Ripple Simulation Based on OpenGL

Parallel Programming Course Project, Team 10

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Outline

- Introduction
- Data Structure
- Experimentation
- Conclusions

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Ripple Simulation

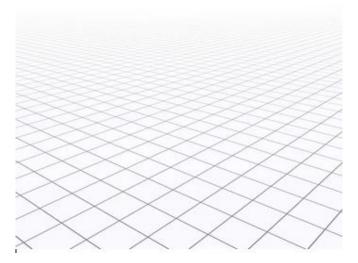
A ripple effect occurs when an initial disturbance propagates due to initial vibration to disturb an increasingly larger portion of the system.

This phenomenon results from the system trying to spread out the energy so it can go back to the initial position.



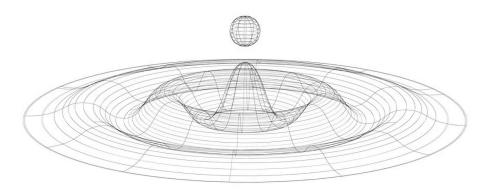
Water Surface Definition

The water surface is a plane in the 3D space, composed of a set of surface_size * surface_size connected vertices, where the vertex has 3 identifiers x, y, and z.



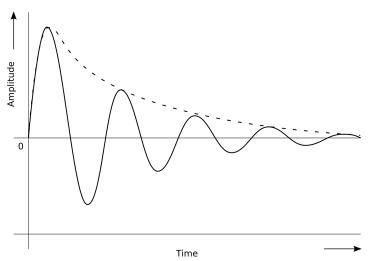
Droplet Simulation

The droplet simulation can be achieved with a particle system. The droplet is a sphere. The acceleration, velocity, and location will be calculated according to the physics laws in the real world. Once the droplet reaches the water surface, it will trigger the isocontour simulation.

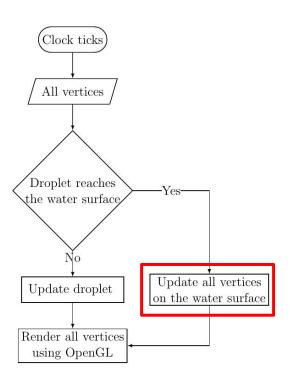


Height Calculation

The isocontour can be seen as a circle with a damped sine amplitude extended outward on a plane of vertices. In world coordinates, each vertex will change its y-coordinate based on its neighbors and a damp coefficient.



System Architecture



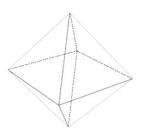
When the clock ticks, the system checks whether the droplet reaches the water surface. The system will update the vertices on the water surface if it reaches the surface.

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Sphere

The proposed sphere consists of an array of vertices and an array of normals. The sphere is constructed by subdivision from a tetrahedron. Each triangle in the tetrahedron is divided into four triangles, and the level of subdivision in our model is 6.



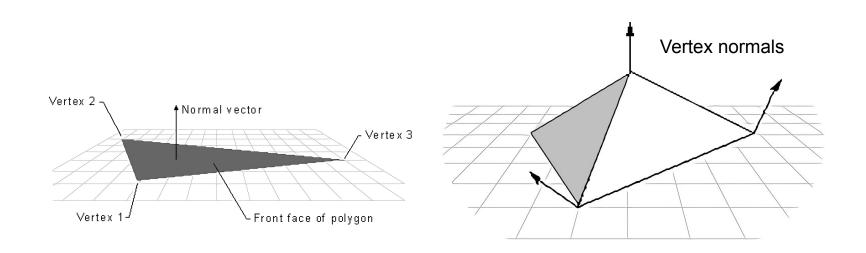






Normal Vector of the Sphere

Because the tetrahedron is centered at the origin of the object coordinate, the normal vector of each point is the same as its normalized coordinate.



Water Surface

There are two arrays containing the vertices. One contains the vertices in the current state, and the other contains the vertices in the next state.

