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High-order cell-centered DG
scheme for Lagrangian hydro-
dynamics

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- 1 Introduction
- 2 2D Lagrangian hydrodynamics
- 3 DG discretization general framework
- 4 Second-order DG scheme
- 5 Third-order DG scheme
- 6 Conclusions and perspectives

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Discontinuous Galerkin (DG)

- Natural extension of Finite Volume method
 - Piecewise polynomial approximation of the solution in the cells
 - High-order scheme to achieve high accuracy
-
- Local variational formulation
 - Choice of the numerical fluxes (global L^2 stability, entropy inequality)
 - Time discretization - TVD multistep Runge-Kutta
 -  C.-W. SHU, *Discontinuous Galerkin methods : General approach and stability*, 2008
 - Limitation - vertex-based hierarchical slope limiters
 -  D. KUZMIN, *A vertex-based hierarchical slope limiter for p-adaptive discontinuous Galerkin methods* J. Comp. Appl. Math., 2009
 -  M. YANG AND Z.J. WANG, *A parameter-free generalized moment limiter for high-order methods on unstructured grids* Adv. Appl. Math. Mech., 2009

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Cell-Centered Lagrangian schemes

Finite volume schemes on moving mesh

- J. K. Dukowicz : CAVEAT scheme

A computer code for fluid dynamics problems with large distortion and internal slip, 1986

- B. Després : GLACE scheme

Lagrangian Gas Dynamics in Two Dimensions and Lagrangian systems, 2005

- P.-H. Maire : EUCLHYD scheme

A cell-centered Lagrangian scheme for two-dimensional compressible flow problems, 2007

- G. Kluth : Hyperelasticity

Discretization of hyperelasticity with a cell-centered Lagrangian scheme, 2010

- S. Del Pino : Curvilinear Finite Volume method

A curvilinear finite-volume method to solve compressible gas dynamics in semi-Lagrangian coordinates, 2010

- P. Hoch : Finite Volume method on unstructured conical meshes

Extension of ALE methodology to unstructured conical meshes, 2011

DG scheme on initial mesh

- R. Loubère : PhD thesis

Une Méthode Particulaire Lagrangienne de type Galerkin Discontinu. Application à la mécanique des Fluides et l'Interaction Laser/Plasma, 2002

Lagrangian and Eulerian descriptions

- Let's have a continuous mapping

$$\mathbf{x} = \Phi(\mathbf{X}, t)$$

- \mathbf{X} the Lagrangian (initial) coordinate
- \mathbf{x} the Eulerian (actual) coordinate
- \mathbf{N} the Lagrangian normal
- \mathbf{n} the Eulerian normal
- $\mathbf{F} = \nabla_{\mathbf{X}} \Phi = \frac{\partial \mathbf{x}}{\partial \mathbf{X}}$ the deformation gradient tensor, $J = \det \mathbf{F}$

- $\rho J = \rho^0$
- $dV = J dV$
- $d\mathbf{x} = \mathbf{F} d\mathbf{X}$
- $J \mathbf{F}^{-t} \mathbf{N} dS = \mathbf{n} ds$ Nanson formula
- $\nabla_{\mathbf{x}} P = \frac{1}{J} \nabla_{\mathbf{X}} \cdot (P \mathbf{J} \mathbf{F}^{-t})$
- $\nabla_{\mathbf{x}} \cdot \mathbf{U} = \frac{1}{J} \nabla_{\mathbf{X}} \cdot (\mathbf{J} \mathbf{F}^{-1} \mathbf{U})$

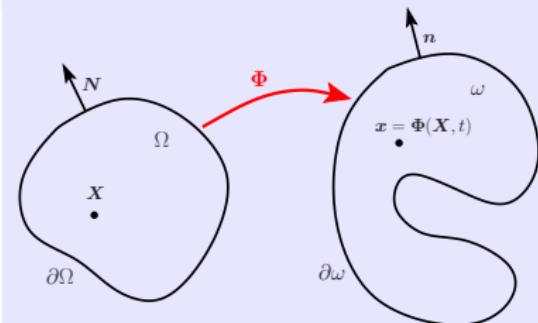


FIG.: Notation for the flow map.

Lagrangian and Eulerian descriptions

- $G = JF^{-t}$ the cofactor matrix of F
- $\nabla_x \cdot G = \mathbf{0}$ Piola compatibility condition

$$\int_{\Omega_c} \nabla_x \cdot G \, dV = \int_{\partial\Omega_c} G \mathbf{N} \, dS = \int_{\partial\omega_c} \mathbf{n} \, ds = \mathbf{0}$$

- Gas dynamics system in Lagrangian formalism

$$\frac{dF}{dt} = \nabla_X \mathbf{U}$$

$$\rho^0 \frac{d}{dt} \left(\frac{1}{\rho} \right) - \nabla_X \cdot (G^t \mathbf{U}) = 0$$

$$\rho^0 \frac{d \mathbf{U}}{dt} + \nabla_X \cdot (P G) = \mathbf{0}$$

$$\rho^0 \frac{d E}{dt} + \nabla_X \cdot (G^t P \mathbf{U}) = 0$$

- Thermodynamical closure EOS : $P = P(\rho, \varepsilon)$ where $\varepsilon = E - \frac{1}{2} \mathbf{U}^2$

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DG discretization general framework

DG discretization

- Let $\{\Omega_c\}_c$ be a partition of the domain Ω into polygonal cells
- $\{\sigma_k^c\}_{k=0 \dots K}$ basis of $\mathbb{P}^\alpha(\Omega_c)$
- $\phi_h^c(\mathbf{X}, t) = \sum_{k=0}^K \phi_k^c(t) \sigma_k^c(\mathbf{X})$ approximate function of $\phi(\mathbf{X}, t)$ on Ω_c

Definitions

- Center of mass $\Xi_c = (\Xi_c^X, \Xi_c^Y)^t = \frac{1}{m_c} \int_{\Omega_c} \rho^0(\mathbf{X}) \mathbf{X} dV$,
where m_c is the constant mass of the cell Ω_c
- The mean value $\langle \phi \rangle_c = \frac{1}{m_c} \int_{\Omega_c} \rho^0(\mathbf{X}) \phi(\mathbf{X}) dV$
of the function ϕ over the cell Ω_c
- The associated scalar product $\langle \phi, \psi \rangle_c = \int_{\Omega_c} \rho^0(\mathbf{X}) \phi(\mathbf{X}) \psi(\mathbf{X}) dV$

Polynomial Taylor basis

- Taylor expansion on the cell, located at the center of mass Ξ_c
- $\sigma_0^c = 1$ and going further in space discretization, the $q + 1$ basis functions of degree q , with $0 < q \leq \alpha$, write

$$\sigma_{\frac{q(q+1)}{2}+j}^c = \frac{1}{j!(q-j)!} \left[\left(\frac{X - \Xi_c^X}{\Delta X_c} \right)^{q-j} \left(\frac{Y - \Xi_c^Y}{\Delta Y_c} \right)^j - \left\langle \left(\frac{X - \Xi_c^X}{\Delta X_c} \right)^{q-j} \left(\frac{Y - \Xi_c^Y}{\Delta Y_c} \right)^j \right\rangle_c \right]$$

where $j = 0 \dots q$, $\Delta X_c = \frac{X_{max} - X_{min}}{2}$ and $\Delta Y_c = \frac{Y_{max} - Y_{min}}{2}$ with X_{max} , Y_{max} , X_{min} , Y_{min} the maximum and minimum coordinates in the cell Ω_c

Outcome

- Same basis functions whatever the shape of the cells
- $\langle \sigma_k^c \rangle_c = m_c \delta_{0k}$ where δ_{kl} is the Kronecker symbol
- $\langle \sigma_0^c, \sigma_k^c \rangle_c = 0, \forall k \neq 0$
- $\phi_0^c = \langle \phi \rangle_c$ the mass averaged value of the function ϕ over the cell Ω_c

Local variational formulations

- $$\begin{aligned} \bullet \quad \frac{d}{dt} \int_{\Omega_c} \rho^0 \left(\frac{1}{\rho}\right) \sigma_q^c dV &= \sum_{k=0}^K \frac{d}{dt} \left(\frac{1}{\rho}\right)_k^c \int_{\Omega_c} \rho^0 \sigma_q^c \sigma_k^c dV \\ &= - \int_{\Omega_c} \boldsymbol{U} \cdot \mathbf{G} \nabla_X \sigma_q^c dV + \int_{\partial \Omega_c} \overline{\boldsymbol{U}} \cdot \sigma_q^c \mathbf{G} \mathbf{N} dL \end{aligned}$$

- $$\begin{aligned} \bullet \quad \frac{d}{dt} \int_{\Omega_c} \rho^0 \boldsymbol{U} \sigma_q^c dV &= \sum_{k=0}^K \frac{d \boldsymbol{U}_k^c}{dt} \int_{\Omega_c} \rho^0 \sigma_q^c \sigma_k^c dV \\ &= \int_{\Omega_c} \boldsymbol{P} \mathbf{G} \nabla_X \sigma_q^c dV - \int_{\partial \Omega_c} \overline{\boldsymbol{P}} \sigma_q^c \mathbf{G} \mathbf{N} dL \end{aligned}$$

