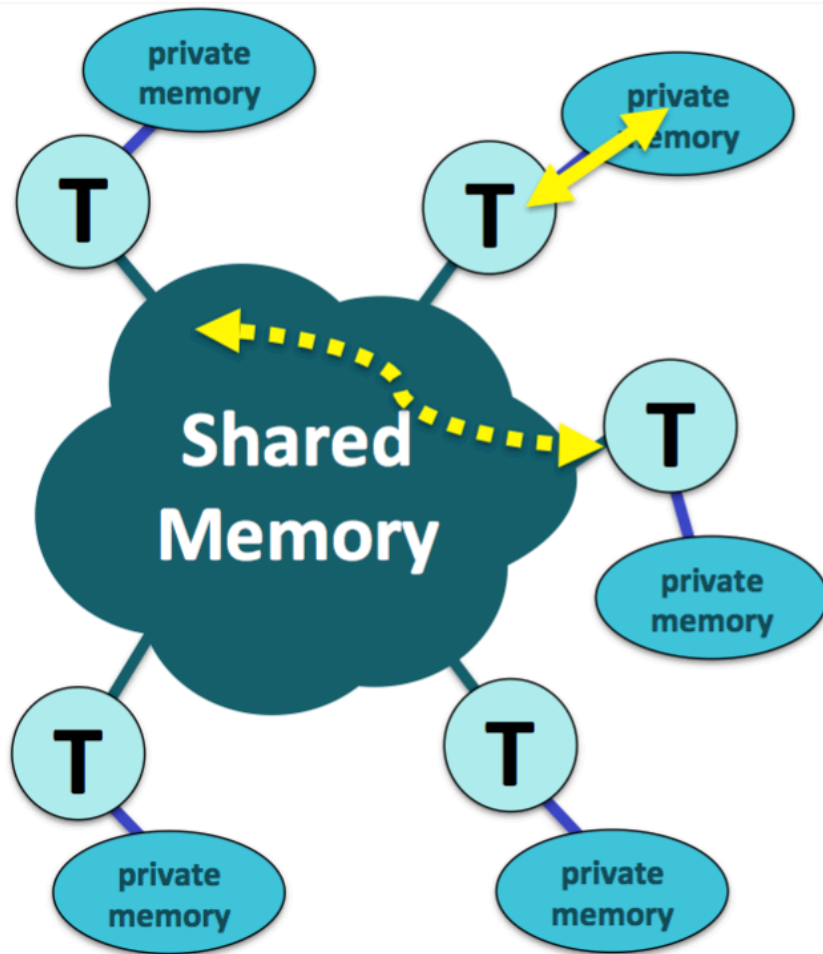
ifort (ou icc) –openmp prog.f	(INTEL)
f90 (ou cc ou CC) –openmp prog.f	(SUN Studio)
gcc/gfortran –fopenmp –std=f95 prog.f	(GNU)
export OMP_NUM_THREADS=2	

./a.out

# ps –eLF

USER	PID	PPID	LWP	C	NLWP	SZ	RSS	PSR ...
------	-----	------	-----	---	------	----	-----	---------

# The OpenMP memory model



- All the threads have access to the same **globally shared** memory
- Each thread has access to **its own private memory area** that can not be accessed by other threads
- Data transfer is performed through shared memory and is **100% transparent to the application**
- The application programmer is responsible for providing the corresponding data-sharing attributes

# Data sharing attributes

- Need to set the visibility of each variable that appears inside an OpenMP parallel region using the following **data-sharing attributes**
- **shared**: the data can be read and written by any thread of the team. All changes are visible to all threads
- **private**: each thread is working on its own version of the data that cannot be accessed by other threads of the team
- **firstprivate**: each thread is working on its own version of the variable. The data is initialized using the value it had before entering the parallel region
- **lastprivate**: each thread is working on its own version of the variable. The value of the last thread leaving the region is copied back to the variable.

# Variable status

The status of a variable in a parallel zone

- SHARED, it's located in the global memory
- PRIVATE, it's located in the thread of each thread. It's value is undefined at the entrance of the zone
- Declaring the variable status
  - # pragma omp parallel private (list)
  - # pragma omp parallel firstprivate (list)
  - # pragma omp parallel shared (list)
- Declaring a default status
  - DEFAULT(PRIVATE|SHARED|NONE) clause

```
program private_var.f
  !$USE OMP_LIB
  integer:: tmp =999
  Call OMP_SET_NUM_THREADS(4)
  !$OMP PARALLEL PRIVATE(tmp)
    print *, tmp
    tmp = OMP_GET_THREAD_NUM()
    print *, OMP_GET_THREAD_NUM(), tmp
  !$OMP END PARALLEL
  print *, tmp
end
```

# Putting Threads to Work: the Worksharing Constructs

```
1 void simple_loop(int N,  
2                 float *a,  
3                 float *b)  
4 {  
5     int i;  
6     // i, N, a and b are shared by  
7     // default  
8     #pragma omp parallel firstprivate(N)  
9     {  
10        // i is private by default  
11        #pragma omp for  
12        for (i = 1; i <= N; i++) {  
13            b[i] = (a[i] + a[i-1]) / 2.0;  
14        }  
15    }
```

- **omp for** : distribute the iterations of a loop over the threads of the parallel region.
- Here, assigns  $N/P$  iterations to each thread,  $P$  being the number of threads of the parallel region.
- **omp for** comes with an implicit **barrier synchronization** at the end of the loop one can remove with the **nowait** keyword.

