

# Parallel Computing

## Exercise 3 (May 30, 2023)

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The result is to be submitted by the deadline stated above via the Moodle interface.  
The submitted result is as a .zip or .tgz file which contains

- a single PDF (.pdf) file with
  - a cover page with the title of the course, your name(s), Matrikelnummer(s), and email-address(es),
  - the source code of the sequential program,
  - the demonstration of a sample solution of the program,
  - the source code of the parallel program,
  - the demonstration of a sample solution of the program,
  - a benchmark of the sequential and of the parallel program.
- the source (.c/.cpp/.java) files of the sequential and of the parallel program.

## Distributed Memory Programming in MPI

The goal of this exercise is to develop in MPI a distributed memory solution to one of the problems specified in Exercise 1. As the base of your parallel programming effort, you may use the sequential program you have developed in Exercise 1; you may also write a new sequential program or ask one of your colleagues for one. Our default assumption is that the programs for this assignment are written in C/C++, using the official MPI binding for the parallel solution.

Having said this, you may also write your sequential program in Java and use for the parallelization one of the MPI bindings for Java provided by

- OpenMPI: <https://www.open-mpi.org/faq/?category=java>
- FastMPJ: <http://gac.udc.es/~rreye/fastmpj/>
- MPJ Express: <http://mpjexpress.org>

OpenMPI is available at the course machine (see `module avail`), without guarantee of a functional Java interface; newer versions of this package respectively the other packages are to be installed on your own.

However, neither do we recommend to solve this assignment in Java nor will we be able or willing to give any support for the use of Java with MPI.

**Benchmarking** Benchmark the programs (both the sequential and the parallel one) as in Exercise 1; you may also use the MPI function `double MPI_Wtime()` which returns the current wall clock time in seconds. Report the results as in Exercise 1.

**Contiguous Matrices** If a matrix  $A$  of dimension  $M \times N$  with values of type  $T$  is to be passed (respectively broadcast/scattered/gathered) among processes, make sure that  $A$  is represented by a contiguous block in memory. This can be achieved either by a global declaration `T A[M][N]` (which allocates the matrix in the data segment of the process,  $M$  and  $N$  have then to be compile-time constants) or by a declaration and initialization `T* A = malloc(M*N*sizeof(T))` (which allocates the matrix on the heap,  $N$  may then be variable; however, the element  $A[i][j]$  is now denoted by the reference `A[i*M+j]`).

**Presentation** Please be prepared to give a short (10 min) presentation of your results on June 6; you will be notified by June 1 whether such a presentation is requested from you.

## Alternative A: Matrix Inversion by Gauss-Jordan Elimination

In the MPI solution to this problem, you may assume that the number  $P$  of processes divides the matrix dimension  $N$  exactly.

- The program starts by distributing matrix  $A$  *row-wise* among the  $P$  processes *in a round-robin fashion* (i.e. process 0 receives rows  $0, P, 2P, \dots$ , process 1 receives rows  $1, P + 1, 2P + 1, \dots$ , and so on). In such a way we generally ensure that the work load of a process is not influenced by a particular distribution of data in the matrix.

To distribute  $A$ , process 0 constructs a correspondingly permuted version  $A'$  of the matrix and scatters its values among all processes (by a single call of `MPI_Scatter`).

- For performing the diagonalization, the program runs in  $N$  iterations, where in iteration  $i$  process  $p = i\%P$  broadcasts row  $i$  to all other processes (`MPI_Bcast`). Each process then uses this row to update all the rows of the matrix for which it is responsible.

To simplify the program, you may assume that  $A(i, i)$  is different from 0 (if this should not be the case, you may abort the computation by `MPI_Abort`).

- Finally, process 0 gathers the permuted inverse matrix  $B'$  from all processes (by a single call of `MPI_Gather`) and constructs from this the actual inverse matrix  $B$ .

## Alternative B: Counting the Satisfying Assignments of a Formula

In the MPI solution to this problem, process 0 serves as a manager of a pool of tasks which it distributes among  $P$  additional worker processes (the speedup/efficiency values are to be computed with respect to the number  $P$  of workers):

- The manager holds a pool of partial assignments still to be processed (initially only the empty assignment  $a = []$ ), the number of satisfying assignments reported by the workers so far (initially 0), and, for each worker, the information whether the worker is currently working or waiting for a new assignment to be processed (initially all workers are waiting).
- To each worker that is waiting the manager attempts to send a partial assignment from its pool. However, before the manager gives a worker the last assignment in its pool (such that the pool would become empty), the manager attempts to extend this assignment to  $2^t$  partial assignments of which it forwards one to the worker and keeps the other ones.
- When a worker has processed the partial assignment(s) on its local stack such that this stack becomes empty, it returns the number of satisfying assignments to the server as a result and waits for a new assignment to be processed.
- Whenever a worker returns as a result a number of satisfying assignments, the manager records this information and attempts to send the worker a new assignment. If this is not possible, the manager records the worker as waiting.
- From time to time (e.g., when it has popped a certain number of assignments from its local stack), every worker sends some of its oldest assignments (those with the most variables not yet assigned) from the top of its local stack to the server. Alternatively (or in addition), the manager may send (when its pool gets empty or close to being empty) requests to all workers for new assignments; the workers may regularly (e.g., when popping an assignment from the local stack) poll for such requests and answer them.
- If a server receives from a worker some assignments (rather than a result), it puts them into its pool and attempts (as described above) to send assignments to all waiting workers.
- If the manager detects that all clients are waiting and its pool is empty, it reports the total number of satisfying assignments and informs all workers. The manager and the workers may then terminate.

Please note that in this solution, since each worker may send without request additional assignments to the server, the number of assignments that the manager keeps in its pool is not bounded by  $2^t - 1$ ; thus the pool may have to be dynamically expanded.