Current:	x1	y1	z1	x2	y2	z2	
Next:	x1	y1	z1	x2	y2	z2	

Proposed Methods

- OpenMP
- CUDA

Ripple Simulation

Each time the ripple function is called, it calculates the new height of each vertex from the current state and stores the results into the next state.

0	0	0	0.4	0.4	0.4
0	1	0	0.4	0.8	0.4
0	0	0	0.4	0.4	0.4

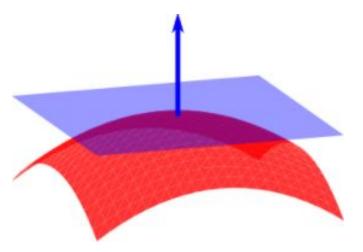
Ripple Simulation (cont.)

```
for (int z = 0; z < surface_size; z++)

for (int x = 0; x < surface_size; x++) {
    new_y = (sum of heights of neighbors in the current state) / coeff;
    new_y -= height at (z, x) in the next state;
    new_y -= new_y / damp;
    height at (z, x) in the next state = new_y;
}</pre>
```

Ripple Simulation (cont.)

After flipping state, it uses the vertices in the current state to update the normals.



Ripple Simulation (cont.)

```
for (int z = 0; z < surface_size; z++)</pre>
          for (int x = 0; x < surface_size; x++) {
             neg z = negative z direction from (z, x);
              pos_z = positive z direction from (z, x);
 4
             neg_x = negative x direction from (z, x);
              pos_x = positive x direction from (z, x);
 6
              new_normal = normalize(cross(neg_z, neg_x) +
 8
                                     cross(neg x, pos z) +
 9
                                     cross(pos_z, pos_x) +
10
                                      cross(pos x, neg z));
11
```

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Environment Settings

CPU: Intel Core i7-9700 (8 cores, 8 threads)

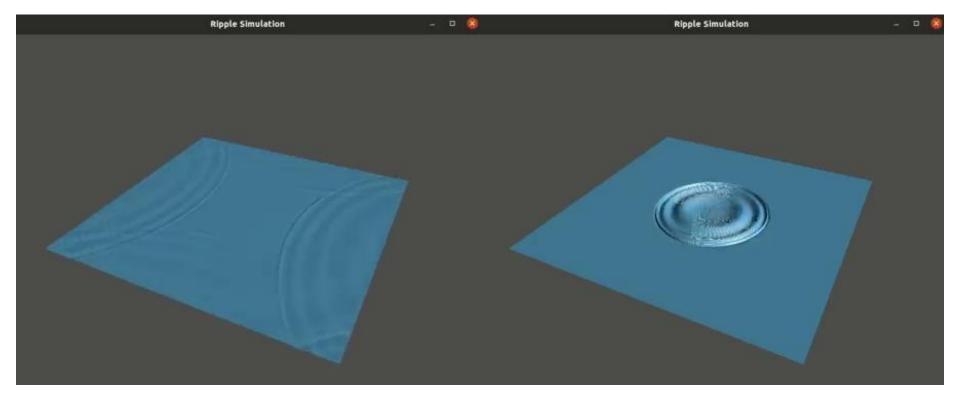
GPU: Nvidia GeForce GTX 1650 SUPER (4 GB GDDR6)

