

Local subcell monolithic DG/FV scheme

François Vilar

Institut Montpelliérain Alexander Grothendieck, Université de Montpellier, France

Séminaire **SISMA**



UNIVERSITÉ DE
MONTPELLIER



- 1 Introduction
- 2 DG as a subcell FV
- 3 Monolithic subcell DG/FV scheme
- 4 Entropy stabilities
- 5 Maximum principles
- 6 Conclusion

Scalar Conservation Law (SCL)

- $\partial_t u(\mathbf{x}, t) + \nabla_x \cdot \mathbf{F}(u(\mathbf{x}, t)) = 0, \quad (\mathbf{x}, t) \in \mathbb{R}^d \times [0, T]$
- $u(\mathbf{x}, 0) = u_0(\mathbf{x}), \quad \mathbf{x} \in \mathbb{R}^d$

Fundamental difficulty

- Even considering smooth flux function $\mathbf{F}(\cdot)$ and initial datum $u_0(\cdot)$ (C^∞ for instance), solution may become discontinuous in finite time
- Standard example: Burgers equation

$$\begin{cases} \partial_t u + \partial_x \left(\frac{1}{2} u^2 \right) = 0, & (x, t) \in [0, 1] \times \mathbb{R}^+, \\ u(x, 0) = \sin(2\pi x), & x \in [0, 1]. \end{cases}$$

Strong solution

- Local in time existence and uniqueness of a solution $u \in C^1(\mathbb{R}^d \times [0, t_c[)$

Formation of a discontinuity in finite time

Weak solution

$$\forall \psi \in \mathcal{C}_0^1(\mathbb{R}^d \times \mathbb{R}^+)$$

- $\int \int_{\mathbb{R}^d \times \mathbb{R}^+} \left(u \partial_t \psi + \mathbf{F}(u) \cdot \nabla_x \psi \right) d\mathbf{x} dt - \int_{\mathbb{R}^d} u_0(\mathbf{x}) \psi(\mathbf{x}, 0) d\mathbf{x} = 0$
- Non-uniqueness of a weak solution $u \in L_{loc}^\infty(\mathbb{R}^d \times \mathbb{R}^+)$

Entropy definition

- Let $u \in \mathcal{C}^1(\mathbb{R}^d \times [0, t_c[)$ be a strong solution
- η a strictly convex function is an entropy if $\exists \phi$ s.t.

$$\partial_t \eta(u) + \nabla_x \cdot \phi(u) = 0$$

Unique entropic weak solution

$$\forall \psi \in \mathcal{C}_0^1(\mathbb{R}^d \times \mathbb{R}^+)$$

- Let $u \in L_{loc}^\infty(\mathbb{R}^d \times \mathbb{R}^+)$ be a weak solution
- u is the unique entropic weak solution if it satisfies, for any pair (η, ϕ)

$$\int \int_{\mathbb{R}^d \times \mathbb{R}^+} \left(\eta(u) \partial_t \psi + \phi(u) \cdot \nabla_x \psi \right) d\mathbf{x} dt - \int_{\mathbb{R}^d} \eta(u_0(\mathbf{x})) \psi(\mathbf{x}, 0) d\mathbf{x} \geq 0$$

Formation of a discontinuity in finite time

Challenges of numerical discretization

- The weak solution we wish to approach may present areas of great regularity as well as discontinuities of great intensity
 - Smooth areas: high accuracy
 - Discontinuities: strong robustness and stability
 - Everywhere: additional mathematical or physical constraints
 - ↪ Maximum principle or positivity
 - ↪ Entropy inequalities
 - ↪ ...

Numerical schemes and discontinuous approximated solution

- Finite Volume FV
- (Weighted) Essentially Non-Oscillatory (W)-ENO
- Discontinuous Galerkin DG

Discontinuous Galerkin scheme

- Introduced by Reed and Hill in 1973 in the frame of the neutron transport
- Major development and improvements by B. Cockburn and C.-W. Shu

Procedure

- Local variational formulation
- Piecewise polynomial approximation of the solution in the cells
- Choice of the numerical fluxes
- Time integration

Advantages

- Natural extension of Finite Volume method
- Excellent analytical properties (L_2 stability, hp -adaptivity, ...)
- Extremely high accuracy (superconvergent for scalar conservation laws)
- Compact stencil (involve only face neighboring cells)

Scalar conservation law

- $\partial_t u(\mathbf{x}, t) + \nabla_x \cdot \mathbf{F}(u(\mathbf{x}, t)) = 0, \quad (\mathbf{x}, t) \in \omega \times [0, T]$
- $u(\mathbf{x}, 0) = u_0(\mathbf{x}), \quad \mathbf{x} \in \omega$

$(k+1)^{\text{th}}$ order semi-discretization

- $\{\omega_c\}_c$ a partition of ω , such that $\omega = \bigcup_c \omega_c$
- $u_h(\mathbf{x}, t)$ the numerical solution, such that $u_{h|\omega_c} = u_h^c \in \mathbb{P}^k(\omega_c)$

$$u_h^c(\mathbf{x}, t) = \sum_{m=1}^{N_k} u_m^c(t) \sigma_m^c(\mathbf{x})$$

- $\{\sigma_m^c\}_{m=1,\dots,N_k}$ a basis of $\mathbb{P}^k(\omega_c)$, with $N_k = \frac{(k+1)(k+2)}{2}$ in 2D.

Local variational formulation on ω_c

- $$\int_{\omega_c} \frac{\partial u_h^c}{\partial t} \psi \, dV = \oint_{\omega_c} \mathbf{F}(u_h^c) \cdot \nabla_x \psi \, dV - \oint_{\partial \omega_c} \psi \mathcal{F}_n \, dS, \quad \forall \psi \in \mathbb{P}^k(\omega_c)$$
- $\mathcal{F}_n = \mathcal{F}(u_h^c, u_h^\nu, \mathbf{n})$ numerical flux

Numerical example: solid body rotation

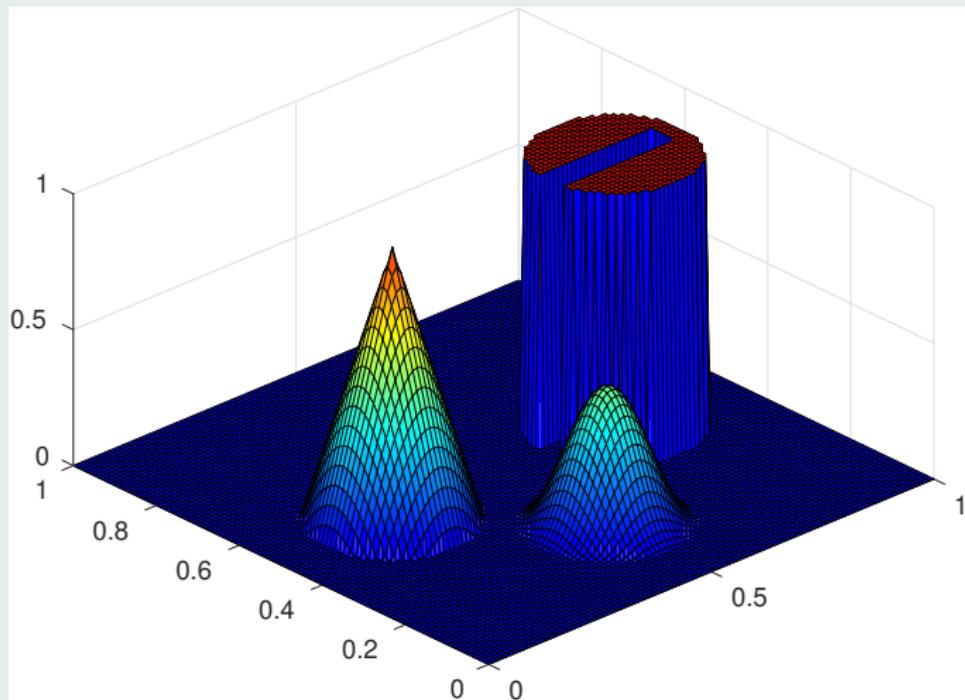
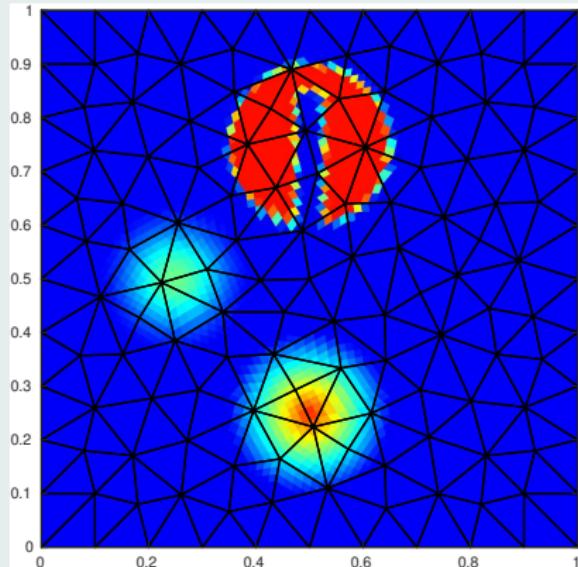
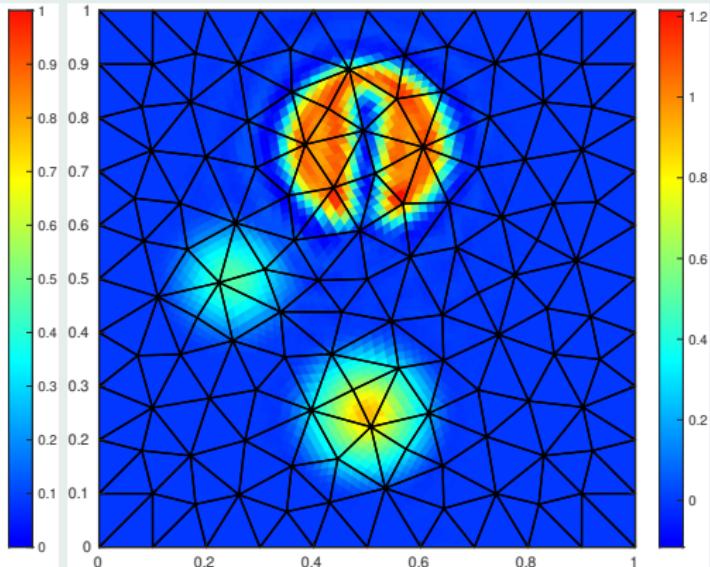


