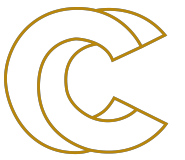


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Leon Kos

University of Ljubljana, FME, LECAD lab

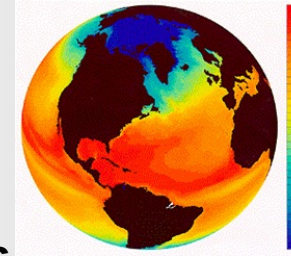
Why supercomputing?



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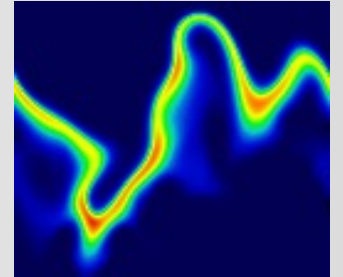
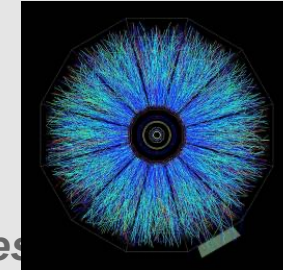
- **Weather, Climatology, Earth Science**

- degree of warming, scenarios for our future climate.
- understand and predict ocean properties and variations
- weather and flood events



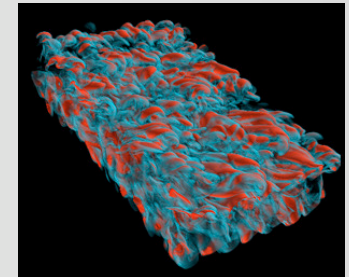
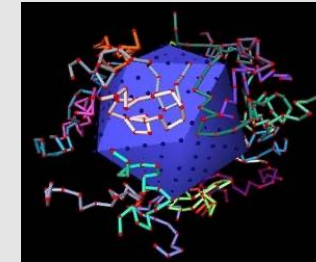
- **Astrophysics, Elementary particle physics, Plasma physics**

- *systems, structures which span a large range of different length and time scales*
- *quantum field theories like QCD, ITER*



- **Material Science, Chemistry, Nanoscience**

- *understanding complex materials, complex chemistry, nanoscience*
- *the determination of electronic and transport properties*

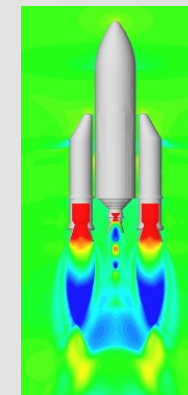


- **Life Science**

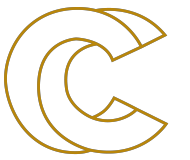
- *system biology, chromatin dynamics, large scale protein dynamics, protein association and aggregation, supramolecular systems, medicine*

- **Engineering**

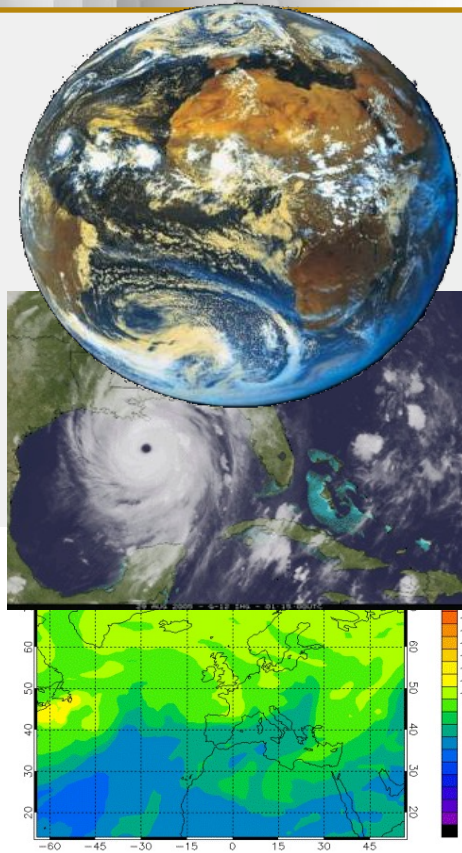
- *complex helicopter simulation, biomedical flows, gas turbines and internal combustion engines, forest fires, green aircraft,*
- *virtual power plant*



