Project Team INRIA: HiePACS Author of the post- doctoral research subject: O. Coulaud, L. Giraud

Title of the post- doctoral research subject: Fast tensor arithmetic for large scale computing and applications

Scientific priorities (cf strategic plan): Computing the future: models, software and digital systems

Applications: materials physics, uncertainty quantification, biodiversity, machine learning

Scientific Research context (10-15 lines):

High-dimensional problems are encountered in many areas of practical interest like in stochastic equations, uncertainty quantification problems, quantum and vibrational chemistry, optimization, machine learning, ... Such problems have led to the introduction of numerical techniques such as variable reduction, reduced modeling, and randomized algorithms. However, the typical dimension of these problems makes unaffordable the existing standard techniques. In the last decades, the theoretical study [1, 2] of low-rank tensor methods has developed and reached a satisfactory level of maturity to be applied on such domains. A new approach particularly well suited is to seek a hierarchical low-rank tensor format that captures the algebraic structure of the system. Formats such as hierarchical Tucker [2] and tensor train [1] have become the convention, and proved to provide efficient and stable data sparse representation for extremely high-dimensional systems. However, accessible and general high-performance parallel software has not yet emerged despite the obvious and its crucial need. We wish to develop new computational tools designed to achieve high-performance for high dimension problems on supercomputing platforms. This entails the development of specialized high-performance, more scalable, robust and further accurate algorithms than contemporary designs.

Post- doctoral researcher work description (10-15 lines):

The first stage will be dedicated to identify and implement tools to compute fast matrix-vector products with the tensor arithmetic in the context of iterative solvers such as eigensolvers or Krylov methods. Many questions have to be addressed namely 1) which hierarchical tensor format should be used 2) how we compute the rank of tensor and 3) what is the best stable algorithm to project the resulting vector in the hierarchical format.

In the second step we will develop new parallel algorithms to reduce the complexity and boost their performance. The test bed for the numerical validation and benchmark of these kernels will be the computation of vibrational spectra of molecules in dimension's space 12 ([4]) and 20. Other domains like biodiversity or machine learning will likely be considered as second application domain. This research action falls within the upcoming research agenda of the HiePACS team with initiatives at the national and the European scale to foster collaboration and exchanges.

Required Knowledge and background:

Candidates should have a strong background in scientific computing with a PhD in computational sciences (applied mathematics, numerical linear algebra, numerical algorithms, ...).

References (max 5):

- [1] W. Hackbusch, *Tensor Spaces and Numerical Tensor Calculus*, vol. 42. Berlin, Heidelberg: Springer Berlin Heidelberg, 2012.
- [2] T. G. Kolda and B. W. Bader, "Tensor Decompositions and Applications," *SIAM Review*, vol. 51, no. 3, pp. 455–500, 2009.
- [3] J. Ballani, L. Grasedyck, and M. Kluge, "A Review on Adaptive Low-Rank Approximation Techniques in the Hierarchical Tensor Format," in *Extraction of Quantifiable Information from Complex Systems*, vol. 102, 2014, pp. 195–210.
- [4] P. S. Thomas and T. Carrington, "Using Nested Contractions and a Hierarchical Tensor Format To Compute Vibrational Spectra of Molecules with Seven Atoms," *The Journal of Physical Chemistry A*, vol. 119, no. 52, pp. 13074–13091, 2015.
- [5] T. Huckle, K. Waldherr, and T. Schulte-Herbrüggen, "Computations in quantum tensor networks," *Linear Algebra and Its Applications*, vol. 438, pp. 750–781, 2013.

Keywords (max 5-6):

Tensor decomposition, high-dimensional, low rank approximation, eigensolver, matric vector products.

Duration:

12 months or 16 months