3D relativistic hydrodynamic computations on graphics processing units

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- Motivation,
- NVIDIA CUDA Framework,
- Hydrodynamics on the GPU,
- Future plans

Problem formulation

Fast and efficient code for hydrodynamics would enable us to study collisions event–by–event in full 3+1D.

For this a huge amount of processing power is required.

GPUs (Graphics Processing Units) seem to be a promising and adequate solution.

Our project aims to implement hydrodynamic algorithms using NVIDIA GPGPU (General Purpose Computing on Graphics Processing Units) solution, namely the CUDA framework.

Early stage of developement.

Motivation

Main reasons to invest in GPGPU:

- High performance leading to big speedups in parallel problems,
- Lower power consumption per FLOPS,
- Low price per FLOPS (Floating point OPeration per Second) typically 0.2–0.3 USD per theoretical GFLOPS for NVIDIA cards 5–15 USD per GFLOPS for Intel processors

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CUDA framework

Programming in a C++—like language.

A cluster of 128–512 processors which execute the *kernel* in parallel.

Up to 4 GPUs in a single unit.

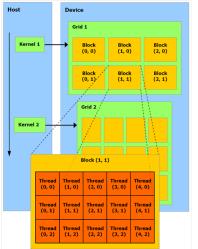
Memory types:

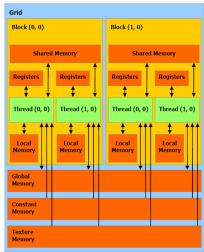
global memory big, slow, common to all threads, accessible by CPU shared memory small, relatively fast, accessed by threads in a block

registers the fastest, only several per thread

Threads are divided into a grid of blocks, and are executed in parallel within a warp.

CUDA framework





images from NVIDIA CUDA Programming Guide

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Hydrodynamic calculations

To solve the hydrodynamics equations:

