

OpenMP tutorials

## 1 – Introduction

The tutorials will take place on the Vargas computer (IBM Power6 with 3586 cores). They are located in the `$WORKDIR/OpenMP_tp` directory of each student account. They consist of 9 independent tutorials ; each of them has its own directory which includes a `Makefile` for compilation, a `batch.sh` file to launch a dedicated batch execution and, of course, one or several `Fortan source files` that you'll have to edit and modify (according to the instructions).

In each tutorial directory, you can find the following sub-directories :

- ☞ `sans_indications_openmp` has the plain sequential source files.
- ☞ `avec_indications_openmp` has source files that include clues about the modifications you have to work on.
- ☞ `solution` has the source file that you're supposed to look at when you're done, to check if your work is correct.

## General instructions

- ☞ type **gmake mono** to build the sequential executable (the compiler sees the OpenMP directives as comments).
- ☞ type **gmake para** to build the parallel executable, with the compiler interpreting the OpenMP directives.
- ☞ type **gmake clean** to wipe out object files build by a previous compilation and core files resulting from a previous execution ; type **gmake cleanall** to erase previous executable files too.
- ☞ type **llsubmit batch.sh** to submit a job in the batch system. The job will run on dedicated nodes, first a sequential execution, then parallel ones with 2, 4, 6, 8 threads. The executables must have been build before (see previous gmake commands).
- ☞ The **Qstat** command lists the submitted/running jobs. Once the batch job is done, the results is available in a **.res** file in the current directory.

Good work!

## 2 – tp1 : matrix product

The `prod_mat.f90` source file computes a matrix product :

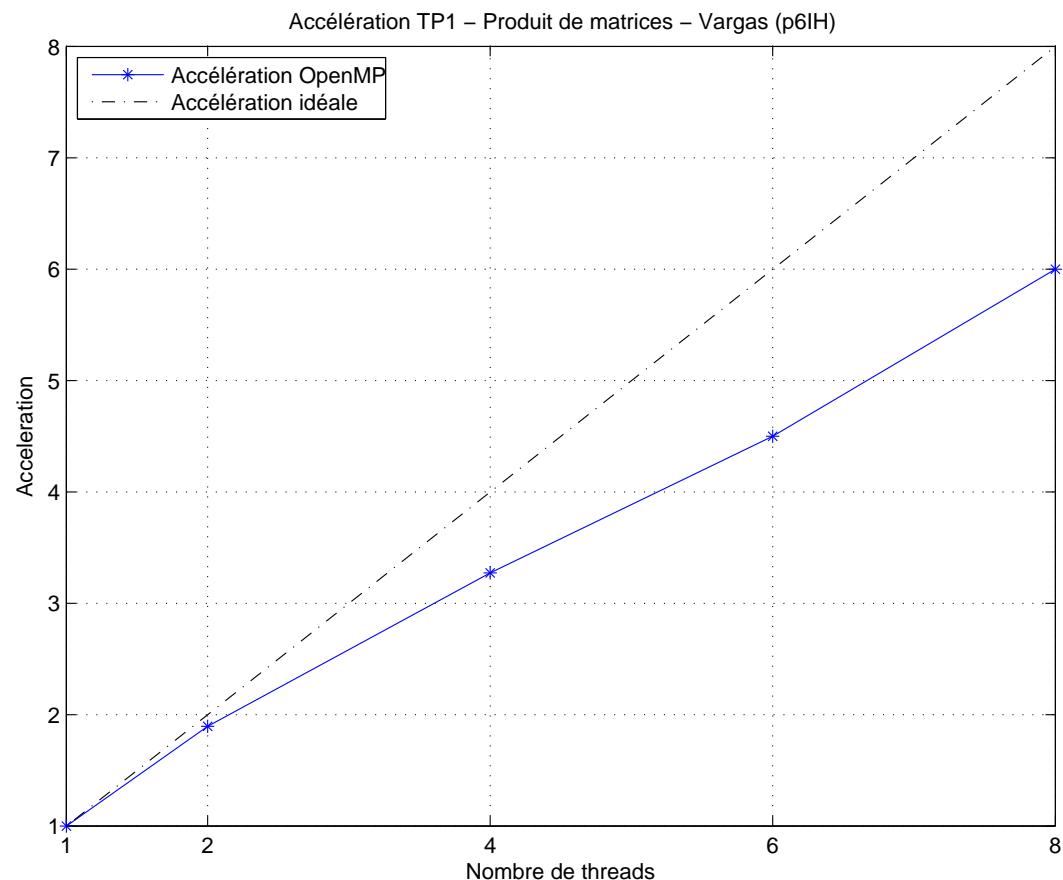
$$C = A \times B$$

In this basic tutorial, you're asked you to :

1. Insert proper OpenMP directives.
2. Analyse the parallel performance on 2, 4, 6, 8 threads comparing to a sequential execution (please do a dedicated execution by submitting a *batch* job). You can test the various loop distribution modes (**STATIC**, **DYNAMIC**, **GUIDED**) and test various chunk sizes.
3. Plot the acceleration curves.

