A Finite Volume Approximation for a Two-Temperature Plasma Fusion Model

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Strasbourg, November 17th, 2016

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- Framework
- A Two-Temperature Fusion Plasma Model
- **③** Finite Volume Approximation
- O Numerical Tests
- **o** Conclusions and Perspectives

1. Framework: A Fortunate Meeting of Programs

BORDEAUX: C. Berthon¹, B. Dubroca², A. S.

then E. Estibals, D. Arégba, J. Breil, S. Brull, ...

2 NICE: H. Guillard, B. Nkonga, A. S.

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Assumptions:

- Unmagnetized quasineutral totally ionized plasma
- \bullet Particles undergoing the electric field ${\bf E}$ given by the Ohm's law:

 $c_i \operatorname{grad} p_e - c_e \operatorname{grad} p_i = n_e q_e \mathbf{E}$ where: $c_e = \rho_e / \rho$ and $c_i = \rho_i / \rho$ The Model:

$$\begin{cases} \partial_t \rho + \operatorname{div}(\rho \mathbf{u}) = \mathbf{0}, \\ \partial_t(\rho \mathbf{u}) + \operatorname{div}(\rho \mathbf{u} \otimes \mathbf{u} + (p_e + p_i)\mathbf{I}) = \mathbf{0} \\ \partial_t(\rho_e \varepsilon_e + \frac{1}{2}\rho_e \mathbf{u} \cdot \mathbf{u}) + \operatorname{div}((\rho_e \varepsilon_e + \frac{1}{2}\rho_e \mathbf{u} \cdot \mathbf{u} + p_e)\mathbf{u}) \\ - (c_i \operatorname{grad} p_e - c_e \operatorname{grad} p_i) \cdot \mathbf{u} = \nu_{ei}^{\mathscr{E}} (T_i - T_e) \\ \partial_t(\rho_i \varepsilon_i + \frac{1}{2}\rho_i \mathbf{u} \cdot \mathbf{u}) + \operatorname{div}((\rho_i \varepsilon_i + \frac{1}{2}\rho_i \mathbf{u} \cdot \mathbf{u} + p_i)\mathbf{u}) \\ + (c_i \operatorname{grad} p_e - c_e \operatorname{grad} p_i) \cdot \mathbf{u} = -\nu_{ei}^{\mathscr{E}} (T_i - T_e) \end{cases}$$

The mathematical properties of the model

- The model is non-conservative
- The seminal work of *Coquel* and *Marmignon* ³ make it conservative under: assumption of **null jump of entropy across shocks**
- \bullet The same transformation was recently rederived in the internship of $\textit{Estibals}^4$

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F. COQUEL, C. MARMIGNON, Astro. Space Sc., 260, 1-2, 15-27 (1998)

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D. ARÉGBA, J. BREIL, S. BRULL, B. DUBROCA, E. ESTIBALS, submitted for publication

The mathematical properties of the model

• The following conservative system is then obtained:

$$\begin{cases} \partial_t \rho + \operatorname{div}(\rho \mathbf{u}) &= 0\\ \partial_t(\rho \mathbf{u}) + \operatorname{div}(\rho \mathbf{u} \otimes \mathbf{u} + (p_e + p_i)\mathbf{I}) &= 0\\ \partial_t(\rho \mathscr{E}) + \operatorname{div}((\rho \mathscr{E} + p_i + p_e)\mathbf{u}) &= 0\\ \partial_t(\rho_e s_e) + \operatorname{div}(\rho_e s_e \mathbf{u}) &= \nu_{ei}^{\mathscr{E}} \rho_e^{1-\gamma_e}(T_i - T_e) \end{cases}$$

where: $\rho \mathscr{E} = \rho_i \varepsilon_i + \frac{1}{2} \rho_i \mathbf{u} \cdot \mathbf{u} + \rho_e \varepsilon_e + \frac{1}{2} \rho_e \mathbf{u} \cdot \mathbf{u}$ is the total energy, $s_e = p_e \rho_e^{-\gamma_e}$ is the electron's entropy

The mathematical properties of the model

• The following conservative system is then obtained:

$$\begin{cases} \partial_t \rho + \operatorname{div}(\rho \mathbf{u}) &= 0\\ \partial_t(\rho \mathbf{u}) + \operatorname{div}(\rho \mathbf{u} \otimes \mathbf{u} + (p_e + p_i)\mathbf{l}) &= 0\\ \partial_t(\rho \mathscr{E}) + \operatorname{div}((\rho \mathscr{E} + p_i + p_e)\mathbf{u}) &= 0\\ \partial_t(\rho s_e) + \operatorname{div}(\rho s_e \mathbf{u}) &= \nu_{ei}^{\mathscr{E}} c_e^{-\gamma e} \rho^{1-\gamma_e}(T_i - T_e) \end{cases}$$