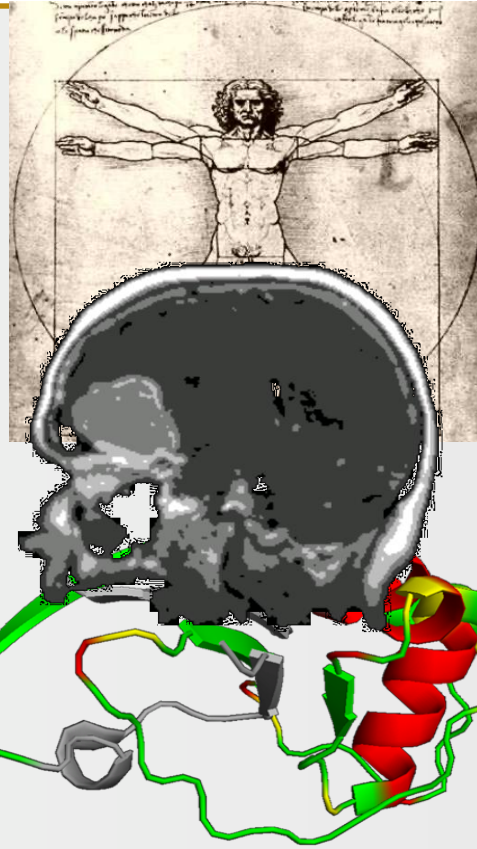
Supercomputing drives science with simulations



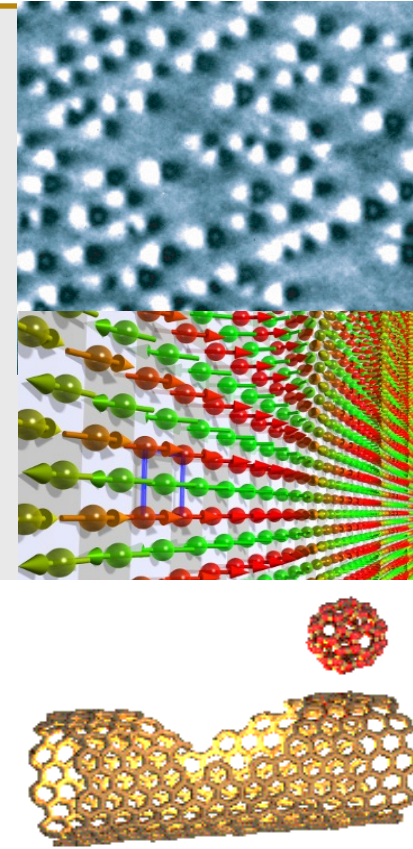
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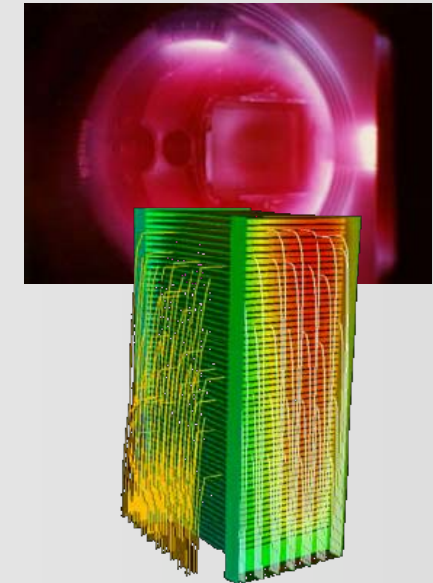
Environment
Weather/ Climatology
Pollution / Ozone Hole



