

Tecniche di programmazione in chimica computazionale

Optimization and parallelization

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- Check the performances (**profiling**)

Optimization options

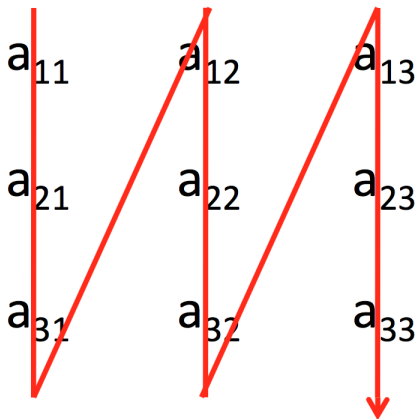
- Use **optimization options** of the **compiler**: -Ok (k=0,1, 2, 3)
(type *man ifort* and search “Specifies the code optimization for applications.”)
- **Automatic** optimization improving code performance
- Example with **profiling.f90**
- Compile with -O0, -O1, -O2 and -O3 (n=100 and m=1000 as input)
- Run the executable by typing **time ./profiling.x**

Swap indexes in matrices

- Methods to linearly store multidimensional arrays in RAM
- Fortran memory management: **column-major order**

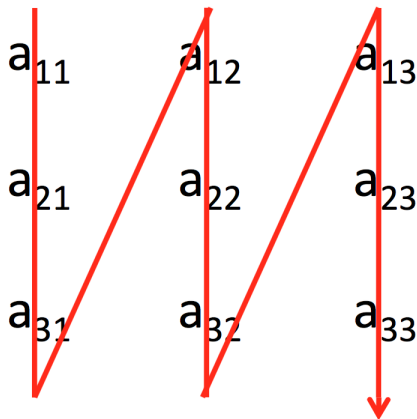
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- Example **matrix_swap.f90** (compile using **-O0**)

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- RS/6000: Practical MPI Programming

Parallelization: where

- RS/6000: Practical MPI Programming
- OpenMP Application Program Interface
- <https://www.openmp.org/resources/tutorials-articles/>

MPI parallelization

- MPI: Message Passing Interface protocol

MPI parallelization

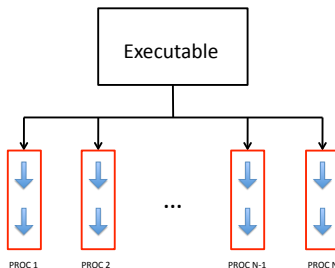
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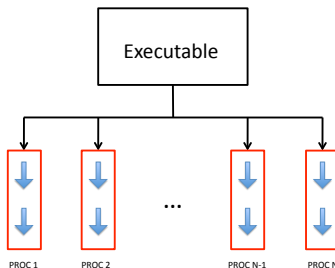
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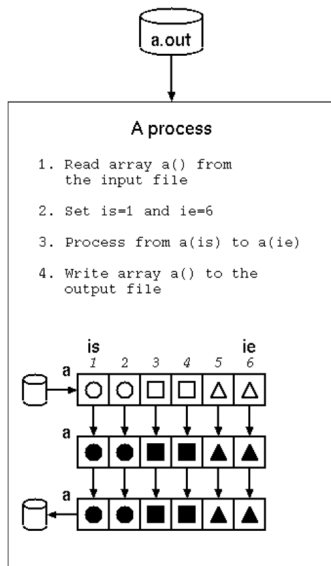
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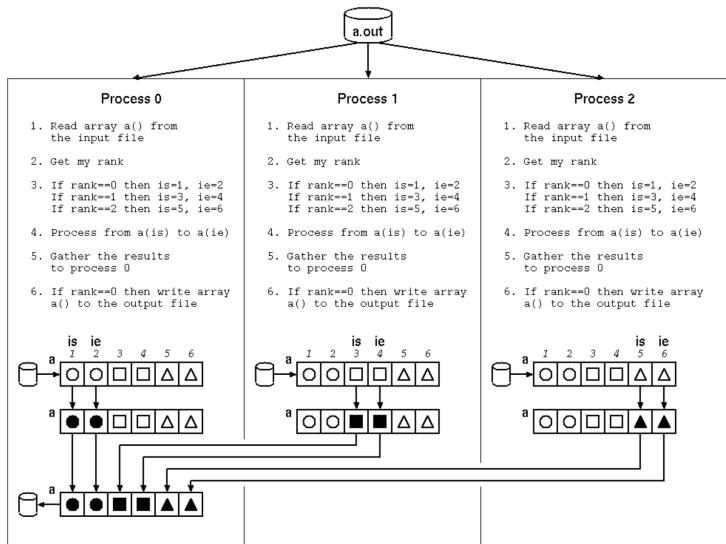


- Each process has a **unique** identifier (**rank**)

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- Example `hello_world.f90`

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- Example: **simd.f90**