Figure: Rotation of composite signal: initial solution

Rotation of composite signal



(a) Initial solution



(b) Solution after one rotation

Figure: 6th-order DG scheme on 242 cells

Rotation of composite signal

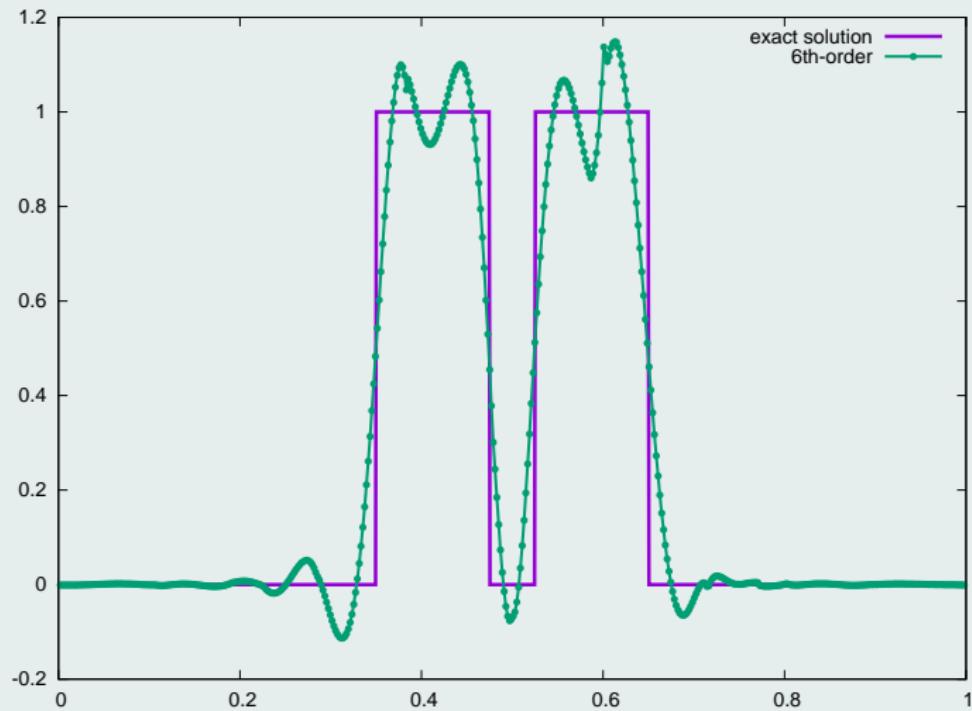


Figure: 6th-order DG scheme on 242 cells: profiles for $y = 0.75$

Spurious oscillations, aliasing and non-entropic behavior

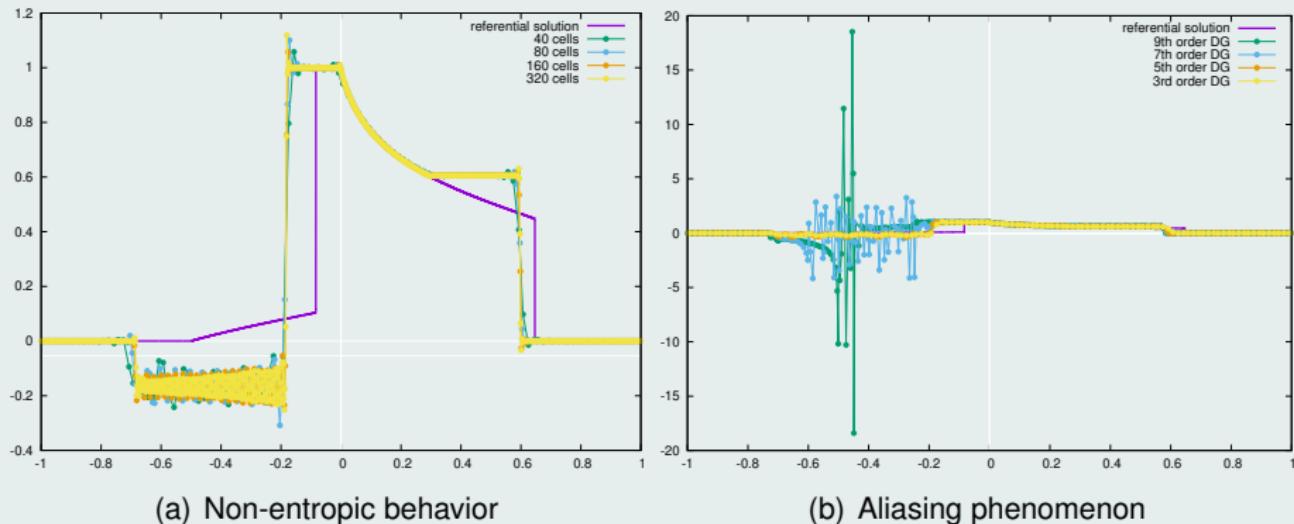


Figure : DG solutions for the Buckley non-convex flux case

Gibbs phenomenon and non-admissible solution

- High-order schemes leads to spurious oscillations near discontinuities
- Non-admissible solution potentially leading to a crash
- Vast literature of how prevent this phenomenon to happen:
 \implies *a priori* and *a posteriori* limitations

A priori limitation

- Artificial viscosity
- Slope/moment/hierarchical limiter
- ENO/WENO limiter
- Flux limiter and FCT schemes
- **Monolithic HO/LO schemes**

A posteriori limitation

- MOOD (“Multi-dimensional Optimal Order Detection”)
- A posteriori subcell correction

Admissible numerical solution

- Maximum principle / positivity preserving
- Limit the apparition of spurious oscillations
- Ensure a correct entropic behavior

Preserving high-accuracy and subcell resolution

Reduce the characteristic length of action

Methodology

Blend, at the subcell scale, high-order DG and 1st-order FV

-  *F.V., A Posteriori Correction of High-Order DG Scheme through Subcell Finite Volume Formulation and Flux Reconstruction.* JCP, 2018.
-  *F.V., Local subcell monolithic DG/FV convex property preserving scheme on unstructured grids and entropy consideration.* JCP, 2024.

1 Introduction

2 DG as a subcell FV

3 Monolithic subcell DG/FV scheme

4 Entropy stabilities

5 Maximum principles

6 Conclusion

DG as a subcell Finite Volume

- Rewrite DG scheme as a FV-like scheme on a subgrid

Cell subdivision into N_S subcells

$$N_S \geq N_k \equiv \#\mathbb{P}^k$$

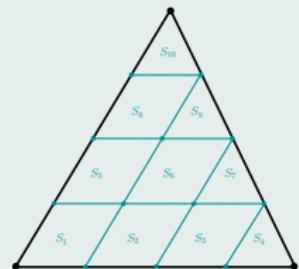
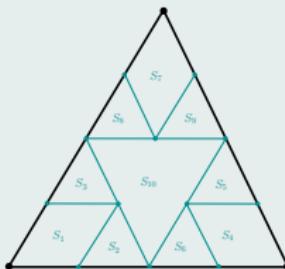
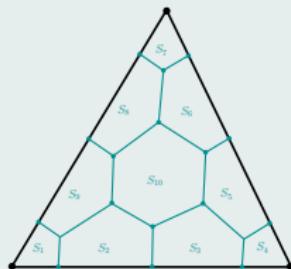


Figure : Examples of $N_S = N_k$ subdivision for \mathbb{P}^3 DG scheme on a triangle

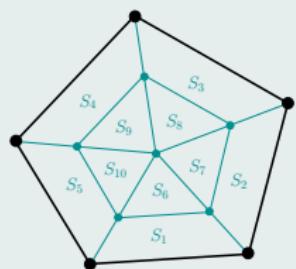
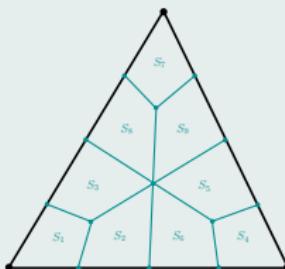
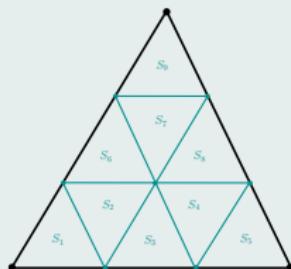