# Description of tp1

Nb. of threads	Elapsed time	Speedup
seq.		
1		
2		
4		
6		
8		



## 3 – tp2 : Jacobi Method

The jacobi.f90 source file contains the resolution of a linear system

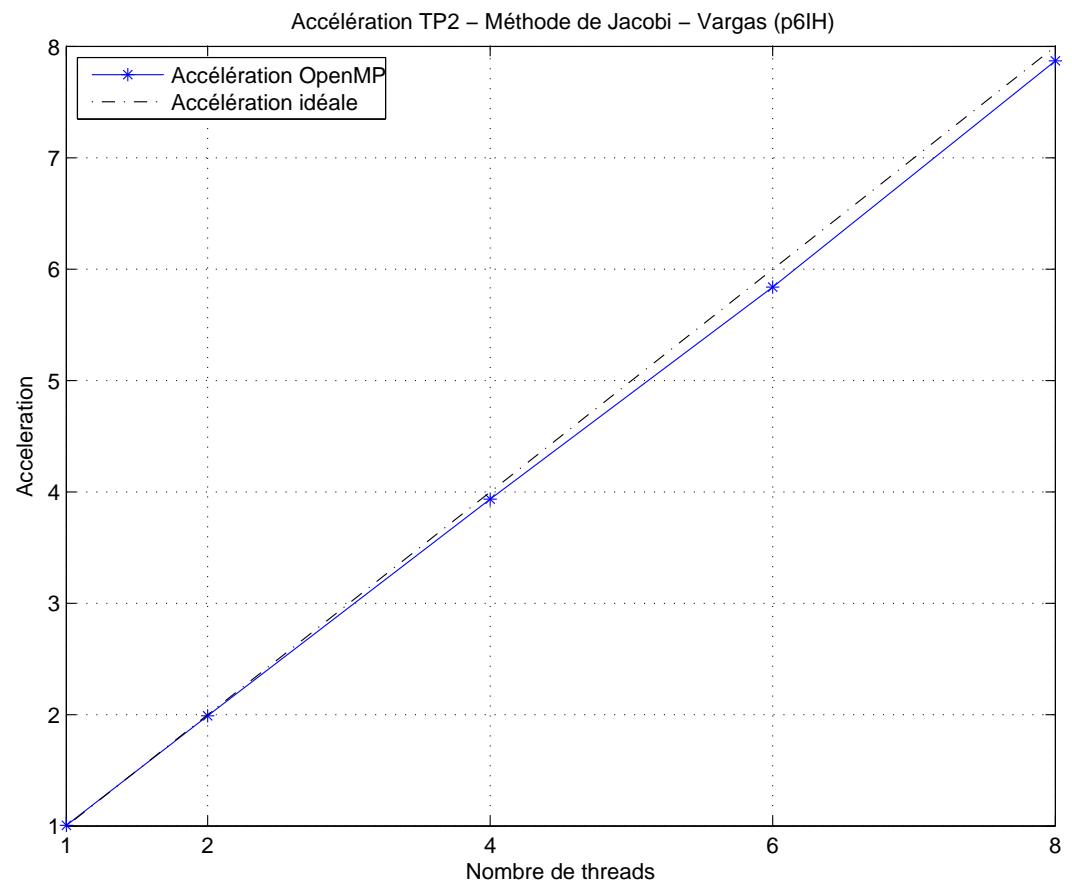
$$A \times x = b$$

with the Jacobi iterative method.

1. Analyze the variable data-sharing attributes and insert the appropriate OpenMP directives in the **jacobi.f90** source file.
2. Analyze the parallel performance with 2, 4, 6, 8 threads referring to a sequential execution (please do a dedicated execution by submitting a batch job).
3. Plot the acceleration curve.