• This system can be written in the following compact form:

$$\partial_t \mathscr{U} + \operatorname{div} \mathscr{F}(\mathscr{U}) = \mathscr{S}(\mathscr{U})$$

with:

$$\mathscr{U} = \begin{pmatrix} \rho \\ \rho \mathbf{u} \\ \rho \mathscr{E} \\ \rho \mathscr{S}_{e} \end{pmatrix} \quad \mathscr{F}(\mathscr{U}) = \begin{pmatrix} \rho \mathbf{u} \\ \rho \mathbf{u} \otimes \mathbf{u} + (p_{i} + p_{e}) \mathbf{I} \\ (\rho \mathscr{E} + p_{i} + p_{e}) \mathbf{u} \\ \rho \mathscr{S}_{e} \mathbf{u} \end{pmatrix} \quad \mathscr{I}(\mathscr{U}) = \begin{pmatrix} 0 \\ \mathbf{0} \\ 0 \\ \mathscr{S}_{e} \end{pmatrix}$$
$$\mathscr{S}_{e} = \nu_{e_{i}}^{\mathscr{E}} c_{e}^{-\gamma_{e}} \rho^{1-\gamma_{e}}(T_{i} - T_{e})$$

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The mathematical properties of the model

• The solution \mathscr{U} of the model belongs to the set of *physically admissible states* \mathscr{O} defined by:

$$\mathscr{O} = \left\{ \mathscr{U} = \left(\rho, \rho \mathbf{u}, \rho \mathscr{E}, \rho s_{e}\right)^{\mathsf{T}} \in \mathbb{R}^{6}, \quad \rho > 0, \quad \mathscr{E} - \frac{1}{2}\mathbf{u} \cdot \mathbf{u} > 0, \quad s_{e} > 0 \right\}$$

• A useful Lemma:

Lemma

Let $\mathscr{U} = (\rho, \rho \mathbf{u}, \rho \mathscr{E}, \rho s_e)^T$ be a solution of the model. Then the following systems are equivalent:

$$\begin{cases} \partial_t \rho = \partial_t \mathbf{u} = \partial_t \mathscr{E} = 0 \\ \partial_t s_e = \nu_{ei}^{\mathscr{E}} c_e^{-\gamma_e} \rho^{-\gamma_e} (T_i - T_e) \\ \\ \partial_t T_e = \nu_{ei}^{\mathscr{E}} (T_i - T_e) \\ \partial_t T_i = -\nu_{ei}^{\mathscr{E}} (T_i - T_e) \end{cases}$$

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The mathematical properties of the model

- The model is Galilean invariant
- Hyperbolicity

Theorem

The 1D version of the model without the source term \mathscr{S}_e is hyperbolic. The eigenvalues are given by the set

$$\Xi = \{u - c_{ei}, u, u, u, u + c_{ei}\}$$

where $c_{ei} = \sqrt{rac{\partial(p_i + p_e)}{\partial
ho}}.$

The characteristic fields associated to the eigenvalues $u \pm c_{ei}$ are genuinely nonlinear while the characteristic fields associated to the eigenvalue u are linearly degenerated.

A numerical strategy to approximate the model based on our work reported in:

- A. BONNEMENT, et al., *ESAIM Proceedings*, **32**, 163-176 (2011)
- M. BILANCERI, et al., *ESAIM Proceedings*, **43**, 164-179 (2013)
- Toroidal geometry and cylindrical coordinates for a torus
- Tessellation:

 \Box unstructured mesh composed of triangles in polar planes

 \Box structured mesh in the toroidal direction

- *i.e.* curved primastic elements in the toroidal direction
- Our finite volume approximation in curvilinear coordinates:

 The divergence of the momentum equation is kept in local cylindrical coordinates

Our finite volume approximation in curvilinear coordinates:

 The divergence of the momentum equation is kept in local cylindrical coordinates

 □ integration of this divergence form on control cells
 □ definition of adequate discrete cylindrical base and projection of the result of the integration step on this base

This yields the following generic FV approximation:

The details:

• Toroidal geometry and cylindrical coordinates

$$\begin{cases} x = R \cos \varphi \\ y = R \sin \varphi \\ z = Z \end{cases}$$

 $(\mathbf{e}_R, \, \mathbf{e}_arphi, \, \mathbf{e}_Z)$ is the cylindrical local base

