

SIMULATEQCD - a Simple Multi-GPU Lattice code for QCD calculations

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Introduction

Increasing GPU power across a competitive market of various GPU manufacturers and GPU-based supercomputers pushes lattice programmers to develop code usable for multiple APIs. In this poster we showcase SIMULATEQCD [1], a Simple Multi-GPU Lattice code for QCD calculations, developed and used by the HotQCD collaboration for large-scale projects on both NVIDIA and AMD GPUs. Our code has been made publicly available on GitHub [1]. We explain our design strategy, give a list of available features and modules, and provide our most recent benchmarks on state-of-the-art supercomputers.

Design strategy

SIMULATEQCD is a multi-GPU, multi-node lattice code written using C++17 and utilizing the OOP paradigm and modern C++ features.

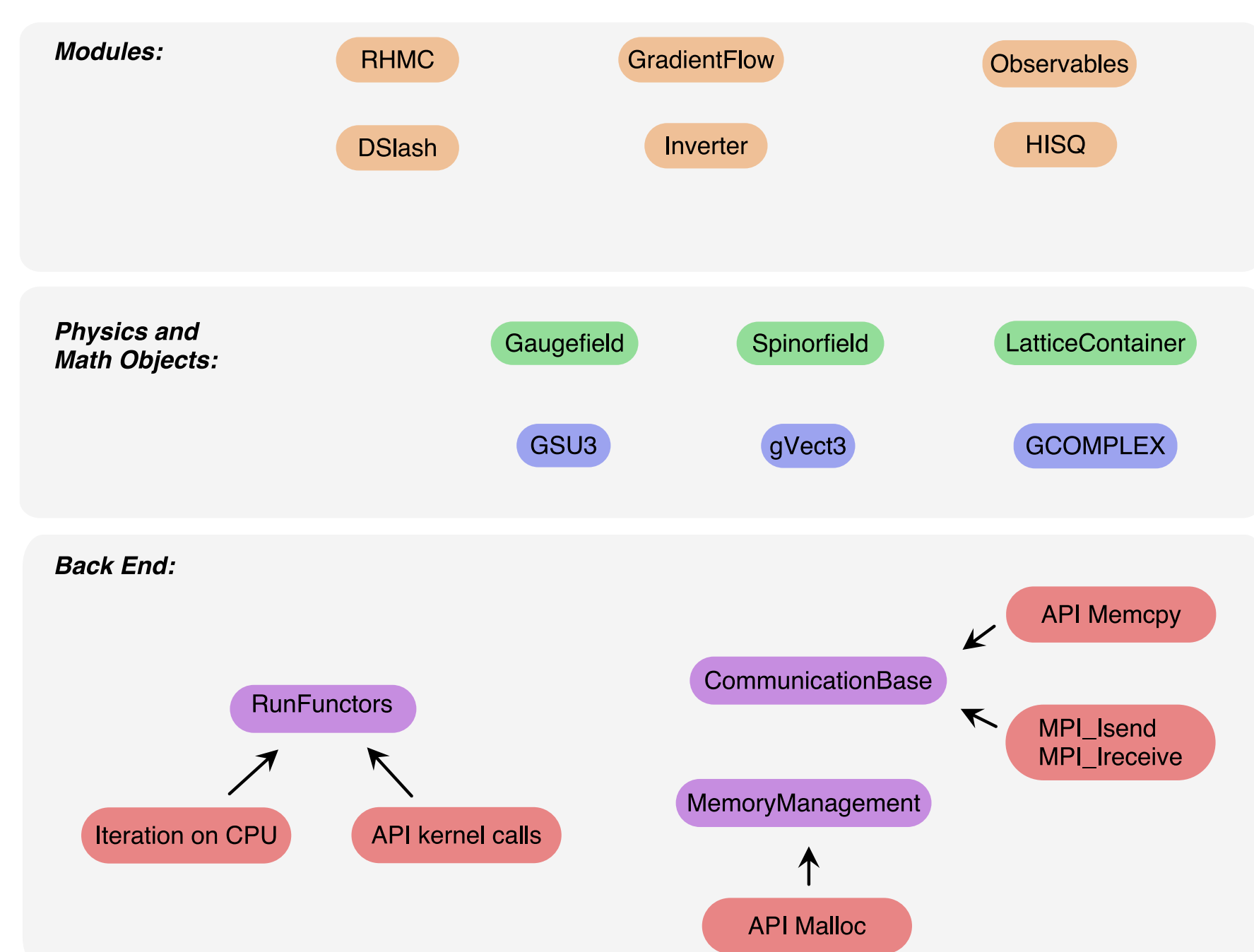


Figure 1: Sketch of inheritance hierarchy. Modules inherit from physics and math objects, which in turn inherit from the back end. Image taken from [3].