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\partial_t \mathbf{U} + \nabla \cdot \mathbf{F}(\mathbf{U}) = \mathbf{0}
```

after discretization (in 1+1D):

$$U_i^{n+1} = U_i^n + \frac{\Delta t}{\Delta x} \left(F_{i-\frac{1}{2}} - F_{i+\frac{1}{2}} \right)$$

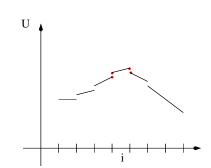
we use MUSCL w/ slope limiting + Musta–Force algorithm, which gives second order accuracy in time and space. In order to apply this scheme, each cell must access itself in the

preceding timestep, and 2 neighbouring cells on each side.

Hydrodynamic calculations

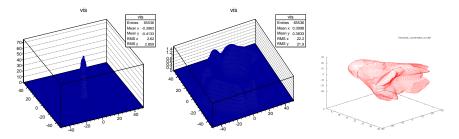
MUSCL makes a linearization inside the cells, edges (red dots) are propagated 1/2 of a timestep and given to Musta–Force, which computes final inter–cell values.

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Preliminary results for 2D with UrQMD generated initial conditions

Plots show initial energy density, energy density after freezeout, and freezeout coordinates.



GPU implementation

We have to map cells \rightarrow threads.

Each thread corresponds to a point in XY plane. The kernel loops over Z axis.

In order to minimize redundant global memory reads, data is cached in shared memory and in registers.

XY planes are cached in shared memory, neighbours in Z direction in registers.

Blocks of threads must overlap by 4 threads in XY plane, as a border condition of size 2 is always required.

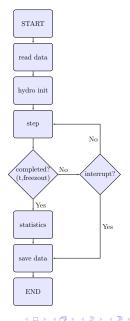
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Future plans

Encapsulate in a flexible, object–oriented C++ interface that allows easy integration with other models and software as a library or stand–alone program.



Draft UML diagram

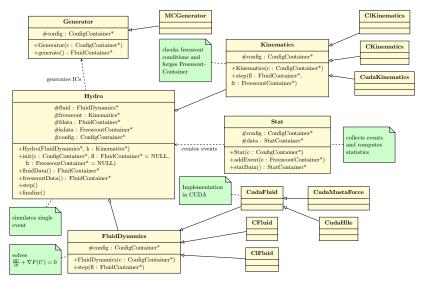


Figure 1: Class hierarchy

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Thank you

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Equations Hydrodynamics equations:

$$\partial_t E + \nabla \cdot \left[(E+p) \, \vec{\nu} \right] = 0 \tag{1}$$

$$\partial_t \vec{M} + \nabla \cdot \left[\vec{M} \vec{v} + p \hat{I} \right] = 0$$
(2)

$$\partial_t R + \nabla \cdot [R \vec{\nu}] = 0 \tag{3}$$

where:

$$E = (e+p)\gamma^2 - p$$
(4)

$$\vec{M} = (e+p)\gamma^2 \vec{\nu}$$
(5)

$$R = n\gamma$$
(6)

$$\partial_t \mathbf{U} + \nabla \cdot \mathbf{F}(\mathbf{U}) = \mathbf{0} \tag{7}$$

LAB frame variables: E, \vec{M}, R, \vec{v} . Fluid element frame variables: e, p, n.

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Integration algorithm

The scheme for equation

 $\partial_t \mathbf{U} + \nabla \cdot \mathbf{F}(\mathbf{U}) = \mathbf{0}$

is

$$U_{i}^{n+1} = U_{i}^{n} + \frac{\Delta t}{\Delta x} \left(F_{i-\frac{1}{2}} - F_{i+\frac{1}{2}} \right)$$
(8)

Image: A matrix

We need to know $F_{i+\frac{1}{2}}$ and $F_{i-\frac{1}{2}}$.

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Computing $F_{i+\frac{1}{2}}$ and $F_{i-\frac{1}{2}}$ — Musta–Force method

In first iteration $U_L = U_i$ and $U_R = U_{i+1}$. Step 1:

$$F_L = F(U_L), F_R = F(U_R)$$

$$U_M = \frac{1}{2} [U_L + U_R] - \frac{1}{2} \frac{\Delta t}{\Delta x} [F_L - F_R]$$
(9)

$$F_M = F(U_M) \tag{10}$$

$$F(U) = \begin{pmatrix} (E+p)\vec{v} \\ \vec{M}\vec{v} + p\hat{I} \\ R\vec{v} \end{pmatrix}$$
(11)

$$F_{i+\frac{1}{2}} = \frac{1}{4} \left[F_L + 2F_M + F_R - \frac{\Delta x}{\Delta t} \left(U_R - U_L \right) \right]$$
(12) (13)

To compute F(U) we additionally need to know \vec{v} and p.

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Computing $F_{i+\frac{1}{2}}$ and $F_{i-\frac{1}{2}}$ — Musta–Force method

Step 2:

$$U_L^{new} = U_L - \frac{\Delta t}{\Delta x} \left[F_{i+\frac{1}{2}} - F_L \right]$$
(14)
$$U_R^{new} = U_R - \frac{\Delta t}{\Delta x} \left[F_R - F_{i+\frac{1}{2}} \right]$$
(15)

We then substitute new $U_{L,R}$ in step 1.

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SLIC/MUSCL scheme

In this method we create a linear approximation of U(x) inside $\left[x_{i-\frac{1}{2}}, x_{i+\frac{1}{2}}\right]$, so that on limit points:

$$U_L = U_i - \frac{1}{2}\Delta_i$$
(16)
$$U_R = U_i + \frac{1}{2}\Delta_i$$
(17)

where Δ_i is the slope vector. Next step is to evolve $U_{L,R}$ by a half step:

$$\bar{U}_{L,R} = U_{L,R} + \frac{1}{2} \frac{\Delta t}{\Delta x} \left[F(U_L) - F(U_R) \right]$$
 (18)

Then we use Musta–Force, using U_R in i–th cell and U_L in i+1–th cell as initial conditions.

SLIC/MUSCL scehem — choosing Δ_i

The formula is:

$$\Delta_i = \frac{1}{2}(1+\omega)(U_i - U_{i-1}) + \frac{1}{2}(1-\omega)(U_{i+1} - U_i), \quad \omega \in [-1,1]$$
(19)

To avoid oscillations we use *slope limiters* — we change $\Delta_i \rightarrow \overline{\Delta}_i = \xi(r)\Delta_i$, where $r = \frac{U_i - U_{i-1}}{U_{i+1} - U_i}$ One of choices for ξ is called MINBEE/MINMOD:

$$\xi(r) = \max[0, \min(1, r)] \tag{20}$$

Getting \vec{v} and pUsing

$$M = (E+p)v$$

we have:

$$\nu = \frac{M}{E+p} \tag{21}$$

We obtain:

$$e = E - Mv \tag{22}$$

$$n = R\sqrt{1-\nu^2} \tag{23}$$

Pressure *p* is computed using eos p = p(e, n):

$$\nu = \frac{M}{E + p(E - M\nu, R\sqrt{1 - \nu^2})}$$
 (24)

$$\vec{\nu} = \nu \frac{\vec{M}}{M} \tag{25}$$