Figure : Examples of $N_S \geq N_k$ subdivision

DG schemes through residuals

- $\sum_{m=1}^{N_k} \frac{d u_m^c}{dt} \int_{\omega_c} \sigma_m \sigma_p dV = \oint_{\omega_c} \mathbf{F}(u_h^c) \cdot \nabla_x \sigma_p dV - \oint_{\partial \omega_c} \sigma_p \mathcal{F}_n dS, \quad \forall p \in [1, N_k]$

$$\implies M_c \frac{d \mathbf{U}_c}{dt} = \Phi_c$$

- $(\mathbf{U}_c)_m = u_m^c$ Solution moments
- $(M_c)_{mp} = \int_{\omega_c} \sigma_m \sigma_p dV$ Mass matrix
- $(\Phi_c)_m = \oint_{\omega_c} \mathbf{F}(u_h^c) \cdot \nabla_x \sigma_m dV - \oint_{\partial \omega_c} \sigma_m \mathcal{F}_n dS$ DG residuals

Subdivision and definition

- ω_c is subdivided into N_s subcells S_m^c
- Let us define $\bar{\psi}_m^c = \frac{1}{|S_m^c|} \int_{S_m^c} \psi dV$ the subcell mean value

Submean values

- $\bar{u}_m^c = \frac{1}{|S_m^c|} \sum_{q=1}^{N_k} u_q^c \int_{S_m^c} \sigma_q \, dV \quad \Rightarrow \quad \boxed{\bar{U}_c = P_c U_c}$

- $(\bar{U}_c)_m = \bar{u}_m^c$ Submean values

- $(P_c)_{mp} = \frac{1}{|S_m^c|} \int_{S_m^c} \sigma_p \, dV$ Projection matrix

$$\Rightarrow \boxed{\frac{d\bar{U}_c}{dt} = P_c M_c^{-1} \Phi_c}$$

Admissibility of the cell sub-partition into subcells

- $P_c^t P_c$ has to be non-singular

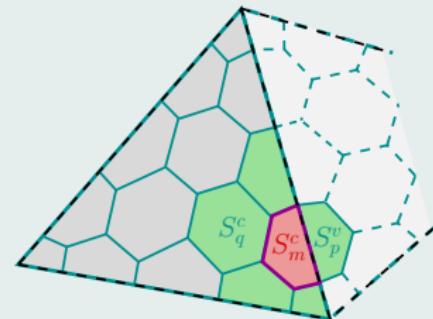
$$\Rightarrow \boxed{U_c = (P_c^t P_c)^{-1} P_c^t \bar{U}_c}$$
 Least square procedure

- If $N_s = N_k$, $\bar{U}_c = P_c U_c \iff U_c = P_c^{-1} \bar{U}_c$

Subcell Finite Volume: reconstructed fluxes

- Let us introduce the **reconstructed fluxes**

$$\frac{d \bar{U}_m^c}{dt} = -\frac{1}{|S_m^c|} \sum_{S_p^v \in \mathcal{V}_m^c} I_{mp} \widehat{F}_{pm}$$



- \mathcal{V}_m^c the set of face neighboring subcells of S_m^c
- We impose that on the boundary of cell ω_c , so for $S_p^v \notin \omega_c$

$$I_{mp} \widehat{F}_{pm} = \oint_{f_{mp}^c} \mathcal{F}_n dS \equiv \oint_{f_{mp}^c} \mathcal{F}(u_h^c, u_h^v, \mathbf{n}_{mp}^c) dS$$

- $\widetilde{\mathcal{V}}_m^c$ the set of face neighboring subcells of S_m^c belonging to ω_c
- Let A_c be the adjacency matrix such that

$$(A_c)_{mp} = \begin{cases} 1 & \text{if } S_p^v \in \widetilde{\mathcal{V}}_m^c \text{ with } m < p, \\ -1 & \text{if } S_p^v \in \widetilde{\mathcal{V}}_m^c \text{ with } m > p, \\ 0 & \text{if } S_p^v \notin \widetilde{\mathcal{V}}_m^c. \end{cases}$$

Subcell Finite Volume: reconstructed fluxes

- Let us introduce $D_c = \text{diag}(|S_1^c|, \dots, |S_{N_k}^c|)$ and $(B_c)_m = \int_{\partial S_m^c \cap \partial \omega_c} \mathcal{F}_n \, dS$
- Let \widehat{F}_c be the vector containing all the interior faces reconstructed fluxes

$$-A_c \widehat{F}_c = D_c P_c M_c^{-1} \Phi_c + B_c$$

Graph Laplacian technique

- $A_c \in \mathcal{M}_{N_s \times N_f^c}$ with N_f^c the number of interior faces
- $(L_c)_{mp} := (A_c A_c^\dagger)_{mp} = \begin{cases} |\widetilde{\mathcal{V}}_m^c| & \text{if } m = p, \\ -1 & \text{if } S_p^v \in \widetilde{\mathcal{V}}_m^c, \\ 0 & \text{otherwise.} \end{cases}$
- $L_c \mathbf{1} = \mathbf{0}$ where $\mathbf{1} = (1, \dots, 1)^t \in \mathbb{R}^{N_s}$
- $\Pi = \frac{1}{N_s} (\mathbf{1} \otimes \mathbf{1}) \in \mathcal{M}_{N_s}$

Graph Laplacian technique

- Let \mathcal{L}_c^{-1} be the pseudo-inverse of L_c such that

$$\mathcal{L}_c^{-1} = (L_c + \lambda \Pi)^{-1} - \frac{1}{\lambda} \Pi \quad \forall \lambda \neq 0$$

- Then, $\widehat{\mathbf{F}}_c$ is uniquely defined as following

$$\widehat{\mathbf{F}}_c = -A_c^t \mathcal{L}_c^{-1} (D_c P_c M_c^{-1} \Phi_c + \mathbf{B}_c)$$

- The only terms depending on the time are Φ_c and \mathbf{B}_c**
- Equivalently, the polynomial solution governing equation is given by

$$\frac{d \mathbf{U}_c}{dt} = -P_c^{-1} D_c^{-1} (A_c \widehat{\mathbf{F}}_c + \mathbf{B}_c)$$

remark

- This unique solution does exit since $(D_c P_c M_c^{-1} \Phi_c + \mathbf{B}_c) \cdot \mathbf{1} = 0$

Different cell subdivisions

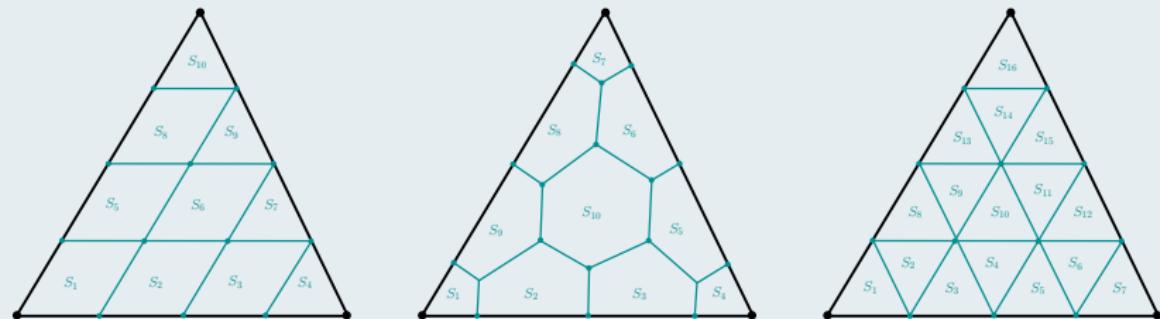
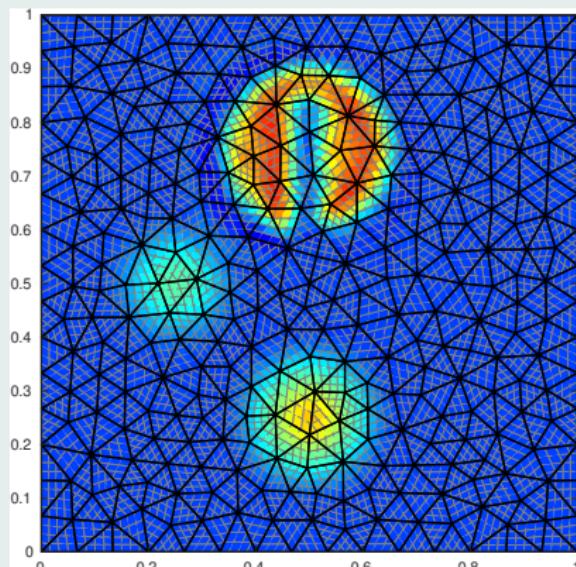