RAM: 32 GB DDR4 2666MHz

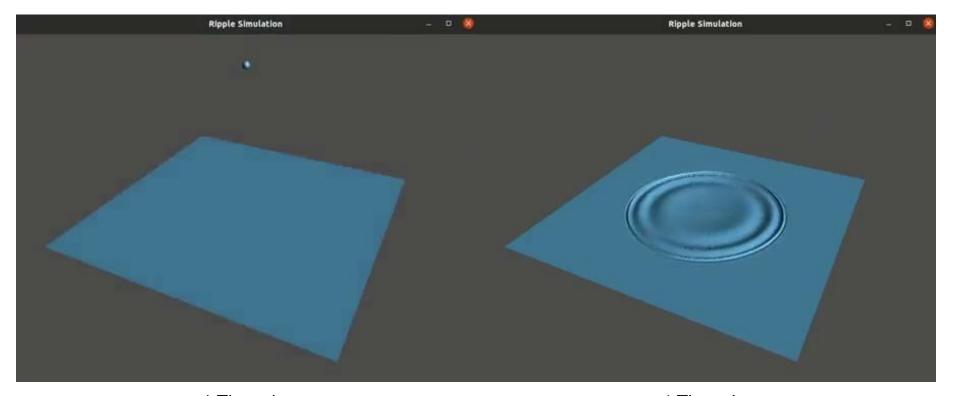
OS: Ubuntu 20.04

CUDA: 11.5

```
Experiment Variables
    OpenMP
        # of Threads: 1, 2, 4, 6, 8
        Surface Size: 400, 800, 1600
        Iteration: 100, 200
    CUDA
        Surface Size: 400, 800, 1600, 3200, 6400
        Block Dim: 25, 50, 100, 200
        # of cudaStream: 1, 2, 4, 8
```



Surface Size 400 Surface Size 1600



1 Thread 4 Threads 22

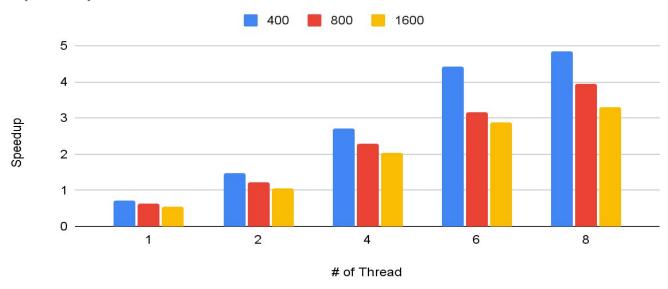
OpenMP Implementation

```
#pragma omp parallel
                                Update Surface
#pragma omp for collapse (2)
        for (x = 1; x < surface size - 1; x++) {...}
#pragma omp for
        for (x = 1; x < surface size - 1; x++) {...}
#pragma omp for
#pragma omp sections
#pragma omp section {...}
#pragma omp section { . . . }
#pragma omp section {...]
#pragma omp section { . . }
```

```
Synchronization
    state = 1 - state;
#pragma omp parallel
                            Update Normals
        // Update normals
#pragma omp for collapse(2)
        for (z = 1; z < surface size - 1; z++) {...}
#pragma omp for
        for (x = 1; x < surface size - 1; x++) {...}
#pragma omp for
        for (z = 1; z < surface size - 1; z++) {...}
#pragma omp for
        for (x = 1; x < surface size - 1; x++) {...}
#pragma omp for
        for (z = 1; z < surface size - 1; z++) {...}
#pragma omp sections {...}
```

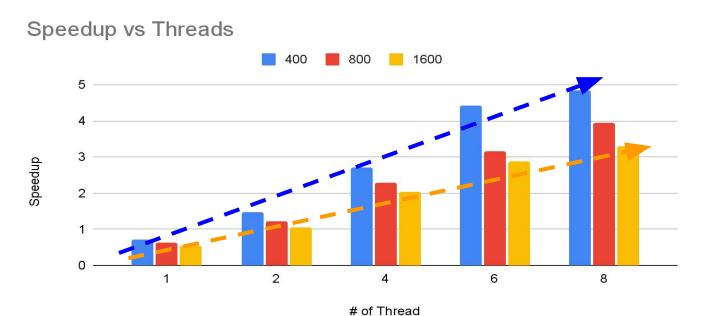
Experiment: OpenMP

Speedup vs Threads



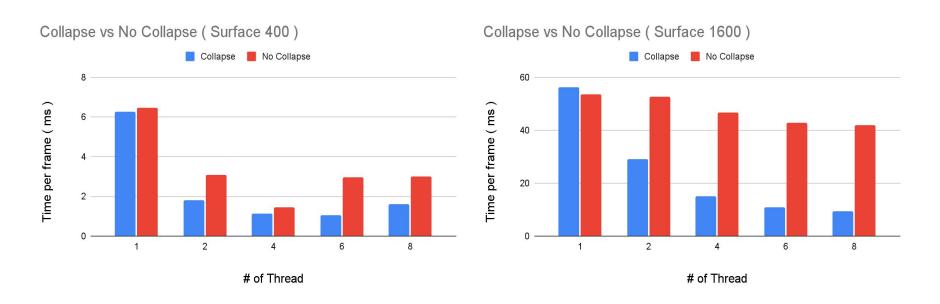
Measured in 100 iterations, best speedup per frame Note that the Time Complixity is $O(\ S^2\)$