Aging Society
Medicine
Biology



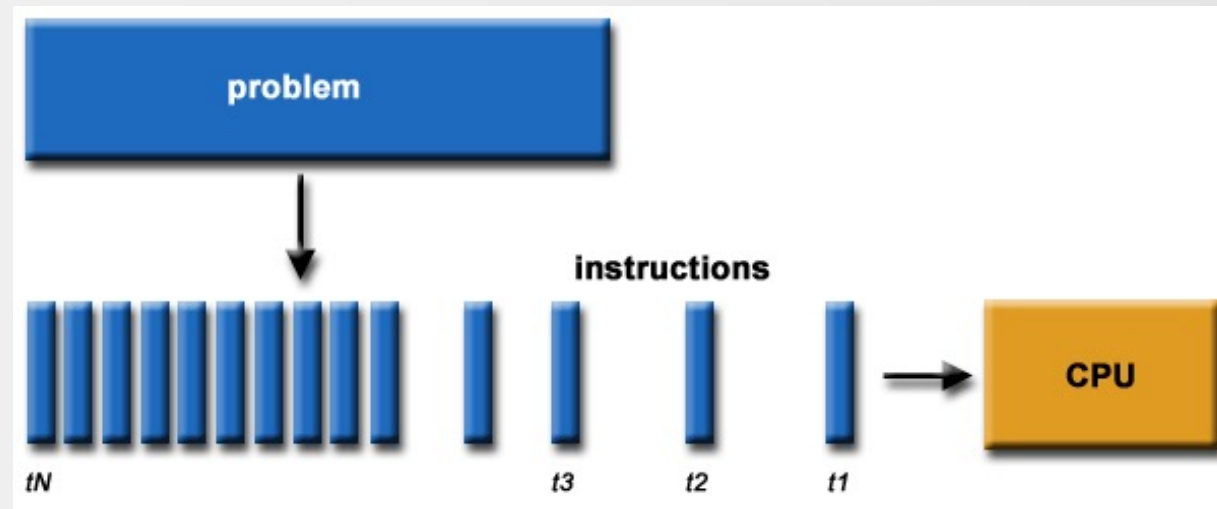
Materials/ Inf. Tech
Spintronics
Nano-science



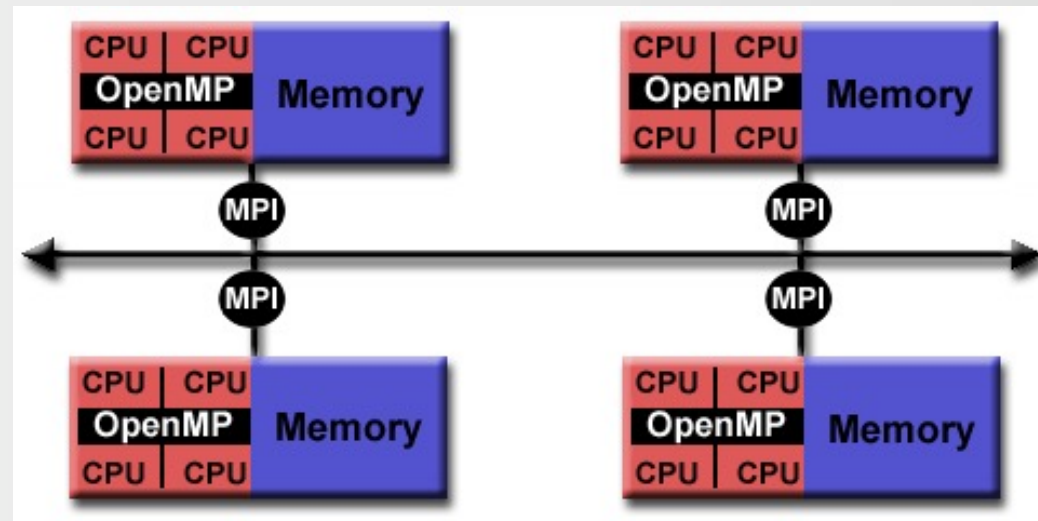
Energy
Plasma Physics
Fuel Cells

Introduction to parallel computing

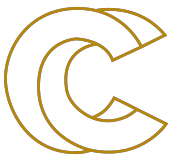
- ▶ Usually is the program written for serial execution on one processor
- ▶ We divide the problem into series of commands that can be executed in parallel
- ▶ Only one command at a time can be executed on one CPU



- ▶ Threading
- ▶ **OpenMP** – *automatic parallelization*
- ▶ Distributed memory model = **Message Passing Interface (MPI)** – *manual parallelization needed*
- ▶ **Hybrid model OpenMP/MPI**