# Work sharing

- Distributing a loop between threads (// loop)
- Distribution of several sections of code between threads, one section of code per thread (// sections)
- Running a portion of code on a single thread
- Execution of several occurrences of the same function by different threads
- Execution by different threads of different work units, tasks

# Work sharing: parallel loop

## DO Directive in Fortran, **for** in C

Parallelism by distribution of iterations of a loop

- The way in which the iterations can be distributed can be specified in the **SCHEDULE** clause (coded in the program or by an environment variable)
- A global synchronization is performed at the end of construction **END DO** (unless **NOWAIT**)
- Possibility to have several **DO** constructions in a parallel region
- The loop indices are integers and private
- Infinite loops and **do while** are not parallelizable

# DO and PARALLEL DO Directives

```
Program loop
implicit none
integer, parameter :: n=1024
integer          :: i, j
real, dimension(n, n) :: tab
!$OMP PARALLEL
...              ! Replicated code
!$OMP DO
do j=1, n        ! Shared loop
do i=1, n        ! Replicated loop
    tab(i, j) = i*j
end do
end do
!$OMP END DO
!$OMP END PARALLEL
end program loop
```

```
Program parallelloop
implicit none
integer, parameter :: n=1024
integer          :: i, j
real, dimension(n, n) :: tab
!$OMP PARALLEL DO
do j=1 n         ! Shared loop
do i=1, n        ! Replicated loop
    tab(i, j) = i*j
end do
end do
!$OMP END PARALLEL DO
end program parallelloop
```

PARALLEL DO is a fusion of 2 directives

Beware: END PARALLEL DO includes a synchronization barrier!

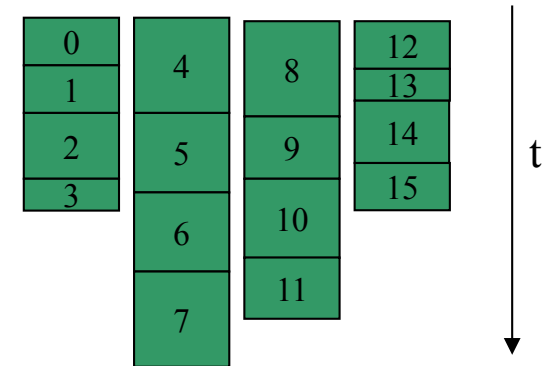


# Work sharing: SCHEDULE

**!\$OMP DO SCHEDULE(STATIC, packet-size)**

By default packet-size =  $\#\_iterations / \#\_threads$

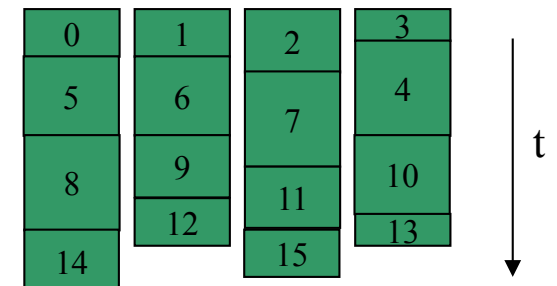
Ex: 16 iterations (0 to 15), 4 threads: packet size by default is 4



**!\$OMP DO SCHEDULE(DYNAMIC, packet-size)**

Packets are distributed to free threads in a dynamic way

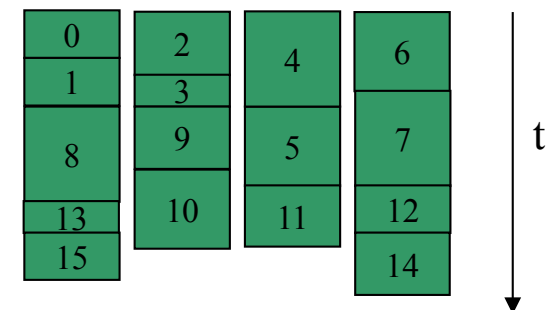
All the packets have the same size (except maybe the last one), by default the packet size is one



**!\$OMP DO SCHEDULE(GUIDED, packet-size)**

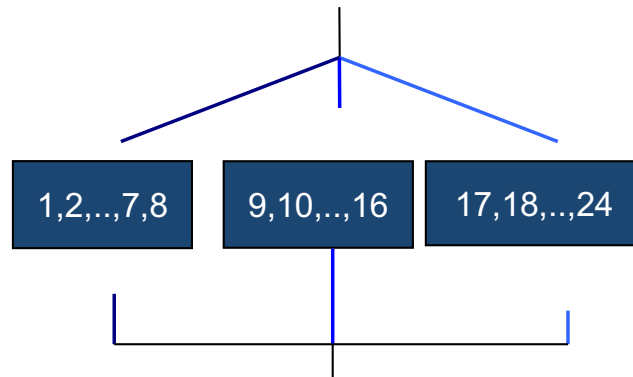
Packet-size: minimal packet size (1 by default) except the last one

Maximal packet size at the beginning of the loop (here 2) then decrease to balance the load