• The conservative form of the model in cylindrical coordinates is given by:

$$\begin{cases} \partial_t(R\rho) + \partial_{\xi_k}(R\rho \mathbf{u} \cdot \mathbf{e}^k) &= 0\\ \partial_t(R\rho \mathbf{u}) + \partial_{\xi_k}(R\mathbf{T} \cdot \mathbf{e}^k) &= 0\\ \partial_t(R\rho \mathscr{E}) + \partial_{\xi_k}(R(\rho \mathscr{E} + p_i + p_e)\mathbf{u} \cdot \mathbf{e}^k) &= 0\\ \partial_t(R\rho s_e) + \partial_{\xi_k}(R\rho s_e \mathbf{u} \cdot \mathbf{e}^k) &= R \mathscr{S}_e \end{cases}$$

where: $e_k \in \{\mathbf{e}_R, \mathbf{e}_{\varphi}, \mathbf{e}_Z\}$; $\mathbf{e}_k \cdot \mathbf{e}^l = \delta_k^l$; $\mathbf{T} = (\rho u_k u_l + (p_i + p_e)\delta_k^l) \mathbf{e}_k \otimes \mathbf{e}_l$ $\mathscr{S}_e = \nu_{ei}^{\mathscr{E}} c_e^{-\gamma_e} \rho^{1-\gamma_e}(T_i - T_e)$

• Tessellation:

 \Box triangles T_{β} in (R, Z)-coordinates to mesh polar planes \Box interval of angles $(\varphi_k, \varphi_{k+1})$, where $k \in \{1, \dots, N_{plan}\}$ to mesh the computational domain in the toroidal direction

This leads to curved primastic elements in the toroidal direction to partition the computational domain

 \Box INRIA (R, Z)-coordinates 2D control cells Ω_{α} leading to the 3D control cells Ω_{α}^{3D} associated to each node α of the mesh of the computational domain

 \Box The boundary of each control cell Ω^{3D}_{α} is composed of poloidal surfaces and toroidal surfaces

The details:

• Our finite volume approximation in curvilinear coordinates:

◊ scalar equations: continuity, energy, entropy

vectorial equation: momentum

 $\hfill\square$ Our procedure applied to scalar equations is same as the well-known FV scheme

□ For vectorial equations, our strategy proceeds as follows: \diamond Integration of momentum equation over the control cell Ω^{3D}_{α} :

$$|\Omega_{\alpha}^{3D}| \ \partial_t \left(\frac{1}{|\Omega_{\alpha}^{3D}|} \int_{\Omega_{\alpha}^{3D}} R\rho \mathbf{u} \, d\Omega \right) + \int_{\Omega_{\alpha}^{3D}} \partial_{\xi_k} (R\mathbf{T} \cdot \mathbf{e}^k) \, d\Omega = 0$$

 $\diamond \mbox{ Crucial choice of components of the vector } \\ \frac{1}{|\Omega_{\alpha}^{3D}|} \int_{\Omega_{\alpha}^{3D}} R\rho \mathbf{u} \, d\Omega \mbox{ to be stored in order to represent it } \\ \hline \end{tabular}$

The details:

- Our finite volume approximation in curvilinear coordinates:
 - $\hfill\square$ For vectorial equations, our strategy proceeds as follows:
 - Crucial choice of components of the vector

 $\frac{1}{|\Omega_{\alpha}^{3D}|}\int_{\Omega_{\alpha}^{3D}}R\rho\mathbf{u}\,d\Omega\text{ to be stored in order to represent it }$

♦ We chose to store the components of the vector \mathbf{u}_{α} with respect to the local basis ($\mathbf{e}_{R}(\alpha)$, $\mathbf{e}_{Z}(\alpha)$, $\mathbf{e}_{\varphi}(\alpha)$) of the control cell Ω_{α}^{3D}

♦ This automatically leads to:

$$\frac{1}{|\Omega_{\alpha}^{3D}|} \int_{\Omega_{\alpha}^{3D}} R\rho \mathbf{u} \, d\Omega = \rho_{\alpha} \left(\eta_{\alpha} u_{R,\alpha} \mathbf{e}_{R}(\alpha) + u_{Z,\alpha} \mathbf{e}_{Z}(\alpha) + \eta_{\alpha} u_{\varphi,\alpha} \mathbf{e}_{\varphi}(\alpha) \right)$$
with:

$$\eta_{\alpha} = \frac{\sin\left(\frac{\varphi_{\alpha+1/2} - \varphi_{\alpha-1/2}}{2}\right)}{\frac{\varphi_{\alpha+1/2} - \varphi_{\alpha-1/2}}{2}}, \, \mathbf{u}_{\alpha} = u_{R,\alpha} \mathbf{e}_{R}(\alpha) + u_{Z,\alpha} \mathbf{e}_{Z}(\alpha) + u_{\varphi,\alpha} \mathbf{e}_{\varphi}(\alpha)$$

$$= \frac{\varphi_{\alpha+1/2} - \varphi_{\alpha-1/2}}{2}$$

The details:

