



# *Computação em Larga Escala*

*Introduction to High-Performance Computing*

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# *Summary*

- *High-performance computing*
- *Architectural basics of a parallel machine*
- *Parallel decomposition*
- *Amdahl's Law*
- *Tools to be used to code parallel applications*
- *Suggested reading*

# *High-performance computing - 1*

The area of *High-performance computing* (HPC) is always changing as new technologies and processes become established. In general, it pertains to the use of multiple tightly-coupled processors or computer clusters to run concurrently computational-intensive tasks with high throughput and efficiency. It is common to include in the HPC concept not only the computer architecture, but also a set of elements such as hardware systems, software tools, programming platforms and parallel programming paradigms.

Over the last decade, HPC has evolved significantly, namely due to the emergence of CPU-GPU heterogeneous architectures, which has led to a fundamental paradigm shift in parallel programming.

# High-performance computing - 2

## Top supercomputer sites (as in November 2021)

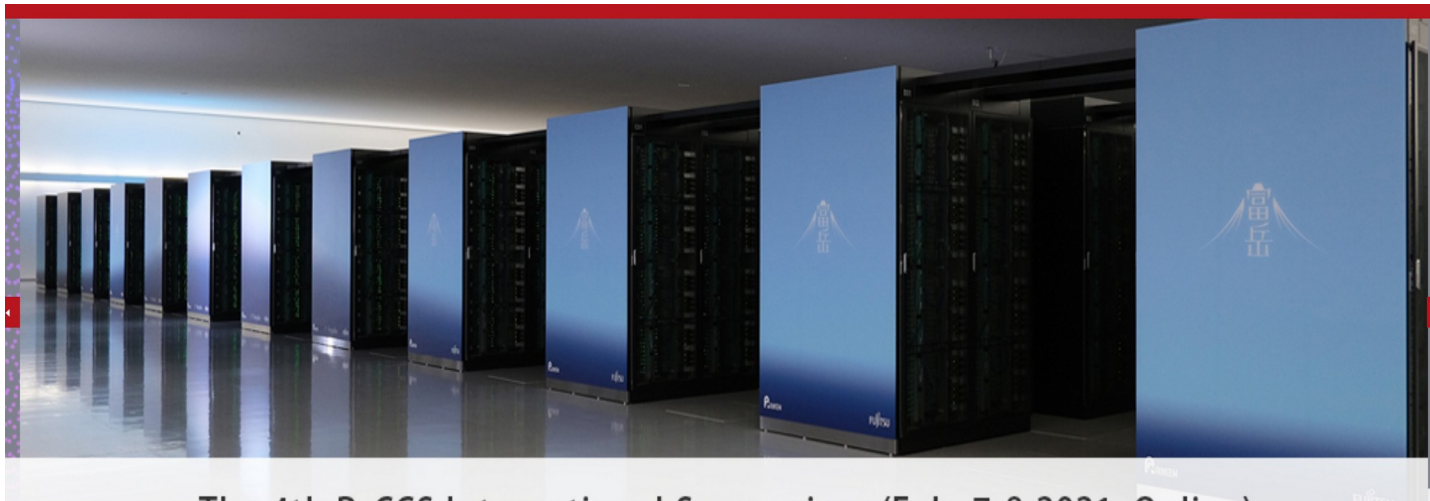
Adapted from: <https://www.top500.org/lists/2021/11/>