Experiment: OpenMP (cont.)



The result shows this application using multithreading, leads to poor scalability w.r.t. data size!

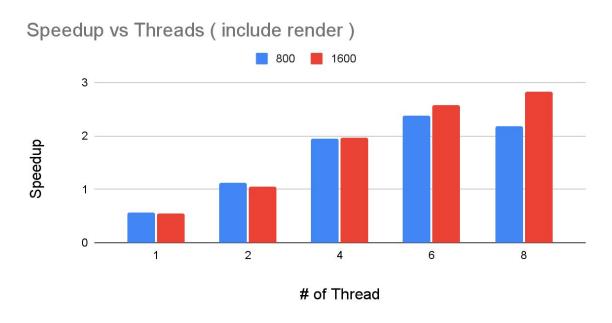
Experiment: OpenMP (for collapse)



Measured in 100 iterations, average execution time per frame

This shows the benefit of collapse for loop during mesh computation

Experiment: OpenMP (OpenGL performance pitfall)



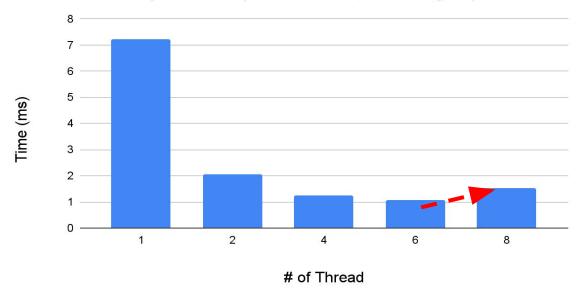
Measured in 200 iterations, average speedup per iteration

Experiment: OpenMP (OpenGL performance pitfall)

Flat p	orofile:					
Each sample counts as 0.01 seconds.					gprof shows 56.37% of program is	
%	cumulative	self		self	total	running OpenGL Mathematics
time	seconds	seconds	calls	ms/call	ms/call	name
43.63	2.43	2.43	100	24.30	54.88	ripple_omp(Surface*, int&)
12.57	3.13	0.70	8606493	0.00	0.00	glm::detail::compute_cross <float, (glm::qualifier)0,="" false=""></float,>
9.87	3.68	0.55	9277987	0.00	0.00	glm::vec<3, float, (glm::qualifier)0> glm::operator- <float,< td=""></float,<>
7.18	4.08	0.40	67281668	0.00	0.00	std::vector <float, std::allocator<float=""> >::operator[](unsi</float,>
7.09	4.48	0.40	6336606	0.00	0.00	glm::vec<3, float, (glm::qualifier)0> glm::operator+ <float,< td=""></float,<>
6.46	4.84	0.36	41015456	0.00	0.00	glm::vec<3, float, (glm::qualifier)0>::vec(float, float, fl
3.41	5.03	0.19	8379948	0.00	0.00	glm::vec<3, float, (glm::qualifier)0> glm::cross <float, (gl<="" td=""></float,>
3.23	5.21	0.18	2338992	0.00	0.00	glm::vec<3, float, (glm::qualifier)0> glm::operator* <float,< td=""></float,<>
2.33	5.34	0.13	2367126	0.00	0.00	glm::vec<3, float, (glm::qualifier)0> glm::operator* <float,< td=""></float,<>
1.44	5.42	0.08	2185603	0.00	0.00	<pre>glm::detail::compute_normalize<3, float, (glm::qualifier)0,</pre>
0.90	5.47	0.05	2334716	0.00	0.00	glm::detail::compute_dot <glm::vec<3, (glm::qualifier<="" float,="" td=""></glm::vec<3,>
0.72	5.51	0.04	204	0.20	0.20	std::vector <float, std::allocator<float=""> >::data() const</float,>