# Description of tp2

Nb. of threads	Elapsed time	Speedup
seq.		
1		
2		
4		
6		
8		



## 4 – tp3 : computation of $\pi$

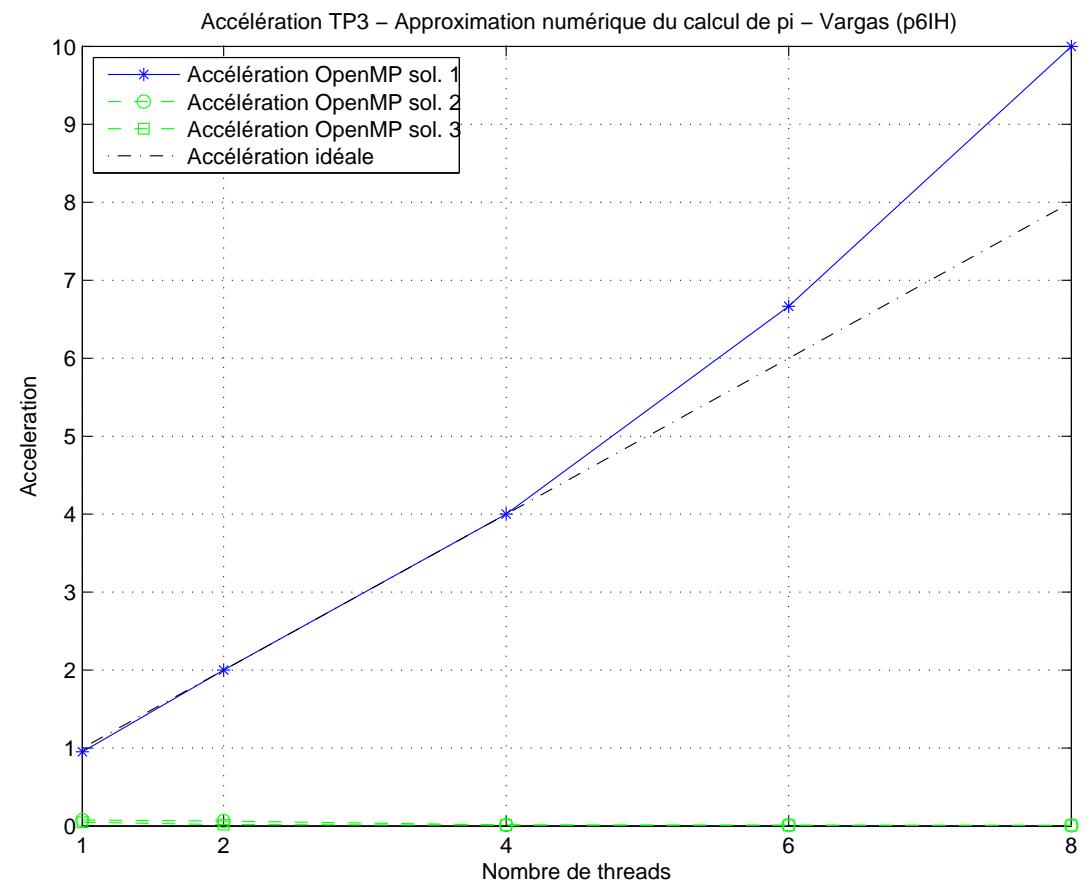
The aim of this exercice is to compute  $\pi$  by numerical integration. In the [pi.f90](#) source file you can find the program that computes  $\pi$  using the rectangle method (mean point). Let  $f(x) = \frac{4}{1+x^2}$  the function to integrate and  $N$ ,  $h = \frac{1}{N}$  the number of points and the discretization interval  $[0, 1]$ , respectively.

1. Analyze the variable data-sharing attributes and insert the appropriate OpenMP directives in the [pi.f90](#) source file. This tutorial's point is to use three different methods (three different OpenMP directives, one for each version) to parallelize the code.
2. Analyze the parallel performance with 2, 4, 6, 8 threads referring to a sequential execution (please do a dedicated execution by submitting a *batch* job).
3. Optimize the 2 less performing versions, without changing the OpenMP directives that were used, to reach the same performance for the three versions.
4. Plot the acceleration curves.

# Description of tp3

10

Nb. of threads	Elapsed time	Speedup
seq.		
1		
2		
4		
6		
8		



## 5 – tp4 : Conjugate Gradient (*CG*) Method

The `gradient_conjugue.f90` source file contains a preconditionned conjugate gradient method to solve a symmetric linear system.

