

Calcul Parallèle avec OpenMP

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LIMOS - Cours ED-SPI

16-17/01/19

OpenMP Overview

OpenMP (Open specifications for MultiProcessing) API for multithreaded applications

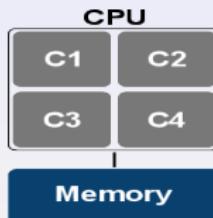
- A set of compiler directives and library routines for parallel application programmers
- Greatly simplifies writing multi-threaded (MT) programs in Fortran, C and C++
- Standardizes last 20 years of Symmetric Multi-Processing (SMP) practice

Suitable architectures

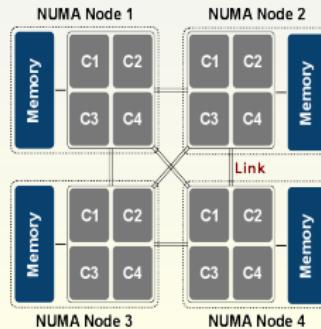
- Multi-threading processors
- Multi-core machines
- Shared-memory machines

Shared memory architecture for SMP

Architecture **UMA** (Uniform Memory Access) : Equal access time



Architecture **NUMA** (Non-Uniform Memory Access) : Different access time



OpenMP Syntax

- OpenMP compiler directives

```
#pragma omp construct[clause/]...
```

Example : #pragma omp parallel for

- Function prototypes and types in the file :

```
#include “omp.h”
```

- OpenMP constructs apply to a structured block :
 - a block of one or more statements with one point of entry at the top and one point of exit at the bottom.
 - a structured block can contain an exit()

Check if your environment works

hello_omp.c

```
#include <stdio.h>
#include "omp.h"
int main()
{
    int mytid=-1, nb_ths=0;
    #pragma omp parallel
    {
        nb_ths=omp_get_num_threads(); mytid=omp_get_thread_num();
        printf("Hello World from tid %d %d \n", mytid, nb_ths);
    }
    return 0;
}
```

Compile : gcc -fopenmp hello_omp.c

Set the number of threads : export OMP_NUM_THREADS=8

Run : ./a.out

Check if your environment works

hello_omp.c 2

```
#include <stdio.h>
#include "omp.h"
int main()
{
    int mytid=-1, nb_ths=0;
    omp_set_num_threads(8);
    #pragma omp parallel
    {
        nb_ths=omp_get_num_threads(); mytid=omp_get_thread_num();
        printf("Hello World from tid %d %d \n", mytid, nb_ths);
    }
    return 0;
}
```

Compile : gcc -fopenmp hello_omp.c
Run : ./a.out

Fixing the number of threads

Environment variable

```
export OMP_NUM_THREADS=8
```

Runtime library

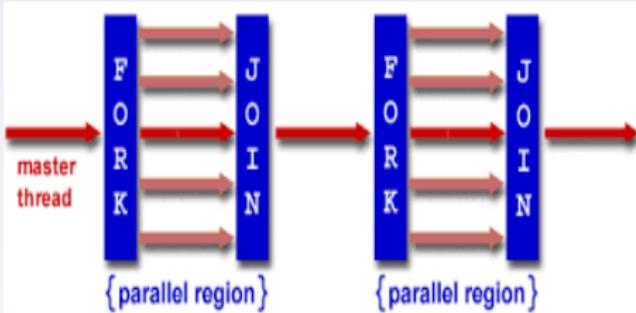
```
omp_set_num_threads(8);
```

Clause of parallel region

```
#pragma omp parallel num_threads(8)
```

Threads interaction

Fork-Join model



OpenMP is a multi-threading, shared address model

- Threads communicate by sharing variables
- Sharing of data causes data conflicts
- Synchronization protects data conflicts
- Synchronization is expensive

Parallel sections

```
#pragma omp parallel sections (+clauses)
{
    #pragma omp section
    {
        block 1
    }
    ...
    #pragma omp section
    {
        block n
    }
} /* synchronization */
```

The blocks defined by `#pragma omp section` are performed in parallel by the threads in the team.

