High Performance Computing

Time Complexity and Parallisation

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Given a program or algorithm (assuming a suitable machine model) that depends on a parameter n, or input of size n, we say that it has time complexity (or runtime) $\mathcal{O}(f(n))$ for a function $f : \mathbb{Z}_{\geq n_0} \to \mathbb{R}$, defined on a $\mathbb{Z}_{\geq n_0} := \{n \in \mathbb{Z} : n \geq n_0\}$ for some integer n_0 , if there is a constant time span c and an integer $n_1 \geq n_0$ such that

runtime for parameter/input size $n \le c \cdot f(n)$ for all $n \ge n_1$.

Alternative, one writes $\mathcal{O}(f(n))$, but this does not coincide with the definition of $\mathcal{O}(\cdot)$ in mathematics.

- For example, summing the first *n* integers naively has time complexity $\mathcal{O}(n)$.
- Summing the first *n* integers via the formula n(n+1)/2 time complexity $\mathcal{O}(\log(n)^*)$.

There are algorithms to solve the travelling salesperson problem in time complexity $\mathcal{O}(n^2 2^n)$. It is not known whether there is an algorithm that solves in $\mathcal{O}(p(n))$ for some polynomial function p.

Parallelization gives rise to a new parameter r, the number of compute instances (compute nodes/processes/threads).

Contributions to the runtime TIME(n; r) split up into three fundamentally different terms:

sequential $TIME_{seq}(n)$ parallelizable $TIME_{par}(n)$, and communication/synchronization $TIME_{sync}(n)$.

A naive approach gives

$$TIME(n; r) = TIME_{seq}(n) + \frac{1}{r}TIME_{par}(n) + rTIME_{sync}(n).$$

$$TIME(n; r) = TIME_{seq}(n) + \frac{1}{r}TIME_{par}(n) + rTIME_{sync}(n)$$

attains its minimum at

$$r = \sqrt{\frac{\text{TIME}_{\text{par}}(n)}{\text{TIME}_{\text{sync}}(n)}}$$

and the estimated optimal runtime is

 $\text{TIME}_{\text{seq}}(n) + 2\sqrt{\text{TIME}_{\text{par}}(n)\text{TIME}_{\text{sync}}(n)}.$

Due to the effort of synchronization the best possible number of compute instance is limited. In many problems, $TIME_{sync}(n)$ is so marginal, that effectively this bound is not reached:

 $TIME_{seq}(n) + 2\sqrt{TIME_{par}(n)TIME_{sync}(n)} \longrightarrow TIME_{seq}(n)$ as $TIME_{sync}(n) \rightarrow 0.$

Sub-optimally parallelized programs

In many real-world attempts to "scalar up" a system, however, the impact of $TIME_{sync}(n)$ deteriorates efforts.

Assuming naively that an implementation that is not parallelized has runtime

 $TIME_{seq}(n) + TIME_{par}(n)$,

we have to require that

$$\text{TIME}_{\text{seq}}(n) \leq \frac{\text{TIME}_{\text{par}}(n)}{\sqrt{2}}$$

in order to beat it with a parallelized one.

But close to this cut-off the optimal number of computing instances is $\approx 1/\sqrt{2} < 1$, while in reality a systems "in the cloud" run on hundreds of nodes.

For given r, to achieve

$$TIME(n; r) = TIME_{seq}(n) + \frac{1}{r}TIME_{par}(n) + rTIME_{sync}(n)$$

$$\leq TIME_{seq}(n) + TIME_{par}(n)$$

we have to have

$$\frac{\text{TIME}_{\text{sync}}(n)}{\text{TIME}_{\text{par}}(n)} \leq \frac{1}{r} \Big(1 - \frac{1}{r} \Big) \to \frac{1}{r} \quad \text{as } r \to \infty.$$

If your big data computation on 1000 nodes in the cloud consists of matching ten words in a text, you might violate this bound.

It is also legitimate to analyze TIME(n; r) for $n \to \infty$, but this requires estimates or expressions for the contributions $\text{TIME}_{\text{seq}}(n)$, $\text{TIME}_{\text{par}}(n)$, and $\text{TIME}_{\text{sync}}(n)$.

The runtime of a matrix-vector multiplication of size n on a single node is approximately $2n^2t$, where t denotes the time consumed for one floating point operation.

A parallelized variant in which the matrix gets sliced could, for example, achieve

$$\frac{2n^2t}{r} + (r-1)\Big(l + \frac{nb}{r}\Big),$$

where l and b denote the latency and bandwidth of an underlying communication system.

As $n \to \infty$ is asymptotically $2n^2 t/r$.

Instead of time complexity one can consider speedups. The following two approaches to speedups put emphasis on different aspects.

Focusing on the problem:

SPEEDUP(r) = $\frac{\text{best possible runtime achieved on 1 node}}{\text{runtime achieved with } r \text{ nodes}}$

Focusing on algorithm or implementation:

SPEEDUP(r) = $\frac{\text{time achieved on 1 node}}{\text{time achieved on } r \text{ nodes}}$.

Effects like memory locality can yield

SPEEDUP(r) > 1.

For instance, L1 cache is usually attached to each core, therefore there is more L1 cache available in total when parallelizing a program.