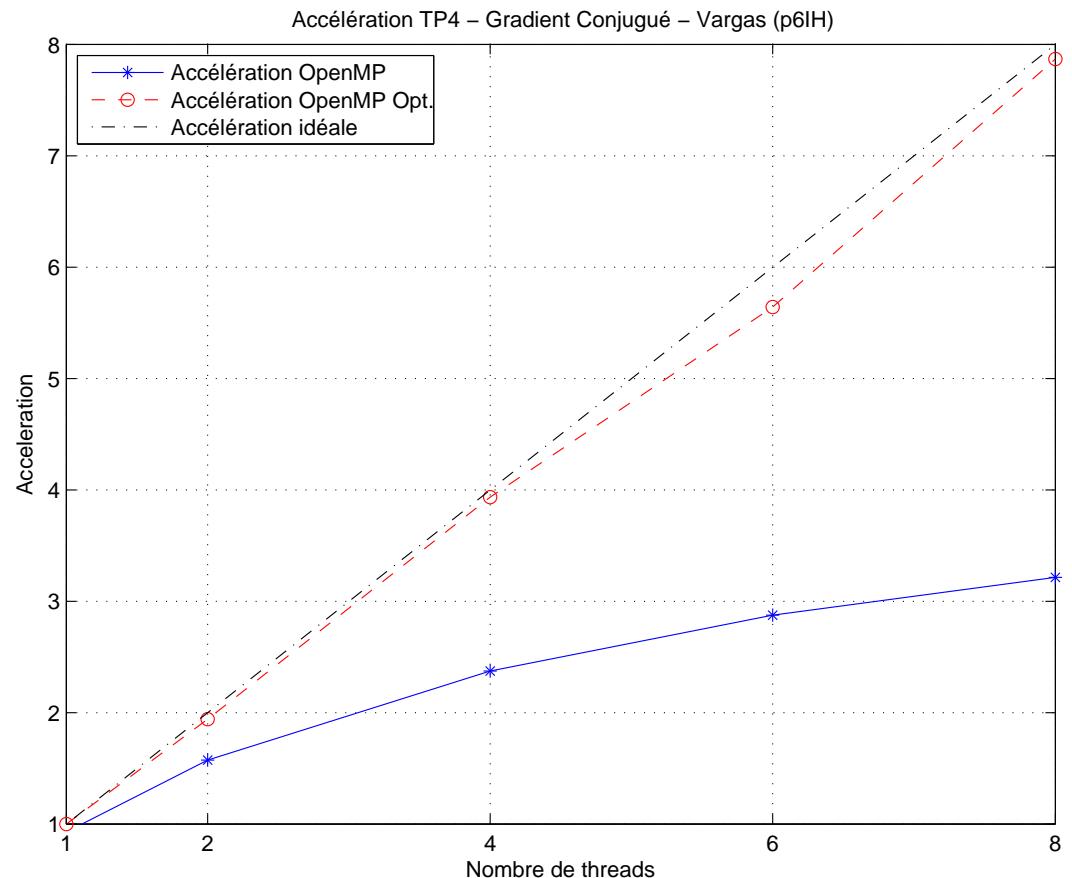
$$A \times x = b$$

1. Analyze the variable data-sharing attribute and insert the appropriate OpenMP directives in the `gradient_conjugue.f90` source file.
2. Analyze the parallel performance with 2, 4, 6, 8 threads referring to a sequential execution (please do a dedicated execution by submitting a *batch* job). What are your conclusions about the **WORKSHARE** directive efficiency ?
3. Optimize the parallel version by slightly modifying the original source code to avoid the only one **WORKSHARE** construct that is inefficient.
4. Plot the appropriate acceleration curves.

# Description of tp4

12

Nb. of threads	Elapsed time	Speedup
seq.		
1		
2		
4		
6		
8		



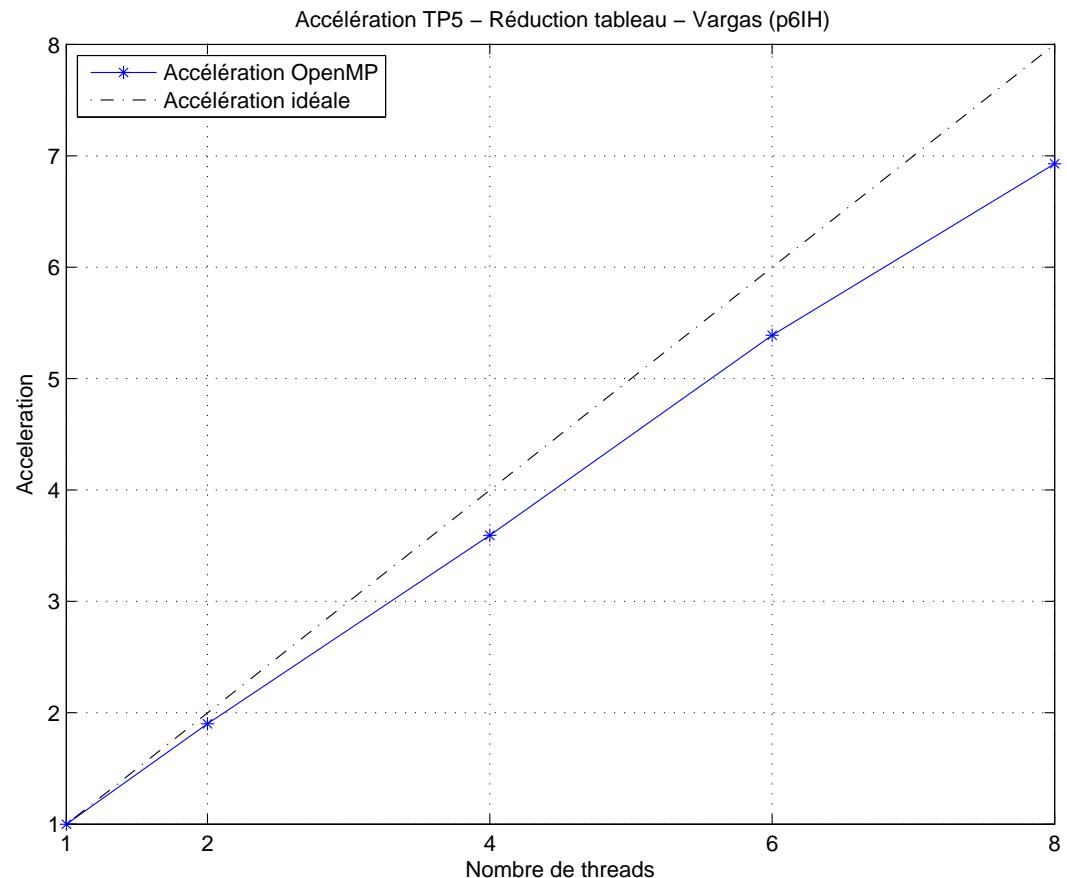
## 6 – tp5 : array reduction

The `reduction_tab.f90` source code comes from a chemistry code ; it computes the reduction of a 3D array into a vector. The aim of this exercice is to use three different ways to parallelize this code.