- $$\begin{aligned} \bullet \quad \frac{d}{dt} \int_{\Omega_c} \rho^0 E \sigma_q^c dV &= \sum_{k=0}^K \frac{d E_k^c}{dt} \int_{\Omega_c} \rho^0 \sigma_q^c \sigma_k^c dV \\ &= \int_{\Omega_c} \boldsymbol{P} \boldsymbol{U} \cdot \mathbf{G} \nabla_X \sigma_q^c dV - \int_{\partial \Omega_c} \overline{\boldsymbol{P} \boldsymbol{U}} \cdot \sigma_q^c \mathbf{G} \mathbf{N} dL \end{aligned}$$

Local variational formulations

- $\int_{\Omega_c} \rho^0 \sigma_q^c \sigma_k^c dV = \langle \sigma_q^c, \sigma_k^c \rangle_c$ generic coefficient of the symmetric positive definite mass matrix
- $\int_{\Omega_c} \rho^0 \sigma_k^c dV = 0, \forall k \neq 0$ implies that the equations corresponding to mass averaged values are independent of the other polynomial basis components equations

- $\int_{\Omega_c} \mathbf{U} \cdot \mathbf{G} \nabla_X \sigma_q^c dV, \int_{\Omega_c} P \mathbf{G} \nabla_X \sigma_q^c dV$ and $\int_{\Omega_c} P \mathbf{U} \cdot \mathbf{G} \nabla_X \sigma_q^c dV$ are evaluated through the use of a two-dimensional quadrature rule
- $\int_{\partial\Omega_c} \overline{\mathbf{U}} \cdot \sigma_q^c \mathbf{G} \mathbf{N} dL, \int_{\partial\Omega_c} \overline{P} \sigma_q^c \mathbf{G} \mathbf{N} dL$ and $\int_{\partial\Omega_c} \overline{P} \mathbf{U} \cdot \sigma_q^c \mathbf{G} \mathbf{N} dL$ required a specific treatment to ensure the GCL

Entropic analysis

Entropic semi-discrete equation

- Fundamental assumption $\overline{P} \overline{\mathbf{U}} = \overline{P} \overline{\mathbf{U}}$
- The use of variational formulations and Piola condition leads to

$$\int_{\Omega_c} \rho^0 \theta \frac{d\eta}{dt} dV = \int_{\partial\Omega_c} (\overline{P} - P)(\mathbf{U} - \overline{\mathbf{U}}) \cdot \mathbf{G}\mathbf{N} dL,$$

where η is the specific entropy and θ the absolute temperature defined by means of the Gibbs identity

Entropic semi-discrete equation

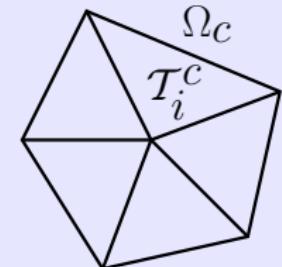
- A sufficient condition to satisfy $\int_{\Omega_c} \rho^0 \theta \frac{d\eta}{dt} dV \geq 0$ is

$$\overline{P} - P = -Z(\overline{\mathbf{U}} - \mathbf{U}) \cdot \frac{\mathbf{G}\mathbf{N}}{\|\mathbf{G}\mathbf{N}\|} = -Z(\overline{\mathbf{U}} - \mathbf{U}) \cdot \mathbf{n}, \quad (2)$$

where $Z \geq 0$ has the physical dimension of a density times a velocity

Deformation gradient tensor discretization

- Requirements : Piola compatibility condition and geometry continuity
- Triangular decomposition $\Omega_c = \bigcup_{i=1}^{ntri} \mathcal{T}_i^c$
- F discretization by means of a mapping defined on triangular cells
- We develop Φ on the Finite Element basis functions λ_p



$$\Phi_h^i(\mathbf{X}, t) = \sum_p \lambda_p(\mathbf{X}) \Phi_p(t),$$

where the points p are control points including vertices in \mathcal{T}_i

- $\Phi_p(t) = \Phi(\mathbf{X}_p, t) = \mathbf{x}_p$
- $\mathbf{F} = \nabla_{\mathbf{X}}\Phi \implies \mathbf{F}_i(\mathbf{X}, t) = \sum_p \Phi_p(t) \otimes \nabla_{\mathbf{X}}\lambda_p(\mathbf{X})$
- $\frac{d\Phi_p}{dt} = \mathbf{U}_p \implies \frac{d}{dt}\mathbf{F}_i(\mathbf{X}, t) = \sum_p \mathbf{U}_p(t) \otimes \nabla_{\mathbf{X}}\lambda_p(\mathbf{X})$

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P_1 deformation gradient tensor discretization

- The chosen linear basis functions are the P_1 barycentric coordinate basis functions which write in a generic triangle \mathcal{T}_i

$$\lambda_p(\mathbf{X}) = \frac{1}{2|\mathcal{T}_i|} [X(Y_{p^+} - Y_{p^-}) - Y(X_{p^+} - X_{p^-}) + X_{p^+}Y_{p^-} - X_{p^-}Y_{p^+}],$$

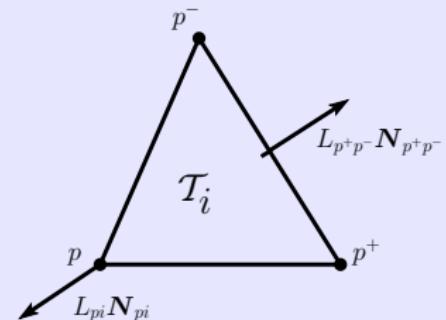
where p , p^+ and p^- are the counterclockwise ordered triangle nodes and $|\mathcal{T}_i|$ the triangle volume

- $\mathbf{F}_i(t) = \frac{1}{|\mathcal{T}_i|} \sum_{p \in \mathcal{P}(\mathcal{T}_i)} \mathbf{x}_p(t) \otimes L_{pi} \mathbf{N}_{pi},$

$$\frac{d}{dt} \mathbf{F}_i(t) = \frac{1}{|\mathcal{T}_i|} \sum_{p \in \mathcal{P}(\mathcal{T}_i)} \mathbf{U}_p(t) \otimes L_{pi} \mathbf{N}_{pi},$$

where $\mathcal{P}(\mathcal{T}_i)$ is the node set of \mathcal{T}_i

- $L_{pi} \mathbf{N}_{pi} = \frac{1}{2}(L_{p^-p} \mathbf{N}_{p^-p} + L_{pp^+} \mathbf{N}_{pp^+})$ the corner normal at node p in the initial configuration



Local boundary terms integration

Numerical fluxes linear approximation

- On face f_{pp^+} , $\bar{\psi}_{|_{pp^+}}^c(\zeta) = \psi_{pc}^+(1 - \zeta) + \psi_{p^+c}^- \zeta$,
where $\zeta \in [0, 1]$ is the curvilinear abscissa
- Hence

$$\bar{\mathbf{U}}_{|_{pp^+}}(\zeta) = \mathbf{U}_p(1 - \zeta) + \mathbf{U}_{p^+} \zeta$$

$$\bar{\mathbf{P}}_{|_{pp^+}}^c(\zeta) = \mathbf{P}_{pc}^+(1 - \zeta) + \mathbf{P}_{p^+c}^- \zeta$$

$$\bar{\mathbf{P}}\bar{\mathbf{U}}_{|_{pp^+}}^c(\zeta) = (\mathbf{P}\mathbf{U})_{pc}^+(1 - \zeta) + (\mathbf{P}\mathbf{U})_{p^+c}^- \zeta$$

- The basis function σ_q^c being linear over Ω_c

$$\sigma_{q|_{pp^+}}^c(\zeta) = \sigma_q^c(\mathbf{X}_p)(1 - \zeta) + \sigma_q^c(\mathbf{X}_{p^+}) \zeta$$

Fundamental assumption

- $\bar{\mathbf{P}}\bar{\mathbf{U}} = \bar{\mathbf{P}}\bar{\mathbf{U}} \implies (\mathbf{P}\mathbf{U})_{pc}^\pm = \mathbf{P}_{pc}^\pm \mathbf{U}_p$