| <i>System name</i>   | <i>Node characteristics / Interconnect</i>  | <i>Location</i>   | <i>Number of nodes / cores<br/>Total memory</i>            | <i>High performance Linpack mark<br/>PFlops</i> | <i>Power (MW)</i> | <i>Operating System</i> |
|--|---|---|--|---|-------------------|-------------------------|
| <b><i>Supercomputer Fugaku</i></b><br>Fujitsu A64FX 48C                                  | A64FX 48C 2.2GHz<br>Tofu Interconnect D   | RIKEN Center for Computational Science<br>Kobe, Japan<br><a href="https://www.r-ccs.riken.jp/en/">https://www.r-ccs.riken.jp/en/</a>                                  | 158 976 / 7 630 848<br>48 + 2 core / node<br>5,09 PB       | 442,01  | 29,90             | RedHat Linux            |
| <b><i>Summit</i></b><br>IBM Power System AC922   | IBM POWER9 22C 3.07GHz<br>NVIDIA Volta GV100<br>Dual-rail Mellanox EDR Infiniband | Oak Ridge National Laboratory<br>Oak Ridge, United States<br><a href="https://www.ornl.gov">https://www.ornl.gov</a>  | 4 608 / 2 397 824<br>2 (22 core) + 6 GPU / node<br>2,80 PB | 143,50  | 9,78              | RedHat Linux            |
| <b><i>Sierra</i></b><br>IBM Power System S922LC  | IBM POWER9 22C 3.07GHz<br>NVIDIA Volta GV100<br>Dual-rail Mellanox EDR Infiniband | Lawrence Livermore National Laboratory<br>Livermore, United States<br><a href="http://www.llnl.gov">http://www.llnl.gov</a>   | 3 022 / 1 572 480<br>2 (22 core) + 4 GPU / node<br>1,38 PB | 94,64   | 7,44              | RedHat Linux            |
| <b><i>Sunway TaihuLight</i></b><br>(Divine Power, the light of Taihu Lake)<br>Sunway MPP | Sunway SW26010 260C 1.45GHz<br>Sunway   | National Supercomputing Center<br>Wuxi, China<br><a href="http://www.nscwx.cn">http://www.nscwx.cn</a>  | 40 960 / 10 649 600<br>4 (1+64 core) / node<br>1,31 PB     | 93,01   | 15,37             | Proprietary Linux-based |
| <b><i>Perlmutter</i></b><br>HPE Cray EX235n  | AMD EPYC 7763 64C 2.45GHz<br>NVIDIA A100 SXM4<br>Slingshot-10                     | Lawrence Berkeley National Laboratory<br>Berkeley, United States<br><a href="https://www.nersc.gov/systems/perlmutter/">https://www.nersc.gov/systems/perlmutter/</a> | 4500 / 761 856<br>1 / 2 core + 4 GPU / node<br>0,42 PB     | 70,87   | 2,59              | HPE Cray OS             |

## *High-performance computing - 3*

### **Supercomputer Fugaku**

From: <https://www.r-ccs.riken.jp/en/>



# High-performance computing - 4

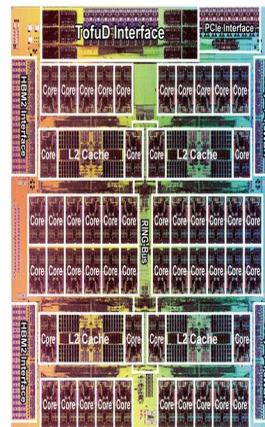
## Fujitsu A64FX 48C

From: <https://www.r-ccs.riken.jp/en/fugaku/about/>

|            |             |
|------------|-------------|
| Technology | 7 nm FinFET |
|------------|-------------|

\*1 Cache performance is with the CPU clock speed of 2 GHz

\*2 Please refer to [GitHub](#)  for details



CPU-Die (Image courtesy of Fujitsu)