$$\begin{split} |\Omega_{\alpha}^{3D}| \partial_{t}(\rho_{\alpha}\eta_{\alpha}\mathbf{u}_{\alpha}) + \int_{\Omega_{\alpha}^{3D}} \partial_{\xi_{k}}(R\mathbf{T}\cdot\mathbf{e}^{k}) d\Omega &= 0 \\ \text{where: } \eta_{\alpha}\mathbf{u}_{\alpha} &= \eta_{\alpha}u_{R,\alpha}\mathbf{e}_{R}(\alpha) + u_{Z,\alpha}\mathbf{e}_{Z}(\alpha) + \eta_{\alpha}u_{\varphi,\alpha}\mathbf{e}_{\varphi}(\alpha) \\ \bullet \text{ This yields the following generic FV approximation:} \\ \left|\Omega_{\alpha}^{3D}\right| \partial_{t} \begin{pmatrix} \rho_{\alpha}\\ \rho_{\alpha}\eta_{\alpha}\mathbf{u}_{\alpha}\\ \rho_{\alpha}\mathscr{E}_{\alpha}\\ \rho_{\alpha}\mathscr{E}_{\alpha} \end{pmatrix} + \sum_{S_{\alpha\beta}\in\mathscr{S}^{pol}} \int_{S_{\alpha\beta}} \mathbf{F}\left(\mathscr{U}_{\alpha},\mathscr{U}_{\beta},\mathbf{n}_{\alpha\beta}\right) d\partial\Omega \\ &+ \sum_{S_{\alpha\beta}\in\mathscr{S}^{tor}} \int_{S_{\alpha\beta}} \mathbf{F}\left(\mathscr{U}_{\alpha},\mathscr{U}_{\beta},\mathbf{n}_{\alpha\beta}\right) d\partial\Omega = \int_{\Omega_{\alpha}^{3D}} \begin{pmatrix} 0\\ 0\\ 0\\ R\mathscr{I}_{\alpha} \end{pmatrix} d\Omega \\ &= 0 \\ 16/4 \end{split}$$

The details:

• This yields the following generic FV approximation:

$$\begin{split} \left|\Omega_{\alpha}^{3D}\right| \partial_{t} \begin{pmatrix} \rho_{\alpha} \eta_{\alpha} \mathbf{u}_{\alpha} \\ \rho_{\alpha} \mathcal{E}_{\alpha} \\ \rho_{\alpha} \mathcal{S}_{e\alpha} \end{pmatrix} + \sum_{S_{\alpha\beta} \in \mathscr{S}^{pol}} \int_{S_{\alpha\beta}} \mathbf{F} \left(\mathscr{U}_{\alpha}, \mathscr{U}_{\beta}, \mathbf{n}_{\alpha\beta}\right) \, d\partial\Omega \\ &+ \sum_{S_{\alpha\beta} \in \mathscr{S}^{tor}} \int_{S_{\alpha\beta}} \mathbf{F} \left(\mathscr{U}_{\alpha}, \mathscr{U}_{\beta}, \mathbf{n}_{\alpha\beta}\right) \, d\partial\Omega = \int_{\Omega_{\alpha}^{3D}} \begin{pmatrix} \mathbf{0} \\ \mathbf{0} \\ \mathbf{0} \\ R\mathscr{S}_{e} \end{pmatrix} \, d\Omega \end{split}$$

where: \mathscr{S}^{pol} are poloidal boundaries of Ω^{3D}_{α} \mathscr{S}^{tor} are toroidal ones

• The fluxes $\int_{S_{\alpha\beta}} F(\mathscr{U}_{\alpha}, \mathscr{U}_{\beta}, \mathbf{n}_{\alpha\beta}) d\partial\Omega$ are computed with a new relaxation scheme

The details: • The fluxes $\int_{S_{\alpha\beta}} \mathbf{F}(\mathscr{U}_{\alpha}, \mathscr{U}_{\beta}, \mathbf{n}_{\alpha\beta}) d\partial\Omega$ are computed with a new relaxation scheme based on C. Berthon, B. Dubroca, A. S., *SINUM*, **50**, 468-491 (2012) C. Berthon, B. Dubroca, A. S., *CMS*, **13**, 2119-2154 (2015) and derived in D. Arégba, J. BREIL, S. BRULL, B. DUBROCA, E. ESTIBALS, submitted for publication

 \Box The 1D model without source term

$$\begin{cases} \partial_t \rho + \partial_x (\rho u) &= 0\\ \partial_t (\rho u) + \partial_x (\rho u^2 + (p_e + p_i)) &= 0\\ \partial_t (\rho v) + \partial_x (\rho v u) &= 0\\ \partial_t (\rho w) + \partial_x (\rho w u) &= 0\\ \partial_t (\rho \mathscr{E}) + \partial_x ((\rho \mathscr{E} + p_i + p_e)u) &= 0\\ \partial_t (\rho s_e) + \partial_x (\rho s_e u) &= 0 \end{cases}$$

