Numerical Approximation of a two-fluid two-pressure diphasic model

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We are interested in the computation of two-phase flows, **using a two-fluid approach.**

Generally speaking, two-fluid models have several drawbacks :

- they are generally non conservative
- they may not be hyperbolic in a large class of conditions
- they may not satify the maximum principle property on the void fraction α

Today, we will focus on the first point only.

The governing equations are the following :

$$\begin{cases} \partial_t \alpha_1 + u_l \partial_x \alpha_1 = 0\\ \partial_t \alpha_1 \rho_1 + \partial_x \alpha_1 \rho_1 u_1 = 0\\ \partial_t (\alpha_1 \rho_1 u_1) + \partial_x (\alpha_1 \rho_1 u_1^2 + \alpha_1 p_1(\rho_1)) - p_l \partial_x \alpha_1 = 0\\ \partial_t \alpha_2 \rho_2 + \partial_x \alpha_2 \rho_2 u_2 = 0\\ \partial_t (\alpha_2 \rho_2 u_2) + \partial_x (\alpha_2 \rho_2 u_2^2 + \alpha_2 p_2(\rho_2)) + p_l \partial_x \alpha_1 = 0 \end{cases}$$

Note first that this model is barotropic and focuses on convective effects only.

It is unconditionally hyperbolic (not strictly) with real eigenvalues given by

$$\lambda_0 = u_1, \lambda_1 = u_1 - c_1, \ \lambda_2 = u_1 + c_1, \lambda_3 = u_2 - c_2, \ \lambda_4 = u_2 + c_2$$

The last four characteristic fields are GNL.

The first characteristic field is LD under the condition

$$u_I = u_2, \quad p_I = p_1 \quad (\text{or} \quad u_I = u_1, \quad p_I = p_2)$$

To summarize, the model is

- non conservative
- hyperbolic
- associated with a pure transport of the void fraction α_1

The literature is large on this subject : see for instance the works by Gallouët, Hérard and Seguin, Ransom and Hicks, Baer and Nunziato, Kapila *et al*, Gavrilyuk and Saurel, Saurel and Abgrall, ...

The Riemann problem is difficult to solve for this model because of

- the large number of waves
- the non linear pressure terms
- the non conservative products



FIG.: General structure of a Riemann problem

Except of course if $\partial_x \alpha_1 = 0$ since we get two classical barotropic Euler systems in this case...

A RELAXATION APPROACH (JOINT WITH A. AMBROSO, T. GALLÉ AND F. COQUEL) A « SHARP INTERFACE » APPROACH WITH RANDOM SAMPLING (STILL IN PROGRESS) NUMERICAL ILLUSTRATIONS

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A RELAXATION APPROACH (JOINT WITH A. AMBROSO, T. GALIÉ AND F. COQUEL)



Introducing the condensed form

$$\partial_t \mathbf{u} + \partial_x \mathbf{f}(\mathbf{u}) + \mathbf{c}(\mathbf{u})\partial_x \mathbf{u} = 0, \ t > 0, \ x \in \mathbb{R},$$

the general idea is to propose a larger but simpler relaxation model with source term

$$\partial_t \mathbf{v} + \partial_x \mathbf{g}(\mathbf{v}) + \mathbf{d}(\mathbf{v}) \partial_x \mathbf{v} = \lambda \mathcal{R}(\mathbf{v}), \ t > 0, \ x \in \mathbb{R}.$$

This system is expected to be such that

$$\lim_{\lambda\to\infty}\mathbf{v}^{\lambda}=\mathbf{u}.$$

Here, simpler means that the Riemann problem of the convective part is easier to solve.

As a consequence, the numerical strategy for solving the equilibrium system

$$\partial_t \mathbf{u} + \partial_x \mathbf{f}(\mathbf{u}) + \mathbf{c}(\mathbf{u})\partial_x \mathbf{u} = 0, \ t > 0, \ x \in \mathbb{R},$$

is based on a splitting strategy on the relaxation model :

• First step : solve by a Godunov scheme the convective part

$$\partial_t \mathbf{v} + \partial_x \mathbf{g}(\mathbf{v}) + \mathbf{d}(\mathbf{v})\partial_x \mathbf{v} = 0$$

• Second step : solve in the regime $\lambda \to \infty$

$$\partial_t \mathbf{v} = \lambda \mathcal{R}(\mathbf{v})$$

that is, impose on each cell the relation

$$\mathcal{R}(\mathbf{v}) = 0$$

How to designe the relaxation model

$$\partial_t \mathbf{v} + \partial_x \mathbf{g}(\mathbf{v}) + \mathbf{d}(\mathbf{v}) \partial_x \mathbf{v} = \lambda \mathcal{R}(\mathbf{v}), \ t > 0, \ x \in \mathbb{R}.$$

such that

- $\lim_{\lambda\to\infty} \mathbf{v}^{\lambda} = \mathbf{u}$
- the model is larger but simpler to solve

Recall that the Riemann problem is difficult to solve for this model because of

- the large number of waves
- the non linear pressure terms
- the non conservative products

Here, our relaxation process will concern the last two points only.