### Tofu Interconnect D

The 6D mesh/torus interconnect is used for communication



# High-performance computing - 5

## SI prefixes for orders of magnitude in units

| <i>Prefix</i> |    | <i>Order of magnitude</i> |          | <i>Computing Performance (FLOPS)</i> | <i>Memory Size (B)</i> |     |
|---------------|----|---------------------------|----------|--------------------------------------|------------------------|-----|
| K (kilo)      | Ki | $10^3$                    | $2^{10}$ | KFLOPS                               | KB                     | KiB |
| M (mega)      | Mi | $10^6$                    | $2^{20}$ | MFLOPS                               | MB                     | MiB |
| G (giga)      | Gi | $10^9$                    | $2^{30}$ | GFLOPS                               | GB                     | GiB |
| T (tera)      | Ti | $10^{12}$                 | $2^{40}$ | TFLOPS                               | TB                     | TiB |
| P (peta)      | Pi | $10^{15}$                 | $2^{50}$ | PFLOPS                               | PB                     | PiB |
| E (exa)       | Eo | $10^{18}$                 | $2^{60}$ | EFLOPS                               | EB                     | EoB |
| Z (zeta)      | Zo | $10^{21}$                 | $2^{70}$ | ZFLOPS                               | ZB                     | ZoB |
| Y (yota)      | Yi | $10^{24}$                 | $2^{80}$ | YFLOPS                               | YB                     | YiB |

Supercomputing community is aiming to reach EFlops by 2020 and ZFLOPS by 2030.

# *High-performance computing - 6*

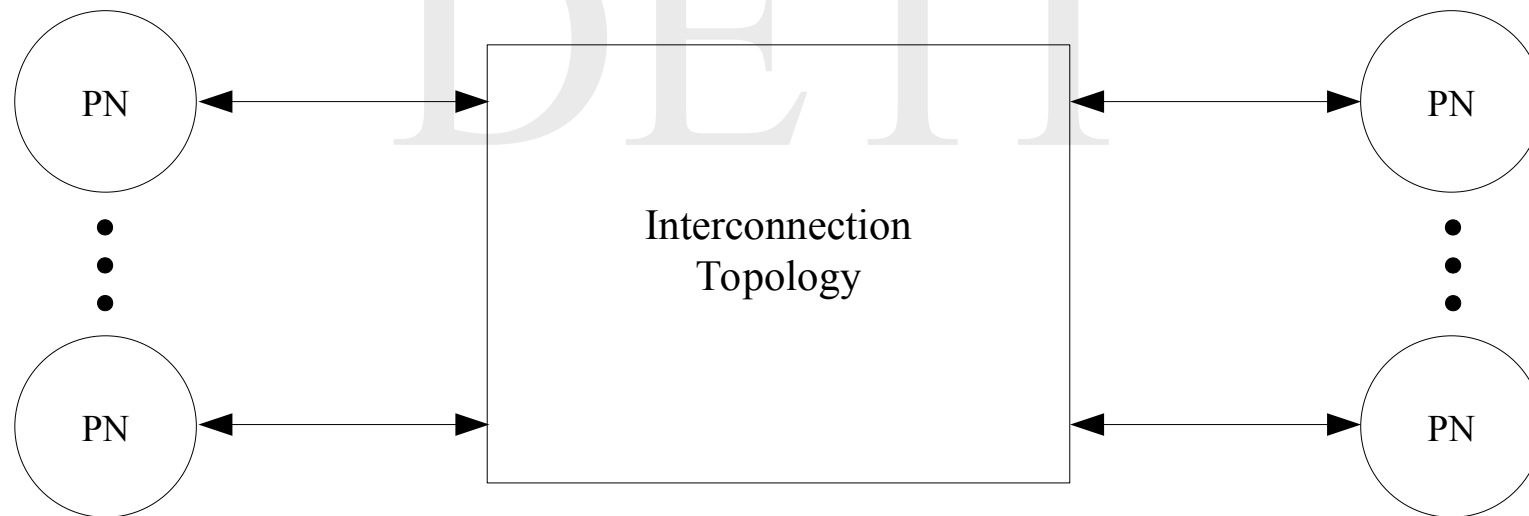
## **Major supercomputing application areas**

- cosmology, astrophysics and astronomy
- computational chemistry, biology and engineering
- computer science
- earth sciences and materials
- weather forecasting
- geographic information science and technology
- global security
- nuclear fusion
- weapons and complex integration