1. Version 1 : analyze the variable data-sharing attributes and insert the appropriate OpenMP directives in the `reduction_tab.f90` source file. Do not make any modification to the original sequential code.
2. Version 2 : analyze the variable data-sharing attributes and modify the original loop order, so that one **PARALLEL DO** directive applies to the outermost loop.
3. Version 3 : without changing the loop order ( $k, j, i$ ), modify the source code to parallelize the  $k$  outermost loop.
4. For each version, analyze the parallel performance with 2, 4, 6, 8 threads referring to a sequential execution (please do a dedicated execution by submitting a batch job). Plot the acceleration curves. How can version 2's bad performance be explained ?

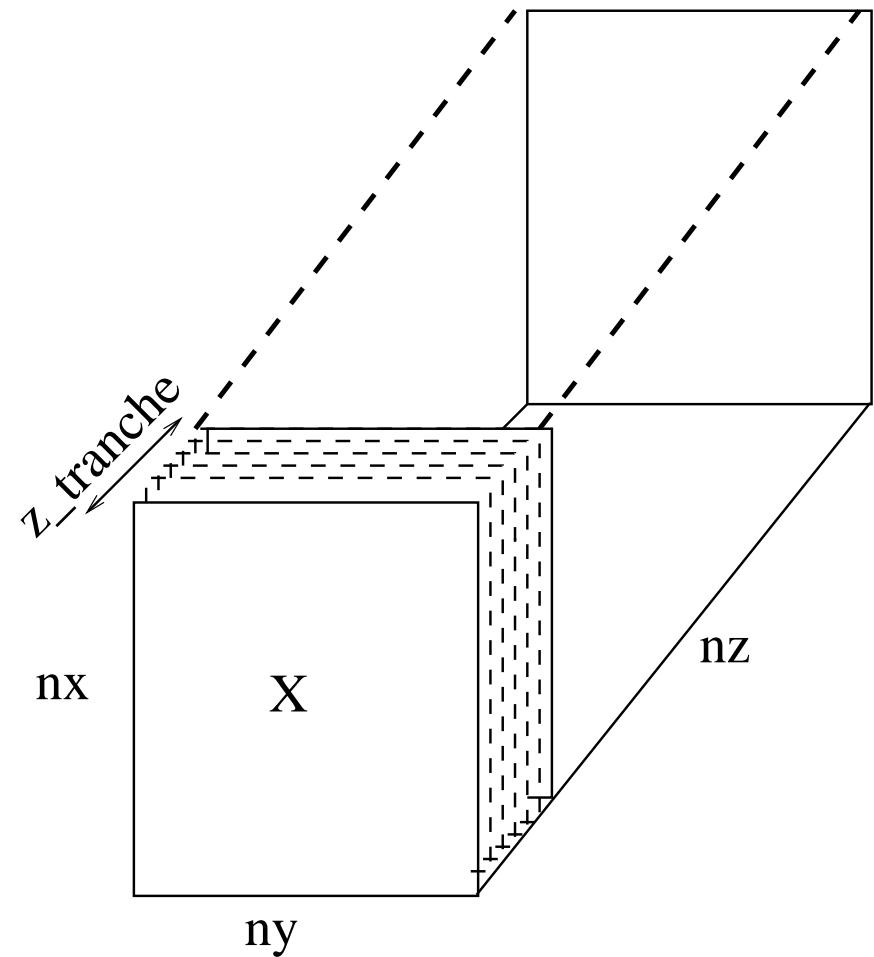
# Description of tp5

Nb. of threads	Elapsed time	Speedup
seq.		
1		
2		
4		
6		
8		



## 7 – tp6 : multiple Fast Fourier Transform

The `fft.f90` source file contains a 3D matrix direct real-complex forward and reverse FFT. Parallelization consists in explicitly distributing the work by cutting the 3D matrix following the 3rd dimension, with as many slices as threads : each thread computes the FFT on its 2D array.



