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Optimizing Hydrodynamic Simulations of Quantum Fluids

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**Introduction**

The lattice Boltzmann method was first popularized for its success in mainstream computational fluid dynamics, but has recently also emerged as an important tool in the fields of physics, biology, and chemistry for understanding complex fluids. Unfortunately, the range of physical problems which can be solved with the algorithm is currently severely limited by its high memory and computational requirements. The capacity of the lattice Boltzmann method to solve scientific problems has therefore become an important problem in high performance scientific computing. The proposed project intends to tackle this problem directly by optimizing an existing scientific code that simulates the evolution of a quantum fluid for more efficient and scalable overall performance.

**Problem**

The naïve implementation of the lattice Boltzmann method which has been previously implemented is overly restrictive on the simulation sizes achievable with available resources, due at least in part to lack of parallelism and inefficient data and memory handling. We propose to remedy these problems by implementing the following:

1. **Single-Core Optimization:** We will vectorize computations and adopt blocking wherever these result in improved performance and reevaluate data structures to improve memory-usage.
2. **OpenMP:** We will parallelize data-independent sections of the single-core optimized code using OpenMP. The final code will combine hybrid MPI and OpenMP model. We need to decide inter-node bandwidth. The problem we need to figure out is although MPI process is multi-threaded, only master thread will make MPI called.
3. **MPI:** We will implement MPI for running OpenMP-implemented functions in parallel and handling intercommunication.
4. **Data Handling:** We will port IO to HDF5 to improve data write times, implement data compression, and improve data processing.

The input data for the program are the grid size, lattice resolution, time duration, and time resolution of the simulation, as well as a prescribed set of initial conditions which depend on the problem to be solved. The output data are the macroscopic variables (for example, local density and fluid velocity) at each time step.

**Analysis**

We implement these techniques and methods with the goal of producing an output code which can solve scientific problems that the input code could not. The performance of each code version will be compared to the original code performance so as to evaluate exactly how effective our changes are and provide us with insight as to what could still be improved. By adopting this incremental development methodology, we aim to be able to correct mistakes and gain foresight into future changes in order to maximize both efficiency and code development. We intend to demonstrate our achievement by producing the following:

1. **Benchmarks** indicating the performance enhancement due to solving each significant sub-problem. We plan to do benchmarking at the beginning of the project and after each new implementation. Performance improvement will be gauged by running the code on INTEL Phi.
2. **Detailed profiling analysis** will be done at the beginning of the project and after the completion of each major project task. Profiling will be done using PerfSuite and Tau.
3. **Scientific Results.** We will finally demonstrate our code’s scalability and scientific capacity by simulating decaying turbulence, a simple-to-implement but computationally-intensive and interesting physical system.

**References**

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