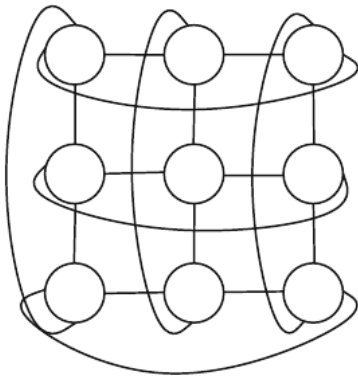
## *Architectural basics of a parallel machine - 1*

Present day high performance computers are at the top level distributed memory parallel machines. They may be thought of as vast clusters of processing nodes (PNs) interconnected by some network topology. The reason for the fact is *scalability*, the ability for system performance to increase as new nodes are attached to it.

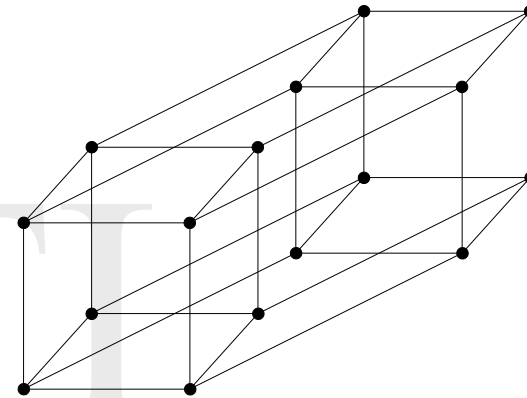


## *Architectural basics of a parallel machine - 2*

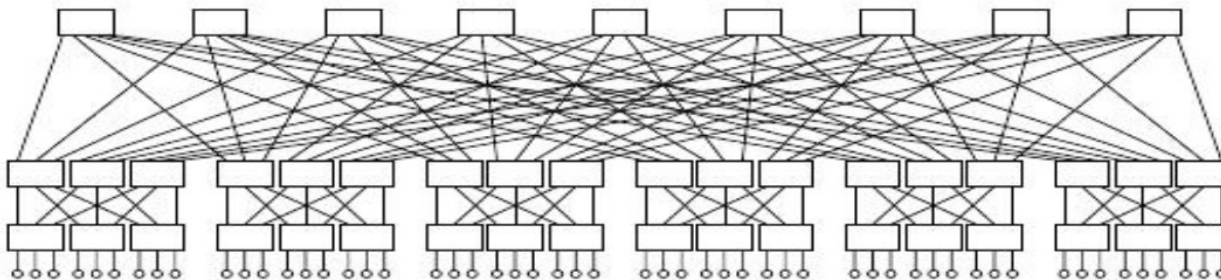
### Common interconnection topologies



*torus mesh*



*hipercube*



*fat tree*

## *Architectural basics of a parallel machine - 3*

hipercube dimension  $n$  = distance  $n$

The key concerns on the interconnection topology are twofold

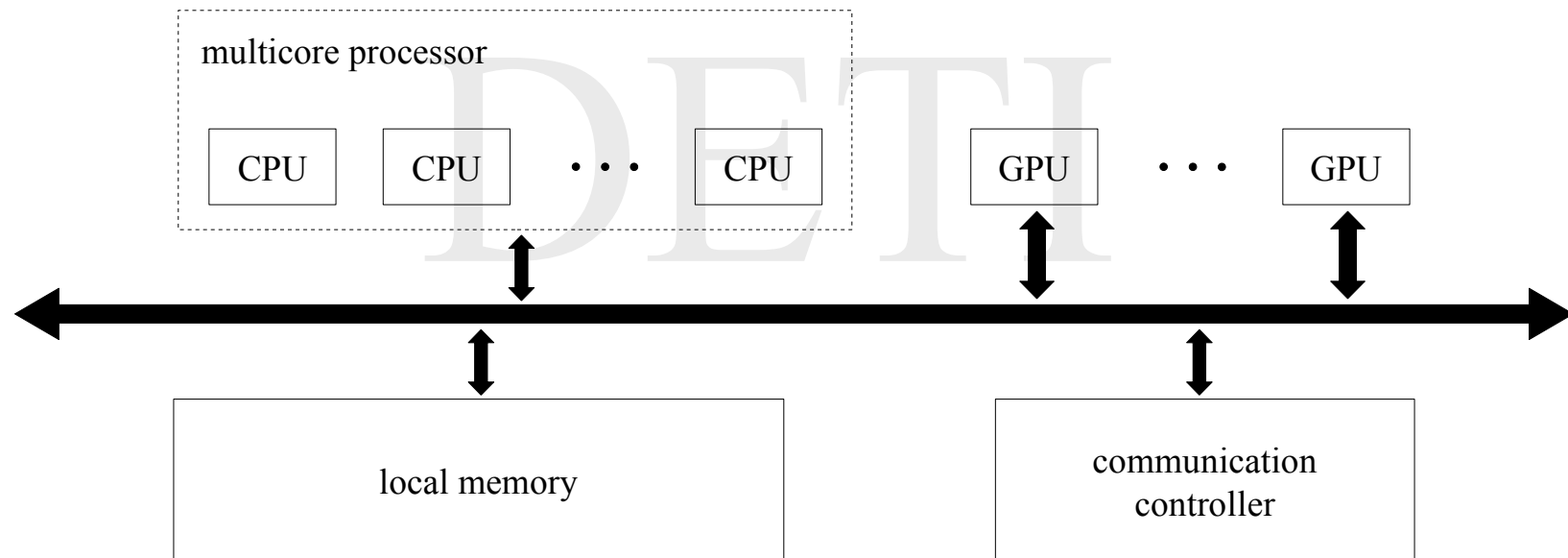
- to keep the number of connections per node small as the number of processing nodes in the cluster increase
- to keep communication time and bandwidth constant as the number of processing nodes in the cluster increase.