Figure: Examples of easily generalizable subdivisions for a triangle cell

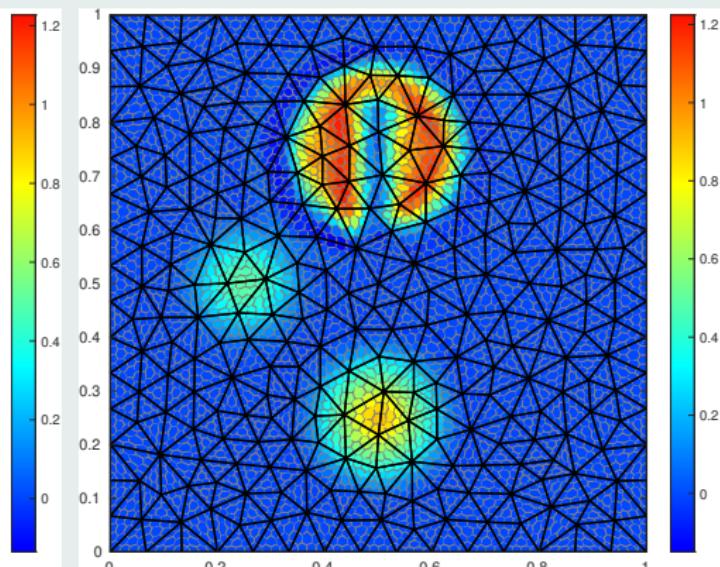
DG is DG

- Only the functional space matters
- The cell subdivision has no influence on the resulting scheme
- Even in the case where $N_s > N_k$

Rotation of a composite signal after one full rotation



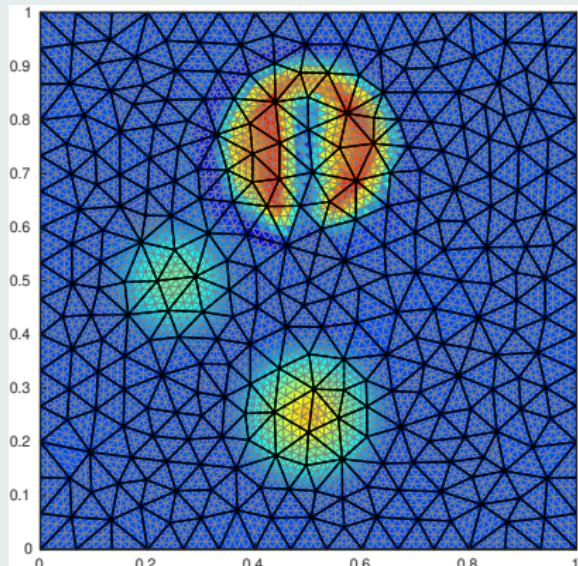
(a) Cartesian subdivision



(b) Polygonal subdivision

Figure: \mathbb{P}^3 reconstructed flux FV schemes on 576 cells

Rotation of a composite signal after one full rotation



(c) Triangular subdivision

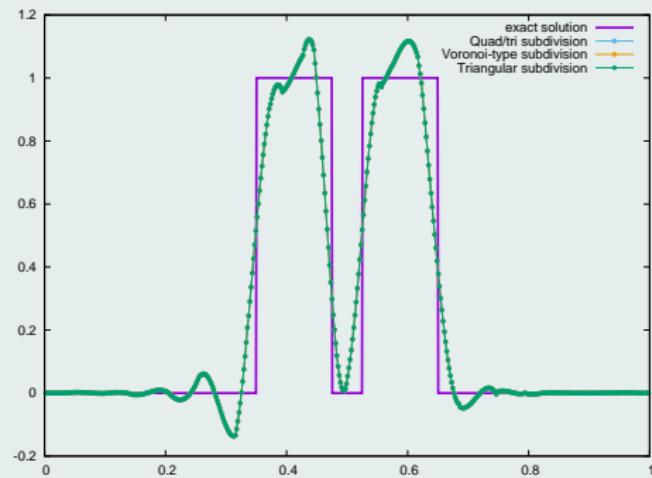
(d) Solution profiles for $y = 0.75$

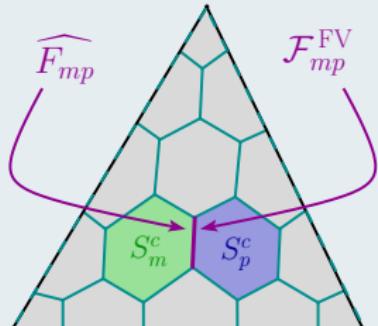
Figure: \mathbb{P}^3 reconstructed flux FV schemes on 576 cells

- 1 Introduction
- 2 DG as a subcell FV
- 3 Monolithic subcell DG/FV scheme
- 4 Entropy stabilities
- 5 Maximum principles
- 6 Conclusion

Blending low and high order fluxes

- Each face f_{mp}^c of each subcell S_m^c will be assigned two fluxes

$$\widetilde{F}_{mp} = \mathcal{F}_{mp}^{\text{FV}} + \underbrace{\theta_{mp}}_{\in [0,1]} \underbrace{\left(\widehat{F}_{mp} - \mathcal{F}_{mp}^{\text{FV}} \right)}_{\Delta F_{mp}}$$



- $\mathcal{F}_{mp}^{\text{FV}} := \mathcal{F}(\bar{u}_m^c, \bar{u}_p^v, \mathbf{n}_{mp})$ first-order subcell numerical flux
- \widehat{F}_{mp} high-order DG reconstructed flux
- The local subcell monolithic DG/FV then writes as follows**

$$\frac{d \bar{u}_m^c}{dt} = -\frac{1}{|S_m^c|} \sum_{S_p^v \in \mathcal{V}_m^c} I_{mp} \widetilde{F}_{mp}$$

Numerical flux

E-flux

$$\mathcal{F}(u^-, u^+, \mathbf{n}) = \frac{(\mathbf{F}(u^-) + \mathbf{F}(u^+))}{2} \cdot \mathbf{n} - \frac{\gamma(u^-, u^+, \mathbf{n})}{2} (u^+ - u^-)$$

- $\gamma(u^-, u^+, \mathbf{n}) \geq \max_{w \in I(u^-, u^+)} (|\mathbf{F}'(w) \cdot \mathbf{n}|)$
- $I(a, b) = [\min(a, b), \max(a, b)]$
- In the system case, we make use of either Rusanov, HLL, HLL-C, ...

Time integration

- For sake of simplicity, we focus on forward Euler (FE) time stepping, as SSP Runge-Kutta can be formulated as convex combinations of FE
- The semi-discrete scheme provided with FE time integration writes

$$\bar{u}_m^{c,n+1} = \bar{u}_m^{c,n} - \frac{\Delta t}{|S_m^c|} \sum_{S_p^v \in \mathcal{V}_m^c} l_{mp} \widetilde{F}_{mp}$$

Reformulation of the monolithic subcell scheme

- $\gamma_{mp} := \gamma(\bar{u}_m^{c,n}, \bar{u}_p^{v,n}, \mathbf{n}_{mp})$

1st-order FV dissipation coefficient

- $$\bar{u}_m^{c,n+1} = \bar{u}_m^{c,n} - \frac{\Delta t}{|S_m^c|} \left(\sum_{S_p^v \in \mathcal{V}_m^c} I_{mp} \widetilde{F_{mp}} \pm \sum_{S_p^v \in \mathcal{V}_m^c} I_{mp} \gamma_{mp} \bar{u}_m^{c,n} + \mathbf{F}(\bar{u}_m^{c,n}) \cdot \sum_{S_p^v \in \mathcal{V}_m^c} I_{mp} \mathbf{n}_{mp} \right)$$

$$= \left(1 - \frac{\Delta t}{|S_m^c|} \sum_{S_p^v \in \mathcal{V}_m^c} I_{mp} \gamma_{mp} \right) \bar{u}_m^{c,n}$$

$$+ \frac{\Delta t}{|S_m^c|} \sum_{S_p^v \in \mathcal{V}_m^c} I_{mp} \gamma_{mp} \left(\bar{u}_m^{c,n} - \frac{\widetilde{F_{mp}} - \mathbf{F}(\bar{u}_m^{c,n}) \cdot \mathbf{n}_{mp}}{\gamma_{mp}} \right)$$

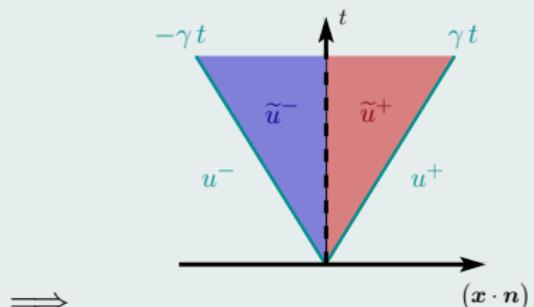
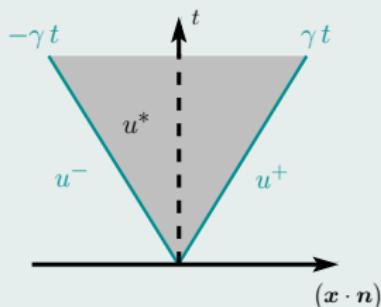
- $$\boxed{\bar{u}_m^{c,n+1} = \left(1 - \frac{\Delta t}{|S_m^c|} \sum_{S_p^v \in \mathcal{V}_m^c} I_{mp} \gamma_{mp} \right) \bar{u}_m^{c,n} + \frac{\Delta t}{|S_m^c|} \sum_{S_p^v \in \mathcal{V}_m^c} I_{mp} \gamma_{mp} \widetilde{U_{mp}} -}$$