Local boundary terms integration

Analytical integration

- $\int_{\partial\Omega_c} \bar{\psi} \sigma_q^c \mathbf{G} \mathbf{N} dL = \sum_{p \in \mathcal{P}(c)} \left(\int_0^1 \bar{\psi}|_{pp^+}(\zeta) \sigma_q^c|_{pp^+}(\zeta) d\zeta \right) \mathbf{G}|_{pp^+} L_{pp^+} \mathbf{N}_{pp^+}$,
where $\mathbf{G}|_{pp^+}$ is the constant value of tensor \mathbf{G} on face f_{pp^+}
- $\mathbf{G}|_{pp^+} L_{pp^+} \mathbf{N}_{pp^+} = l_{pp^+} \mathbf{n}_{pp^+}$ Eulerian normal of face f_{pp^+}
- $$\begin{aligned} \int_{\partial\Omega_c} \bar{\psi} \sigma_q^c \mathbf{G} \mathbf{N} dL &= \sum_{p \in \mathcal{P}(c)} \left(\int_0^1 \bar{\psi} \sigma_q^c|_{pp^+}(\zeta) d\zeta \right) l_{pp^+} \mathbf{n}_{pp^+} \\ &= \sum_{p \in \mathcal{P}(c)} \frac{1}{6} [\psi_{pc}^+(2\sigma_q^c(\mathbf{X}_p) + \sigma_q^c(\mathbf{X}_{p^+}))l_{pp^+} \mathbf{n}_{pp^+} \\ &\quad + \psi_{p^+c}^-(2\sigma_q^c(\mathbf{X}_{p^+}) + \sigma_q^c(\mathbf{X}_p))l_{pp^+} \mathbf{n}_{pp^+}] \\ &= \sum_{p \in \mathcal{P}(c)} \frac{1}{6} [\psi_{pc}^+(2\sigma_q^c(\mathbf{X}_p) + \sigma_q^c(\mathbf{X}_{p^+}))l_{pp^+} \mathbf{n}_{pp^+} \\ &\quad + \psi_{pc}^-(2\sigma_q^c(\mathbf{X}_p) + \sigma_q^c(\mathbf{X}_{p^-}))l_{p-p} \mathbf{n}_{p-p}] \end{aligned}$$

Semi-discrete equations

- Half-Left and half-right corner normals $I_{pc}^{\pm} \mathbf{n}_{pc}^{\pm}$

$$I_{pc}^- \mathbf{n}_{pc}^- = \frac{1}{2} I_{p-p}^- \mathbf{n}_{p-p}^- \quad \text{and} \quad I_{pc}^+ \mathbf{n}_{pc}^+ = \frac{1}{2} I_{pp^+} \mathbf{n}_{pp^+}$$

- The weighted corner normals

$$I_{pc}^q \mathbf{n}_{pc}^q = I_{pc}^{-,q} \mathbf{n}_{pc}^- + I_{pc}^{+,q} \mathbf{n}_{pc}^+ \quad \text{where} \quad I_{pc}^{\pm,q} = \frac{1}{3} (2\sigma_q^c(\mathbf{X}_p) + \sigma_q^c(\mathbf{X}_{p^\pm})) I_{pc}^{\pm}$$

- The q^{th} moment of the subcell forces

$$\mathbf{F}_{pc}^q = P_{pc}^- I_{pc}^{-,q} \mathbf{n}_{pc}^- + P_{pc}^+ I_{pc}^{+,q} \mathbf{n}_{pc}^+$$

- The semi-discrete equations on the specific volume, momentum and total energy successive moments, in respect with the GCL, write

$$\int_{\Omega_c} \rho^0 \frac{d}{dt} \left(\frac{1}{\rho} \right) \sigma_q^c dV = - \sum_{i=1}^{ntri} \mathbf{G}_i^c \nabla_X \sigma_q^c \cdot \int_{T_i^c} \mathbf{U} dV + \sum_{p \in \mathcal{P}(c)} \mathbf{U}_p \cdot I_{pc}^q \mathbf{n}_{pc}^q$$

$$\int_{\Omega_c} \rho^0 \frac{d}{dt} \sigma_q^c dV = \sum_{i=1}^{ntri} \mathbf{G}_i^c \nabla_X \sigma_q^c \int_{T_i^c} P dV - \sum_{p \in \mathcal{P}(c)} \mathbf{F}_{pc}^q$$

$$\int_{\Omega_c} \rho^0 \frac{d}{dt} \sigma_q^c dV = \sum_{i=1}^{ntri} \mathbf{G}_i^c \nabla_X \sigma_q^c \cdot \int_{T_i^c} P \mathbf{U} dV - \sum_{p \in \mathcal{P}(c)} \mathbf{U}_p \cdot \mathbf{F}_{pc}^q$$

Nodal solvers

- Fundamental identity on the cell Ω_c

$$\bar{P} = P_h^c - Z_c (\bar{\mathbf{U}} - \mathbf{U}_h^c) \cdot \mathbf{n},$$

where $Z_c = \rho_c a_c$ is the acoustic impedance

- Using this expression to calculate \mathbf{F}_{pc}^q leads to

$$\begin{aligned}\mathbf{F}_{pc}^q &= P_{pc}^- I_{pc}^{-,q} \mathbf{n}_{pc}^- + P_{pc}^+ I_{pc}^{+,q} \mathbf{n}_{pc}^+ \\ &= P_h^c(\mathbf{X}_p, t) I_{pc}^q \mathbf{n}_{pc}^q - Z_c [(\mathbf{U}_p - \mathbf{U}_h^c(\mathbf{X}_p, t)) \cdot \mathbf{n}_{pc}^-] I_{pc}^{-,q} \mathbf{n}_{pc}^- \\ &\quad - Z_c [(\mathbf{U}_p - \mathbf{U}_h^c(\mathbf{X}_p, t)) \cdot \mathbf{n}_{pc}^+] I_{pc}^{+,q} \mathbf{n}_{pc}^+\end{aligned}$$

- Finally, the q^{th} moment of the subcell force writes

$$\mathbf{F}_{pc}^q = P_h^c(\mathbf{X}_p, t) I_{pc}^q \mathbf{n}_{pc}^q - \mathbf{M}_{pc}^q (\mathbf{U}_p - \mathbf{U}_h^c(\mathbf{X}_p, t)),$$

where $\mathbf{M}_{pc}^q = Z_c (I_{pc}^{-,q} \mathbf{n}_{pc}^- \otimes \mathbf{n}_{pc}^- + I_{pc}^{+,q} \mathbf{n}_{pc}^+ \otimes \mathbf{n}_{pc}^+)$

Nodal solvers

- To be conservative in total energy and momentum over the whole domain, $\sum_{c \in \mathcal{C}(p)} \mathbf{F}_{pc} = \mathbf{0}$ and thus

$$\left(\sum_{c \in \mathcal{C}(p)} M_{pc} \right) \mathbf{U}_p = \sum_{c \in \mathcal{C}(p)} \left[P_h^c(\mathbf{X}_p, t) I_{pc} \mathbf{n}_{pc} + M_{pc} \mathbf{U}_h^c(\mathbf{X}_p, t) \right],$$

where $M_{pc} = Z_c (I_{pc}^+ \mathbf{n}_{pc}^+ \otimes \mathbf{n}_{pc}^+ + I_{pc}^- \mathbf{n}_{pc}^- \otimes \mathbf{n}_{pc}^-)$ are positive semi-definite matrices with a physical dimension of a density times a velocity

First moment equations

- $m_c \frac{d}{dt} \left(\frac{1}{\rho} \right)_0^c = \sum_{p \in \mathcal{P}(c)} \mathbf{U}_p \cdot I_{pc} \mathbf{n}_{pc}$

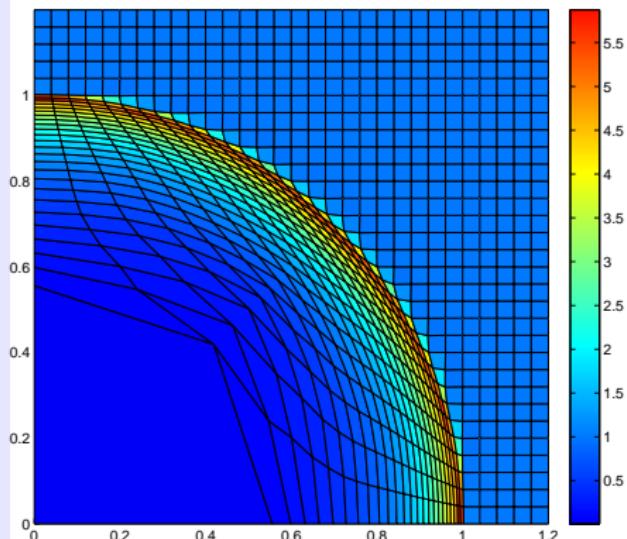
- $m_c \frac{d \mathbf{U}_0^c}{dt} = - \sum_{p \in \mathcal{P}(c)} \mathbf{F}_{pc}$

- $m_c \frac{d E_0^c}{dt} = - \sum_{p \in \mathcal{P}(c)} \mathbf{U}_p \cdot \mathbf{F}_{pc}$

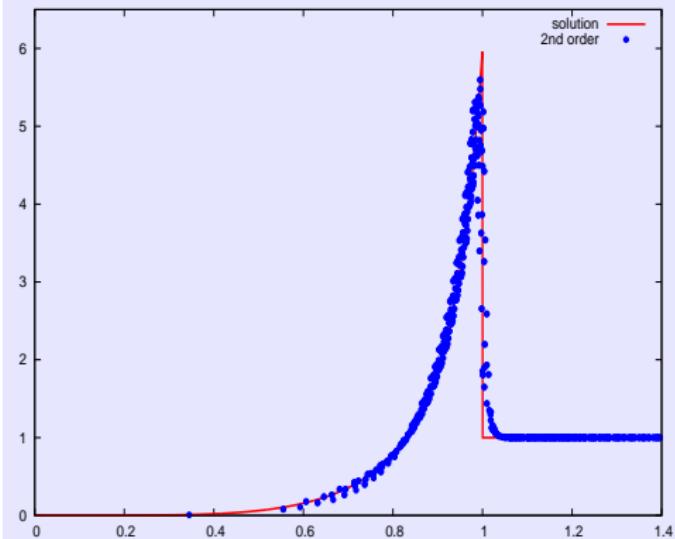
We recover the EUCLHYD scheme

Numerical results

Sedov point blast problem on a Cartesian grid



(a) Second-order scheme with limitation.