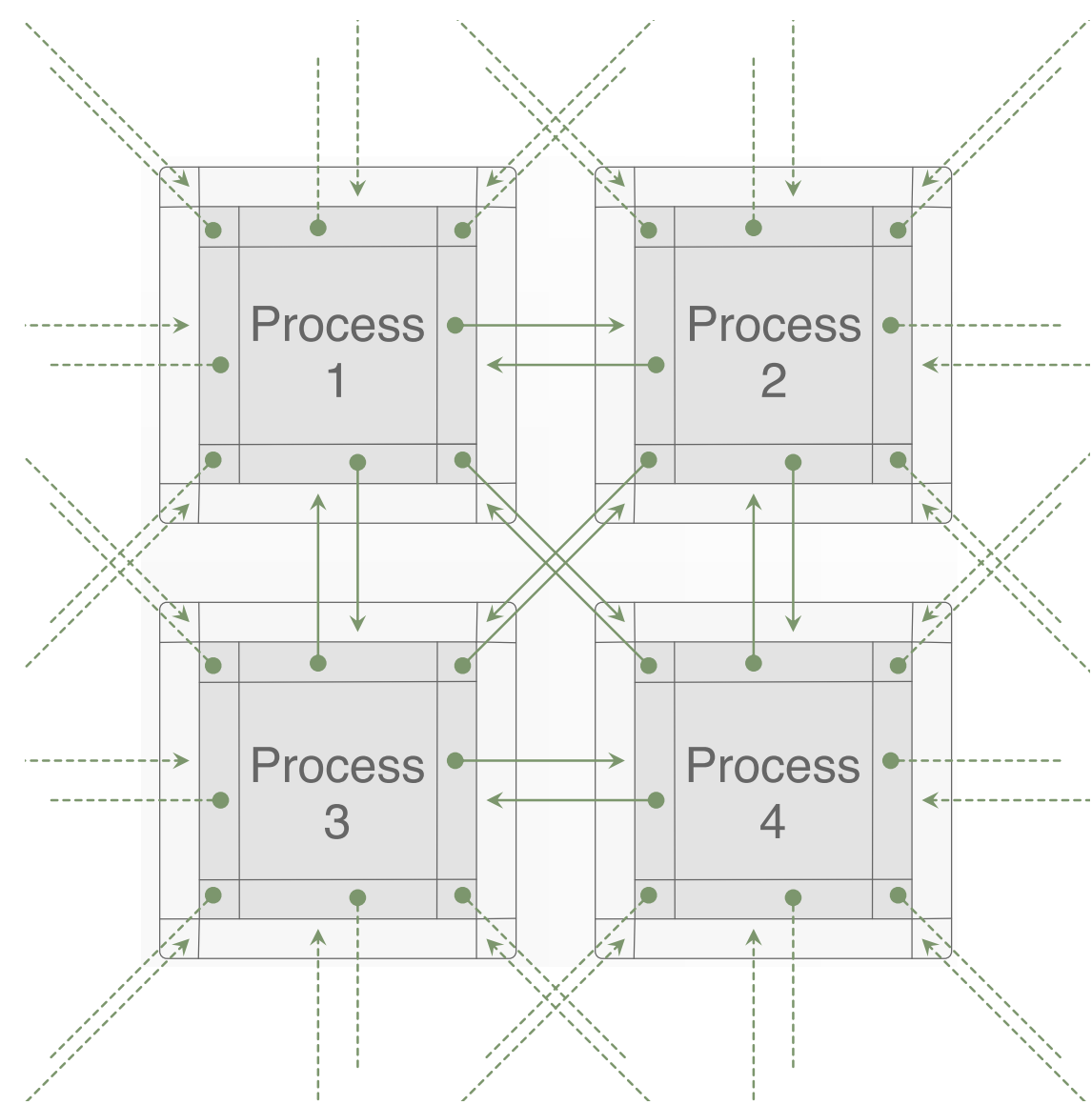


Figure 2: Schematic halo exchange for four processes in two dimensions, each process containing a sublattice. The bulk is indicated by the dark gray squares, and the halo is indicated by light grey squares. Image taken from [2].

Benchmarks

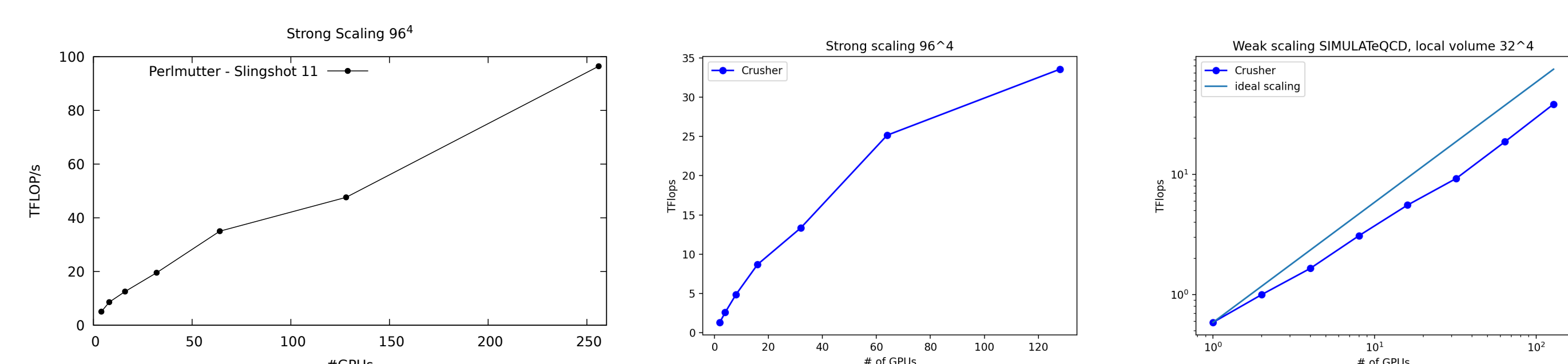


Figure 3: Strong scaling results for the HISQ \mathcal{D} -operator on Perlmutter (left) and Crusher (middle) and weak scaling results on Crusher (right).

Perlmutter: NVIDIA A100 (4 GPUs per Node), slingshot interconnect.

Crusher: AMD MI250X (4 GPUs/8 GPDs per Node), slingshot interconnect.

References

- [1] SIMULATEQCD public code repository, <https://github.com/LatticeQCD/SIMULATEQCD>.
- [2] L. Mazur, Topological Aspects in Lattice QCD, Ph.D. thesis, Bielefeld U. (2021). doi:10.4119/unibi/2956493.
- [3] D. Bollweg, L. Altenkort, D. A. Clarke, O. Kaczmarek, L. Mazur, C. Schmidt, P. Scior, H.-T. Shu, HotQCD on multi-GPU Systems, PoS LATTICE2021 (2022) 196, arXiv:2111.10354.