• Convex combination under CFL condition

Blended Riemann intermediate states

- $$\widetilde{u_{mp}^-} = \bar{u}_m^{c,n} - \frac{\mathcal{F}_{mp}^{\text{FV}} - \mathbf{F}(\bar{u}_m^{c,n}) \cdot \mathbf{n}_{mp}}{\gamma_{mp}} - \theta_{mp} \frac{\Delta F_{mp}}{\gamma_{mp}}$$

$$= \underbrace{\frac{\bar{u}_m^{c,n} + \bar{u}_p^{v,n}}{2} - \frac{(\mathbf{F}(\bar{u}_p^{v,n}) - \mathbf{F}(\bar{u}_m^{c,n})) \cdot \mathbf{n}_{mp}}{2\gamma_{mp}}}_{u_{mp}^{*, \text{FV}}} - \theta_{mp} \frac{\Delta F_{mp}}{\gamma_{mp}}$$
- $$\widetilde{u_{mp}^\pm} = u_{mp}^{*, \text{FV}} \pm \theta_{mp} \frac{\Delta F_{mp}}{\gamma_{mp}} \implies u_{mp}^{*, \text{FV}} = \frac{1}{2} \left(\widetilde{u_{mp}^-} + \widetilde{u_{mp}^+} \right)$$



Admissible solution

Find the correct θ_{mp}

- G a convex admissible set where the solution has to remain in

$$\bullet \quad G = \left\{ u_0 \in [\alpha, \beta] \implies u \in [\alpha, \beta] \right\}$$

SCL

$$\bullet \quad G = \left\{ u = \begin{pmatrix} \rho \\ q \\ E \end{pmatrix}, \quad \rho > 0, p(u) > 0 \right\}$$

Euler

- A sufficient condition to ensure $\bar{u}_m^{c,n+1} \in G$ is that

$$\forall S_p^v \in \mathcal{V}_m^c, \quad \widetilde{u_{mp}}^\pm = u_{mp}^{*, \text{FV}} \pm \theta_{mp} \frac{\Delta F_{mp}}{\gamma_{mp}} \in G$$

- We want to prevent the code from crashing (apparition of NaN, ...)
- We want to prevent the apparition of spurious oscillations
- **We want to ensure discrete entropy inequalities (?)**
- We apply a local blending coefficients smoothening to avoid too stiff transition from high to low orders

1 Introduction

2 DG as a subcell FV

3 Monolithic subcell DG/FV scheme

4 Entropy stabilities

5 Maximum principles

6 Conclusion

Questions regarding entropy

- Can we find the θ_{mp} coefficients ensuring an entropy inequality?
- What do we mean by entropy inequality?
 - for one or any entropy?
 - at the discrete or semi-discrete time level?
 - at the cells or subcells space level?
- If we manage to ensure an entropy inequality, is it worth the effort?
 - in terms of accuracy
 - in terms of other critical properties to ensure, as positivity for instance
- Do we really need an entropy inequality to practically capture the entropic weak solution?

Definitions

- (η, ϕ) entropy - entropy flux
- $v(u) = \eta'(u)$ entropy variable
- $\psi(u) = v(u) \mathbf{F}(u) - \phi(u)$ entropy potential flux

Subcell entropy stability at the discrete level

for all (η, ϕ)

- if $\Delta F_{mp} \cdot (\bar{u}_p^{v,n} - \bar{u}_m^{c,n}) > 0$,

$$\theta_{mp} \leq \min \left(1, \frac{(\gamma_{mp} - \gamma_{\max}) (\bar{u}_p^{v,n} - \bar{u}_m^{c,n})}{2 \Delta F_{mp}} \right)$$

- $\gamma_{\max} := \max_{w \in I(\bar{u}_m^{c,n}, \bar{u}_p^{v,n})} (|\mathbf{F}'(w) \cdot \mathbf{n}_{mp}|)$

\implies **1st order scheme!**

Semi-discrete subcell entropy dissipation

for a given (η, ϕ)

- if $\Delta F_{mp} \cdot \left(v(\bar{u}_p^{v,n}) - v(\bar{u}_m^{c,n}) \right) > 0,$

$$\theta_{mp} \leq \min \left(1, \frac{\left(\frac{\psi(\bar{u}_p^{v,n}) - \psi(\bar{u}_m^{c,n})}{v(\bar{u}_p^{v,n}) - v(\bar{u}_m^{c,n})} \right) \cdot \mathbf{n}_{mp} - \mathcal{F}_{mp}^{\text{FV}}}{\Delta F_{mp}} \right)$$

\implies 2nd order scheme!



A. RUEDA-RAMÍREZ, B. BOLM, D. KUZMIN AND G. GASSNER, *Monolithic Convex Limiting for Legendre-Gauss-Lobatto Discontinuous Galerkin Spectral Element Methods*. Commun. Appl. Math. Comput., 2024.

Sub-resolution basis functions

$N_s = N_k$

- Let $\{\varphi_m^c\}_m$ be the sub-resolution basis function such that, $\forall \psi \in \mathbb{P}^k(\omega_c)$

$$\int_{\omega_c} \varphi_m \psi \, dV = \int_{S_m^c} \psi \, dV$$

- Then, given $v_h^c \in \mathbb{P}^k(\omega_c)$, it writes

$$v_h^c = \sum_{m=1}^{N_k} \underline{v}_m^c \varphi_m^c$$

Semi-discrete cell entropy dissipation

for a given (η, ϕ)

- $\Delta \eta_c := \frac{d}{dt} \oint_{\omega_c} \eta(u_h^c) \, dV = \oint_{\omega_c} v(u_h^c) \partial_t u_h^c \, dV = \int_{\omega_c} v_h^c \partial_t u_h^c \, dV$
- $v_h^c = \sum_{m=1}^{N_k} \underline{v}_m^c \varphi_m^c$ L^2 projection of $v(u_h^c)$ onto \mathbb{P}^k
- $\Delta \eta_c = \sum_{m=1}^{N_k} \underline{v}_m^c \int_{\omega_c} \varphi_m^c \partial_t u_h^c \, dV = \sum_{m=1}^{N_k} |S_m^c| \underline{v}_m^c \frac{d \bar{u}_m^c}{dt} = - \sum_{m=1}^{N_k} \underline{v}_m^c \sum_{S_p^v \in \mathcal{V}_m^c} I_{mp} \widetilde{F}_{mp}$

Continuous Knapsack problem

- The sufficient condition rewrites as

$$\mathbf{C}_c \cdot \boldsymbol{\Theta}_c \leq D_c$$

- $\boldsymbol{\Theta}$ contains all the subcells' faces, \mathbf{C}_c the vector defined as

$$\mathbf{C}_{mp}^c = \begin{cases} \oint_{f_{mp}} \left((\underline{v}(u_h^c) - \underline{v}_m^c) \mathcal{F}_n - (\psi(u_h^c) - \psi(u(\underline{v}_m^c))) \cdot \mathbf{n}_{mp} \right) dS, & \forall f_{mp} \subset \partial\omega_c \\ I_{mp} (\underline{v}_p^c - \underline{v}_m^c) \Delta F_{mp}, & \text{otherwise} \end{cases}$$

- $D_c = \sum_{f_{mp} \in \hat{\mathfrak{f}}_c} I_{mp} \psi(u(\underline{v}_m^c)) \cdot \mathbf{n}_{mp} - \sum_{f_{mp} \in \check{\mathfrak{f}}_c} I_{mp} (\underline{v}_p^c - \underline{v}_m^c) \mathcal{F}_{mp}^{FV}$
- $\mathcal{F}_{mp}^{FV} = \mathcal{F}(u(\underline{v}_m^c), u(\underline{v}_p^c), \mathbf{n}_{mp})$ **modified FV numerical flux**
- Because $D_c \geq 0$ the Knapsack problem is indeed solvable**



Y. LIN AND J. CHAN, High order entropy stable discontinuous Galerkin spectral element methods through subcell limiting. JCP, 2024.

Greedy algorithm

- Find $\mathbf{0} \leq \Theta \leq \Theta^e \leq \mathbf{1}$ maximizing $\sum \theta_{mp}$ such that
- θ_{mp}^e being any given supplementary constraint

$$\mathbf{C}_c \cdot \Theta_c \leq D_c$$

High-order accuracy preservation

- Substituting u_h by a smooth solution u
- h_c is the diameter of cell ω_c
- Then, we have that

$$\mathbf{C}_c \cdot \mathbf{1} - D_c = |\omega_c| \mathcal{O}(h_c^{k+1})$$

- This finally implies that

$$\widetilde{F}_{mp} = \widehat{F}_{mp} + (\theta_{mp} - 1) \Delta F_{mp} = \widehat{F}_{mp} + \mathcal{O}(h_c^{k+2})$$