(b) Density profiles comparison.

FIG.: Point blast Sedov problem on a Cartesian grid made of 30×30 cells : density.

Numerical results

Sedov point blast problem on a polygonal grid

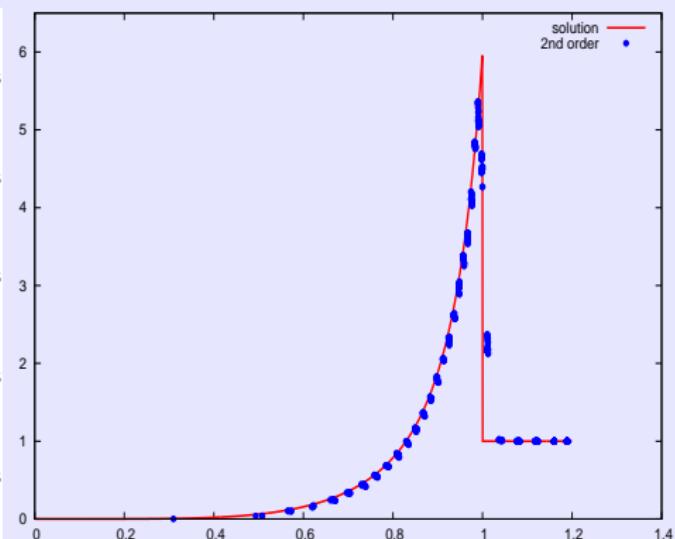
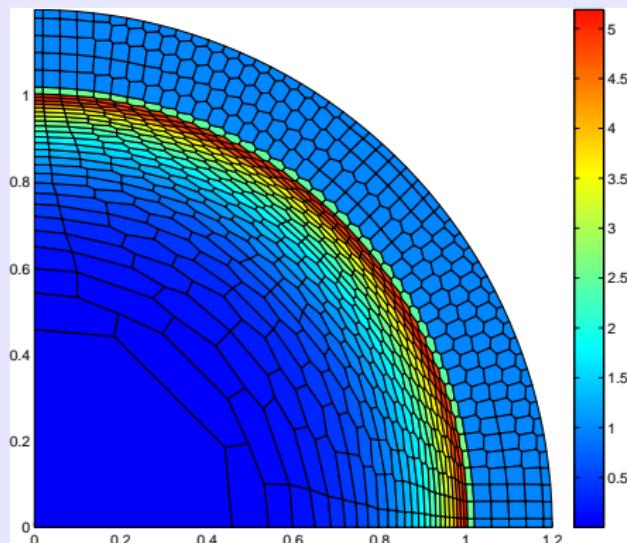
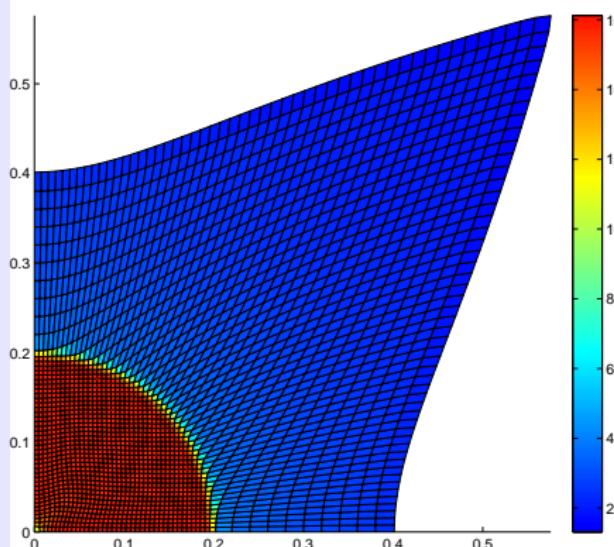


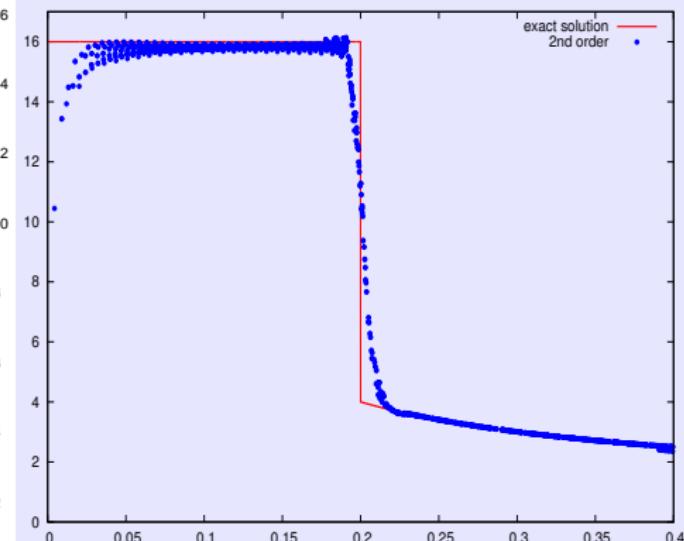
FIG.: Point blast Sedov problem on a unstructured grid made of 775 polygonal cells : density map.

Numerical results

Noh problem



(a) Second-order scheme with limitation.

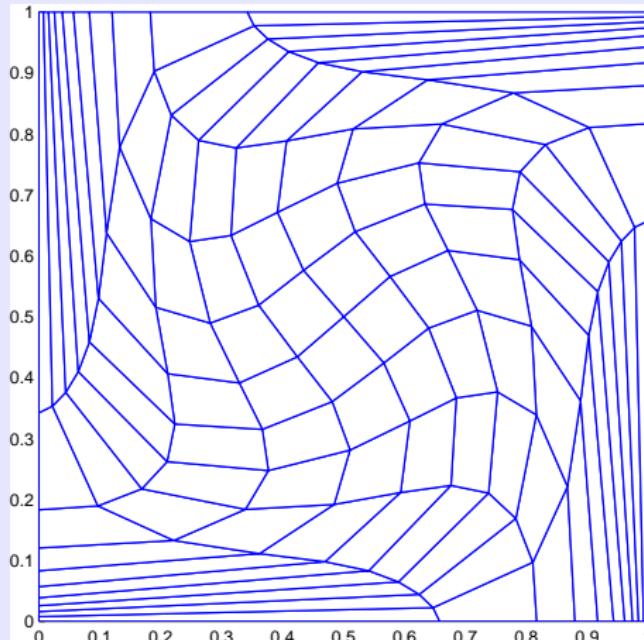


(b) Density profiles comparison.

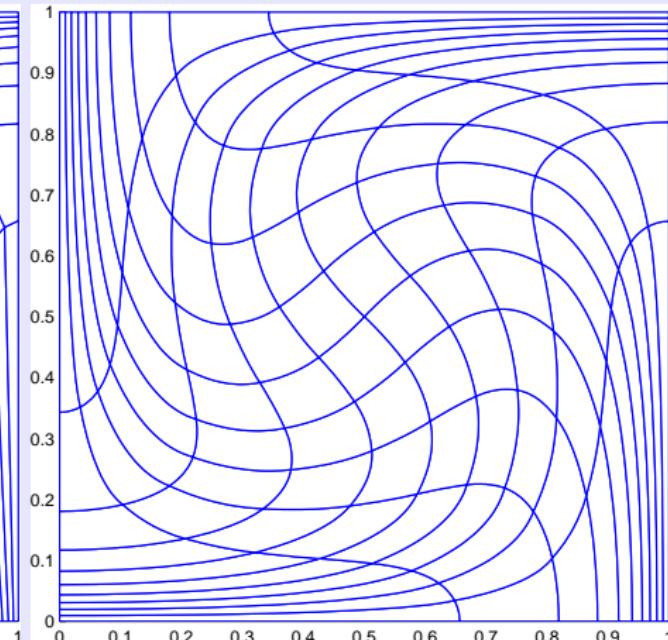
FIG.: Noh problem on a Cartesian grid made of 50×50 cells : density.

Numerical results

Taylor-Green vortex problem



(a) Second-order scheme.



(b) Exact solution.

FIG.: Motion of a 10×10 Cartesian mesh through a T.-G. vortex, at $t = 0.75$.