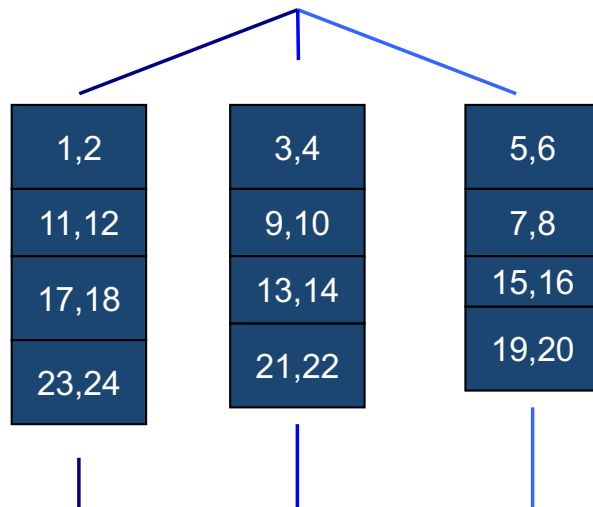


# Work sharing: SCHEDULE

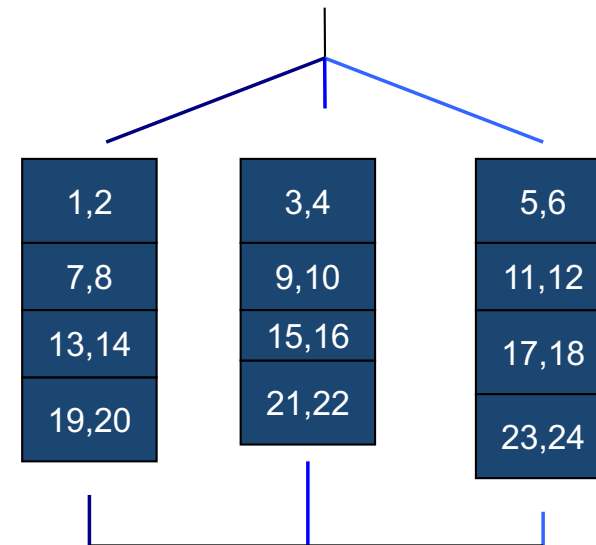
Ex: 24 iterations, 3 threads



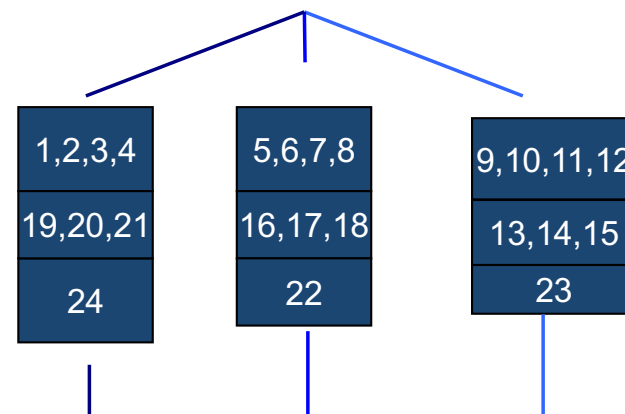
static mode with  
Packet size = # iterations/# threads



Greedy: dynamic



Cyclic: static



Greedy: guided

# Work sharing: SCHEDULE

The choice of the repartition mode can be delayed at the execution time using SCHEDULE(RUNTIME)

Taking into account the environment variable OMP\_SCHEDULE

- Ex

```
export OMP_SCHEDULE="DYNAMIC,400"
```

# A first example to illustrate OpenMP capabilities

Parallelize this simple code using OpenMP

```
f = 1.0
```

```
for (i = 0; i < N; i++)  
    z[i] = x[i] + y[i];
```

```
for (i = 0; i < M; i++)  
    a[i] = b[i] + c[i];
```

```
...
```

```
scale = sum (a, 0, m) + sum (z, 0, n) + f;
```

```
...
```

# A first example to illustrate OpenMP capabilities

First create the parallel region and define the data-sharing attributes


```
#pragma omp parallel default (none) shared (z, x, y, a, b, c, n, m)
private (f, i, scale)
{
    f = 1.0

    for (i = 0; i < n; i++)
        z[i] = x[i] + y[i];

    for (i = 0; i < m; i++)
        a[i] = b[i] + c[i];

    ...

    scale = sum (a, 0, m) + sum (z, 0, n) + f;
    ...
} /* End of OpenMP parallel region */
```

A blue L-shaped arrow is positioned to the right of the code. The vertical part of the arrow points upwards from the closing brace of the parallel region to the opening brace, with the text "parallel region" written vertically along its side. The horizontal part of the arrow points left from the vertical part to the closing brace of the parallel region. Another horizontal arrow points left from the vertical part to the opening brace of the parallel region.

# A first example to illustrate OpenMP capabilities

```
#pragma omp parallel default (none) shared (z, x, y, a, b, c, n, m)
private (f, i, scale)
{
    f = 1.0

    for (i = 0; i < n; i++)
        z[i] = x[i] + y[i];

    for (i = 0; i < m; i++)
        a[i] = b[i] + c[i];

    ...

    scale = sum (a, 0, m) + sum (z, 0, n) + f;
    ...
} /* End of OpenMP parallel region */
```

Statements executed by all the threads of the parallel region !

parallel region

At this point, all the threads execute the whole program (you won't get any speed-up from this!)

# A first example to illustrate OpenMP capabilities

Now distribute the loop iterations over the threads using omp for

```
#pragma omp parallel default (none) shared (z, x, y, a, b, c, n, m)
private (f, i, scale)
{
    f = 1.0
    #pragma omp for
    for (i = 0; i < n; i++)
        z[i] = x[i] + y[i];
    #pragma omp for
    for (i = 0; i < m; i++)
        a[i] = b[i] + c[i];
    ...
    scale = sum (a, 0, m) + sum (z, 0, n) + f;
    ...
} /* End of OpenMP parallel region */
```

Statements executed by all the threads of the parallel region

parallel loop (work is distributed)

parallel loop (work is distributed)

parallel region

Statements executed by all the threads of the parallel region

# Optimization #1: Remove Unnecessary Synchronizations

There are no dependencies between the two parallel loops, we remove the implicit barrier between the two

```
#pragma omp parallel default (none) shared (z, x, y, a, b, c, n, m)
private (f, i, scale)
{
    f = 1.0

    #pragma omp for nowait
    for (i = 0; i < n; i++)
        z[i] = x[i] + y[i];

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

    ...

    #pragma omp barrier
    scale = sum (a, 0, m) + sum (z, 0, n) + f;
    ...
} /* End of OpenMP parallel region */
```

parallel region



## Optimization #2: Don't Go Parallel if the Workload is Small

We don't want to pay the price of thread management if the workload is too small to be computed in parallel

```
#pragma omp parallel default (none) shared (z, x, y, a, b, c, n, m)
private (f, i, scale) if (n > some_threshold && m > some_threshold)
{
    f = 1.0

    #pragma omp for nowait
    for (i = 0; i < n; i++)
        z[i] = x[i] + y[i];

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

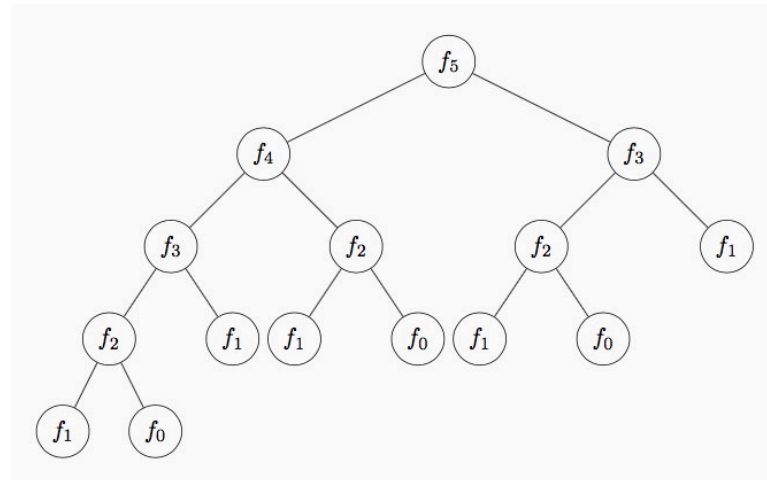
    ...

    #pragma omp barrier
    scale = sum (a, 0, m) + sum (z, 0, n) + f;
    ...
} /* End of OpenMP parallel region */
```