Both in the torus mesh and the hipercube, all connections are point-to-point and, as such, have a fixed bandwidth. The number of connections per node is always four in the former case and  $\log_2 n$  in the latter case, where  $n = 2^k$  is the number of processing nodes in the cluster. The communication time, however, depends on the location of the communication nodes, being at the most  $\sqrt{n}$  and  $\log_2 n$ , respectively, of the equivalent communication time between two adjacent nodes.

A fat tree, on the other hand, is a hierarchical network that tries to keep the same bandwidth at all bisections. All processing nodes transmit at the line speed if the packets are uniformly distributed along the available paths. Since a single connection per node is required and, by using  $k$ -port switches,  $k^3/4$  processing nodes may be attached to it, it presents good scalability properties.

## *Architectural basics of a parallel machine - 4*

### Processing node



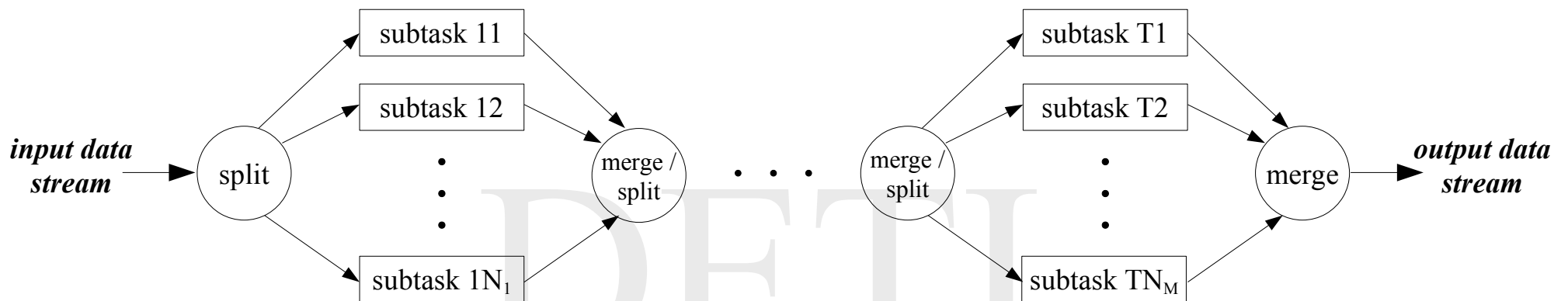
## *Architectural basics of a parallel machine - 5*