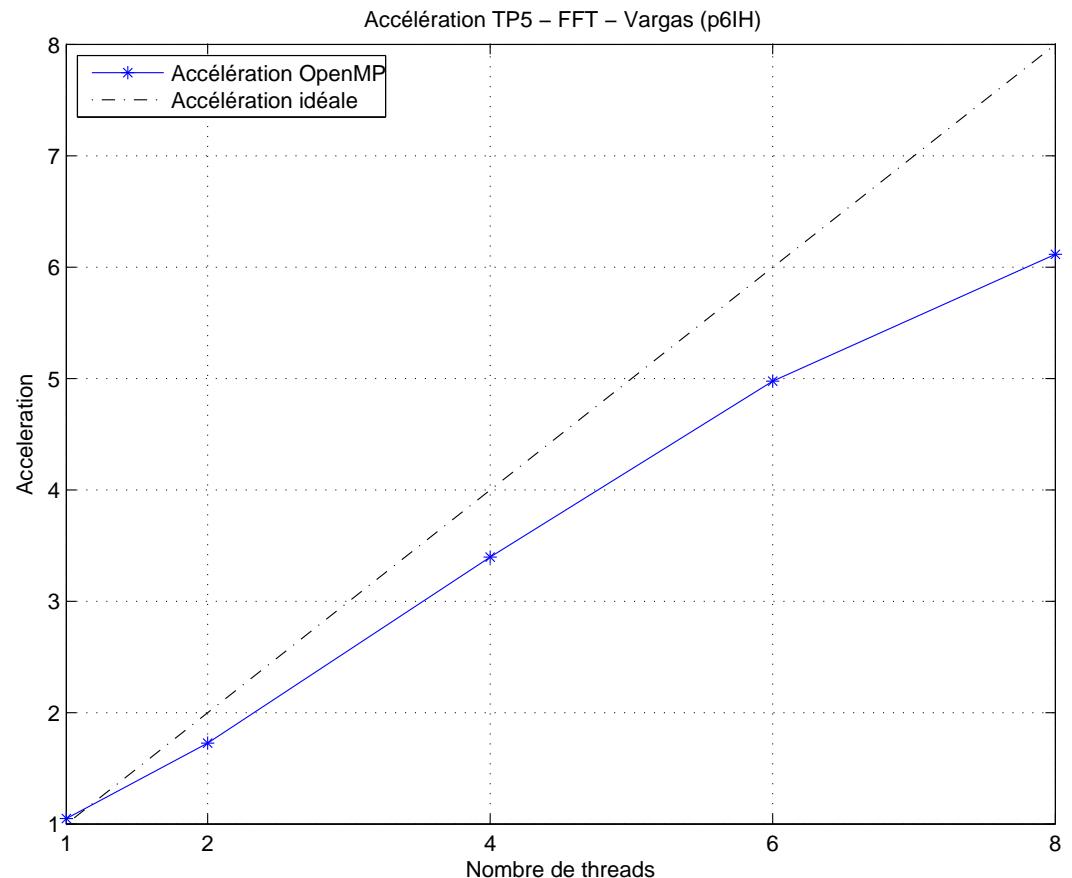
1. Analyze the variable data-sharing attributes and insert the appropriate OpenMP directives in the `fft.f90` source file. In this tutorial, use conditional execution to take care of sequential execution.
2. Analyze the parallel performance with 2, 4, 6, 8 threads referring to a sequential execution (please do a dedicated execution by submitting a *batch* job).
3. Plot the acceleration curve.

**Please note :** the jmfft.a FFT library file should not be modified.

It includes the references to the `scfftm` and `csfftm` subroutines that are called within `fft.f90`.

# Description of tp6

Nb. of threads	Elapsed time	Speedup
seq.		
1		
2		
4		
6		
8		



## 8 – tp7 : Biconjugate gradient stabilized method

The `principal.f90` source file contains the main program, which calls the `bi-cgstab` subprogram (`bi-cgstab.f90` source file) to solve a linear system with multiple right-hand vectors :