Linear advection of a sinus

$$\eta(u) = \frac{1}{2} u^2$$

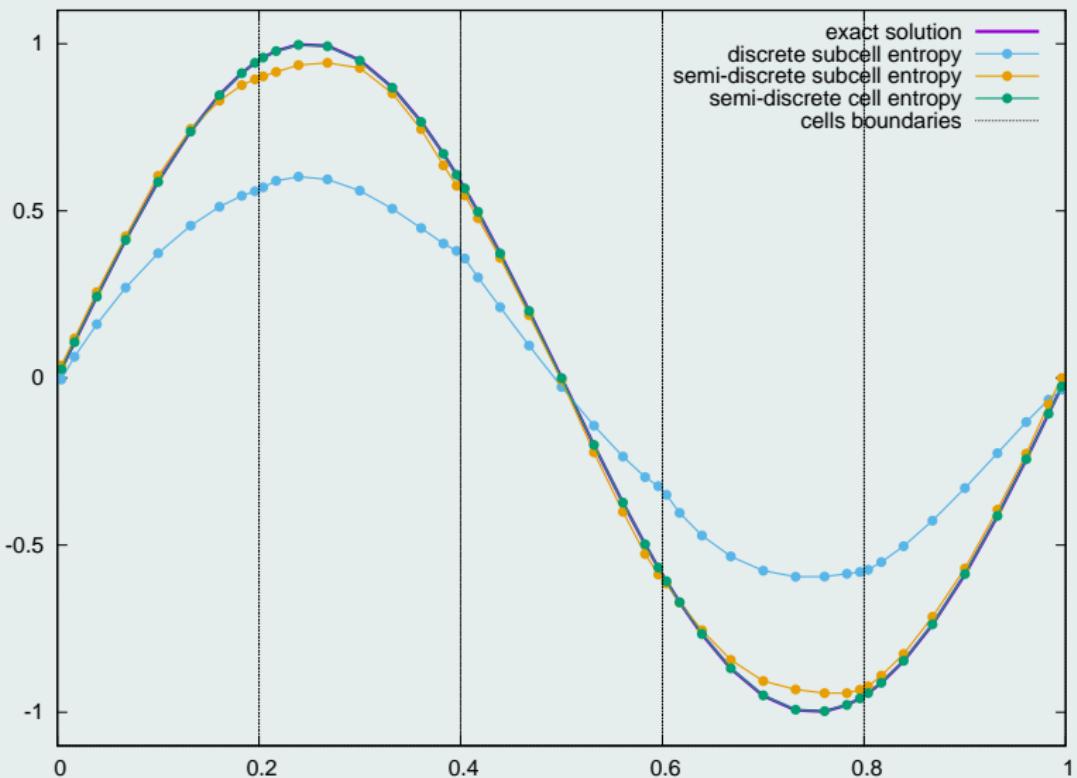


Figure: \mathbb{P}^8 -DG/FV with entropy stabilities on 5 cells

Linear advection of a composite signal

$$\eta(u) = \frac{1}{2} u^2$$

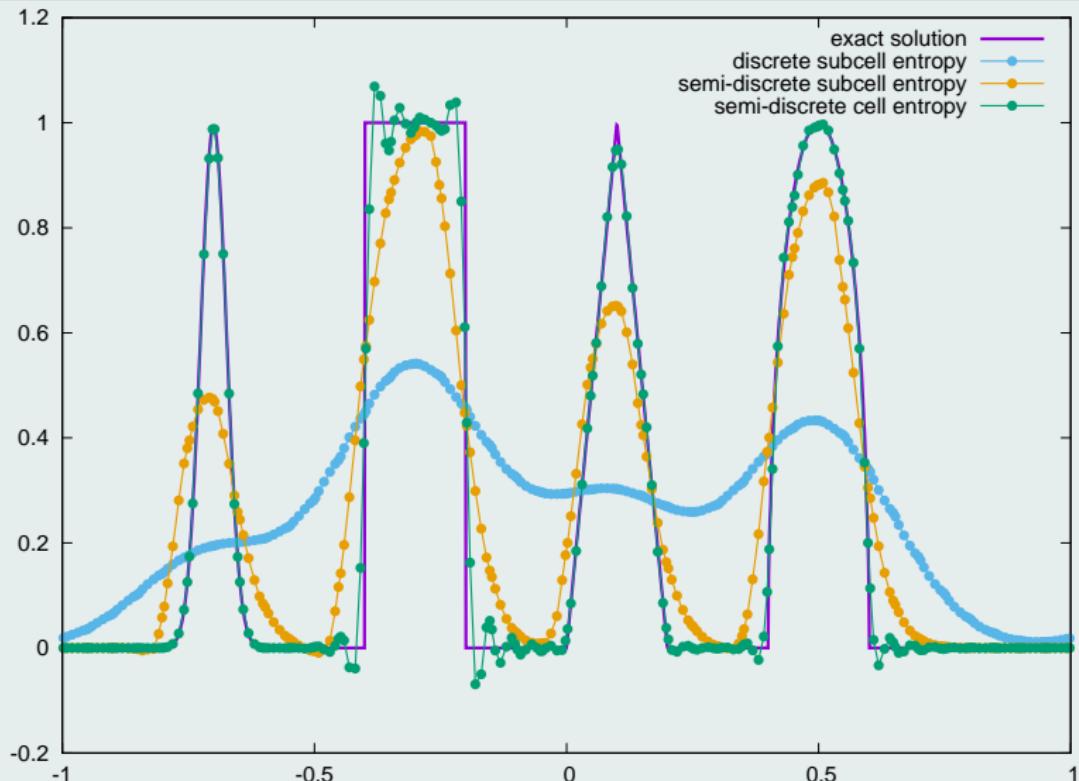


Figure: \mathbb{P}^5 -DG/FV with entropy stabilities on 40 cells

Linear advection of a composite signal $\eta(u) = |u - k_e|^{1+\epsilon} / (1 + \epsilon)$

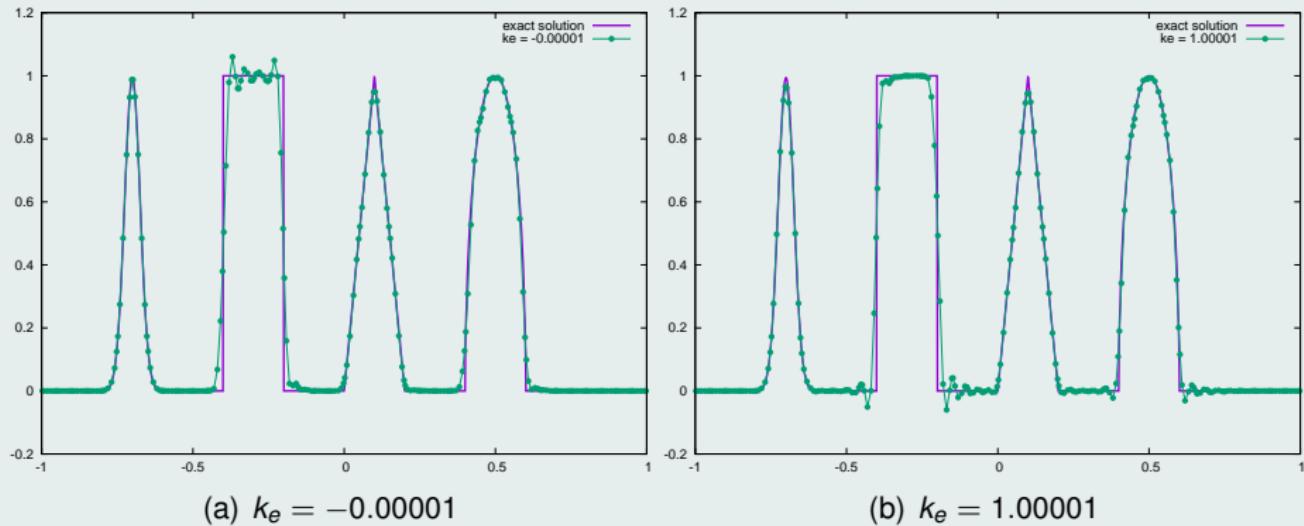


Figure: \mathbb{P}^5 -DG/FV submean values on 40 cells: $\epsilon = 0.25$

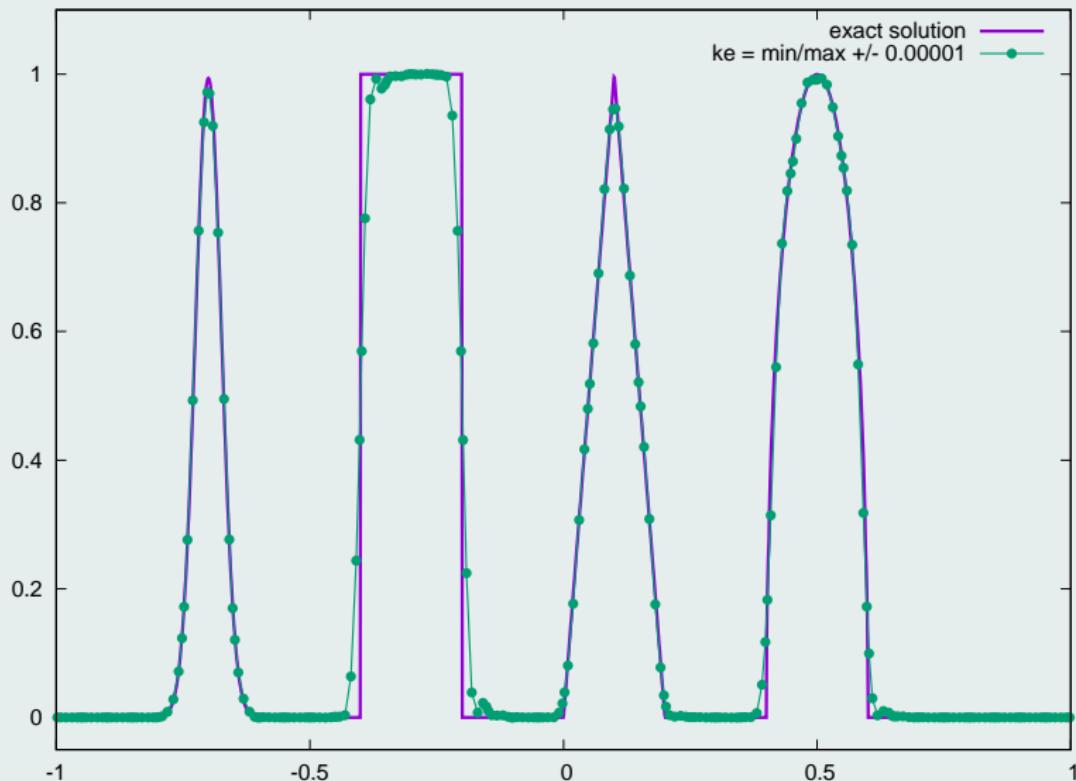
Linear advection of a composite signal $\eta(u) = |u - k_e|^{1+\epsilon} \backslash (1 + \epsilon)$ 