Rate of convergence computed on the pressure in the case of the Taylor-Green vortex

	L_1		L_2		L_∞	
h	$E_{L_1}^h$	$q_{L_1}^h$	$E_{L_2}^h$	$q_{L_2}^h$	$E_{L_\infty}^h$	$q_{L_\infty}^h$
$\frac{1}{20}$	8.98E-3	1.88	1.51E-2	1.75	6.73E-2	1.27
$\frac{1}{40}$	2.44E-3	1.94	4.48E-3	1.95	2.79E-2	1.68
$\frac{1}{80}$	6.36E-4	2.00	1.16E-3	2.00	8.68E-3	1.95
$\frac{1}{160}$	1.59E-4	2.01	2.90E-4	2.01	2.24E-3	2.01
$\frac{1}{320}$	3.94E-5	-	7.18E-5	-	5.54E-4	-

TAB.: Second-order DG scheme without limitation at time $t = 0.6$.

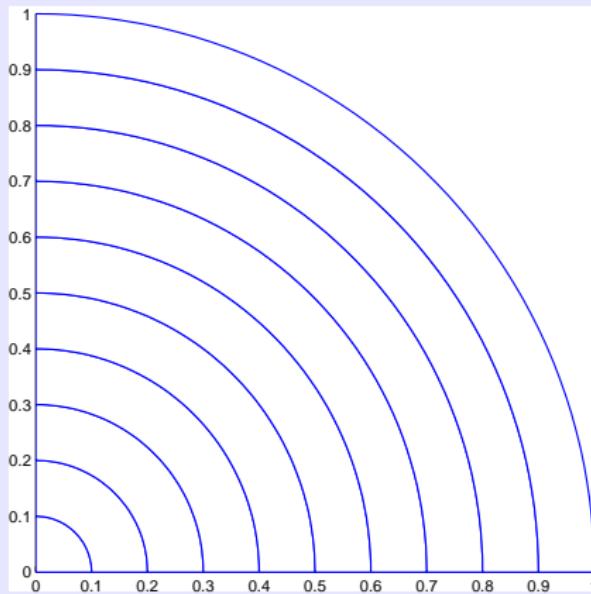
	L_1		L_2		L_∞	
h	$E_{L_1}^h$	$q_{L_1}^h$	$E_{L_2}^h$	$q_{L_2}^h$	$E_{L_\infty}^h$	$q_{L_\infty}^h$
$\frac{1}{20}$	1.99E-2	2.33	2.92E-2	2.03	8.27E-2	1.34
$\frac{1}{40}$	3.96E-3	2.25	7.16E-3	2.20	3.26E-2	1.61
$\frac{1}{80}$	8.31E-4	2.17	1.56E-3	2.15	1.07E-2	1.52
$\frac{1}{160}$	1.85E-4	2.11	3.52E-4	2.14	3.73E-3	2.41
$\frac{1}{320}$	4.28E-5	-	8.01E-5	-	7.01E-4	-

TAB.: Second-order DG scheme with limitation at time $t = 0.6$.

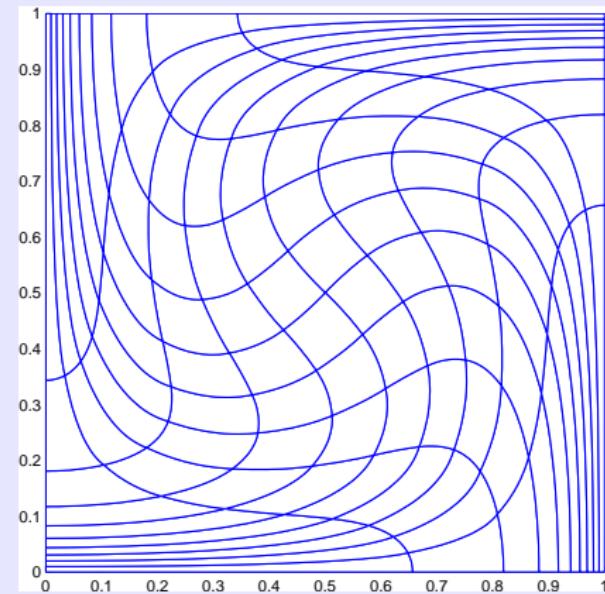
- 1 Introduction
- 2 2D Lagrangian hydrodynamics
- 3 DG discretization general framework
- 4 Second-order DG scheme
- 5 Third-order DG scheme
- 6 Conclusions and perspectives

Curvilinear elements motivation

Circular polar grid : 10×1 cells



Taylor-Green exact motion



V. DOBREV, T. ELLIS, T. KOLEV AND R. RIEBEN, *High Order Curvilinear Finite Elements for Lagrangian Hydrodynamics. Part I : General Framework*, 2010. Presentation available at
<https://computation.llnl.gov/casc/blast/blast.html>

Geometric consideration

- The P_2 quadratic mapping function writes

$$\mathbf{x} = \Phi(\mathbf{X}, t) = \sum_p \mathbf{x}_p(t) \mu_p(\mathbf{X}),$$

where the points p are the triangular nodes and the control points Q of the Bezier edges, and the P_2 barycentric coordinate functions μ_p write

$$\begin{aligned}\mu_p &= (\lambda_p)^2, \quad \mu_{p^+} = (\lambda_{p^+})^2, \quad \mu_{p^-} = (\lambda_{p^-})^2, \\ \mu_Q &= 2\lambda_p\lambda_{p^+}, \quad \mu_{Q^+} = 2\lambda_{p^+}\lambda_{p^-}, \quad \mu_{Q^-} = 2\lambda_{p^-}\lambda_p,\end{aligned}$$

where the functions λ_l , with $l \in \{p, p^+, p^-\}$, are the P_1 Finite Elements linear basis functions

- Finally, the quadratic mapping expresses as

$$\Phi(\mathbf{X}, t) = \sum_{p \in \mathcal{P}(\mathcal{T}_i)} [\mathbf{x}_p(t) (\lambda_p(\mathbf{X}))^2 + 2\mathbf{x}_Q(t) \lambda_p(\mathbf{X}) \lambda_p^+(\mathbf{X})]$$

Geometric consideration

$$\mathbf{F}_i(\mathbf{X}, t) = \frac{2}{|\mathcal{T}_c|} \sum_{p \in \mathcal{P}(\mathcal{T}_i)} \lambda_p [\mathbf{x}_p \otimes L_{pc} \mathbf{N}_{pc} + \mathbf{x}_Q \otimes L_{p^+c} \mathbf{N}_{p^+c} + \mathbf{x}_Q^- \otimes L_{p^-c} \mathbf{N}_{p^-c}]$$

$$\frac{d}{dt} \mathbf{F}_i(\mathbf{X}, t) = \frac{2}{|\mathcal{T}_i|} \sum_{p \in \mathcal{P}(\mathcal{T}_i)} \lambda_p [\mathbf{U}_p \otimes L_{pc} \mathbf{N}_{pc} + \mathbf{U}_Q \otimes L_{p^+c} \mathbf{N}_{p^+c} + \mathbf{U}_Q^- \otimes L_{p^-c} \mathbf{N}_{p^-c}]$$

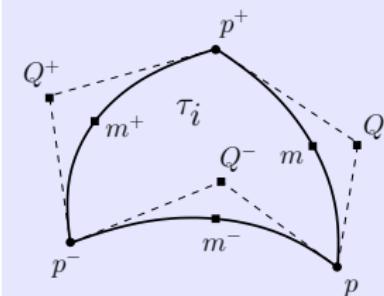
where $\mathbf{U}_Q = \frac{4\mathbf{U}_m - \mathbf{U}_p - \mathbf{U}_{p^+}}{2}$ and $L_{pc} \mathbf{N}_{pc} = \frac{1}{2} (\mathbf{X}_{p^+} - \mathbf{X}_{p^-}) \times \mathbf{e}_Z$

- Given p , Q and p^+ , and ζ in $[0, 1]$, we define the Bezier curve as

$$\mathbf{x}(\zeta) = (1 - \zeta)^2 \mathbf{x}_p + 2\zeta(1 - \zeta) \mathbf{x}_Q + \zeta^2 \mathbf{x}_{p^+}$$

- Midpoint $\mathbf{x}_m = \mathbf{x}(\frac{1}{2}) = \frac{2\mathbf{x}_Q + \mathbf{x}_p + \mathbf{x}_{p^+}}{4}$

- $\mathbf{t} dI = \frac{d\mathbf{x}}{d\zeta} d\zeta = 2((1 - \zeta)(\mathbf{x}_Q - \mathbf{x}_p) + \zeta(\mathbf{x}_{p^+} - \mathbf{x}_Q)) d\zeta$



Local boundary terms integration

Numerical fluxes quadratic approximation

- On face f_{pp^+}

$$\mathbf{U}_{|_{pp^+}}(\zeta) = (1 - \zeta)(1 - 2\zeta)\mathbf{U}_p + 4\zeta(1 - \zeta)\mathbf{U}_m + \zeta(2\zeta - 1)\mathbf{U}_{p^+}$$

$$\overline{P}^c_{|_{pp^+}}(\zeta) = (1 - \zeta)(1 - 2\zeta)P_{pc}^+ + 4\zeta(1 - \zeta)P_{mc} + \zeta(2\zeta - 1)P_{p^+c}^-$$

$$\overline{PU}^c_{|_{pp^+}}(\zeta) = (1 - \zeta)(1 - 2\zeta)(PU)_{pc}^+ + 4\zeta(1 - \zeta)(PU)_{mc} + \zeta(2\zeta - 1)(PU)_{p^+c}^-$$