The details:

- The fluxes $\int_{S_{\alpha\beta}} \mathbf{F}(\mathscr{U}_{\alpha}, \mathscr{U}_{\beta}, \mathbf{n}_{\alpha\beta}) d\partial\Omega$ are computed with a
- new relaxation scheme

 $\hfill\square$ A relaxation approximation of the 1D model without source term

$$\begin{cases}
\frac{\partial_t \rho + \partial_x (\rho u)}{\partial_t (\rho u) + \partial_x (\rho u^2 + (\pi_e + \pi_i))} &= 0 \\
\frac{\partial_t (\rho u) + \partial_x (\rho u v)}{\partial_t (\rho v) + \partial_x (\rho u v)} &= 0 \\
\frac{\partial_t (\rho w) + \partial_x (\rho u w)}{\partial_t (\rho \mathcal{E}) + \partial_x ((\rho \mathcal{E} + \pi_i + \pi_e)u)} &= 0 \\
\frac{\partial_t (\rho \mathcal{E}) + \partial_x (\rho \mathcal{E}eu)}{\partial_t (\rho \mathcal{E}_e) + \partial_x (\rho \pi_e u + c_e a^2 u)} &= \frac{1}{\tau} \rho(p_e - \pi_e) \\
\frac{\partial_t (\rho \pi_i + c_i a^2) + \partial_x (\rho \pi_i u + c_i a^2 u)}{\partial_t (\rho a) + \partial_x (\rho a u)} &= \frac{1}{\tau} \rho(p_i - \pi_i) \\
\frac{\partial_t (\rho a) + \partial_x (\rho a u)}{\partial_t (\rho a)} &= 0
\end{cases}$$

The details:

• The fluxes $\int_{S_{\alpha\beta}} \mathbf{F}(\mathscr{U}_{\alpha}, \mathscr{U}_{\beta}, \mathbf{n}_{\alpha\beta}) \ d\partial\Omega$ are computed with a

new relaxation scheme

 $\hfill\square$ A relaxation approximation of the 1D model without source term

$$\circ \tau \to 0 \Longrightarrow \pi_e \to p_e, \quad \pi_i \to p_i$$

◊ The relaxation system for 1D model turns into:

$$\partial_t \mathbb{U} + \partial_x \mathbb{F}(\mathbb{U}) = rac{1}{ au} \mathbb{T}(\mathbb{U})$$

Theorem

The 1D relaxation system $\partial_t \mathbb{U} + \partial_x \mathbb{F}(\mathbb{U}) = 0$ is hyperbolic. The eigenvalues are given by: $\Lambda = \{u - a/\rho, u, u, u, u, u, u + a/\rho\}$. All the associated characteristic fields are linearly degenerated.

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♦ The Riemann problem can be solved exactely:

Lemma

Assume \mathbb{U}_I and \mathbb{U}_r are constant states and consider

$$\mathbb{U}_0(x) = \left\{ \begin{array}{ll} \mathbb{U}_l & \text{if } x < 0 \\ \mathbb{U}_r & \text{if } x > 0 \end{array} \right.$$

as the initial data for the system $\partial_t \mathbb{U} + \partial_x \mathbb{F}(\mathbb{U}) = 0$. Let a_l and a_r be positive real numbers $a_l > 0$, $a_r > 0$, satisfying:

$$\begin{split} b_{l} &= u_{l} - a_{l}/\rho_{l} < u^{*} < u_{r} + a_{r}/\rho_{r} = b_{r} \\ \text{where: } u^{*} &= \frac{a_{l}u_{l} + a_{r}u_{r}}{a_{l} + a_{r}} - \frac{(\pi_{i,r} + \pi_{e,r}) - (\pi_{i,l} + \pi_{e,l})}{a_{l} + a_{r}} \\ \text{Then the weak solution of system } \partial_{t}\mathbb{U} + \partial_{x}\mathbb{F}(\mathbb{U}) = 0 \text{ with the} \\ \text{initial data } (\mathbb{U}_{l}, \mathbb{U}_{r}) \text{ is given by} \end{split}$$

$$\mathbb{U}_{\mathcal{R}}\left(x/t,\mathbb{U}_{I},\mathbb{U}_{r}\right) = \begin{cases} \mathbb{U}_{I}, & \text{if } b_{I} > x/t \\ \mathbb{U}_{I}^{\star}, & \text{if } b_{I} \leq x/t \leq u^{\star} \\ \mathbb{U}_{r}^{\star}, & \text{if } u^{\star} \leq x/t \leq b_{r} \\ \mathbb{U}_{r}, & \text{if } b_{r} < x/t \end{cases}$$

