



A well-balanced scheme for two-fluid flows in spherical coordinates

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Outline

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- 3 Numerical results
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Model

We consider the Euler equations in spherical coordinates:

$$\partial_t(A\rho) + \partial_x(A\rho u) = 0,$$

$$\partial_t(A\rho u) + \partial_x(A(\rho u^2 + p)) = p\partial_x A,$$

$$\partial_t(A\rho E) + \partial_x(A(\rho E + p)u) = 0,$$

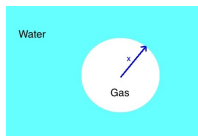
$$\partial_t(A\rho\varphi) + \partial_x(A\rho\varphi u) = 0,$$

where A is the cross-section, ρ the density, u the velocity, E the total energy, φ the fraction of mass of gas and p satisfies the stiffened gas pressure law:

$$p(\rho, e, \varphi) = (\gamma(\varphi) - 1)\rho e - \gamma(\varphi)\pi(\varphi).$$

Applications:

- Spherical bubble:



- Nozzle:



Model

- Let A_Δ a piecewise constant approximation of A :

$$A_\Delta(x) = \sum_i A_i h_{|x_{i-1/2}; x_{i+1/2}|}(x).$$

- The system becomes:

$$\begin{aligned}\partial_t A_\Delta &= 0, \\ \partial_t(A_\Delta \rho) + \partial_x(A_\Delta \rho u) &= 0, \\ \partial_t(A_\Delta \rho u) + \partial_x(A_\Delta(\rho u^2 + p)) &= p \partial_x A_\Delta, \\ \partial_t(A_\Delta \rho E) + \partial_x(A_\Delta(\rho E + p)u) &= 0, \\ \partial_t(A_\Delta \rho \varphi) + \partial_x(A_\Delta \rho \varphi u) &= 0.\end{aligned}$$

Model

We can write the system in the non-conservative form:

$$\partial_t W + \partial_x F(W) - S(W) \partial_x A_\Delta = 0,$$

where the vector of conservative variables is:

$$W = (A_\Delta, A_\Delta \rho, A_\Delta \rho u, A_\Delta \rho E, A_\Delta \rho \varphi)^T,$$

the conservative flux is:

$$F(W) = (0, A_\Delta \rho u, A_\Delta (\rho u^2 + p), A_\Delta (\rho E + p) u, A_\Delta \rho \varphi u)^T,$$

and the non-conservative source term is:

$$S \partial_x A_\Delta = (0, 0, p \partial_x A_\Delta, 0, 0).$$

Hyperbolicity

Theorem

The jacobian matrix of the non-conservative system:

$$\partial_t W + \partial_x F(W) - S(W)\partial_x A_\Delta = 0,$$

admits five real eigenvalues:

- $\lambda_0 = 0(LD)$,
- $\lambda_1 = u - c(GNL)$,
- $\lambda_2 = \lambda_3 = u(LD)$,
- $\lambda_4 = u + c(GNL)$.

Numerical scheme

- We propose a first order finite volume scheme with an Arbitrary Lagrangian Eulerian approach, the boundary $x_{i+1/2}^n$ can move at the speed of the fluid $v_{i+1/2}^n$ between t_n and t_{n+1} :

$$x_{i+1/2}^{n+1,-} = x_{i+1/2}^n + \Delta t v_{i+1/2}^n.$$

- We define an ALE numerical flux:

$$F(W_L, W_R, v^\pm) := F(W(W_L, W_R, v^\pm)) - vW(W_L, W_R, v^\pm)$$

where $W(W_L, W_R, v^\pm)$ is obtained with an approximated Riemann solver described below.

Numerical scheme

The integration of $\partial_t W + \partial_x F(W) - S(W)\partial_x A_\Delta = 0$ on $]t_n; t_{n+1}[\times]x_{i-1/2}(t); x_{i+1/2}(t)[$ gives:

- if $v_{i+1/2}^n \leq 0$ and $v_{i-1/2}^n \geq 0$, the ALE scheme is:

$$\Delta x_i^{n+1} W_i^{n+1,-} - \Delta x_i^n W_i^n + \Delta t \left(F(W_i^n, W_{i+1}^n, v_{i+1/2}^{n,-}) - F(W_{i-1}^n, W_i^n, v_{i-1/2}^{n,+}) \right) = 0.$$

- if $v_{i+1/2}^n > 0$, we have to add the following term on the left side of the equation above:

$$\Delta t \left(F(W_i^n, W_{i+1}^n, 0^-) - F(W_i^n, W_{i+1}^n, 0^+) \right),$$

- if $v_{i-1/2}^n < 0$, we add:

$$\Delta t \left(F(W_{i-1}^n, W_i^n, 0^-) - F(W_{i-1}^n, W_i^n, 0^+) \right).$$

Numerical scheme

To update $W_i^{n+1,-}$, we have to compute $F(W_L, W_R, v^\pm)$. To do that, we need to:

- precise how we choose the ALE velocity v ,
- explain how we compute $F(W_L, W_R, v^\pm)$ for a given ALE velocity v .

