# Optimising and Parallelising d2q9-bgk.c

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February 23, 2022

#### Abstract

d2q9-bgk.c implements the Lattice Boltzmann methods (LBM) to simulate a fluid density on a lattice. This report outlines the techniques I utilised to optimise and parallelise d2q9-bgk.c, as well as a detailed analysis of those techniques. To do so, this report is split into several sections corresponding to different iterations of my code.

# 1 Original Code

I compiled the original d2q9-bgk.c using the GNU Compiler Collection (GCC) with the following command:

Table 1: Execution times of the original code

Test Case Size	Time (s)
$128 \times 128$	29.16
$128 \times 256$	58.71
$256 \times 256$	233.32
$1024 \times 1024$	980.89

Figure 1 contains the total time to initialise, compute and collate each of the test cases when running the ELF file produced. It was important to measure the original code, so that I could quantify the performance improvements of my latter implementations. I measured each of the total times by taking an average of 10 runs on Blue-Crystal Phase 4's (BC4's) compute nodes. Each of BC4's compute nodes is a Lenovo nx360 M5, which contains two 14-core 2.4 GHz Intel E5-2680 v4 (Broadwell) CPUs and 128 GiB of RAM [1]. I took an average of multiple runs because of the variation between runs, which exists due to the inconsistent performance of compute nodes. It is important to note that I measured every execu-

tion time in this report in an identical manner to ensure my results were fair and unskewed.

# 2 Serial Optimisations

# 2.1 Compiler

Table 2: Execution times after compiler changes, and speedup over the original code

Grid Size	Time (s)	Speedup
$128 \times 128$	22.25	1.31
$128\times256$	44.42	1.32
$256\times256$	176.69	1.33
$1024\times1024$	795.41	1.23

I compiled my serial optimised implementation using the Intel C Compiler as opposed to GCC, since it provides better optimised code for Intel processors. Furthermore, I compiled my code with the Ofast option, which set aggressive options to improve the speed of my program, including O3 optimisations and aggressive floating point optimisations [2].

## 2.2 Loop Fusion and Pointer Swap

Table 3: Execution times after loop fusion and pointer swap, and speedup over the original code

Test Case Size	Time (s)	Speedup
$128 \times 128$	19.42	1.50
$128 \times 256$	39.21	1.50
$256 \times 256$	155.64	1.50
$1024\times1024$	635.61	1.54

LBM is a memory bound problem. As a result of this, there was a significant opportunity to optimise d2q9-bgk.c by decreasing the number of memory accesses. One method I utilised to accomplish this was loop fusion. In the original code,

the entire grid was iterated over in four sequential procedures within each timestep: propagate, rebound, collision and av\_velocity. By absorbing these four procedures into the timestep procedure and fusing the four loops, I was able to drastically decrease the number of memory accesses, thereby improving the performance of my program.

Implementing loop fusion offered another significant opportunity to eliminate redundant memory accesses. The original code had a significant quantity of value copying between the cells and tmp\_cells arrays. I was able to eliminate this by writing all new values of cells to a cells\_new array, and simply swapping the pointers of cells\_new and cells at the end of each timestep. I eliminated the tmp\_cells array entirely.

#### 2.3 Arithmetic Improvements

Table 4: Execution times after arithmetic improvements, and speedup over the original code

Test Case Size	Time (s)	Speedup
$128 \times 128$	19.10	1.53
$128 \times 256$	38.49	1.53
$256 \times 256$	153.39	1.52
$1024\times1024$	621.52	1.58

Despite the compiler being able to partially optimise the arithmetic within each timestep without making any changes to the code, there were still some manual improvements that I made to improve the performance of the program. Division is a very slow arithmetic operation relative to multiplication. Therefore, to eliminate a large number of unnecessary division operations I precalculated several values including:

$$\frac{1}{c^2} = 3$$
  $\frac{1}{2c^2} = 1.5$   $\frac{1}{2c^4} = 4.5$ 

where c is the speed of sound. Additionally, I noticed that the number of cells in the grid that were not obstacles tot\_u was recalculated and then divided by each timestep. I eliminated this inefficiency by counting number of cells that were not obstacles only once (during the initialisation phase). I then saved the reciprocol of this value as a parameter num\_non\_obstacles\_r, which I used once per timestep in a multiplicative operation to compute the average velocity.