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♦ The Riemann problem can be solved exactely: Then the star intermediate states U_I^* and U_r^* are given by

$$\mathbb{U}_{g}^{\star} = \begin{pmatrix} \rho_{g}^{\star} & \\ \rho_{g}^{\star} u^{\star} & \\ \rho_{g}^{\star} v^{\star} & \\ \rho_{g}^{\star} v^{\star} & \\ \frac{1}{2} \rho_{g}^{\star} (u^{\star})^{2} + \rho_{g}^{\star} \varepsilon_{i,g}^{\star} + \rho_{g}^{\star} \varepsilon_{e,g}^{\star} \\ \rho_{g}^{\star} s_{e,g}^{\star} & \\ \rho_{g}^{\star} \pi_{i,g}^{\star} + c_{i} (a_{g}^{\star})^{2} \\ \rho_{g}^{\star} \pi_{e,g}^{\star} + c_{e} (a_{g}^{\star})^{2} \\ \rho_{g}^{\star} \pi_{e,g}^{\star} & \\ \rho_{g}^{\star} a_{g}^{\star} & \end{pmatrix}$$

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• Practical implementation of our approximation

At time t^n , we consider a piecewise constant approximation of the solution of the initial model given by,

$$\mathscr{U}^{\Delta}(\mathsf{R},\mathsf{Z},arphi,\ t^{n})=\mathscr{U}^{n}_{lpha},\ (\mathsf{R},\mathsf{Z},arphi)\in\Omega^{3D}_{lpha}$$

where

$$\mathscr{U}_{\alpha}^{n} = \begin{pmatrix} \rho_{\alpha}^{n} & \\ \rho_{\alpha}^{n} \mathbf{u}_{\alpha}^{n} & \\ \frac{1}{2} \rho_{\alpha}^{n} \mathbf{u}_{\alpha}^{n} \cdot \mathbf{u}_{\alpha}^{n} + \rho_{\alpha}^{n} \varepsilon_{i,\alpha}^{n} + \rho_{\alpha}^{n} \varepsilon_{e,\alpha}^{n} \\ \rho_{\alpha}^{n} \mathbf{s}_{e,\alpha}^{n} & \end{pmatrix}$$

with
$$\rho_{\alpha}^{n} \varepsilon_{i,\alpha}^{n} = \frac{p_{i,\alpha}^{n}}{\gamma_{i}-1}$$
, $\rho_{\alpha}^{n} \varepsilon_{e,\alpha}^{n} = \frac{p_{e,\alpha}^{n}}{\gamma_{e}-1}$, $s_{e,\alpha}^{n} = p_{e,\alpha}^{n} (c_{e} \rho_{\alpha}^{n})^{-\gamma_{e}}$

To evolve in time this approximation, we proceed in two steps:

To evolve in time this approximation, we proceed in two steps:

 \Box *First step: Evolution step.* We set the relaxation state as:

$$\mathbb{U}_{\alpha}^{n} = \begin{pmatrix} \rho_{\alpha}^{n} \\ \rho_{\alpha}^{n} \mathbf{u}_{\alpha}^{n} \\ \frac{1}{2} \rho_{\alpha}^{n} \mathbf{u}_{\alpha}^{n} \cdot \mathbf{u}_{\alpha}^{n} + \rho_{\alpha}^{n} \varepsilon_{i,\alpha}^{n} + \rho_{\alpha}^{n} \varepsilon_{e,\alpha}^{n} \\ \rho_{\alpha}^{n} s_{e,\alpha}^{n} \\ \rho_{\alpha}^{n} p_{i,\alpha}^{n} + c_{i} (a_{\alpha}^{n})^{2} \\ \rho_{\alpha}^{n} p_{e,\alpha}^{n} + c_{e} (a_{\alpha}^{n})^{2} \\ \rho_{\alpha}^{n} a_{\alpha}^{n} \end{pmatrix}$$

where the positive real numbers a^n_{α} satisfy a subcharacteristic condition

To evolve in time this approximation, we proceed in two steps:

□ First step: Evolution step.

 \diamond The relaxation scheme leads to the updated relaxation state $\widetilde{\mathbb{U}}_{\alpha}^{n}$ along the normal $\mathbf{n}_{\alpha\beta}$ of the interface $\mathcal{S}_{\alpha\beta}$

 \diamond The fluxe **F** ($\mathscr{U}_{\alpha}, \mathscr{U}_{\beta}, \mathbf{n}_{\alpha\beta}$) is reconstructed as:

 $\mathsf{F}\left(\mathscr{U}_{\alpha},\mathscr{U}_{\beta},\mathsf{n}_{\alpha\beta}\right)=\mathscr{F}\left(\mathscr{N}\widetilde{\mathbb{U}}_{\alpha}^{n}\right)$

where: \mathscr{F} is the exact physical flux,

 $\ensuremath{\mathcal{N}}$ is the projection operator of relaxation state space to real state space.