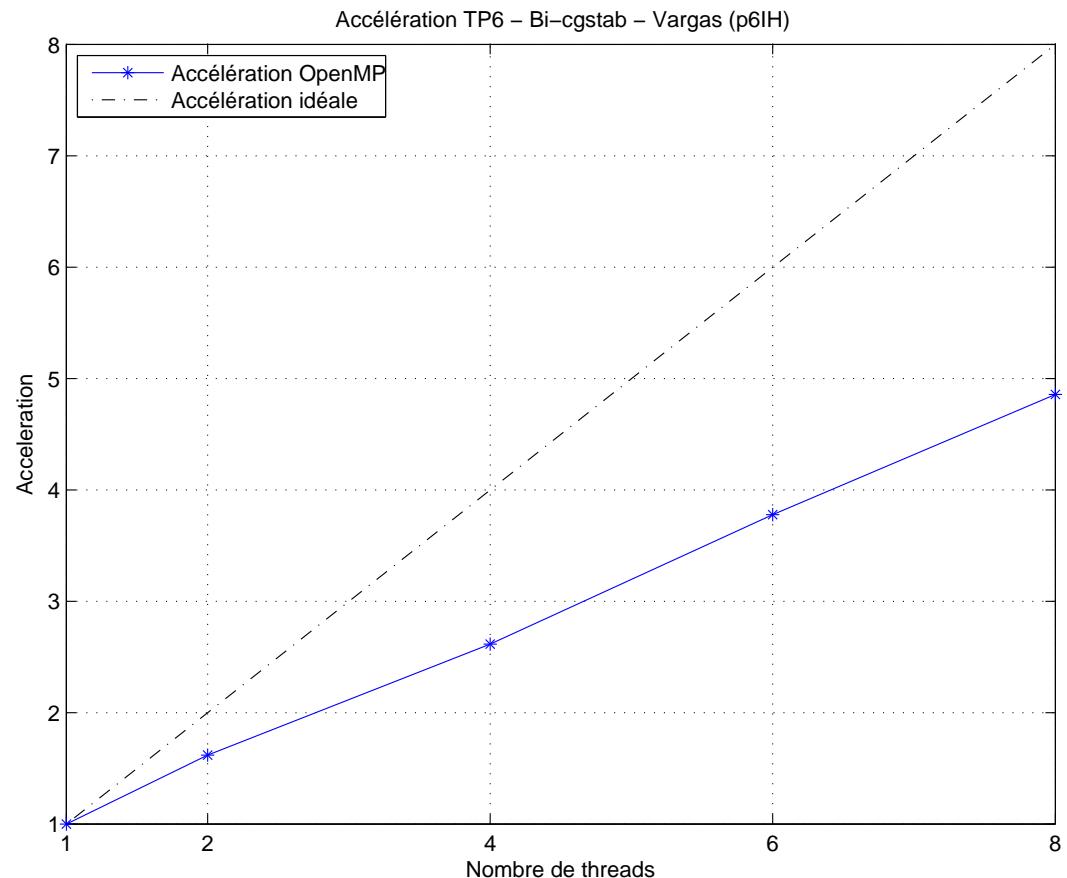
$$A \times x = B$$

with the biconjugate gradient stabilized method (BiCGSTAB)

1. Analyze the variable data-sharing attributes and insert the appropriate OpenMP directives in the `principal.f90` and `bi-cgstab.f90` source file. The `bi-cgstab` subprogram should be modified to behave as an orphaned procedure.
2. Analyze the parallel performance with 2, 4, 6, 8 threads referring to a sequential execution (please do a dedicated execution by submitting a *batch* job).
3. Plot the acceleration curve.

# Description of tp7

Nb. of threads	Elapsed time	Speedup
seq.		
1		
2		
4		
6		
8		



## 9 – tp8 : resolution equation of Poisson

The `poisson.f90` and `gradient_conjugue.f90` source files are used to solve the POISSON problem (1). The analytic solution  $u_a(x, y)$  is given by :

$$u_a(x, y) = \cos \pi x \times \sin \pi y \quad ; \quad (x, y) \in [0, 1] \times [0, 1]$$

$$\left\{ \begin{array}{l} -\frac{\partial^2 u}{\partial x^2} - \frac{\partial^2 u}{\partial y^2} = b(x, y) \\ u(0, y) = u_a(0, y) \\ u(1, y) = u_a(1, y) \\ u(x, 0) = u_a(x, 0) \\ u(x, 1) = u_a(x, 1) \end{array} \right. \quad (1)$$

We will use a center finite difference method in the  $x$  direction followed by an FFT (sinus) in the  $y$  direction. Let be  $\tilde{u}$  and  $\tilde{b}$  the result of the FFT of  $u$  and  $b$  respectively and apply the FFT to POISSON (1) equation :

$$-\frac{\partial^2 \tilde{u}}{\partial x^2} - \widetilde{\frac{\partial^2 u}{\partial y^2}} = \tilde{b}(x, y)$$

The sinus transformation  $\widetilde{\frac{\partial^2 u}{\partial y^2}}$  of  $\frac{\partial^2 u}{\partial y^2}$  operator is diagonal, each element representing an eigenvalue of the operator. These eigenvalues can be calculated analytically. Let  $j = 1, \dots, N_j$  represents the indice of discretization and  $h_y$  the spatial discretization in the  $y$  direction, then the eigenvalues are :

$$\text{vp}_j = \frac{4}{h_y^2} \sin^2 \frac{\pi(j-1)}{2(N_j-1)} \quad ; \quad j = 2, \dots, N_j - 1$$

Hence, in the basis composed of the eigenvectors, the POISSON problem turns out to be a resolution of  $N_j - 2$  symetric tridiagonal linear systems (use Conjugate Gradient method to solve them) of size  $(N_i - 2) \times (N_i - 2)$  each, where  $N_i$  represents the number

of discretization points in the  $x$  direction :

$$\begin{pmatrix} d_j & -c_x & 0 & \dots & \dots & 0 \\ -c_x & d_j & -c_x & 0 & \dots & \vdots \\ 0 & \ddots & \ddots & \ddots & \ddots & \vdots \\ \vdots & \ddots & \ddots & \ddots & \ddots & 0 \\ \vdots & 0 & -c_x & d_j & -c_x & \\ 0 & \dots & \dots & 0 & -c_x & d_j \end{pmatrix} \begin{pmatrix} \tilde{u}_{2,j} \\ \tilde{u}_{3,j} \\ \vdots \\ \vdots \\ \tilde{u}_{N_i-2,j} \\ \tilde{u}_{N_i-1,j} \end{pmatrix} = \begin{pmatrix} \tilde{b}_{2,j} + \text{CL} \\ \tilde{b}_{3,j} \\ \vdots \\ \vdots \\ \tilde{b}_{N_i-2,j} \\ \tilde{b}_{N_i-1,j} + \text{CL} \end{pmatrix}$$

where  $c_x = \frac{1}{h_x^2}$ ,  $c_y = \frac{1}{h_y^2}$ ,  $d_j = 2c_x + \text{vp}_j$  and  $h_x$  is the spatial discretization in the  $x$  direction. The term CL deals with boundary conditions.