Experiment: OpenMP (OpenGL performance pitfall)

Execution Time per frame (Surface 400, include glm)



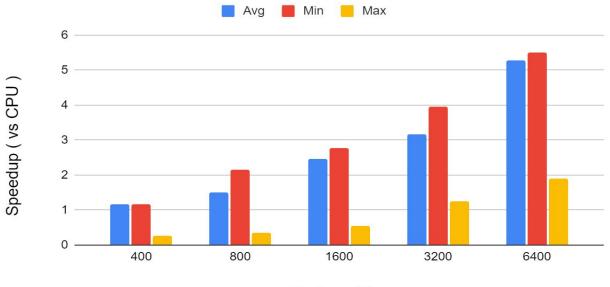
In early implementation, execution time increase if the # of thread == # of CPU cores,

OpenGL Mathematics (glm::cross, etc) requires a dedicated processor

This will result in context switching from frame to frame

Experiment: CUDA (Surface)

CUDA Speedup vs Surface Size

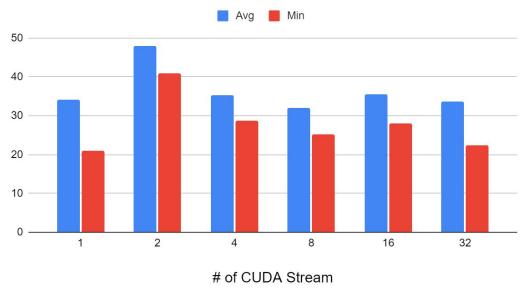


Surface Size

Measured in 100 iterations, only parallel parts Block Dim 25x25, with 8 CUDA streams

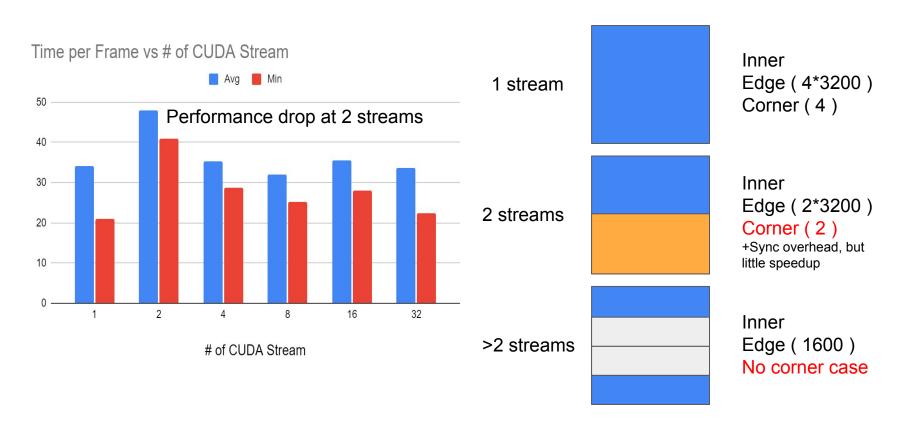
Experiment: CUDA (Stream)





Measured in 100 iterations, only parallel parts Block Dim 25x25, with 8 CUDA streams

Experiment: CUDA (Branching Pitfall)

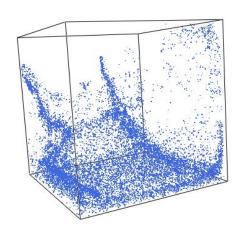


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Conclusions

In this project, a novel thread parallelism-based model is proposed to improve the simulation of the water surface. The future work also includes developing other forms of water simulation such as water fountain, a glass of water, droplets of water on a glass surface, water waves inside a tank.





source: 3D Real-Time Fluid Simulation by a1ex90 https://github.com/a1ex90/Fluids3D.git

Q & A