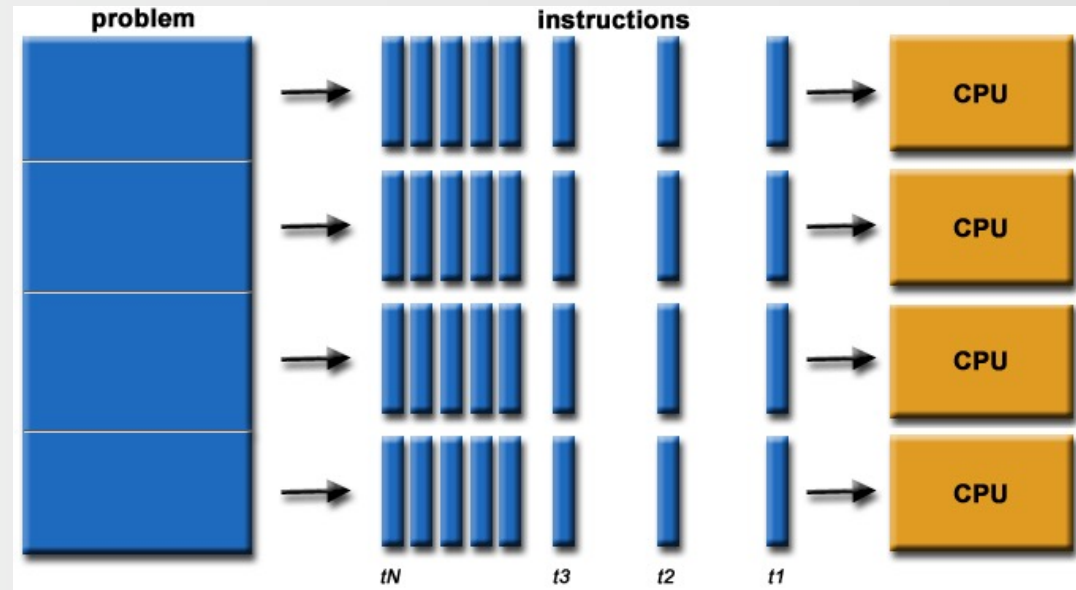


Embarrassingly simple parallel processing

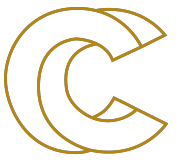


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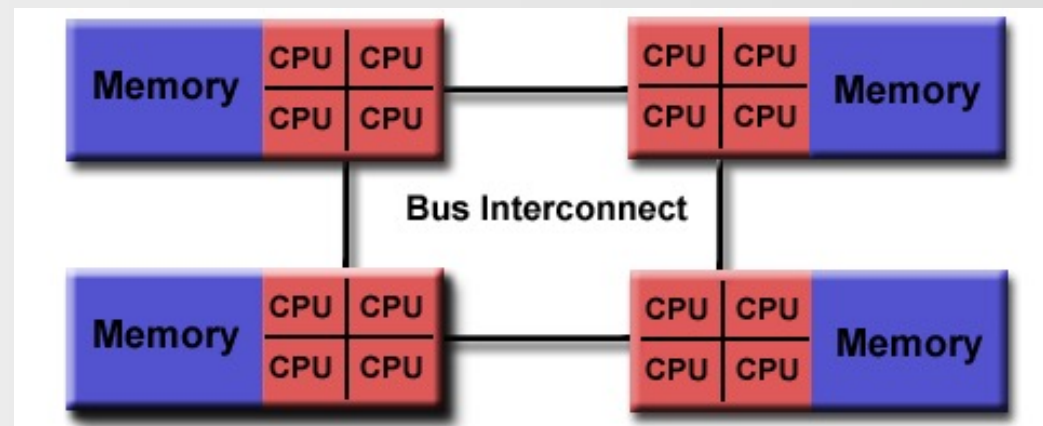
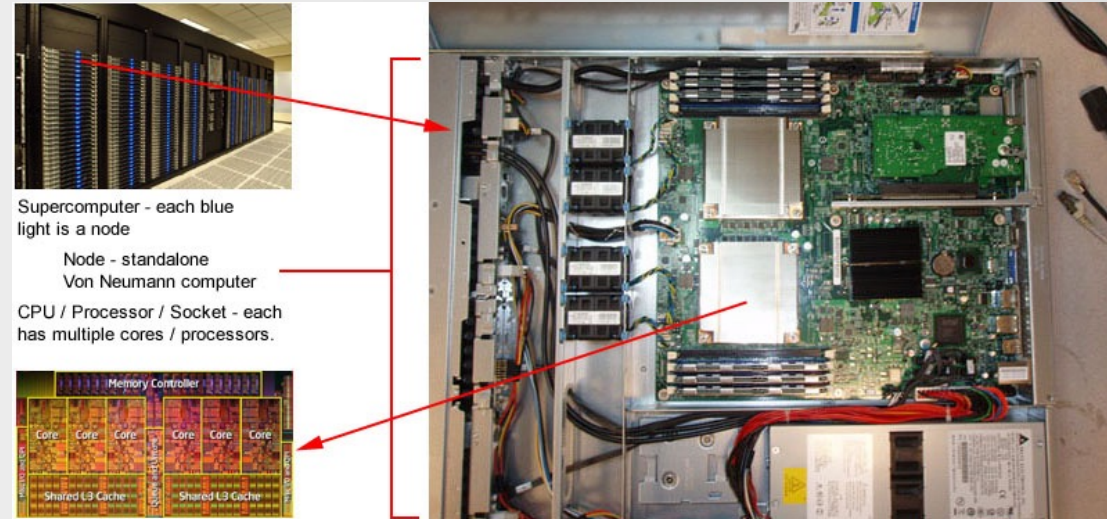
- ▶ Parallel processing of the same subproblems on multiple processors
- ▶ No communication is needed between processes