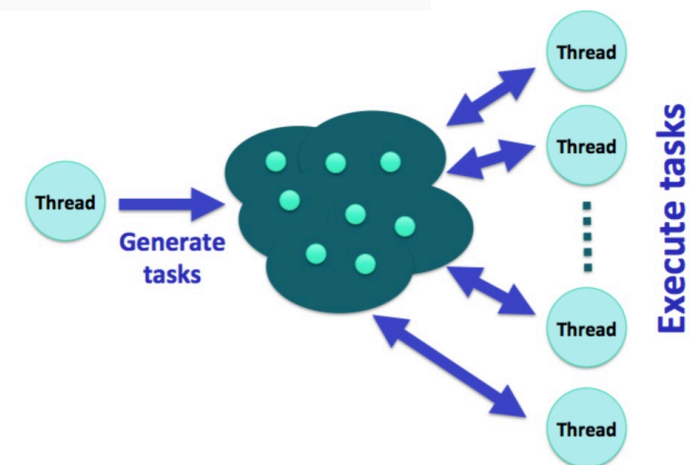
# Extending the Scope of OpenMP with Task Parallelism

- **omp for** has made OpenMP popular and remains for most users its central feature. But what if my application was not written in a loop-based fashion?

```
1 int fib(int n) {  
2     int i, j;  
3     if (n < 2) {  
4         return n;  
5     } else {  
6         i = fib(n - 1)  
7         ;  
8         j = fib(n - 2)  
9         ;  
10        return i + j;  
11    }
```



- The OpenMP tasking concept : tasks generated by one OpenMP thread can be executed by any of the threads of the parallel region



# Tasking in OpenMP: Basic Concept (cont'd)

- The application programmer specifies regions of code to be executed in a task with the **#pragma omp task** construct
- All tasks can be executed *independently*
- When any thread encounters a task construct, a task is generated
- Tasks are executed **asynchronously** by any thread of the parallel region
- Completion of the tasks can be guaranteed using the **taskwait** synchronization construct

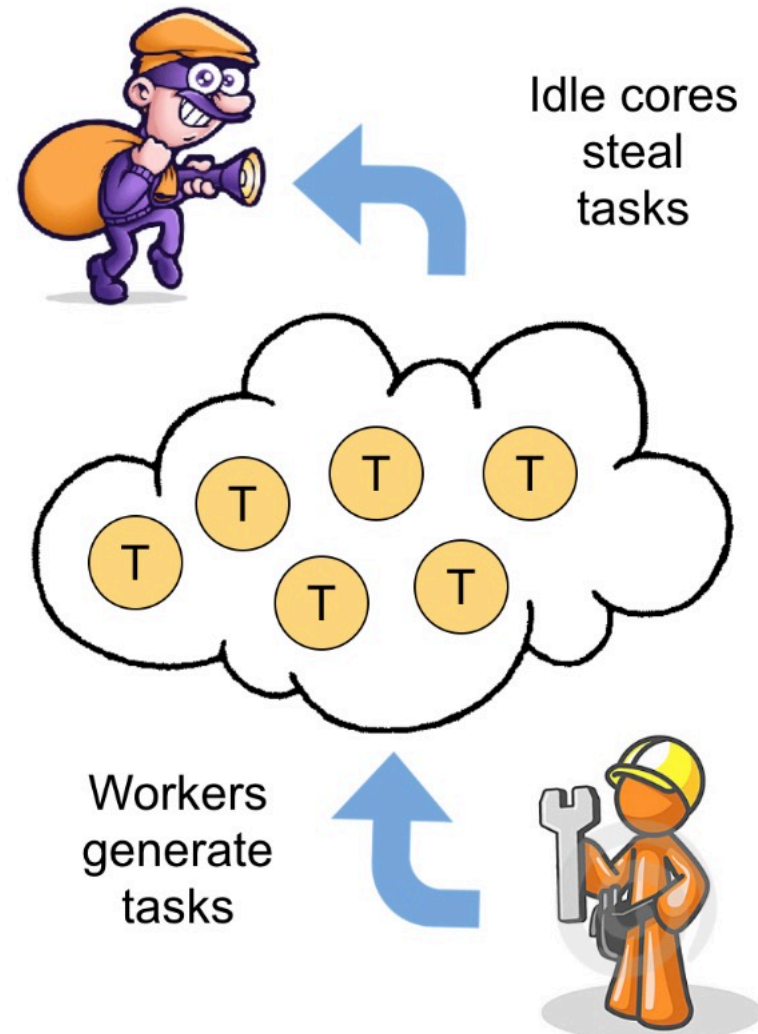
```
1  int main(void) {
2      ...
3      #pragma omp parallel
4      {
5          #pragma omp single
6          res = fib(50);
7      }
8      ...
9  }
10
11 int fib(int n) {
12     int i, j;
13     if (n < 2) {
14         return n;
15     } else {
16         #pragma omp task
17         i = fib(n - 1);
18         #pragma omp task
19         j = fib(n - 2);
20         #pragma omp taskwait
21         return i + j;
22     }
23 }
```

# Tasking in OpenMP: Execution Model

## The Work-Stealing execution model

- Each thread has its own *task queue*
- Entering an omp task construct pushes a task to the thread's local queue
- When a thread's local queue is empty, it steals tasks from other queues

Tasks are well suited to applications with irregular workload.



