Scaling performance of the SAGECal calibration package: from LOFAR to SKA

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Abstract

This decade, the Square Kilometre Array (SKA) will perform its first observations. Preparations for building dishes, antennas, correlators and infrastructure are well underway. Concurrently, software for the processing of SKA observations is developed at a number of levels. At a more basic level there are the telescope monitoring and control systems and also the correlator software. On top of that, in order to deliver science-ready data products, software pipelines are needed for radio frequency interference (RFI) mitigation, averaging, calibration and imaging. Here, we focus on the SAGECal calibration package, in particular on the times needed to obtain calibration solutions. This is an important aspect, since this package is now used for the Epoch of Reionization (EoR) Key Science Project of LOFAR, but will also have to run optimally on SKA1 LOW. In terms of number of stations used for observing this amounts to a factor 10 increase, from 51 to 512 stations. Consequently, the disk space needed to store an observation will increase by a factor 100, provided the number of frequency channels remains the same. In this paper we investigate the scaling behaviour of SAGECal, whose runtimes should ideally scale linearly with the number of stations. We also explain the algorithms inside SAGECal and use them to explain its scaling behaviour.

1 Introduction

A number of radio telescopes with many low-gain elements have recently been deployed, designed and/or funded. However, the history of such telescopes goes back to the Mills Cross Telescopes in Australia. Concurrently, software for the processing of SKA observations is developed at a number of levels. At a more basic level there are the telescope monitoring and control systems and also the correlator software. On top of that, in order to deliver science-ready data products, software pipelines are needed for radio frequency interference (RFI) mitigation, averaging, calibration and imaging. Here, we focus on the SAGECal calibration package, in particular on the times needed to obtain calibration solutions. This is an important aspect, since this package is now used for the Epoch of Reionization (EoR) Key Science Project of LOFAR, but will also have to run optimally on SKA1 LOW. In terms of number of stations used for observing this amounts to a factor 10 increase, from 51 to 512 stations. Consequently, the disk space needed to store an observation will increase by a factor 100, provided the number of frequency channels remains the same. In this paper we investigate the scaling behaviour of SAGECal, whose runtimes should ideally scale linearly with the number of stations. We also explain the algorithms inside SAGECal and use them to explain its scaling behaviour.

2 Description of the SAGECal calibration package

SAGECal is an amalgamation of several key algorithms for optimal performance and accuracy in calibration. We briefly discuss these algorithms in the following and highlight their unique characteristics.

- Expectation maximization (EM) and space alternating
generalized expectation maximization (SAGE) algorithms [4]: These algorithms enable us to decompose calibration along multiple directions in the sky into a set of uni-directional calibration problems [5]. Therefore, we get linear scaling with the number of directions being calibrated.

- Expectation-conditional maximization either (ECME) algorithm [6]: We use this algorithm to minimize a Student’s T cost function [7] in robust calibration.

- Limited memory, Broyden Fletcher Goldfarb Shanno (LBFGS) algorithm [8]: By approximating the (inverse) Hessian using a set of vectors, we can calibrate with a large number of unknowns, without running out of memory. Moreover, by using a stochastic LBFGS algorithm [9], we can calibrate data at the highest resolution.

- Riemannian trust region (RTR) algorithm [10]: Combined with the truncated conjugate gradient method [11, 12], we get a solver that scales linearly in memory usage with the number of stations [13] unlike quadratic scaling of other Newton methods. Moreover, because we calibrate on a Riemannian quotient manifold, we get faster convergence compared to Euclidean space calibration.

- Alternating direction method of multipliers (ADMM) [14]: We can calibrate data taken over a wide bandwidth and also data that are stored across a compute cluster by formulating it as a distributed consensus optimization problem [15, 16].

All the above algorithms have been accelerated using GPUs [17, 18]. We show the scaling of the workhorses of SAGECal, i.e., the optimization algorithms RTR and LBFGS, using simulated data. We simulate mock radio telescopes having the number of stations \( N \) from 64 (LOFAR) to 1024 (super SKA). The calibration is performed along one direction only because the linear scaling with more than one direction is well illustrated in our previous work [17, 18].

For the comparison, we use a 12 core Intel Xeon E5-2670 CPU and an NVIDIA Tesla K40c GPU. We plot the time taken by RTR and LBFGS as well as the total time taken to calibrate 10 time samples. We plot the time taken (normalized by the GPU time for \( N = 64 \)) against \( N^2 \) (instead of \( N \)) in Figs. 1 and 2. This is to illustrate the scaling with the number of baselines which is proportional to \( N^2 \).

From Fig. 1 we see that the performance of RTR algorithm is similar in both GPU and CPU versions. The CPU is not overwhelmed while running the RTR algorithm and can keep up with the GPU. In contrast, we see a significant overhead in the CPU version of LBFGS in Fig. 2. The total time in Fig. 2 is dominated by the LBFGS algorithm. Overall, we clearly see the scaling that is dependent on the data, which is in turn dependent on the number of baselines.

Using mini-batches of data as in [9] will improve this and we pursue this as future work. Note also that we have only shown the scaling of the optimization algorithms and for full calibration, additional time is required for model computation and data reading and writing [18].

3 Conclusions

We have shown the scaling of the core optimization algorithms used in SAGECal. With GPU acceleration, we require about \( \times20 \) more computations when we scale from LOFAR to SKA. Considering that there is \( \times8 \) increase in the number of stations and \( \times64 \) increase in the number of baselines, we see that the scaling is superlinear.

4 Acknowledgements

This work is supported by Netherlands eScience Center (project DIRAC, grant 27016G05).
Figure 3. The scaling of both RTR and LBFGS algorithms with $N^2$. Compared to LOFAR, SKA will require about $\times 20$ ($\times 50$ for CPU) more time.

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