Computing the interface speed v

When initial datas satisfy $\varphi \in \{0, 1\}$, the algorithm verifies

- if $\varphi_L = \varphi_R$, we take $v = 0$.
- if $\varphi_L \neq \varphi_R$, we use an exact Riemann solver in $(\varrho, u, p, \varphi)^T$ to compute u^* and p^* . We take $v = u^*$ and the ALE flux takes the Lagragian following form:

$$F(W_L, W_R, v^\pm) = (0, A^* p^*, A^* u^* p^*, 0, -A^* u^*)^T,$$

where

$$A^* = \begin{cases} A_L, & \text{if } v < 0, \\ A_R, & \text{if } v > 0. \end{cases}$$

Computing $W(W_L, W_R, 0^\pm)$

Proposition

The Riemann invariants associated to the stationary wave $\lambda_0 = 0$ are:

- $\varphi,$
- $Q = \rho Au,$
- $s = (p + \pi(\varphi))\rho^{-\gamma(\varphi)},$
- $H = E + \frac{p}{\rho}.$

Considering the set of variables $Z = (A, \varphi, s, Q, H)^T$, the system $\partial_t W + \partial_x F(W) - S(W)\partial_x A_\Delta = 0$ becomes:

$$\partial_t Z + C(Z)\partial_x Z = 0.$$

VFRoe-ncv solver

We want to solve

$$\partial_t Z + C(Z)\partial_x Z = 0,$$

with the initial datas

$$Z(x, 0) = \begin{cases} Z(W_L) & , x < 0 \\ Z(W_R) & , x > 0 \end{cases}.$$

- We linearize $C(Z)$ and we solve the linear Riemann problem:

$$\partial_t Z + C(\hat{Z})\partial_x Z = 0,$$

with the same initial datas.

- We obtain $Z(W_L, W_R, 0^\pm)$.

Property of $Z(W_L, W_R, 0^\pm)$

Proposition

The interface Riemann solver computes intermediate states $Z(W_i, W_{i+1}, 0^\pm)$ which verify:

- $A_{i+1/2}^- = A_i,$
- $\varphi_{i+1/2}^- = \varphi_{i+1/2}^+,$
- $s_{i+1/2}^- = s_{i+1/2}^+,$
- $A_{i+1/2}^+ = A_{i+1},$
- $Q_{i+1/2}^- = Q_{i+1/2}^+,$
- $H_{i+1/2}^- = H_{i+1/2}^+.$

$$Z(W_L, W_R, 0^\pm) \rightarrow W(W_L, W_R, 0^\pm)$$

Problem

The map $W \mapsto Z$ is not bijective.

To obtain $\rho_{i+1/2}^-$, we need to solve:

$$L(\rho) = \frac{\gamma(\varphi_{i+1/2})}{\gamma(\varphi_{i+1/2}) - 1} s_{i+1/2} \rho^{\gamma(\varphi_{i+1/2}) - 1} + \frac{Q_{i+1/2}^2}{2\rho^2 (A_{i+1/2}^-)^2} - H_{i+1/2} = 0.$$

and we take:

- $u_{i+1/2}^- = \frac{Q_{i+1/2}}{\rho_{i+1/2}^- A_{i+1/2}^-},$
- $p_{i+1/2}^- = s_{i+1/2} (\rho_{i+1/2}^-)^{\gamma(\varphi_{i+1/2})} - \pi(\varphi_{i+1/2}).$

$\implies W(W_L, W_R, 0^\pm)$ is computed.

