CUDA Point-Plane Collisions

The goal is to maximize particles possible in a physics simulation with 6 plane checks and basic gravity while minimizing kernel and thus convergence time.

Linux

Dependencies

For code that is run on the Linux machine for access to a 4090, this is what was necessary:

1. To get the VM CUDA to run, append

```
export PATH=/usr/local/cuda-11.7/bin$PATH
```

to ~/.profile and source ~/.bashrc . To know it works, run nvcc --version .

- 2. Install GLM somewhere I added an environment variable

 export GLM_INCLUDE_DIR=~/packages/glm to my .profile and got the code from here. Feel free to
 just modify CMakeLists.txt to point to your glm installation if its local to the project, or elsewhere.
- 3. CMakeLists.txt needed to be modified to find the header files in /usr/ using this.
- 4. To run ncu meaningfully, you need root access. https://developer.nvidia.com/nvidia-development-tools-solutions-err nvgpuctrperm-permission-issue-performance-counters

Helpful Commands

- cat /etc/os-release shows the distro and device architecture.
- nvidia-smi provides GPU information (assuming it is a NVIDIA gpu).
- ncu --target-processes all -o <report name> <executable> [executable args] will generate a
 compute report over all kernels given some executable and its args.

Notes

1. Benchmarks were run on a NVIDIA GeForce RTX 4090 with CUDA 12.2.

- 2. The simulation was refactored to correctly run without rendering and over Linux. New logic was added to iterate until last particle velocity reaches a nonzero threshold ≈ 0 , which is the point of "convergence". This is possible with a cudaMemcpy of sizeof(uint32_t), which is assumed to be a very minimal cost per cycle, and an atomicAdd to a counter of "dead particles".
- 3. The original timestep (#define DT_SIMULATION (1.0f / 20.0f)) was updated to (1 / 60.0f) or once every second. That is a pretty smooth DT value.
 - i. Regarding just how quickly the simulation "converges" even for n=1 that is a subjective rate based on my modeling, it does render the base code was stripped of rendering capability.
- 4. Thread size and particle size vary and are tested as hyperparameters. Block size is according to the following:

```
size_t problem_sz = g_particles.h_maxParticles;
g_blocksPerGrid = dim3((problem_sz + g_threadsPerBlock.x - 1) / g_threadsPerBlock.x);
```

- 5. test/... is where many of the .sh scripts were made to iterate over various particle sizes and thread counts.
 - i. The files are labeled in cout/{particle ct}/.
- 6. GPU Memory:
 - i. Global Memory (~25 GB, of which only ~70 MB is storable in L2 cache)
 - a. I am allocating vec3 pos, vec3 vel, and float radius, which is

$$2*sizeof(vec3) + sizeof(float) = 28$$
 bytes per particle

b. The **theoretical limit** to particle allocations on the GPU was just under **1 billion** particles:

$$\frac{\text{device memory}}{size of (\text{particle})} = \frac{25393692672 \text{ bytes}}{28 \text{ bytes}} \approx 906917595 \text{ particles}$$

That said, out of concern for potential margin of error (and other users...), I made the upper bound 7,500,000 particles. My maximum particle runtime allocation was **20.4GB**.

- c. There are of course some other overhead allocations, i.e. d_deadParticles (isn't ideal for global, but can't be in __constant__) either.
- ii. Constant Memory (~64 KB, cached in L1)
 - a. I allocated 4 __constant__ members due to their constant nature, and frequent or "hot" use across all particles for collision checks and whatever else. These were size_t d_maxParticles, size_t d_numPlanes, vec3 d_planeP[6], and vec3 d_planeN[6].
 - b. These may seem small, but GPU constant space is very limited, around about **64 KB**.

7. NCU Concerns

i. It is clear that while global accesses are necessary in this case, compute reports "60% excessive sectors" - less memory is used than is transferred. For example look at this function:

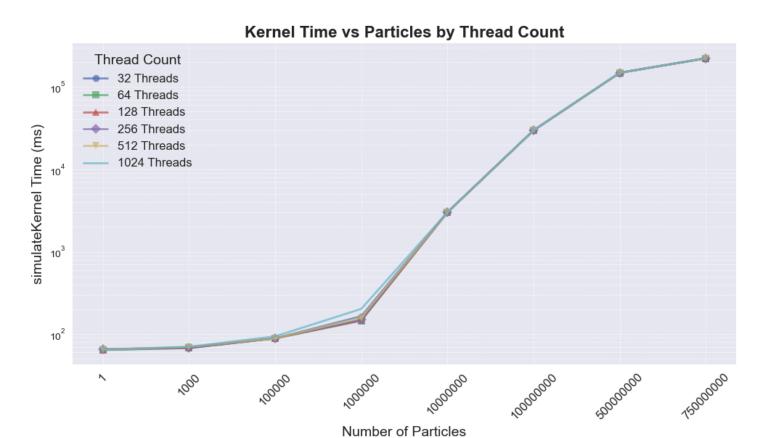
- It is not necessary to make a retrieval from global memory 6 times per plane for the same position.
- Even though we only ask for 12 bytes of our float3, memory boundaries will chunk global memory into sectors, which are usually ~32 bytes per chunk. The data is then retrieved in sectors, meaning when we ask for pos[0], we are perhaps getting pos[1], and **some** of pos[2].
- The reason **some** of pos[2] is problematic is that we now have to do an entirely additional fetch for just 4 bytes of pos[2]. Therefore, it is ideal to align your "object" into even chunks of the sector size.

Improvements

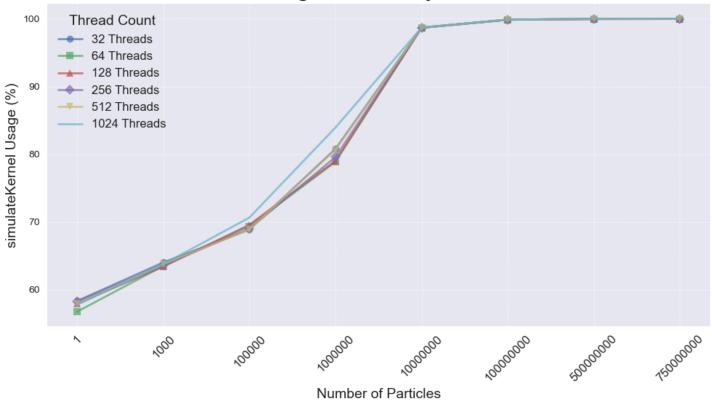
- 1. Padding was added, moving vec3 to vec4 for both pos and vel. This update brought sizeof(particle) = 2 * sizeof(vec4) + sizeof(float) = 28.
 - i. This was done due to the uncoalesced global memory access, which was a major bottleneck. The idea was that sector size, at 32 bytes, is better stored with two vec4s, meaning no request would ever need to use another fetch request or cycle to get the necessary memory.
- 2. Any redundant fetches from global memory were refactored
 - i. const glm::vec3& x(pos[idx]), v(vel[idx]); was moved outside of the loop iterating 6 times per solveConstraints iteration. This doesn't sound like a lot, but when it happens 5 times too many for each particle, it scales quickly.

3. For the most part, the results of this were nonexistent. I would like to run more conclusive tests.

Plots & Images



Kernel Usage vs Particles by Thread Count



Lowest Kernel Time vs Number of Particles (with Corresponding Thread Count)