- The basis function σ_q^c being quadratic over Ω_c

$$\sigma_q^c|_{pp^+}(\zeta) = (1 - \zeta)(1 - 2\zeta)\sigma_q^c(\mathbf{X}_p) + 4\zeta(1 - \zeta)\sigma_q^c(\mathbf{X}_m) + \zeta(2\zeta - 1)\sigma_q^c(\mathbf{X}_{p^+})$$

Fundamental assumption

- $\overline{PU} = \overline{P} \overline{U} \implies (PU)_{pc}^\pm = P_{pc}^\pm \mathbf{U}_p \text{ and } (PU)_{mc} = P_{mc} \mathbf{U}_m$

Normal and subcell forces definitions

$$I_{pc}^q \mathbf{n}_{pc}^q = I_{pc}^{-,q} \mathbf{n}_{pc}^{-,q} + I_{pc}^{+,q} \mathbf{n}_{pc}^{+,q}$$

$$I_{pc}^{-,q} \mathbf{n}_{pc}^{-,q} = \frac{1}{10} [(6\sigma_q^c(\mathbf{X}_p) + 4\sigma_q^c(\mathbf{X}_{m-})) I_{Q-p} \mathbf{n}_{Q-p} + (\sigma_q^c(\mathbf{X}_p) - \sigma_q^c(\mathbf{X}_{p-})) I_{p-p} \mathbf{n}_{p-p}]$$

$$I_{pc}^{+,q} \mathbf{n}_{pc}^{+,q} = \frac{1}{10} [(6\sigma_q^c(\mathbf{X}_p) + 4\sigma_q^c(\mathbf{X}_m)) I_{pQ} \mathbf{n}_{pQ} + (\sigma_q^c(\mathbf{X}_p) - \sigma_q^c(\mathbf{X}_{p+})) I_{pp^+} \mathbf{n}_{pp^+}]$$

$$\mathbf{F}_{pc}^q = P_{pc}^- I_{pc}^{-,q} \mathbf{n}_{pc}^{-,q} + P_{pc}^+ I_{pc}^{+,q} \mathbf{n}_{pc}^{+,q}$$

$$I_{mc}^q \mathbf{n}_{mc}^q = I_{mc}^{-,q} \mathbf{n}_{mc}^{-,q} + I_{mc}^{+,q} \mathbf{n}_{mc}^{+,q}$$

$$I_{mc}^{-,q} \mathbf{n}_{mc}^{-,q} = \frac{1}{5} (4\sigma_q^c(\mathbf{X}_m) + \sigma_q^c(\mathbf{X}_p)) I_{pQ} \mathbf{n}_{pQ}$$

$$I_{mc}^{+,q} \mathbf{n}_{mc}^{+,q} = \frac{1}{5} (4\sigma_q^c(\mathbf{X}_m) + \sigma_q^c(\mathbf{X}_{p+})) I_{Qp^+} \mathbf{n}_{Qp^+}$$

$$\mathbf{F}_{mc}^q = P_{mc} I_{mc}^q \mathbf{n}_{mc}^q$$

Case of $q = 0$

$$I_{pc} \mathbf{n}_{pc} = I_{Q-Q} \mathbf{n}_{Q-Q}$$

$$I_{mc} \mathbf{n}_{mc} = I_{pp^+} \mathbf{n}_{pp^+}$$

- $\int_{\partial\Omega_c} \overline{\mathbf{U}} \cdot \mathbf{G} \mathbf{N} dL = \sum_{p \in \mathcal{P}(c)} \frac{1}{3} \mathbf{U}_p \cdot I_{Q-Q} \mathbf{n}_{Q-Q} + \sum_{m \in \mathcal{M}(c)} \frac{2}{3} \mathbf{U}_m \cdot I_{pp^+} \mathbf{n}_{pp^+}$

Normal and subcell forces definitions

- The semi-discrete equations on the specific volume, momentum and total energy successive moments, ensuring the GCL, write

$$\int_{\Omega_c} \rho^0 \frac{d}{dt} \left(\frac{1}{\rho} \right) \sigma_q^c dV = - \sum_{i=1}^{ntri} \int_{T_i^c} \mathbf{U} \cdot \mathbf{G} \nabla_X \sigma_q^c dV + \sum_{p \in \mathcal{P}(c)} \frac{1}{3} \mathbf{U}_p \cdot l_{pc}^q \mathbf{n}_{pc}^q + \sum_{m \in \mathcal{M}(c)} \frac{2}{3} \mathbf{U}_m \cdot l_{mc}^q \mathbf{n}_{mc}^q$$

$$\int_{\Omega_c} \rho^0 \frac{d \mathbf{U}}{dt} \sigma_q^c dV = \sum_{i=1}^{ntri} \int_{T_i^c} P \mathbf{G} \nabla_X \sigma_q^c dV - \sum_{p \in \mathcal{P}(c)} \frac{1}{3} \mathbf{F}_{pc}^q - \sum_{m \in \mathcal{M}(c)} \frac{2}{3} \mathbf{F}_{mc}^q$$

$$\int_{\Omega_c} \rho^0 \frac{d E}{dt} \sigma_q^c dV = \sum_{i=1}^{ntri} \int_{T_i^c} P \mathbf{U} \cdot \mathbf{G} \nabla_X \sigma_q^c dV - \sum_{p \in \mathcal{P}(c)} \frac{1}{3} \mathbf{U}_p \cdot \mathbf{F}_{pc}^q - \sum_{m \in \mathcal{M}(c)} \frac{2}{3} \mathbf{U}_m \cdot \mathbf{F}_{mc}^q$$

- $\bar{P} = P_h^c - Z_c (\bar{\mathbf{U}} - \mathbf{U}_h^c) \cdot \mathbf{n}$
- $\mathbf{F}_{pc}^q = P_h^c(\mathbf{X}_p, t) l_{pc}^q \mathbf{n}_{pc}^q - M_{pc}^q (\mathbf{U}_p - \mathbf{U}_h^c(\mathbf{X}_p, t)),$

where $M_{pc}^q = Z_c \left(l_{pc}^{-,q} \mathbf{n}_{pc}^{-,q} \otimes \mathbf{n}_{pc}^- + l_{pc}^{+,q} \mathbf{n}_{pc}^{+,q} \otimes \mathbf{n}_{pc}^+ \right)$

- $\mathbf{F}_{mc}^q = P_h^c(\mathbf{X}_m, t) l_{mc}^q \mathbf{n}_{mc}^q - M_{mc}^q (\mathbf{U}_m - \mathbf{U}_h^c(\mathbf{X}_m, t)),$
- where $M_{mc}^q = Z_c l_{mc}^q \mathbf{n}_{mc}^q \otimes \mathbf{n}_{mc}$

Nodal and midpoint solvers

- Fundamental identity on the cell Ω_c

$$\bar{P} = P_h^c - Z_c (\bar{\mathbf{U}} - \mathbf{U}_h^c) \cdot \mathbf{n},$$

where $Z_c = \rho_c a_c$ is the acoustic impedance

- Using this expression to calculate \mathbf{F}_{pc}^q leads to

$$\mathbf{F}_{pc}^q = P_h^c(\mathbf{X}_p, t) I_{pc}^q \mathbf{n}_{pc}^q - \mathbf{M}_{pc}^q (\mathbf{U}_p - \mathbf{U}_h^c(\mathbf{X}_p, t)),$$

where $\mathbf{M}_{pc}^q = Z_c \left(I_{pc}^{-,q} \mathbf{n}_{pc}^{-,q} \otimes \mathbf{n}_{pc}^- + I_{pc}^{+,q} \mathbf{n}_{pc}^{+,q} \otimes \mathbf{n}_{pc}^+ \right)$

- Regarding the midpoint subcell forces, \mathbf{F}_{mc}^q writes

$$\mathbf{F}_{mc}^q = P_h^c(\mathbf{X}_m, t) I_{mc}^q \mathbf{n}_{mc}^q - \mathbf{M}_{mc}^q (\mathbf{U}_m - \mathbf{U}_h^c(\mathbf{X}_m, t)),$$

where the \mathbf{M}_{mc}^q matrices are defined as

$$\mathbf{M}_{mc}^q = Z_c I_{mc}^q \mathbf{n}_{mc}^q \otimes \mathbf{n}_{mc}$$

Nodal solver

- To be conservative in total energy and momentum over the whole domain, we set the following sufficient conditions

$$\sum_{c \in \mathcal{C}(p)} \mathbf{F}_{pc} = \mathbf{0} \quad \text{and} \quad \mathbf{F}_{mL} + \mathbf{F}_{mR} = \mathbf{0},$$

where $\mathcal{C}(p)$ is the set of cells surrounding the p node, and Ω_L and Ω_R the two cells surrounding the midpoint m

- Thanks to $\sum_{c \in \mathcal{C}(p)} \mathbf{F}_{pc} = \mathbf{0}$, we finally have an explicit expression of \mathbf{U}_p

$$\left(\sum_{c \in \mathcal{C}(p)} M_{pc} \right) \mathbf{U}_p = \sum_{c \in \mathcal{C}(p)} [P_h^c(\mathbf{X}_p, t) I_{pc} \mathbf{n}_{pc} + M_{pc} \mathbf{U}_h^c(\mathbf{X}_p, t)]$$

where $M_{pc} = Z_c (I_{pc}^+ \mathbf{n}_{pc}^+ \otimes \mathbf{n}_{pc}^+ + I_{pc}^- \mathbf{n}_{pc}^- \otimes \mathbf{n}_{pc}^-)$ are positive semi-definite matrices with a physical dimension of a density times a velocity.