#### 2.4 Vectorization

Vectorization is the process of converting a scalar implementation to a vector implementation, which enables the compiler to make use of additional registers to perform multiple operations in a single instruction [3]. I utilised several techniques to enforce single-instruction-multiple-data (SIMD) vectorization of the inner loop within each timestep.

Firstly, I converted the t\_speed structure holding cell speeds from an array of structures (AoS) to a structure of arrays (SoA). Previously, the grid was represented with an array of t\_speed structures, whereby each structure contained nine vectors. I altered this such that the grid was represented by one t\_speed structure containing nine pointers, each to an array of floats. Each array of floats contained the values of one vector for each cell within the grid. The SoA format greatly suited vectorization of the inner loop, since it kept memory accesses contiguous when vectorization was performed over structure instances [4].

Having altered the data layout to suit vectorization, I utilised several other techniques to enforce vectorization. Within my code, I implemented the #pragma omp simd pragma to vectorise the inner loop within each timestep. This pragma indicated to the compiler that SIMD instructions could be used to execute iterations of the loop concurrently. The qopenmp-simd option is enabled by default at optimisation levels of O2 or higher to enable OpenMP SIMD compilation [2].

Additionally, I used the restrict keyword to assert that the memory referenced by pointers was not aliased and compiled my program with the restrict to enable pointer disambiguation. This provided a performanced advantage as it prevented the compiler from performing a runtime test for aliasing.

Processors are designed to efficiently move data located on specific byte boundaries, and the compiler is able to perform optimisations when data access is known to be aligned by 64 bytes [5]. To align the cells, cells\_new and obstacles variables, I replaced calls to the malloc and free procedures with the alignment specific replacements: \_mm\_malloc and \_mm\_free, respectively. I used the \_\_assume\_\_aligned procedure and the statement \_\_assume(params.nx % 16 == 0) to inform the compiler that the dynamically allocated variables were aligned. Doing so prevented the compiler from generating conservative code, which would have been detrimental to performance.

Once I had utilised these techniques to enforce

efficient vectorization of the inner loop, I compiled d2q9-bgk.c with the xAVX2 option to direct the compiler to optimise for Intel processors that support Advanced Vector Extensions 2 (AVX2), as BC4's compute nodes do [6].

Table 5: Execution times after vectorization, and speedup over the original code

Test Case Size	Time (s)	Speedup
$128 \times 128$	5.77	5.05
$128 \times 256$	11.57	5.07
$256 \times 256$	41.55	5.62
$1024\times1024$	215.52	4.55

Table 6: Execution times after parallelising (run with 28 threads), and speedup over both the original and vectorized code

		Speedup	
Grid Size	Time (s)	Original	Vectorized
$128 \times 128$	0.70	41.65	8.24
$128 \times 256$	0.81	71.48	14.28
$256 \times 256$	2.63	88.71	15.80
$1024\times1024$	13.43	73.04	16.05

## 3.3 Scaling

## 3 Parallelism

#### 3.1 OpenMP

OpenMP implements parallelism by launching a set of threads that execute portions of code concurrently [7]. I utilised OpenMP's #pragma omp parallel pragma to direct the compiler to parallelise the outer loop in the timestep procedure. Furthermore, I compiled my code with the qopenmp option, which enabled the paralellizer to generate multi-threaded code based on OpenMP directives, as that which I defined. Since the tot\_u variable needed to contain the total velocities of each cell, I used the clause reduction(+:tot\_u) to prevent race conditions. This informed the compiler to create a copy of the tot\_u variable for each thread (initialised to zero), and to sum the local results when the outer loop terminates.

# 3.2 Non-Uniform Memory Access (NUMA)

NUMA is a computer memory design in which memory access time depends on the memory location relative to the processor [8]. Memory is allocated to the closest NUMA region to the thread that first touches the data [9]. Since BC4's compute nodes contain two sockets, the memory access time for a given thread primarily depends on whether the memory is connected to the socket the thread resides in or not. I parallelised the initialisation loops for cells and obstacles to ensure that each thread touched the same data in both the initialise and compute procedures. Furthermore, I set the environment variables OMP\_PROC\_BIND=true and OMP\_PLACES=cores to prevent threads from moving cores.

# References

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