We first focus on the non linear pressure terms.

We propose the following relaxation system :

$$\begin{split} &\partial_{l}\alpha_{1} + u_{2}\partial_{x}\alpha_{1} = 0 \\ &\partial_{t}\alpha_{1}\rho_{1} + \partial_{x}\alpha_{1}\rho_{1}u_{1} = 0 \\ &\partial_{t}(\alpha_{1}\rho_{1}u_{1}) + \partial_{x}(\alpha_{1}\rho_{1}u_{1}^{2} + \alpha_{1}\Pi_{1}) - \Pi_{1}\partial_{x}\alpha_{1} = 0 \\ &\partial_{t}\alpha_{2}\rho_{2} + \partial_{x}\alpha_{2}\rho_{2}u_{2} = 0 \\ &\partial_{t}(\alpha_{2}\rho_{2}u_{2}) + \partial_{x}(\alpha_{2}\rho_{2}u_{2}^{2} + \alpha_{2}\Pi_{2}) + \Pi_{1}\partial_{x}\alpha_{1} = 0 \\ &\partial_{t}\alpha_{1}\rho_{1}\Pi_{1} + \partial_{x}\alpha_{1}\rho_{1}\Pi_{1}u_{1} + \alpha_{1}a_{1}^{2}\partial_{x}u_{1} - a_{1}^{2}(u_{1} - u_{I})\partial_{x}\alpha_{1} = \lambda\alpha_{1}\rho_{1}(p_{1} - \Pi_{1}), \\ &\partial_{t}\alpha_{2}\rho_{2}\Pi_{2} + \partial_{x}\alpha_{2}\rho_{2}\Pi_{2}u_{2} + \alpha_{2}a_{2}^{2}\partial_{x}u_{2} - a_{2}^{2}(u_{2} - u_{I})\partial_{x}\alpha_{2} = \lambda\alpha_{2}\rho_{2}(p_{2} - \Pi_{2}), \end{split}$$

avec $a_k > \rho_k c_k$, k = 1, 2. (the so-called Whitham conditions).

This is unconditionally hyperbolic (not strictly) with real eigenvalues given by

$$\begin{split} \lambda_0^r &= u_2, \\ \lambda_1^r &= u_1 - a_1 \tau_1, \ \lambda_2^r &= u_1 + a_1 \tau_1, \\ \lambda_3^r &= u_2 - a_2 \tau_2, \ \lambda_4^r &= u_2 + a_2 \tau_2. \end{split}$$

All the fields are LD.

We then focus on the non conservative products.

We propose the following modified relaxation system :

$$\begin{array}{l} \partial_t \alpha_1 + u_2 \partial_x \alpha_1 = 0 \\ \partial_t \alpha_1 \rho_1 + \partial_x \alpha_1 \rho_1 u_1 = 0 \\ \partial_t (\alpha_1 \rho_1 u_1) + \partial_x (\alpha_1 \rho_1 u_1^2 + \alpha_1 \Pi_1) = \overline{\Pi_1 \partial_x \alpha_1} \delta_{x-u_2^* t} \\ \partial_t \alpha_2 \rho_2 + \partial_x \alpha_2 \rho_2 u_2 = 0 \\ \partial_t (\alpha_2 \rho_2 u_2) + \partial_x (\alpha_2 \rho_2 u_2^2 + \alpha_2 \Pi_2) = -\overline{\Pi_1 \partial_x \alpha_1} \delta_{x-u_2^* t} \\ \partial_t \alpha_1 \rho_1 \Pi_1 + \partial_x \alpha_1 \rho_1 \Pi_1 u_1 + \alpha_1 a_1^2 \partial_x u_1 - a_1^2 (u_1 - u_I) \partial_x \alpha_1 = \lambda \alpha_1 \rho_1 (p_1 - \Pi_1), \\ \partial_t \alpha_2 \rho_2 \Pi_2 + \partial_x \alpha_2 \rho_2 \Pi_2 u_2 + \alpha_2 a_2^2 \partial_x u_2 - a_2^2 (u_2 - u_I) \partial_x \alpha_2 = \lambda \alpha_2 \rho_2 (p_2 - \Pi_2), \end{array}$$

avec $a_k > \rho_k c_k$, k = 1, 2. (the so-called Whitham conditions).

The Riemann problem is now explicitly solved.

How do we guess $\overline{\Pi_I \partial_x \alpha_1}$?