Parallel loop

```
#pragma omp parallel shared(x,y,z)
{
    #pragma omp for private(i)
    for (i=0; i<DIM; i++)
        z[i] = x[i]*y[i];
}
```

Equivalent combined version

```
#pragma omp parallel for shared(x,y,z) private(i)
for (i=0; i<DIM; i++)
    z[i] = x[i]*y[i];
```

Programming example $C = A + B$ serial

```
#include <stdio.h>
#include <stdlib.h>
#define N 4096

int main ()
{
    double A[N][N], B[N][N], C[N][N];
    int i,j;

    for (i = 0; i < N; i++) /* Initialization */
        for (j = 0; j < N; j++){
            A[i][j] = i*N+j; B[i][j] = i+j;
        }

    for (i = 0; i < N; i++) /* Computing */
        for (j = 0; j < N; j++){
            C[i][j] = A[i][j] + B[i][j];
        }

    return 0;
}
```

Programming example $C = A + B$ parallel

```
#include <stdio.h>
#include <stdlib.h>
#include "omp.h"
#define N 4096

int main ()
{
    double A[N][N], B[N][N], C[N][N];
    int i,j;

#pragma omp parallel for shared(A,B) private(i,j)
    for (i = 0; i < N; i++) /* Initialization */
        for (j = 0; j < N; j++){
            A[i][j] = i*N+j; B[i][j] = i+j;
        }

#pragma omp parallel for shared(A,B,C) private(i,j)
    for (i = 0; i < N; i++) /* Computing */
        for (j = 0; j < N; j++){
            C[i][j] = A[i][j] + B[i][j];
        }
    return 0;
}
```

OpenMP Programming summary

Programming

- Incrementally parallelization
- Multi-task/worksharing
- Variable management (shared/private, initial/final values, etc)
- Threads synchronization

Issues

- Reliability of results
- Deadlocks

omp parallel clauses

Clause	Do
<code>num_threads(nt)</code>	the number of threads to use
<code>private(list)</code>	declare private variable to each thread
<code>shared(list)</code>	variables shared by the threads in the team
<code>firstprivate(list)</code>	private + pre-initialization
<code>lastprivate(list)</code>	private + the final value of variable is the last one in the last iteration
<code>reduction(operator:list)</code>	Peform a reduction on all scalar variables in list using the specified operator

omp for clauses

Clause	Do
private(list)	declare private variables to each thread
firstprivate(list)	private + pre-initialization
lastprivate(list)	private + the final value of variable is the last one in the last iteration
reduction(operator:list)	Peform a reduction on all scalar variables in list using the specified operator
schedule(type)	Specify how iterations are divided (static, dynamic, guided)
nowait	avoid the implicit barrier at the end of the for-loop

Dot product $x^\top y = \sum x_i y_i$

```
#include <math.h>
#include <stdio.h>
#include <stdlib.h>
#include "omp.h"
int main(){
    int i,n=10000;
    double u[n], v[n], s,sx, ts, te;

    for (i=0; i<n; i++){
        u[i] = 1.0/3.0; v[i] = 1.0/3.0;
    }
    sx = (double) n/9.0;

    s = 0.0; /* put "#pragma omp single" before in parallel region */
    ts = omp_get_wtime();
#pragma omp parallel for schedule(static) reduction(+:s)
    for (i=0; i<n; i++) s += u[i]*v[i];
    te = omp_get_wtime()-ts;

    printf("s-sx=%15.8e elapsed time = %f \n",s-sx,te);
    return 0;
}
```

Jacobi's method

Solve

$$Ax = b$$

by the Jacobi iteration :

$$\begin{aligned}x_i^{(k+1)} &= \frac{1}{a_{ii}} \left(b_i - \sum_{j=1}^{i-1} a_{ij}x_j^{(k)} - \sum_{j=i+1} a_{ij}x_j^{(k)} \right), \\&= \frac{1}{a_{ii}} \left(b_i + a_{ii}x_i^{(k)} - \sum_{j=1}^n a_{ij}x_j^{(k)} \right)\end{aligned}$$

Convergence if

$$\| Ax^{(k)} - b \| < \varepsilon \| b \|$$

THANK YOU FOR YOUR ATTENTION