To evolve in time this approximation, we proceed in two steps:

□ First step: Evolution step.

 \diamond The discrete scheme

$$\begin{pmatrix} \widehat{\rho}_{\alpha}^{n+1} \\ \widehat{\rho}_{\alpha}^{n+1} \widehat{\eta}_{\alpha} \widehat{\mathbf{u}}_{\alpha}^{n+1} \\ \widehat{\rho}_{\alpha}^{n+1} \widehat{\mathcal{E}}_{\alpha}^{n+1} \\ \widehat{\rho}_{\alpha}^{n+1} \widehat{s}_{e_{\alpha}}^{n+1} \end{pmatrix} = \begin{pmatrix} \rho_{\alpha}^{n} \\ \rho_{\alpha}^{n} \eta_{\alpha} \mathbf{u}_{\alpha}^{n} \\ \rho_{\alpha}^{n} \mathcal{E}_{\alpha}^{n} \\ \rho_{\alpha}^{n} \mathbf{s}_{e_{\alpha}}^{n} \end{pmatrix} - \frac{\Delta t}{|\Omega_{\alpha}^{3D}|} \sum_{S_{\alpha\beta} \in \mathscr{S}^{pol} \cup \mathscr{S}^{tor}} |S_{\alpha\beta}| \mathscr{F} \left(\mathscr{N} \widetilde{\mathbb{U}}_{\alpha}^{n} \right)$$

solves, at time t^{n+1} , with the initial data \mathscr{U}_{α}^{n} , the system:

$$\begin{cases} \partial_t(R\rho) + \partial_{\xi_k}(R\rho \mathbf{u} \cdot \mathbf{e}^k) = 0\\ \partial_t(R\rho \mathbf{u}) + \partial_{\xi_k}(R\mathbf{T} \cdot \mathbf{e}^k) = 0\\ \partial_t(R\rho \mathscr{E}) + \partial_{\xi_k}(R(\rho \mathscr{E} + p_i + p_e)\mathbf{u} \cdot \mathbf{e}^k) = 0\\ \partial_t(R\rho s_e) + \partial_{\xi_k}(R\rho s_e \mathbf{u} \cdot \mathbf{e}^k) = 0 \end{cases}$$

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To evolve in time this approximation, we proceed in two steps:

 \Box Second step: Relaxation. The following system is solved at time t^{n+1} :

$$\begin{cases} \partial_t \rho &= & 0\\ \partial_t \mathbf{u} &= & 0\\ \partial_t \mathcal{E} &= & 0\\ \partial_t s_e &= & \nu_{ei}^{\mathcal{E}} \ c_e^{-\gamma_e} \ \rho^{-\gamma_e} (T_i - T_e) \end{cases}$$

with the data $\widehat{\mathscr{U}}_{\alpha}^{n+1} = \begin{pmatrix} \widehat{\rho}_{\alpha}^{n+1} \\ \widehat{\rho}_{\alpha}^{n+1} \widehat{\eta}_{\alpha} \widehat{\mathbf{u}}_{\alpha}^{n+1} \\ \widehat{\rho}_{\alpha}^{n+1} \widehat{\mathcal{E}}_{\alpha}^{n+1} \\ \widehat{\rho}_{\alpha}^{n+1} \widehat{\mathcal{E}}_{e\alpha}^{n+1} \end{pmatrix}$ at time t^n .

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To evolve in time this approximation, we proceed in two steps:

 \Box Second step: Relaxation. Thanks to the useful Lemma, it amounts to solve at time t^{n+1} the system:

$$\begin{cases} \partial_t \rho = 0, \\ \partial_t \mathbf{u} = 0, \\ \partial_t T_e = \nu_{ei}^{\mathscr{E}} (T_i - T_e), \\ \partial_t T_i = -\nu_{ei}^{\mathscr{E}} (T_i - T_e), \end{cases}$$

with the data $\widehat{\mathscr{U}}_{\alpha}^{n+1}$ at time t^n and temperatures $T_{i,\alpha}^n$ and $T_{e,\alpha}^n$. Solving this system leads to the temperatures $T_{i,\alpha}^{n+1}$ and $T_{e,\alpha}^{n+1}$, and then the energy $\mathscr{E}_{\alpha}^{n+1}$ and entropy state s_e^{n+1} are reconstructed. The state $\mathscr{U}_{\alpha}^{n+1}$ at time t^{n+1} is thus determined.