Figure: \mathbb{P}^5 -DG/FV on 40 cells: $k_e = -0.00001$ and $k_e = 1.00001$

Non-linear non-convex flux Buckley case

80 cells

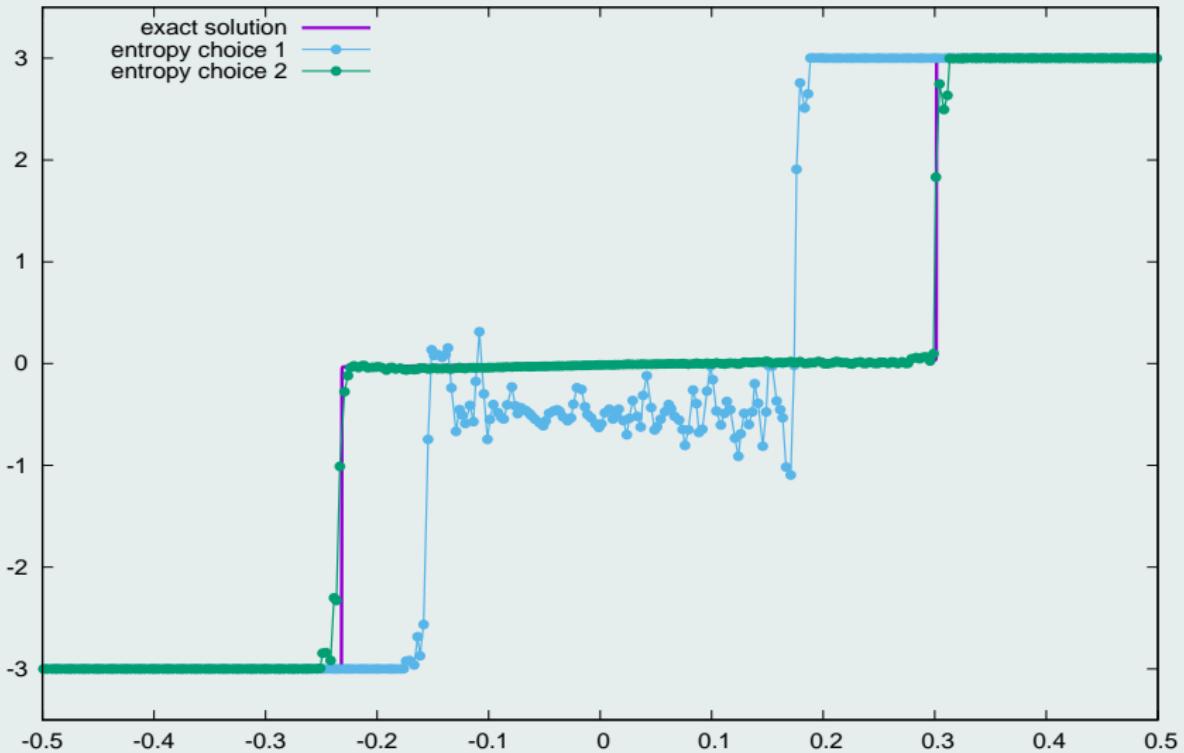


Figure: \mathbb{P}^3 -DG/FV submean values: $\eta_1(u) = \frac{1}{2}u^2$ and $\eta_2(u) = \int \text{atan}(20u) du$

Non-linear non-convex flux Buckley case

80 cells

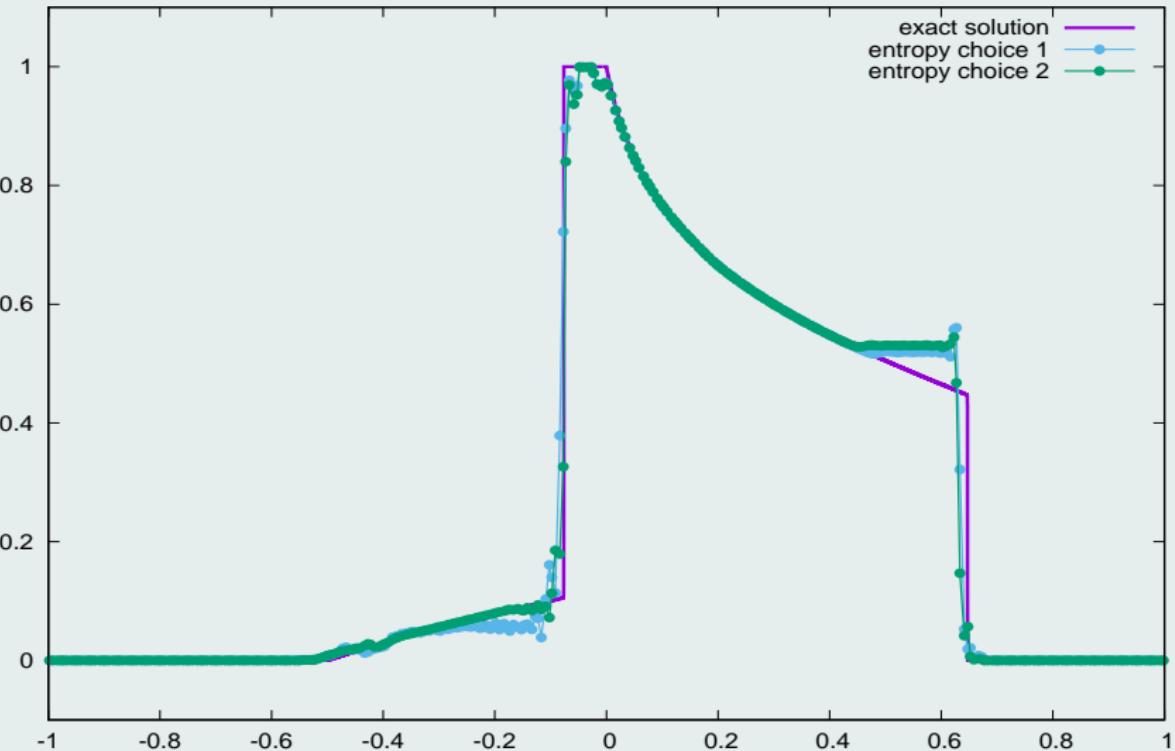


Figure: \mathbb{P}^3 -DG/FV submean values: $\eta_1(u) = \frac{1}{2}u^2$ and $\eta_2(u) = \int \text{atan}(20u) du$

KPP non-convex flux problem

$$\eta(u) = \frac{1}{2}u^2$$

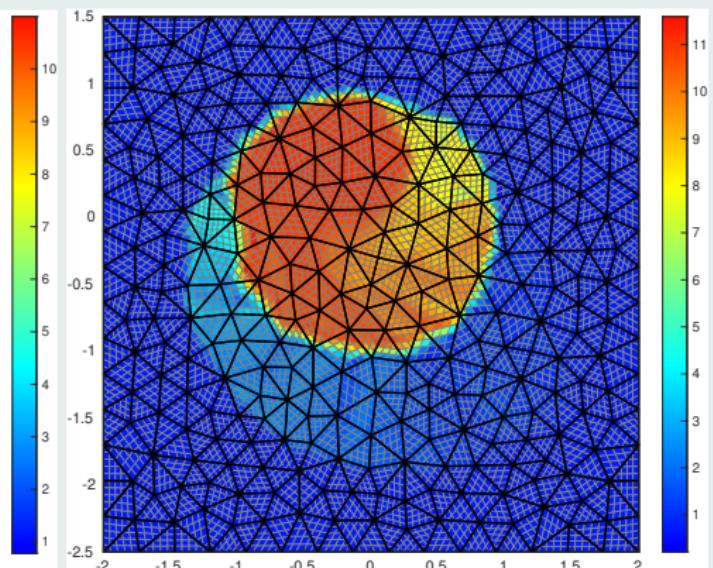
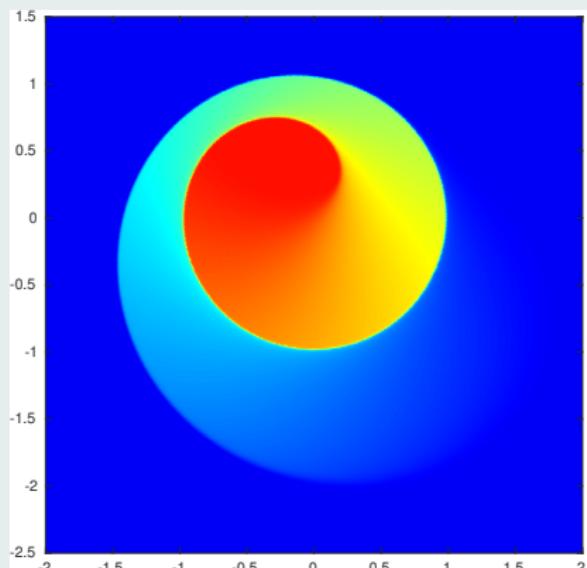