# First OpenMP Tasking Experience

```
1 int main(void)
2 {
3     printf("A ");
4     printf("race ");
5     printf("car ");
6
7     printf("\n");
8     return 0;
9 }
```

Program output:

```
$ OMP_NUM_THREADS=2 ./task-1
```

```
$ A race car
```

- We want to use OpenMP to make this program print either A race car or A car race using tasks.
- Here is a battle plan :
  1. Create the threads that will execute the tasks
  2. Create the tasks and make one of the thread generate them



# First OpenMP Tasking Experience, contd.

```
1 int main(void)
2 {
3     #pragma omp parallel
4     {
5         printf("A ");
6         printf("race ");
7         printf("car ");
8     }
9
10    printf("\n");
11    return 0;
12 }
```

Program output:

```
$ OMP_NUM_THREADS=2 ./task-2
```

```
$ A race A race car car
```

- We want to use OpenMP to make this program print either A race car or A car race using tasks.
- Here is a battle plan :
  1. Create the threads that will execute the tasks
  2. Create the tasks and make one of the thread generate them

# First OpenMP Tasking Experience, contd.

```
1 int main(void)
2 {
3     #pragma omp parallel
4     {
5         #pragma omp single
6         {
7             printf("A ");
8             #pragma omp
9             task
10             printf("race ")
11             ;
12             #pragma omp
13             task
14             printf("car ");
15         }
16     }
17
18     printf("\n");
19     return 0;
20 }
```

- We want to use OpenMP to make this program print either A race car or A car race using tasks.
- Here is a battle plan :
  1. Create the threads that will execute the tasks
  2. Create the tasks and make one of the thread generate them

Program output:

```
$ OMP_NUM_THREADS=2 ./task-3
$ A race car
$ OMP_NUM_THREADS=2 ./task-3
$ A car race
```

# First OpenMP Tasking Experience, contd.

```
1 int main(void)
2 {
3     #pragma omp parallel
4     {
5         #pragma omp single
6         {
7             printf("A ");
8             #pragma omp task
9             printf("race ");
10            #pragma omp task
11            printf("car ");
12
13            printf("is fun ");
14            printf("to watch ");
15        }
16    }
17
18    printf("\n");
19    return 0;
20 }
```

- Now that everything is working as intended, we would like to print is fun to watch at the end of the output string.
- This example illustrates the **asynchronous execution** of tasks.

Program output:

```
$ OMP_NUM_THREADS=2 ./task-4
$ A is fun to watch race car
$ OMP_NUM_THREADS=2 ./task-4
$ A is fun to watch car race
```



# First OpenMP Tasking Experience, contd.

```
1 int main(void)
2 {
3     #pragma omp parallel
4     {
5         #pragma omp single
6         {
7             printf("A ");
8             #pragma omp task
9             printf("race ");
10            #pragma omp task
11            printf("car ");
12            #pragma omp task
13            wait
14            printf("is fun ");
15            printf("to watch ");
16        };
17    }
18    printf("\n");
19    return 0;
20 }
```

- Now that everything is working as intended, we would like to print `is fun to watch` at the end of the output string.
- This example illustrates the **asynchronous execution** of tasks.
- To fix this, you need to explicitly wait for the completion of the tasks with **taskwait** before printing "is fun to watch"

Program output:

```
$ OMP_NUM_THREADS=2 ./task-5
$ A race car is fun to watch
$ OMP_NUM_THREADS=2 ./task-5
$ A car race is fun to watch
```

## What About Tasks with Dependencies on Other Tasks?

- Here, task A is writing some data that will be processed by task C. The same goes for task B and task D.
- The taskwait construct here makes sure task C won't execute before task A and task D before task B.
- As a side effect, task C won't execute until the execution of task B is over, creating some kind of **fake dependency** between task B and C.

```
1 void data_flow_example (void)
2 {
3     type x, y;
4
5     #pragma omp parallel
6     #pragma omp single
7     {
8         #pragma omp task
9         write_data(&x); // Task A
10        #pragma omp task
11        write_data(&y); // Task B
12
13        #pragma omp taskwait
14
15        #pragma omp task
16        print_results(x); // Task
17        C
18        #pragma omp task
19        print_results(y); // Task
20        D
21    }
```

# OpenMP Tasks Dependencies: Rationale

- The **depend** clause allows you to provide information on the way a task will access data.
- It is followed by an access mode that can be **in**, **out** or **inout**.
- Here are some examples of use for the **depend** clause:
  - **depend(in: x, y, z)**: the task will read variables **x**, **y** and **z**
  - **depend(out: res)**: the task will write variable **res**, any previous value of **res** will be ignored and overwritten
  - **depend(inout: k, buffer[0:n])**: the task will both read and write variable **k** and the content of **n** elements of **buffer** starting from index **0**
- The OpenMP runtime system dynamically decides whether a task is ready for execution or not considering its dependencies (there is no need for further user intervention here)

# OpenMP Tasks Dependencies : Some Trivial Example

```
1 void data_flow_example (void)
2 {
3     type x, y;
4
5     #pragma omp parallel
6     #pragma omp single
7     {
8         #pragma omp task depend(out:
9         x)
10        write_data(&x); // Task A
11        #pragma omp task depend(out:
12        y)
13        write_data(&y); // Task B
14
15        #pragma omp task depend(in: x
16        )
17        print_results(x); // Task C
18        #pragma omp task depend(in: y
19        )
20        print_results(y); // Task D
21    }
22 }
```

