# Parallel Solution of Linear Systems 

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#### Abstract

Computational scientists generally seek more accurate results in shorter times, and to achieve this a knowledge of evolving programming paradigms and hardware is important. In particular, optimising solvers for linear systems is a major challenge in scientific computation, and numerical algorithms must be modified or new ones created to fully use the parallel architecture of new computers. Parallel space discretisation solvers for Partial Differential Equations (PDE) such as Domain Decomposition Methods (DDM) are efficient and well documented. At first glance, parallelisation seems to be inconsistent with inherently sequential time evolution, but parallelisation is not limited to space directions. In this article, we present a new and simple method for time parallelisation, based on partial fraction decomposition of the inverse of some special matrices. We discuss its application to the heat equation and some limitations, in associated numerical experiments.


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## 1. Introduction

A major challenge in scientific computing is to decrease the time required in numerical simulations. To achieve this, algorithms and scientific software must be adapted to fit new computer architectures, often involving parallel programming paradigms. Parallelism is not a new topic in the High Performance Computing (HPC) community, but it has become important for all programmers from the beginning of the century when there were maybe 2 cores per processor, for there are typically 12 cores already and some $60-80$ cores are expected in the near future.

The Domain Decomposition Method (DDM) is an efficient approach to parallelisation in spatial directions, where the whole domain is subdivided to produce efficient iterative

[^0]and independent solutions of smaller problems on the resulting subdomains. Correction steps are necessary to propagate information from one subdomain to others. However, the time direction is also a candidate for parallelisation. In 1964, Nievergelt [1] introduced a parallel algorithm based on time decomposition, and a few years later Miranker et al. [2] defined a parallel solver based on a predictor-corrector scheme. Ref. [7] describes the state of the art on time parallel integration.

In Section 2, we present our Partial Differential Equation (PDE) solver involving time parallelisation, where the original linear system is split into uncoupled linear systems to be solved separately. Our method falls in the category of "Direct Parallel Time Integration" cf. Refs. [3,5]. We eventually discuss its application to the two-dimensional heat equation in Section 3 and some limitations of our method in Section 4, and make brief concluding remarks in Section 5.

## 2. Parallel Computation for Linear Systems

Let $\left(A_{i}\right)_{i=1}^{m}$ be a collection of $m$ nonsingular real matrices of size $n \times n$ and set

$$
X=A_{1} \cdots A_{m} .
$$

We are interested in the following problem: how to compute quickly the solution $x \in \mathbb{R}^{n}$ of the linear system

$$
\begin{equation*}
X x=y, \tag{2.1}
\end{equation*}
$$

where $y \in \mathbb{R}^{n}$ is any given vector?

Sequential approach. One should automatically eliminate the computation of the product of all $m$ matrices before solving the linear system, since this is very expensive (requires a large computational time). Let us recall that solving a generic $n \times n$ linear algebraic system by Gauss elimination ( $n$ large) requires $n^{3} / 3+\mathscr{O}\left(n^{2}\right)$ operations. (For simplicity, we only take into account multiplications and divisions, neglecting additions and subtractions.) The product of two matrices of size $n \times n$ involves $n^{3}$ operations, so computation of the product $X$ above requires $m n^{3}$ multiplications. The following algorithm is a sensible sequential computation of the solution:

- compute the vector $x_{1} \in \mathbb{R}^{n}$ such that $A_{1} x_{1}=y$;
- then compute $x_{2}$, the solution of $A_{2} x_{2}=x_{1}$;
- ...
- and ultimately compute $x_{m}=x$, the solution of $A_{m} x_{m}=x_{m-1}$.

Remark 2.1. The computing cost of this algorithm is $m$ times the computing cost of solving one linear system - i.e. $m n^{3} / 3$ operations.


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