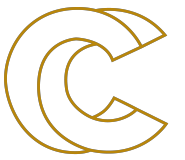
Logical view of a computing node



- ▶ Need to know computer architecture
- ▶ **Interconnect bus for sharing memory between processors (NUMA interconnect)**



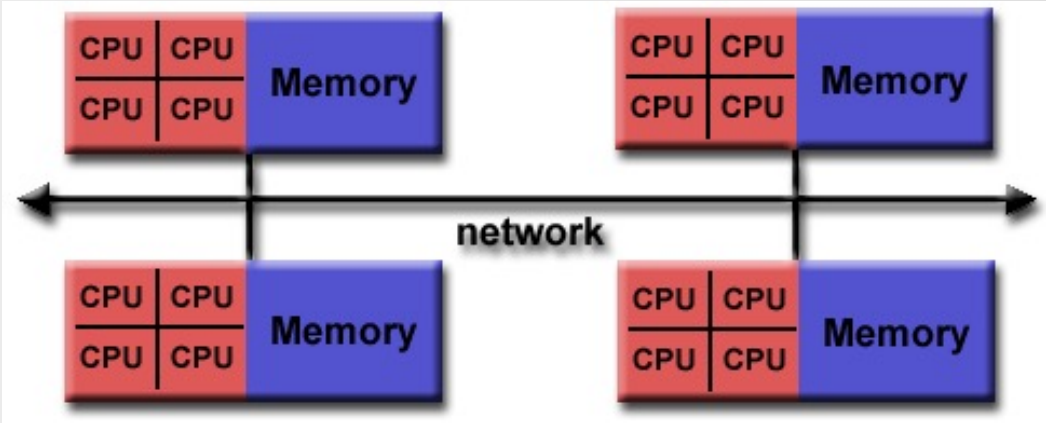
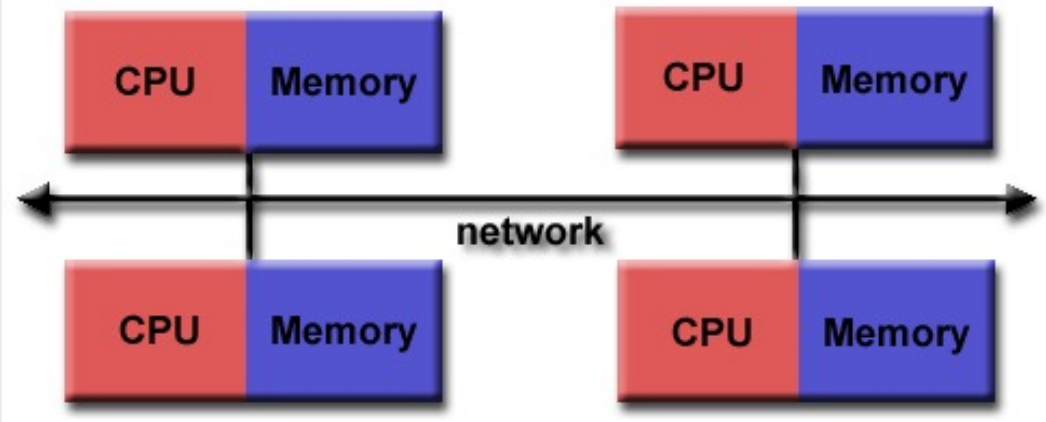
Nodes **interconnect**