A typical processing node consists of one or two multicore CPU sockets and two or more many-core GPUs which gave rise to the name *heterogeneous computing* when referring to this kind of arrangement.

The CPU code is responsible in this context to manage the environment, the code and the data for the GPU, before loading the computation-intensive tasks on the device. GPU computing is not meant to replace CPU computing. CPUs are optimized for dynamic workloads, marked by short sequences of computational operations and unpredictable flow control. On the other hand, GPUs aim at the other hand of the spectrum: workloads that are dominated by computational tasks with simple flow control.

Thus, CPU+GPU heterogeneous parallel computing architectures evolved because the CPU and the GPU have complementary attributes that enable applications to perform best using both types of processors.

## *Parallel decomposition - 1*



Typically, parallel decomposition is data-driven.

Chunks of data of the input stream are fed to a T-stage pipeline of operations. At each stage, data is further split so that operations may be carried out independently in mutual exclusive parts of the chunk being processed. In between the stages, data chunks may undergo reshuffling.

## *Parallel decomposition - 2*

Parallel algorithms may be designed with various degrees of granularity. *Granularity* can be defined as the way how parallel operations are expressed. In this sense, it is not possible to express them without thinking about the hardware platform where the code is going to run.

Parallelism is, thus, organized in three main categories

- *fine-grained parallelism* – parallel operations are expressed at the variable level, it assumes an instruction is executed simultaneously on multiple data sets, a SIMD (single instruction – multiple data) architecture is considered
- *medium-grained parallelism* – parallel operations are expressed at the thread level within a process, a MIMD (multiple instruction – multiple data) architecture of the shared memory type is considered
- *coarse-grained parallelism* – parallel operations are expressed at the process level, a MIMD (multiple instruction – multiple data) architecture of the distributed memory type is considered.

In high performance computing, all three categories of granularity are blended on algorithmic design.

## *Amdahl's Law - 1*

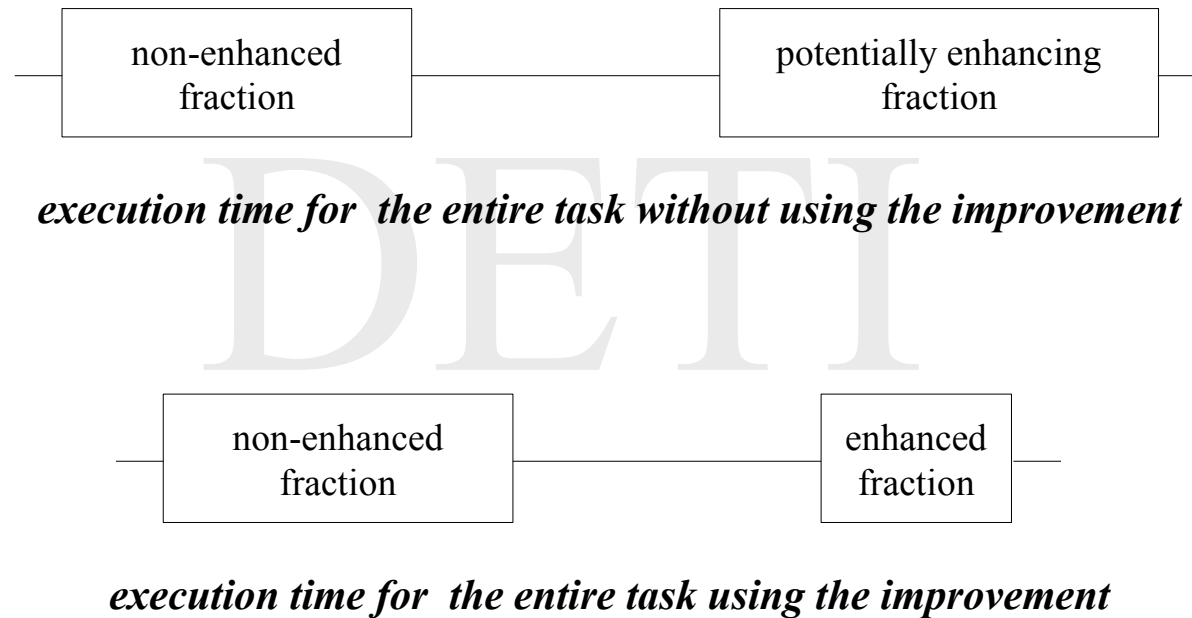
The performance gain that can be obtained by improving some feature of a computer system can be estimated by the *Law of Amdahl*. Amdahl stated in 1967 that *the speed up to be gained from adopting some faster mode of execution is limited by the time fraction of the all operation where the faster mode is used* and is expressed by the formula