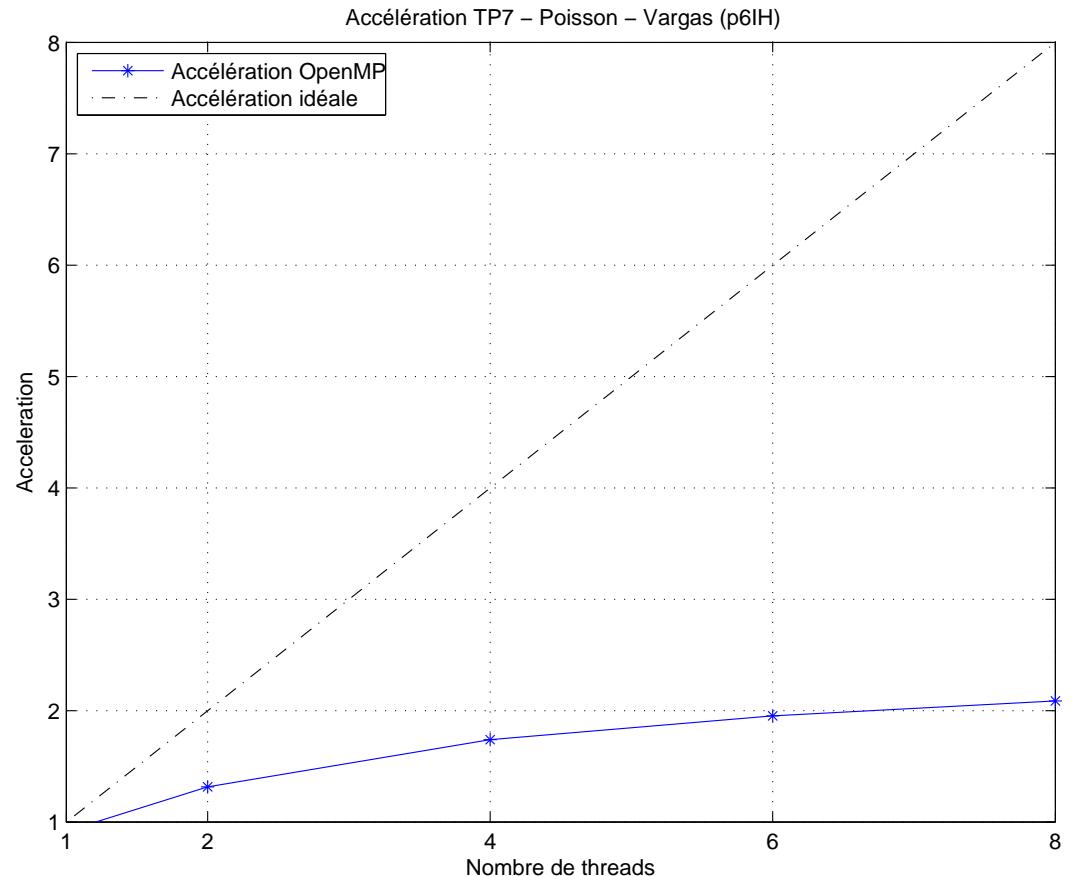
Finally,  $(N_i - 2)$  independent reverse FFT of  $\tilde{u}$  gives the final solution in the canonical basis.

1. Analyze the variable data-sharing attributes and insert the appropriate OpenMP directives in the `poisson.f90` and `gradient_conjugue.f90` source file. The `gradient_conjugue` subprogram should be modified to behave as an orphaned procedure, with only one parallel region.
2. Analyze the parallel performance with 2, 4, 6, 8 threads referring to a sequential execution (please do a dedicated execution by submitting a *batch* job).
3. Plot the acceleration curve.

Please note : the file `c06haf.o` should not be modified. It includes the reference to the subroutine `c06haf` (sinus forward and reverse FFT) called within `poisson.f90`.

# Description of tp8

Nb. of threads	Elapsed time	Speedup
seq.		
1		
2		
4		
6		
8		



## 10 – tp9 : nested loops with dependancies

The source file **dependance.f90** contains two nested loops.

1. Analyse the data dependancy of the nested loops (e.g. are iterations independent one from each other ?). If loop parallelization is done nevertheless, what happens ?
2. Insert the appropriate OpenMP directives in the **dependance.f90** source file. Take care about thread synchronization, to avoid breaking dependancies.
3. Analyze the parallel performance with 2, 4, 6, 8 threads referring to a sequential execution (please do a dedicated execution by submitting a batch job). The execution is validated if the norm printed during execution is 0...
4. Plot the acceleration curve.

# Description of tp9

Nb. of threads	Elapsed time	Speedup
seq.		
1		
2		
4		
6		
8		