The details:

Time-step Δt **issue**:

Consider the 2D contol cell Ω_{α} that generates the 3D control cell Ω_{α}^{3D} . Let T_{β} be any generic triangle that enters in the construction of Ω_{α} . Let $h_{\alpha\beta}$ be the minimum of the heights of the triangle T_{β} . We set:

$$\begin{cases} \widehat{\lambda_{\alpha}} = \max\left\{ |u_{\varphi,\alpha}^{n} \pm c_{ei,\alpha}|, |u_{\varphi,\alpha}^{n} - \hat{c}_{ei,\alpha}|, |u_{\varphi,\alpha next}^{n} + \hat{c}_{ei,\alpha next}| \right\}, \\ \widehat{\lambda_{\alpha\beta}} = \max\left\{ |\mathbf{u}_{\alpha}^{n} \cdot \mathbf{n}_{\alpha\beta} \pm c_{ei,\alpha}|, |\mathbf{u}_{\alpha}^{n} \cdot \mathbf{n}_{\alpha\beta} - \hat{c}_{ei,\alpha}|, |\mathbf{u}_{\beta}^{n} \cdot \mathbf{n}_{\alpha\beta} + \hat{c}_{ei,\beta}| \right\}, \end{cases}$$

$$c_{ei,\alpha} = \sqrt{\frac{\partial(p_i + p_e)}{\partial \rho}} \bigg|_{p_i = p_{i,\alpha}^n, p_e = p_{e,\alpha}^n, \rho = \rho_{\alpha}^n}, \ \hat{c}_{ei,\alpha} = \max\left(\sqrt{\frac{\gamma_i p_{i,\alpha}}{c_i \rho_{\alpha}}}, \sqrt{\frac{\gamma_e p_{e,\alpha}}{c_e \rho_{\alpha}}}\right)$$

Then: $\Delta t = \min_{\alpha} \Delta t_{\alpha}$, with $\Delta t_{\alpha} = \min\left\{\frac{\Delta \varphi_{\alpha}}{\widehat{\lambda_{\alpha}}}, \min_{\beta}\left(\frac{h_{\alpha\beta}}{\widehat{\lambda_{\alpha\beta}}}\right)\right\}$

• Sedov problem in 2D axisymmetric geometry⁵

 \Box Uniform media: $T_i = T_e = 2.901 \times 10^4$ K Hot spot: $T_i = 5.802 \times 10^6$ K, and $T_e = 1.7606 \times 10^7$ K Other uniform parameters: $\rho = 1$ kg.m⁻³,

 $u_{R} = u_{7} = u_{\phi} = 0$

$$\Box$$
 A mesh made of 16384 triangles in (R, Z) -coordinates is used for the simulations

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 P_i , P_e , ρ , at $t = 9.7634 \times 10^{-6}$, and T_e/T_i , T_e , $T_$

• Sedov problem in 3D axisymmetric geometry ⁶ \Box Uniform media: $T_i = T_e = 2.901 \times 10^4$ K Hot spot: $T_i = 5.802 \times 10^6$ K, and $T_e = 1.7606 \times 10^7$ K Other uniform parameters: $\rho = 1$ kg.m⁻³, $u_R = u_Z = u_{\phi} = 0$

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S. GALÉRA, ET AL., J. Comput. Phys., 229, 5755-5787 (2010)



 P_i , P_e , ρ , at $t = 9.7634 \times 10^{-6}$, and Te/Ti, as a second state of 34/41

4. Numerical Tests

• Triple point problem in 2D Cartesian geometry⁷

$$(0,3) \xrightarrow{(1,3)} (7,3)$$

$$(0,3) \xrightarrow{\Omega_2} \rho = 0.125, \ p_e = 0.05, \ p_i = 0.05$$

$$\rho = 1, \ p_e = 0.05, \ p_i = 0.05$$

$$p_i = 0.5 \qquad \rho = 1, \ p_e = 0.05, \ p_i = 0.05$$

$$\rho_i = 0.5 \qquad \Omega_1 \qquad (-1.5,7)$$

$$\rho = 0.125, \ p_e = 0.05, \ p_i = 0.05$$

$$(1,-3) \qquad (-1.5,7) \qquad \Omega_3$$





 T_i , T_e at t = 3.5s



The density ρ at $t = 3.5s^{-1}$

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• Sedov problem in 2D axisymmetric geometry



Figure : The three domain of the triple point problem in the (R, Z) plan.



 T_i , T_e at $t = 1.157410^{-5}s$





 T_i , T_e at $t = 1.157410^{-5}s$



- Presentation of a Two-Temperature model for fusion plasma
- Derivation of A Finite Volume approximation to compute the numerical solutions of this model implemented in *PlaTo*
- Numerical tests have shown the accuracy and robutness of the scheme
- Perspectives:
 - $\hfill\square$ More numerical tests to simulate tokamak physics
 - $\hfill\square$ Extension of the model to include magnetic field
 - $\hfill\square$ Extension of the model to include heat flux and anisotropy

THANK YOU