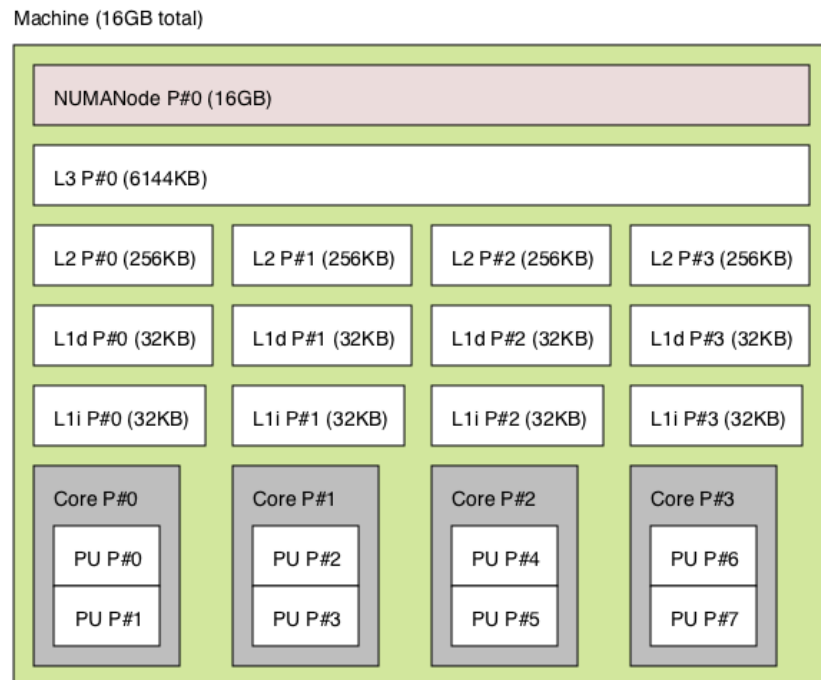
- Here is the previous example program written with tasks dependencies
- The **taskwait** construct is gone
  - The runtime system will rely on data dependencies to choose a ready task to execute
- In this version, task C could be executed before task B, as long as the execution of task A is over
- Expressing dependencies sometimes helps unlocking more parallelism

# Speeding up OpenMP applications

- **Preamble:** Have a closer look at your favorite/target platform
- Improving the execution of a parallel application **requires a good understanding of the target platform architecture**
- In particular, knowing about the following items is always useful:
  - The multicore processor: how many cores are available?
    - Which of them are physical/logical cores (HyperThreading and friends)?
  - The memory hierarchy: what kind of memory is available?
    - How is it organized?
  - The architecture topology: how (multicore) processors are connected together and how do they access memory?

# Getting to Know Your Platform with hwloc

The hwloc library gathers valuable information about your platform and synthesizes it into a generic representation.



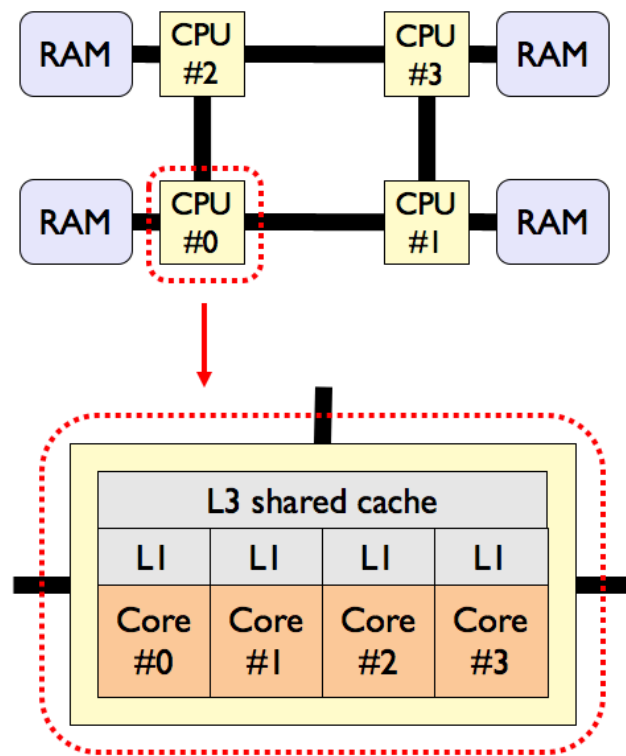
- Provides information about
  - the processing units (logical/physical cores)
  - the cache hierarchy
  - the memory hierarchy (NUMA nodes)
  - However, **hwloc does not provide the entire architecture topology** (the way processors are connected together).

<https://www.open-mpi.org/projects/hwloc/>



# Understanding the Architecture Topology

- The operating system knows about the way processors are connected together to some extent. It can provide a distance table that roughly represent how many crossbars you need to cross to access a specific NUMA node (see the hwloc-distances program)