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- ▶ Distributed computing
- ▶ Many nodes exchange messages on
 - ▶ high speed,
 - ▶ low latency interconnect such as

Infiniband

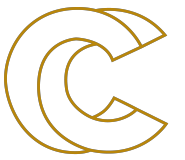


Development of parallel codes

- ▶ Good understanding of the problem being solved in parallel
- ▶ How much of the problem can be run in parallel
- ▶ Bottleneck analysis and profiling gives good picture on scalability of the problem
- ▶ We optimize and parallelize parts that consume most of the computing time
- ▶ Problem needs to be dissected into parts functionally and logically

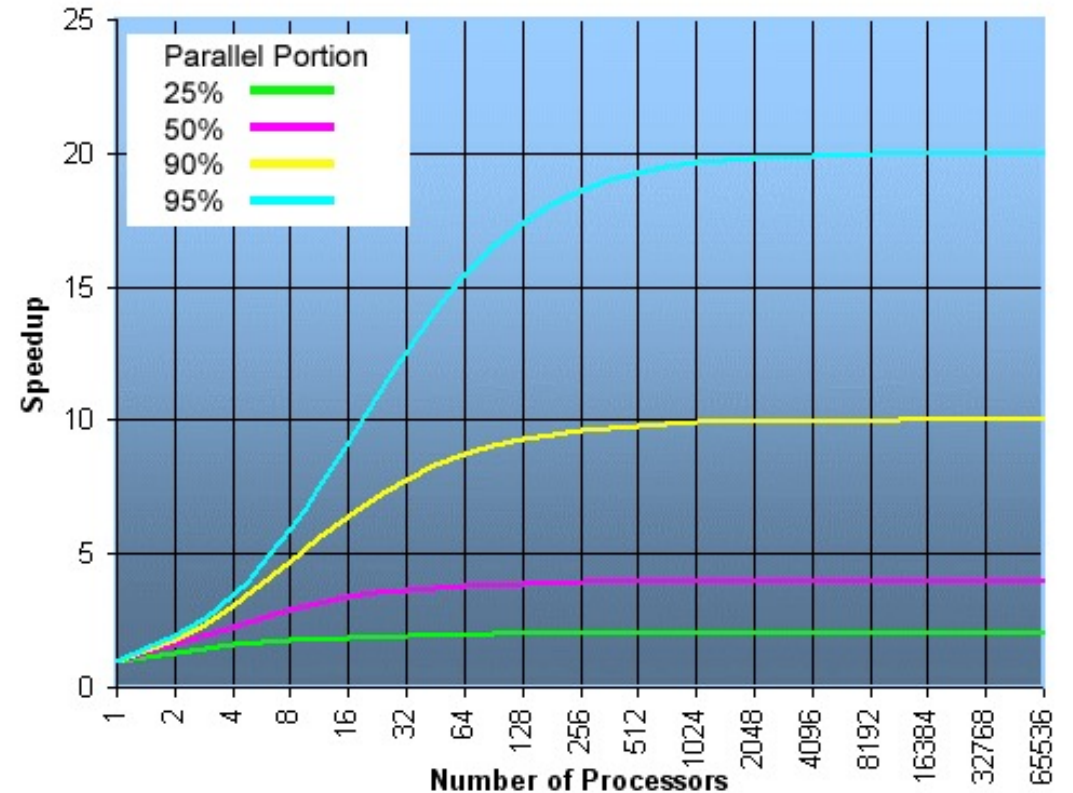
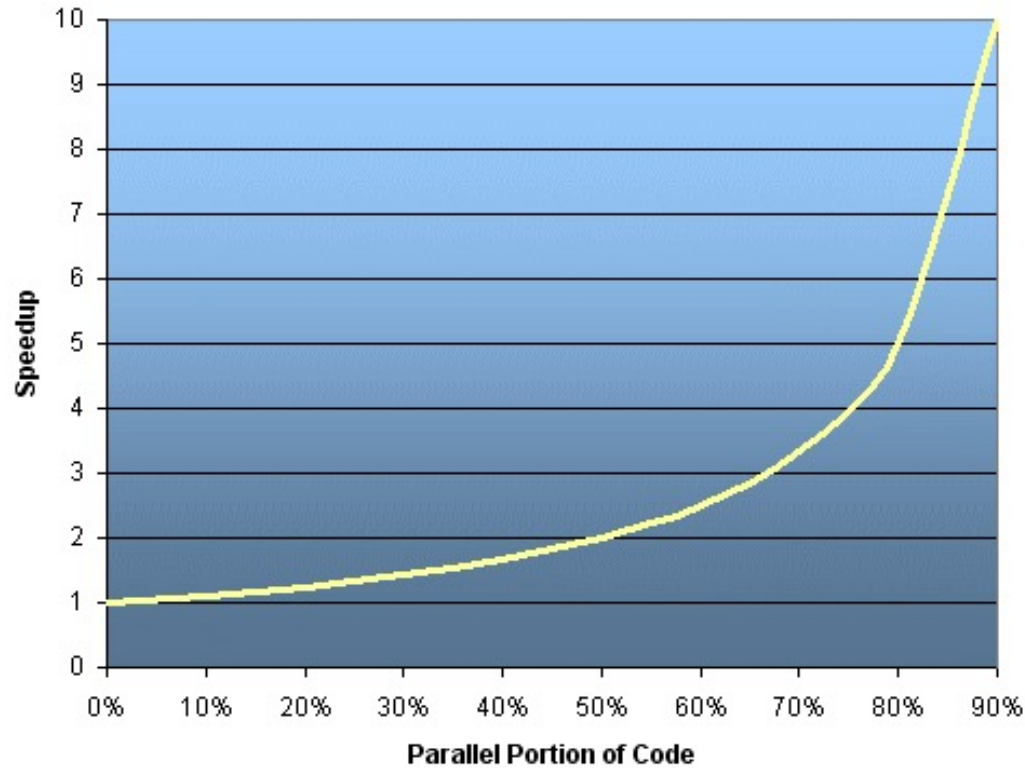
- ▶ Having little and infrequent communication between processes is the best
- ▶ Determining the largest block of code that can run in parallel and still provides scalability
- ▶ Basic properties
 - ▶ *response time*
 - ▶ *transfer speed - bandwidth*
 - ▶ *interconnect capabilities*