Midpoint solver

- The use of the condition $\mathbf{F}_{mL} + \mathbf{F}_{mR} = \mathbf{0}$ leads to

$$\mathbf{M}_m \mathbf{U}_m = \mathbf{M}_m \left(\frac{\mathcal{Z}_L \mathbf{U}_h^L(\mathbf{X}_m) + \mathcal{Z}_R \mathbf{U}_h^R(\mathbf{X}_m)}{\mathcal{Z}_L + \mathcal{Z}_R} \right) - \frac{P_h^R(\mathbf{X}_m) - P_h^L(\mathbf{X}_m)}{\mathcal{Z}_L + \mathcal{Z}_R} I_{pp^+} \mathbf{n}_{pp^+},$$

where the matrix $\mathbf{M}_m = \frac{1}{\mathcal{Z}_L} \mathbf{M}_{mL} = \frac{1}{\mathcal{Z}_R} \mathbf{M}_{mR}$ writes $\mathbf{M}_m = I_{pp^+} \mathbf{n}_{pp^+} \otimes \mathbf{n}_{pp^+}$

Approximate Riemann problem solution

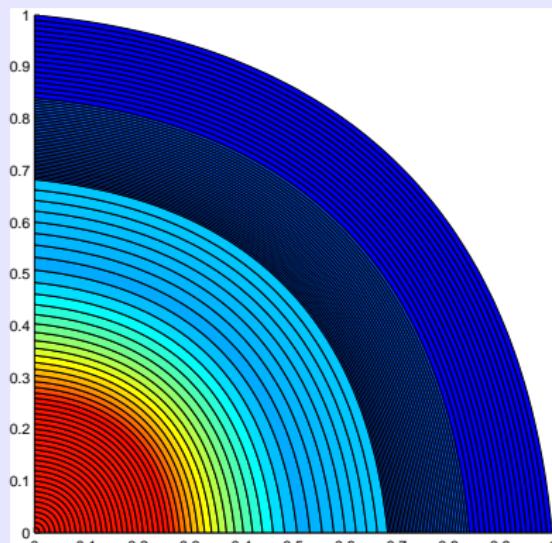
- $(\mathbf{U}_m \cdot \mathbf{n}_{pp^+}) = \left(\frac{\mathcal{Z}_L \mathbf{U}_h^L(\mathbf{X}_m) + \mathcal{Z}_R \mathbf{U}_h^R(\mathbf{X}_m)}{\mathcal{Z}_L + \mathcal{Z}_R} \right) \cdot \mathbf{n}_{pp^+} - \frac{P_h^R(\mathbf{X}_m) - P_h^L(\mathbf{X}_m)}{\mathcal{Z}_L + \mathcal{Z}_R}$

- Regarding the tangential contribution, we make the choice of

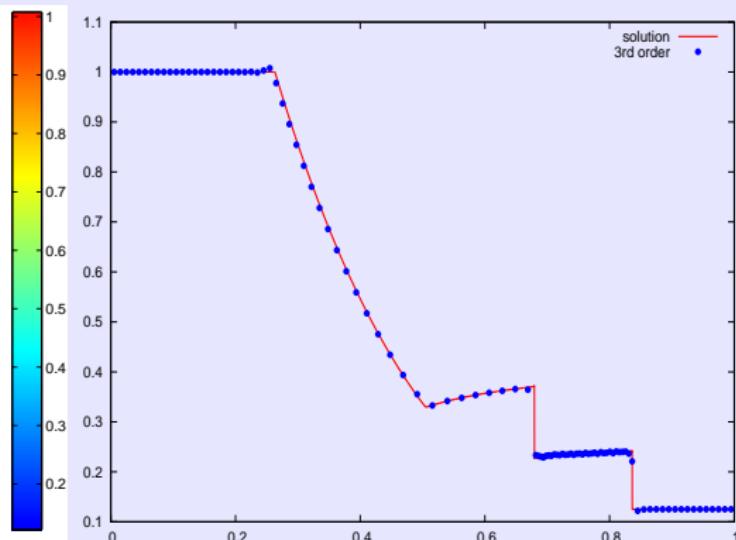
$$(\mathbf{U}_m \cdot \mathbf{t}_{pp^+}) = \left(\frac{\mathcal{Z}_L \mathbf{U}_h^L(\mathbf{X}_m) + \mathcal{Z}_R \mathbf{U}_h^R(\mathbf{X}_m)}{\mathcal{Z}_L + \mathcal{Z}_R} \right) \cdot \mathbf{t}_{pp^+}$$

Numerical results

Polar Sod shock tube problem



(a) Density map.

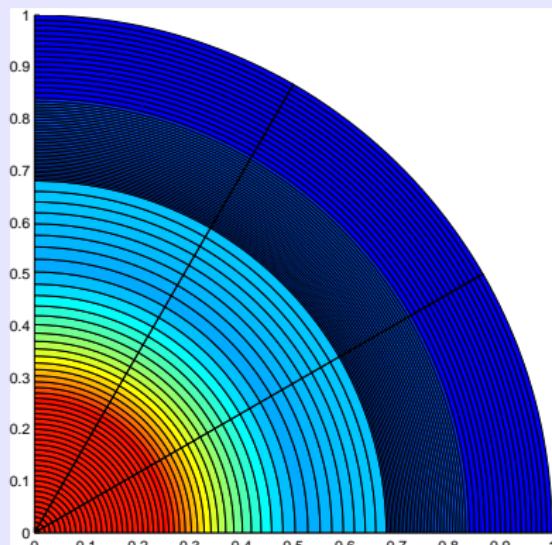


(b) Density profil.

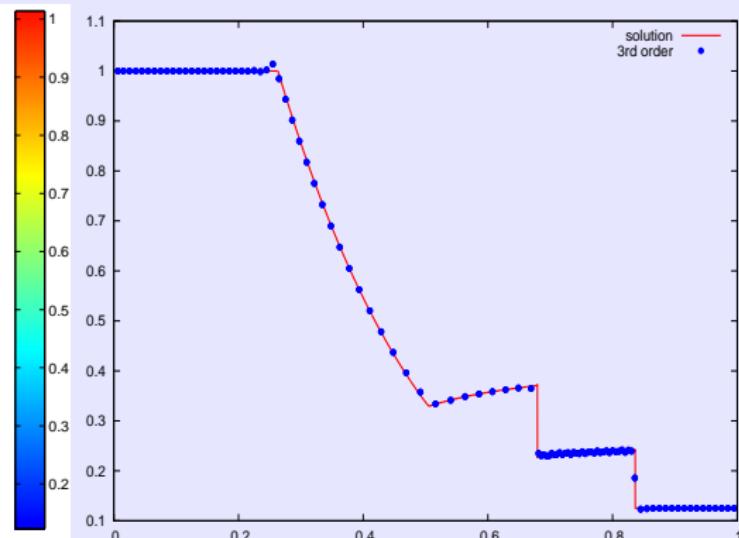
FIG.: Sod shock tube problem on a polar grid made of 100×1 cells.

Numerical results

Polar Sod shock tube problem



(a) Density map.

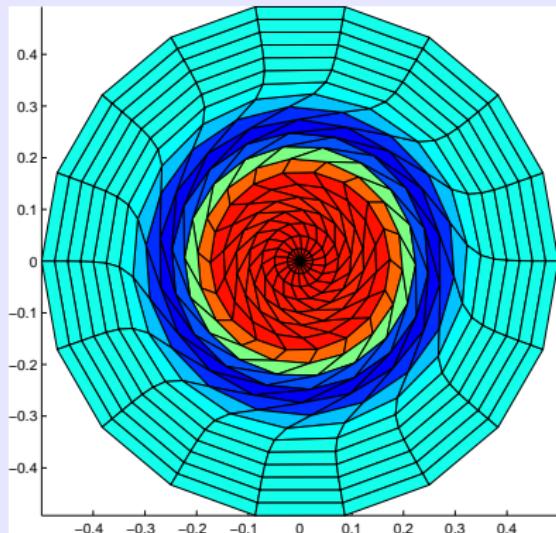


(b) Density profil.