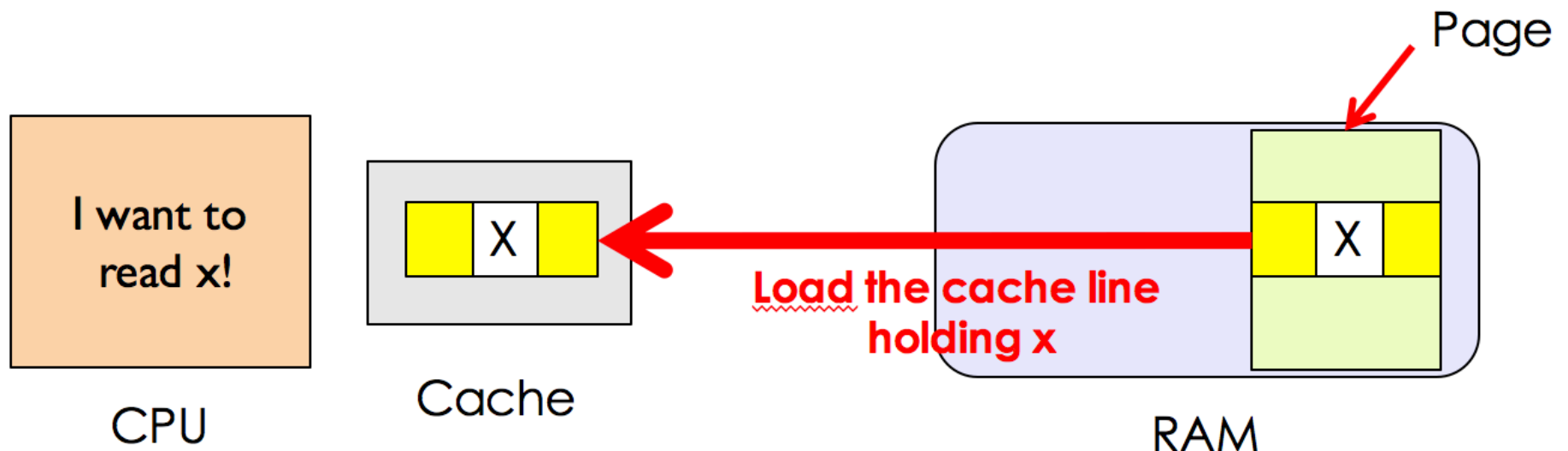


	$N_0$	$N_1$	$N_2$	$N_3$
$N_0$	0	10	10	20
$N_1$	10	0	20	10
$N_2$	10	20	0	10
$N_3$	20	10	10	0

A 4-nodes NUMA machine with the corresponding NUMA distance table

# Cache Memory: Basic Concept

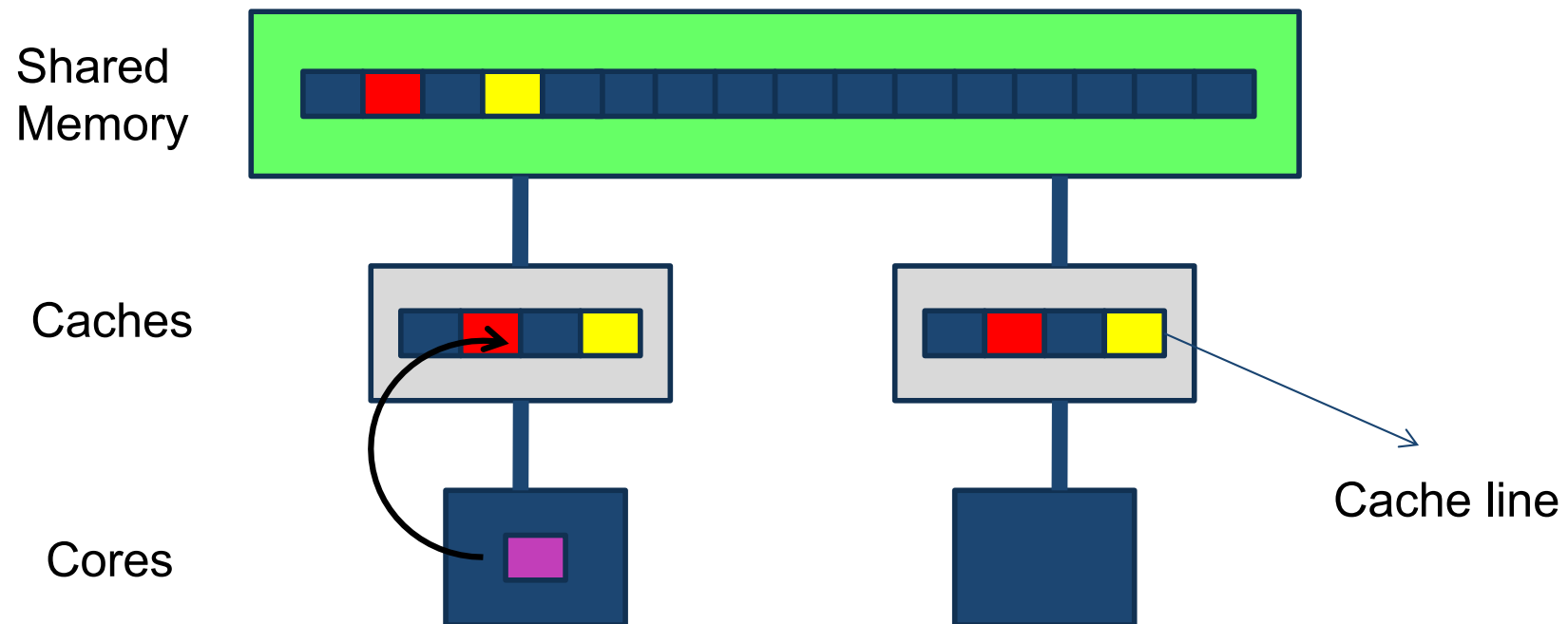
- A **cache** can be seen as a table of **cache lines** holding a predefined amount of memory (64B on most processors)
- Accessing a variable results in a **cache hit** if the corresponding cache line has already been cached (**fast memory access**)
- It can also result in a **cache miss** if the corresponding cache line is not cached yet. The hardware has to load the cache line to the cache before the processor can access it (**longer memory access**).





# Performance: False Sharing effect

Cache coherency and the negative effect of false sharing can have a big impact on performance



Charging a line of a shared cache invalids the other copies of this line

# Performance: False Sharing effect, contd.

**Using data structures in memory may lead to a decrease of performance and a lack of scalability**

- To get performance, use the cache
- If several cores manipulate different data items close in memory, the update of individual elements leads a load of a line of cache (to keep coherency with the main memory)

**False sharing leads to bad performance when the following conditions are met**

- Shared data is modified on # cores
- Several threads on # cores update data which are located on the same cache line
- These updates appear simultaneously and frequently

# Performance: False Sharing effect, contd.

When data are only read, we don't get false sharing

**It can be avoided (or reduced) by**

- Privatizing variables
- Increasing the array size or by using “padding”
- By increasing the packet size (modifying the way loop iterations are shared between threads)

## Performance: False Sharing effect, contd.

```
Integer, dimension(n) :: a
...
!$OMP PARALLEL DO SHARED(nthreads,a) SCHEDULE(static,1)
    DO i=0, nthreads-1
        a(i) = i
    END DO
!$OMP END PARALLEL DO
```

**Nthreads** : # threads executing the loop

If we suppose that every thread possess a copy of **a** in his local cache

Packet size of 1 leads to a false sharing phenomenon for each update

If a cache line can contain **C** elements of vector **a**, we can solve the problem by extending artificially the array dimensions (array padding)

We declare an array **a(C,n)** and replace **a(i)** by **a(1,i)**

## Naive Square Matrix Multiplication Algorithm

We consider a simple matrix multiplication algorithm involving square matrices of double precision floats.

$$\begin{bmatrix} C_0 & C_1 & C_2 & C_3 \\ C_4 & C_5 & C_6 & C_7 \\ C_8 & C_9 & C_{10} & C_{11} \\ C_{12} & C_{13} & C_{14} & C_{15} \end{bmatrix} = \begin{bmatrix} A_0 & A_1 & A_2 & A_3 \\ A_4 & A_5 & A_6 & A_7 \\ A_8 & A_9 & A_{10} & A_{11} \\ A_{12} & A_{13} & A_{14} & A_{15} \end{bmatrix} \times \begin{bmatrix} B_0 & B_1 & B_2 & B_3 \\ B_4 & B_5 & B_6 & B_7 \\ B_8 & B_9 & B_{10} & B_{11} \\ B_{12} & B_{13} & B_{14} & B_{15} \end{bmatrix}$$

where  $C_0 = A_0B_0 + A_1B_4 + A_2B_8 + A_3B_{12}$

Let's implement this using what we've learned about OpenMP!