Parallel portion of the code determines code **scalability**



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▶ Amdahl's law: **Speedup = 1/(1-p)**



Direct Solver or Iterative Solver?

- We are solving a set of matrix equations of the form $[K]\{u\} = \{f\}$. Here $[K]$ is referred to as the stiffness matrix; $\{f\}$ as the force vector and $\{u\}$ as the set of unknowns.
 - Several millions of unknowns
 - Lot of zeros in K
- Direct solvers: Multfront, MUMPS, and LDLT, Pardiso, ...
- Iterative solvers: PETSc and GCPC, ...

Computer Aided Engineering

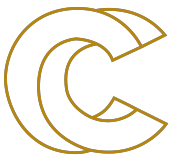
open source tools

- CAD/CAM: **Salome**, **Freecad**, OpenSCAD, LibreCad, Pycam, Camotics, dxf2gcode & Cura
- FEA, CFD & multiphysic simulation: **Salome-Meca** / Code-Aster, SalomeCFD/Code-Saturne, HelyxOs/OpenFOAM, Elmer FEM, **Calculix** with Launcher & CAE GUI, Impact FEM, MBDyn, **FreeFEM**, **MFEM**, **Sparselizard**
- Meshing, pre-post, & visualization: **Salome**, **Paraview**, Helyx-OS, Elmer GUI, VoxelMesher, Tetgen, CGX, GMSH

Questions and practicals on the HPCFS cluster

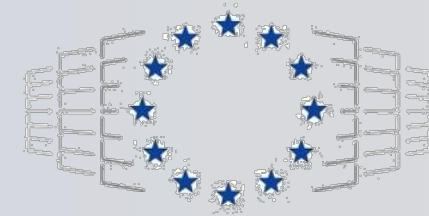


- ▶ Demonstration of the work on the cluster by repeating
- ▶ Access with NX client
- ▶ Learning basic Linux commands
- ▶ SLURM scheduler commands
- ▶ Modules
- ▶ Development with OpenMP and OpenMPI parallel paradigms
- ▶ Exercises and extensions of basic ideas
- ▶ Instructions available at <http://hpc.fs.uni-lj.si/>



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Thanks!



EuroHPC
Joint Undertaking

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