We chose to be exact on contact discontinuity solutions for the equilibrium system



We need to chose one of the two estimates. Do we?

This way, we did construct a Relaxation system such that

- the Riemann solution is explicitly known
- the Riemann solution is exact for isolated contact discontinuities
- the corresponding numerical method is stable and conservative for the mass of each fluid and for the total momentum

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Here, the key point is to remind that the model consists in two classical Euler systems for both phases when $\partial_x \alpha_1 = 0$.

$$\begin{cases} \partial_t \alpha_1 + u_I \partial_x \alpha_1 = 0 \\ \partial_t \alpha_1 \rho_1 + \partial_x \alpha_1 \rho_1 u_1 = 0 \\ \partial_t (\alpha_1 \rho_1 u_1) + \partial_x (\alpha_1 \rho_1 u_1^2 + \alpha_1 p_1(\rho_1)) - p_I \partial_x \alpha_1 = 0 \\ \partial_t \alpha_2 \rho_2 + \partial_x \alpha_2 \rho_2 u_2 = 0 \\ \partial_t (\alpha_2 \rho_2 u_2) + \partial_x (\alpha_2 \rho_2 u_2^2 + \alpha_2 p_2(\rho_2)) + p_I \partial_x \alpha_1 = 0 \end{cases}$$

Our objective is then to design a numerical scheme which

- is based on our « favorite » scheme for the usual Euler equations
- provides in addition sharp contact discontinuities
- obeys stability properties

Our strategy is based on a **Ghost-Fluid** approach coupled with a **Glimm random sampling** strategy.

Let us assume without restriction that α_1 is a step function :



According to the Ghost-Fluid approach, we first split the actual mesh in two α_1 -constant meshes :



We define the reconstructed states such that \mathbf{u}_j and $\tilde{\mathbf{u}}_j$ are joined by an admissible contact discontinuity.

Using a Glimm scheme, we then solve numerically $\partial_t \alpha_1 + u_{2,j+1/2} \partial_x \alpha_1 = 0$

Mesh ±
$$\begin{array}{c} u_{2,j-1/2} \\ \alpha_1 = \alpha_{1+} \\ \mu_{j-3} \\ \mathbf{u}_{j-2} \\ \mathbf{u}_{j-1} \\ \mathbf{u}_{j-1} \\ \mathbf{u}_{j} \\ \mathbf{u}_{j+1} \\ \mathbf{u}_{j+2} \end{array}$$

Being given an equidistributed random sequence $a_n \in (0, 1)$, it amounts to set :

$$\alpha_{1j}^{n+1} = \begin{cases} \alpha_{1+} & \text{if} \quad a_n \le u_{2,j-1/2} \frac{\Delta t}{\Delta x} \\ \alpha_{1-} & \text{if} \quad a_n > u_{2,j-1/2} \frac{\Delta t}{\Delta x} \end{cases}$$

Note that α_1 remains sharp.

With our « favorite » scheme, we numerically solve two classical Euler equations



and we choose on each cell j between the two meshes according to the value $(\alpha_1)_i^{n+1}$.

We immediately get the positivity properties of our « favorite » scheme. Note that the global scheme is not strictly conservative.

For a similar approach in the context of van der Waals fluids, we refer to a work by C. Rohde and C. Merkle which inspired this work (note the presence of a level-set function instead of a random sampling). A RELAXATION APPROACH (JOINT WITH A. AMBROSO, T. GALIÉ AND F. COQUEL) A « SHARP INTERFACE » APPROACH WITH RANDOM SAMPLING (STILL IN PROGRESS) NUMERICAL ILLUSTRATIONS

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 2 A pprox sharp interface pprox approach with random sampling (still in progress)



MOVING CONTACT DISCONTINUITY



FIG.: α_1 and u_2

SHOCK TUBE PROBLEM (1D)



FIG.: α_1 , u_2 and ρ_2

SHOCK TUBE PROBLEM (2D) - 1A



FIG.: Initial data and α_1

SHOCK TUBE PROBLEM (2D) - 1B



FIG.: u_2 and ρ_2

RANSOM FAUCET TEST CASE



FIG.: *u*₂

This work was motivated by a Joint Research Group LJLL-CEA Saclay on multiphase flows and coupling of multiscale models, to which the following persons take part :

LJLL : C. Chalons, F. Coquel, E. Godlewski, F. Lagoutière, N. Seguin, P.-A. Raviart, CEA Saclay : A. Ambroso (now Areva), B. Boutin, T. Galié (now in a service company),

+ some-time participation from EDF : J.-M. Hérard, O. Hurisse.

Advertising :

NTMC'09 (New Trends in Model Coupling)

workshop on coupling problems of multiscale phenomena that are of increasing interest in Applied Mathematics...