Functors: abstracting away the technical

```
template<class floatT, bool onDevice, size_t HaloDepth, CompressionType comp>
struct plaquetteKernel{
    gaugeAccessor<floatT, comp> gAcc;
    plaquetteKernel(Gaugefield<floatT, onDevice, HaloDepth, comp> &gauge) : gAcc(gauge.getAccessor()) {}
    __device__ __host__ floatT operator()(gSite site) {
        typedef GIndexer<All, HaloDepth> GInd;
        GSU3<floatT> temp;
        floatT result = 0;
        for (int nu = 1; nu < 4; nu++) {
            for (int mu = 0; mu < nu; mu++) {
                GSU3<floatT> tmp = gAcc.template getLinkPath<All, HaloDepth>(site, nu, mu, Back(nu));
                result += tr_d(gAcc.template getLinkPath<All, HaloDepth>(site, Back(mu)), tmp);
            }
        }
        return result;
    }
};
```

Listing 1: Functor PlaquetteKernel. Allows arbitrary precision floatT, to run on GPU with onDevice==True, for arbitrary HaloDepth, and arbitrary CompressionType. This functor takes a Gaugefield object as argument, whose elements are accessed in memory through the gaugeAccessor gAcc, set to point to the Gaugefield accessor in the initializer list. The argument of operator() indicates that this functor will be iterated over sites. We indicate with All that we run over both even and odd parity sites. getLinkPath multiplies all links starting at site, following a path in the specified directions.

```
const int halodepth = 0;
const bool useGPU = true;
double plaq;
typedef GIndexer<All, halodepth> GInd;
Gaugefield<double, useGPU, halodepth> gauge
LatticeContainer<useGPU, double> latContainer
latContainer.adjustSize(GInd::getLatData().vol4);
latContainer.template iterateOverBulk<All, halodepth>(plaquetteKernel<floatT, HaloDepth>(gauge));
latContainer.reduce(plaq, GInd::getLatData().vol4);
plaq /= (GInd::getLatData().globalLattice().mult()*18);
```

Listing 2: Using an iterator to calculate the plaquette with the functor given above. We instantiate a LatticeContainer, which is needed for the reduction. Our iterator, iterateOverBulk will assign each gSite in the bulk of a sublattice to a GPU thread.

```
template<typename Accessor, typename Functor, typename CalcReadInd, typename CalcWriteInd>
__global__ void performFunctor(Accessor res, Functor op, CalcReadInd calcReadInd,
                              CalcWriteInd calcWriteInd, const size_t size_x) {
    size_t i = blockDim.x * blockIdx.x + threadIdx.x;
    if (i >= size_x) {
        return;
    }
#ifdef USE_CUDA
    auto site = calcReadInd(blockDim, blockIdx, threadIdx);
#elif defined USE_HIP
    auto site = calcReadInd(dim3(blockDim), GetUint3(dim3(blockIdx)), GetUint3(dim3(threadIdx)));
#endif
    res.setElement(calcWriteInd(site), op(site));
}
template<typename CalcReadInd, typename CalcWriteInd, typename Functor>
void RunFunctors::iterateFunctor(Functor op, CalcReadInd calcReadInd, CalcWriteInd calcWriteInd,
                                const size_t elems, ...) {
    ...
#ifdef USE_CUDA
    performFunctor<<< gridDim, blockDim, ... >>>(getAccessor(), op, calcReadInd, calcWriteInd, elems);
#elif defined USE_HIP
    hipLaunchKernelGGL(performFunctor, gridDim, blockDim, ...,
                      getAccessor(), op, calcReadInd, calcWriteInd, elems);
#endif
    ...
}
```

Listing 3: Sketch of iterator implementation. API-dependent constructions are wrapped inside performFunctor and iterateFunctor methods. The user can decide at compile time whether to use the CUDA or HIP API.

Supported actions and algorithms

- Highly improved staggered quark action (RHMC)
- Standard Wilson gauge action (heatbath and over-relaxation)
- Conjugate gradient
 - multi right-hand side inverter
 - multi shift inverter
- Gradient flow with adaptive step-size
- Multi-level and blocking algorithms
- Gauge fixing (over-relaxation)
- Correlator class
- more to be added ...