$$\begin{aligned}\text{speedup}_{\text{overall}} &= \frac{\text{execution time for the entire task without using the improvement}}{\text{execution time for the entire task using the improvement}} = \\ &= \frac{1}{(1 - \text{frac}_{\text{enhanc}}) + \frac{\text{frac}_{\text{enhanc}}}{\text{speedup}_{\text{enhanc}}}},\end{aligned}$$

where  $\text{frac}_{\text{enhanc}}$  is the time fraction in the original computer system which can be converted to take advantage of the faster mode of execution and  $\text{speedup}_{\text{enhanc}}$  is the speed up to be gained locally by the adoption of the faster mode of execution.



## *Amdahl's Law - 2*



## *Amdahl's Law - 3*

One should point out that, according to Amdahl's Law, there is a well-defined limit to the speedup to be gained. Even if the faster mode of execution reduces the fraction where it is applied to zero, the overall speedup is never larger than the inverse of time fraction where it is not applied.

Thus, parallel decomposition of a problem solution is not usually a straightforward task. Unless the problem lends itself to a self evident partition, one has to look up for decomposition approaches which tend to be quite different from the ones used in the design of the original solution.

When the code is to be run in a distributed memory computer organization, matters become even worse because the parallel algorithm must contemplate the communication time fraction which is not parallelizable. In this latter case, the overall speedup reaches a maximum as the faster mode of execution reduces the fraction where it is applied and then tends to zero

$$\text{speedup}_{\text{overall}} = \frac{1}{(1 - \text{frac}_{\text{enhanc}}) + \text{commOvHead}(\text{speedup}_{\text{enhanc}}) + \frac{\text{frac}_{\text{enhanc}}}{\text{speedup}_{\text{enhanc}}}},$$

## *Tools to be used to code parallel applications*

Parallel applications that are going to be developed, will be written in C Language. Three specific libraries / APIs will be used to implement parallel granularity presented by the algorithms

- *pthread library* – to create multithreaded applications to be run in shared memory architectures (medium-grained parallelism)
- *MPI (message passing interface)* – to create multiprocess applications to be run in distributed memory architectures (coarse-grained parallelism)
- *CUDA C* – to create applications where parallelism is expressed at the variable level, intended to be run in CPU-GPU heterogeneous architectures (fine-grained parallelism).

## *Suggested reading*

- *Introduction to HPC with MPI for Data Science*, Nielsson F., Springer International, 2016
  - Chapter 1: *A Glance at High Performance Computing (HPC)*
- *Programming Massively Parallel Processors: A Hands-on Approach*, Kirk D.B., Hwu W.W., 3rd Edition, Morgan Kaufmann, 2017
  - Chapter 1: *Introduction*