## Speeding-Up OpenMP: Benefit from Cache Memory (2)

```
1 void gemm_omp(double *A, double *B, double *C, int n) {  
2     #pragma omp parallel  
3     {  
4         int i, j, k;  
5         #pragma omp for  
6         for (i=0; i<n; i++) {  
7             for (j=0; j<n; j++) {  
8                 for (k=0; k<n; k++) {  
9                     C[i*n+j] += A[i*n+k]*B[k*n+j];  
10                }  
11            }  
12        }  
13    }  
14 }
```

**On the Intel192 machine**

Serial time : 40.3018250s

Parallel time : 0.270773s

Achieved speed-up: 148

## We Can Do Better : It's All About Being (Cache) Friendly

Assuming we work with 32 bytes-long cache lines and each element of the matrix is 8 bytes long, how many cache lines do I need to compute one element of C?

$$\begin{bmatrix} C_0 & C_1 & C_2 & C_3 \\ C_4 & C_5 & C_6 & C_7 \\ C_8 & C_9 & C_{10} & C_{11} \\ C_{12} & C_{13} & C_{14} & C_{15} \end{bmatrix} = \begin{bmatrix} A_0 & A_1 & A_2 & A_3 \\ A_4 & A_5 & A_6 & A_7 \\ A_8 & A_9 & A_{10} & A_{11} \\ A_{12} & A_{13} & A_{14} & A_{15} \end{bmatrix} \times \begin{bmatrix} B_0 & B_1 & B_2 & B_3 \\ B_4 & B_5 & B_6 & B_7 \\ B_8 & B_9 & B_{10} & B_{11} \\ B_{12} & B_{13} & B_{14} & B_{15} \end{bmatrix}$$

$$\begin{aligned} C_0 &= A_0 B_0 \\ &+ A_1 B_4 \\ &+ A_2 B_8 \\ &+ A_3 B_{12} \end{aligned}$$

## We Can Do Better : It's All About Being (Cache) Friendly

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$$\begin{aligned} C_0 &= A_0 B_0 \\ &+ A_1 B_4 \\ &+ A_2 B_8 \\ &+ A_3 B_{12} \end{aligned}$$



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$$\begin{aligned} C_0 &= A_0 B_0 \\ &+ A_1 B_4 \\ &+ A_2 B_8 \\ &+ A_3 B_{12} \end{aligned}$$

## We Can Do Better : It's All About Being (Cache) Friendly

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$$\begin{aligned} C_0 &= A_0 B_0 \\ &+ A_1 B_4 \\ &+ A_2 B_8 \\ &+ A_3 B_{12} \end{aligned}$$

## We Can Do Better : It's All About Being (Cache) Friendly

Assuming we work with 32 bytes-long cache lines and each element of the matrix is 8 bytes long, how many cache lines do I need to compute one element of C?

$$\begin{bmatrix} C_0 & C_1 & C_2 & C_3 \\ C_4 & C_5 & C_6 & C_7 \\ C_8 & C_9 & C_{10} & C_{11} \\ C_{12} & C_{13} & C_{14} & C_{15} \end{bmatrix} = \begin{bmatrix} A_0 & A_1 & A_2 & A_3 \\ A_4 & A_5 & A_6 & A_7 \\ A_8 & A_9 & A_{10} & A_{11} \\ A_{12} & A_{13} & A_{14} & A_{15} \end{bmatrix} \times \begin{bmatrix} B_0 & B_1 & B_2 & B_3 \\ B_4 & B_5 & B_6 & B_7 \\ B_8 & B_9 & B_{10} & B_{11} \\ B_{12} & B_{13} & B_{14} & B_{15} \end{bmatrix}$$

$$\begin{aligned} C_0 &= A_0 B_0 \\ &+ A_1 B_4 \\ &+ A_2 B_8 \\ &+ A_3 B_{12} \end{aligned}$$

Conclusion:

- Every access to  $A_i$  use the same cache line  $\Rightarrow$  cache friendly
- Every access to  $B_j$  use a different cache line  $\Rightarrow$  poor cache utilization

## Deal with the $B_j$ Situation

To improve cache utilization, we can transpose matrix B to make sure  $B_0$ ,  $B_4$ ,  $B_8$  and  $B_{12}$  are stored on the same cache line

$$\begin{bmatrix} C_0 & C_1 & C_2 & C_3 \\ C_4 & C_5 & C_6 & C_7 \\ C_8 & C_9 & C_{10} & C_{11} \\ C_{12} & C_{13} & C_{14} & C_{15} \end{bmatrix} = \begin{bmatrix} A_0 & A_1 & A_2 & A_3 \\ A_4 & A_5 & A_6 & A_7 \\ A_8 & A_9 & A_{10} & A_{11} \\ A_{12} & A_{13} & A_{14} & A_{15} \end{bmatrix} \times \begin{bmatrix} B_0 & B_4 & B_8 & B_{12} \\ B_1 & B_5 & B_9 & B_{13} \\ B_2 & B_6 & B_{10} & B_{14} \\ B_3 & B_7 & B_{11} & B_{15} \end{bmatrix}$$

$$\begin{aligned} C_0 &= A_0 B_0 \\ &+ A_1 B_4 \\ &+ A_2 B_8 \\ &+ A_3 B_{12} \end{aligned}$$

### Performance evaluation on Intel192:

Serial time: 5.188652s (prev: 40.3018250s)

Parallel time: 0.067657s (prev: 0.270773s)

Achieved speed-up : 77 (rel to prev: 595(!))