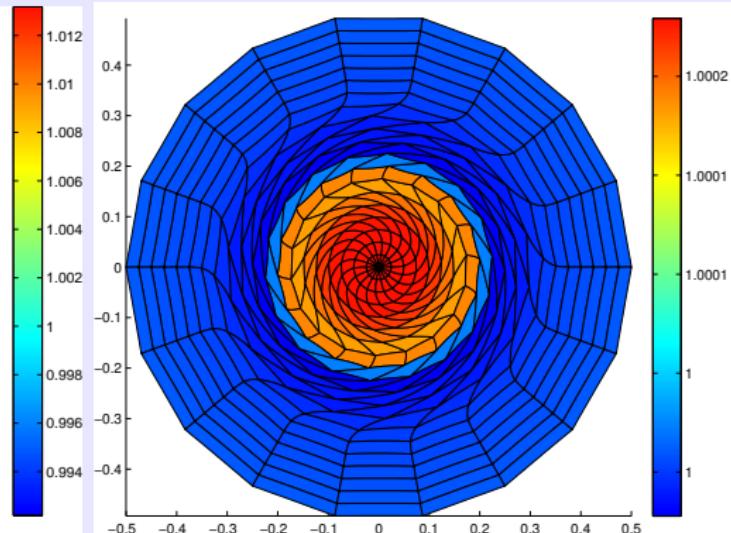
FIG.: Sod shock tube problem on a polar grid made of 100×3 cells.

Numerical results

Variant of the Gresho vortex problem



(a) Second-order scheme.

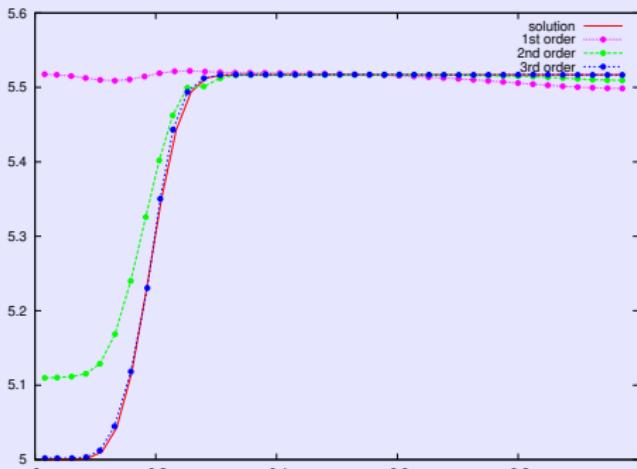


(b) Third-order.

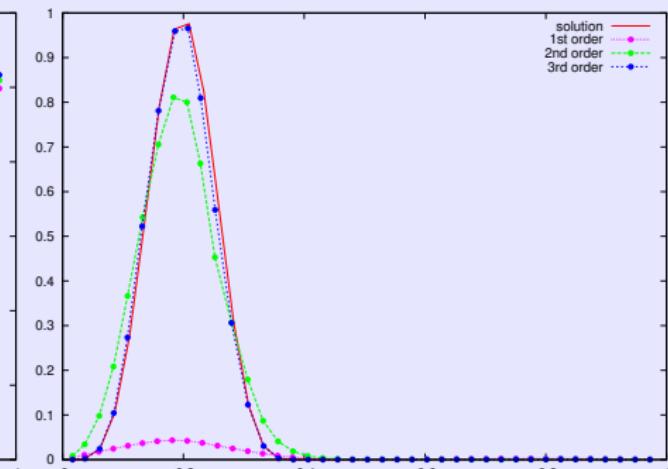
FIG.: Gresho variant problem on a polar grid defined in polar coordinates by $(r, \theta) \in [0, 1] \times [0, 2\pi]$, with 40×18 cells at $t = 0.36$: zoom of the desity map on the zone $(r, \theta) \in [0, 0.5] \times [0, 2\pi]$.

Numerical results

Variant of the Gresho vortex problem



(a) Pressure profil.

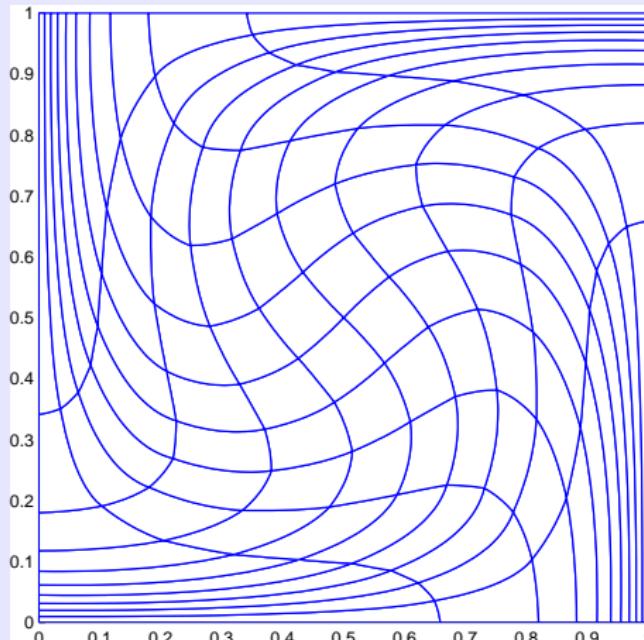


(b) Velocity profil.

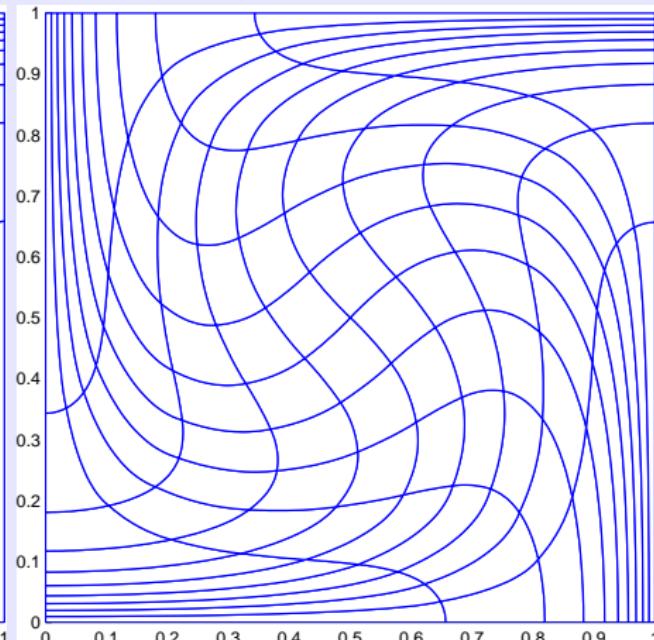
FIG.: Gresho variant problem on a polar grid defined in polar coordinates by $(r, \theta) \in [0, 1] \times [0, 2\pi]$, with 40×18 cells at $t = 0.36$.

Numerical results

Taylor-Green vortex problem



(a) Third-order scheme.



(b) Exact solution.

FIG.: Motion of a 10×10 Cartesian mesh through a T.-G. vortex, at $t = 0.75$.

Rate of convergence computed on the pressure in the case of the Taylor-Green vortex

	L_1		L_2		L_∞	
h	$E_{L_1}^h$	$q_{L_1}^h$	$E_{L_2}^h$	$q_{L_2}^h$	$E_{L_\infty}^h$	$q_{L_\infty}^h$
$\frac{1}{10}$	4.39E-3	3.00	7.73E-3	2.68	3.90E-2	1.93
$\frac{1}{20}$	5.50E-4	3.04	1.21E-3	3.10	1.03E-2	2.98
$\frac{1}{40}$	6.68E-5	2.91	1.40E-4	2.87	1.30E-3	2.66
$\frac{1}{80}$	8.90E-6	2.89	1.92E-5	2.83	2.11E-4	2.74
$\frac{1}{160}$	1.20E-6	-	2.70E-6	-	3.16E-5	-

TAB.: Third-order DG scheme without limitation at time $t = 0.6$.

	L_1		L_2		L_∞	
h	$E_{L_1}^h$	$q_{L_1}^h$	$E_{L_2}^h$	$q_{L_2}^h$	$E_{L_\infty}^h$	$q_{L_\infty}^h$
$\frac{1}{10}$	2.67E-4	2.96	3.36E-7	2.94	1.21E-3	2.86
$\frac{1}{20}$	3.43E-5	2.97	4.36E-5	2.96	1.66E-4	2.93
$\frac{1}{40}$	4.37E-6	2.99	5.59E-6	2.98	2.18E-5	2.96
$\frac{1}{80}$	5.50E-7	2.99	7.06E-7	2.99	2.80E-6	2.99
$\frac{1}{160}$	6.91E-8	-	8.87E-8	-	3.53E-7	-

TAB.: Third-order DG scheme with limitation at time $t = 0.1$.

Conclusions and perspectives

Conclusions

- We developed a 2nd and a 3rd order DG scheme for the 2D gas dynamics system in Lagrangian formalism with particular geometric consideration
- Numerical fluxes study
- Riemann invariants limitation
- GCL and Piola compatibility condition ensured by construction

Prospects

- High-order limitation on curved geometries
- Implementation of a 3rd order DG scheme on moving mesh
- Extension to ALE