Entropy correction

Definition

If $\lambda_k(W_i^n) \leq 0 \leq \lambda_k(W_{i+1}^n)$ in a non linear k-field, then the numerical flux F is replaced by:

$$G(W_i^n, W_{i+1}^n, 0^\pm) = F(W_i^n, W_{i+1}^n, 0^\pm) - \min_k(|\lambda_k(W_i^n)|, |\lambda_k(W_{i+1}^n)|)(W_{i+1}^n - W_i^n)/2.$$

Random sampling remap

We go back to the original grid by the Glimm procedure. We construct a sequence of pseudo-random $\omega_n \in [0, 1[$, and we take:

$$W_i^{n+1} = \begin{cases} W_{i-1}^{n+1,-}, & \text{if } \omega_n < \frac{\Delta t^n}{\Delta x_i} \max(v_{i-1/2}^n, 0), \\ W_i^{n+1,-}, & \text{if } \frac{\Delta t^n}{\Delta x_i} \max(v_{i-1/2}^n, 0) \leq \omega_n \\ & \text{and if } \omega_n \leq 1 + \frac{\Delta t^n}{\Delta x_i} \min(v_{i+1/2}^n, 0), \\ W_{i+1}^{n+1,-}, & \text{if } \omega_n > 1 + \frac{\Delta t^n}{\Delta x_i} \min(v_{i+1/2}^n, 0). \end{cases}$$

Properties of the scheme

The constructed scheme has the following properties:

- it is well-balanced, meaning that **it preserves exactly all stationary states** (i.e. initial datas for which the quantities φ, s, Q, H are constant);
- if the fraction of gas φ takes only the two values 0 or 1, this property is exactly preserved at any time.

Stationary contact

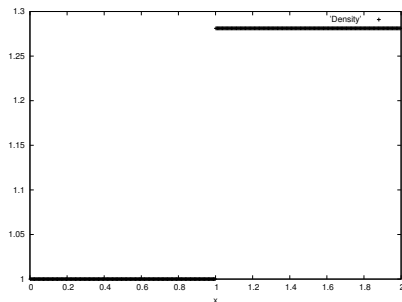
We choose the initial datas:

Quantities	L.	Right
$\rho(\text{kg.m}^{-3})$	1	1.281128576
$u(\text{m.s}^{-1})$	1	0.5203745191
$p(\text{Pa})$	1	1.414587466
φ	1	1
A	1	1.5

such that:

- $\varphi_L = \varphi_R,$
- $Q_L = Q_R,$
- $S_L = S_R,$
- $H_L = H_R.$

We obtain at $t = 0.2s$:

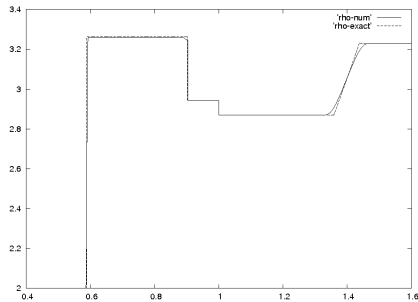


Comparison with an exact solution

The initial datas are:

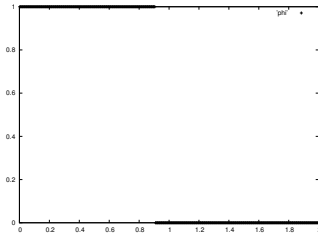
Quantities	L.	Right
$\rho(kg.m^{-3})$	2	3.230672602
$u(m.s^{-1})$	0.5	-0.4442565900
$p(Pa)$	1	12
φ	1	0
A	1.5	1
γ	1.4	1.6
$\pi(Pa)$	0	2

We obtain at $t = 0.2s$:

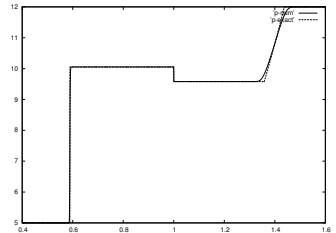


Comparison with an exact solution

- There is no diffusion on φ :



- There is no jump on velocity and pressure:



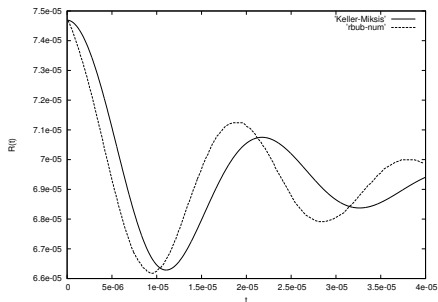
Implosion of a bubble

We simulate the collapse of a spherical bubble of gas in liquid water. Initial datas are:

Quantities	L.	Right
$\rho(kg.m^{-3})$	0.92	1000
$u(m.s^{-1})$	0	0
$p(Pa)$	72567.68	10^5
φ	1	0
A	1.5	1
γ	1.4	3
$\pi(Pa)$	0	733333.33

with $A(x) = 4\pi x^2$.

We plot the radius of the bubble as compared to the EDO model of Keller-Miksis, we obtain:



Conclusion

- We have constructed a new well-balanced scheme which does not diffuse the interface.
- We will pass this algorithm on GPU.
- We will test our algorithm for flow in a nozzle in one or two dimensions.
- We will compare this algorithm with other algorithms.

Thank you for your attention!