Figure: \mathbb{P}^4 -DG/FV entropic scheme: non-entropic solution

1D modified Sod shock tube problem

$$\eta(U) = -\rho \log(p/\rho^\gamma)$$

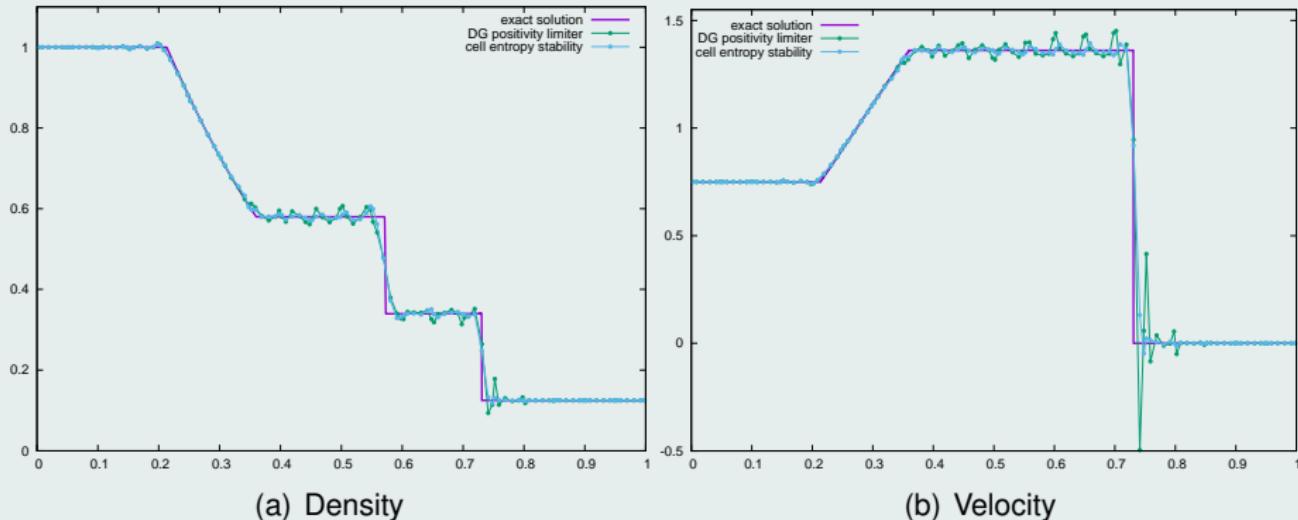


Figure: \mathbb{P}^5 pure DG with positivity limiter and \mathbb{P}^5 -DG/FV monolithic scheme with cell entropy stability on 20 cells

Questions regarding entropy → pieces of answer

- Can we find θ_{mp} coefficients ensuring an entropy inequality?

↪ Yes!

- What do we mean by entropy inequality, and is it worth the effort?

- for any entropy, at the discrete time level and for any subcell

↪ 1st-order

- for a given entropy, at the semi-discrete time level for any subcells

↪ 2nd-order

- for a given entropy, at the semi-discrete time level for any cells

↪ ($k + 1$)th-order

↪ $\mathcal{F}_{mp}^{FV} = \mathcal{F}(u(\underline{v}_m^c), u(\underline{v}_p^v), n_{mp})$



- Do we absolutely need entropy stability while aiming for high order?

↪ Unclear... ⇒ GMP and LMP + relaxation

- 1 Introduction
- 2 DG as a subcell FV
- 3 Monolithic subcell DG/FV scheme
- 4 Entropy stabilities
- 5 Maximum principles
- 6 Conclusion

Global maximum principle

$$\bar{U}_m^{c,n+1} \in [\alpha, \beta]$$

$$\theta_{mp} \leq \min \left(1, \left| \frac{\gamma_{mp}}{\Delta F_{mp}} \right| \min \left(\beta - u_{mp}^{*,\text{FV}}, u_{mp}^{*,\text{FV}} - \alpha \right) \right)$$

Local maximum principle

$$\bar{U}_m^{c,n+1} \in [\alpha_m^c, \beta_m^c]$$

- $\alpha_m^c := \min_{S_q^w \in \mathcal{N}(S_m^c)} (\bar{U}_q^{w,n})$ and $\beta_m^c := \max_{S_q^w \in \mathcal{N}(S_m^c)} (\bar{U}_q^{w,n})$

$$\theta_{mp} \leq \min \left(1, \left| \frac{\gamma_{mp}}{\Delta F_{mp}} \right| \begin{cases} \min \left(\beta_p^v - u_{mp}^{*,\text{FV}}, u_{mp}^{*,\text{FV}} - \alpha_m^c \right) & \text{if } \Delta F_{mp} > 0 \\ \min \left(\beta_m^c - u_{mp}^{*,\text{FV}}, u_{mp}^{*,\text{FV}} - \alpha_p^v \right) & \text{if } \Delta F_{mp} < 0 \end{cases} \right)$$

- The wider set $\mathcal{N}(S_m^c)$ is, the softer this local maximum principle will be
- **Smooth extrema relaxation to preserve accuracy**

Burgers equation

$$u_0(x, y) = \sin(2\pi(x + y))$$

(a) Solution submean values

(b) Blending coefficients

Figure: \mathbb{P}^5 -DG/FV scheme with GMP and relaxed-LMP on 242 cells

KPP non-convex flux problem

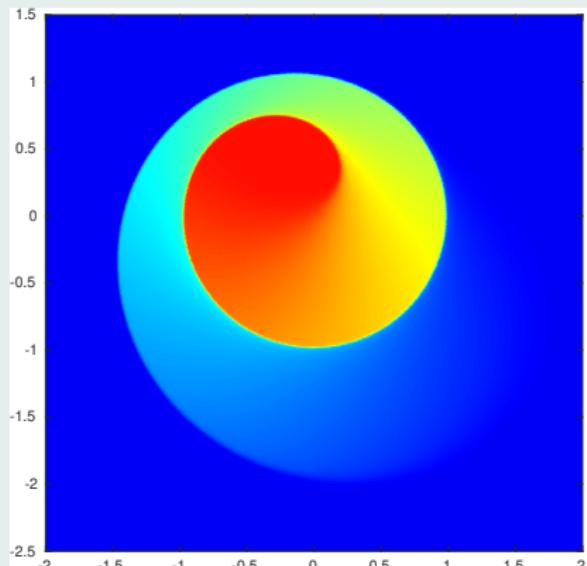
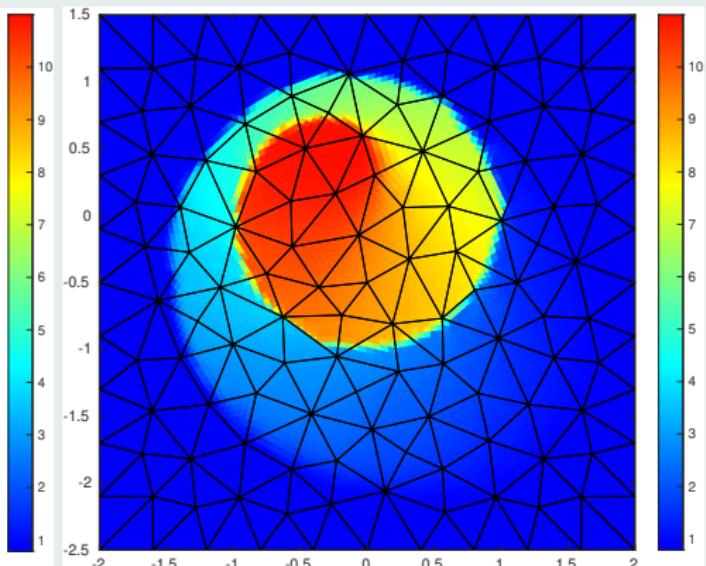
(a) 1th-order FV on 209184 cells(b) \mathbb{P}^7 -DG/FV on 272 cells

Figure: \mathbb{P}^7 -DG/FV scheme with GMP and relaxed-LMP

Non-linear Euler compressible gas dynamics equations

- $\partial_t \mathbf{V} + \nabla_x \cdot \mathbf{F}(\mathbf{V}) = \mathbf{0}$

- $\mathbf{V} = \begin{pmatrix} \rho \\ \mathbf{q} \\ E \end{pmatrix}$ conservative variables

- $\mathbf{F}(\mathbf{V}) = \begin{pmatrix} \mathbf{q} \\ \frac{\mathbf{q} \otimes \mathbf{q}}{\rho} + p I_d \\ (E + p) \frac{\mathbf{q}}{\rho} \end{pmatrix}$ flux function

- $p := p(\mathbf{V}) = (\gamma - 1) \left(E - \frac{1}{2} \frac{\|\mathbf{q}\|^2}{\rho} \right)$ equation of state

Monolithic subcell DG/FV scheme property

- Positivity of the density and internal energy, at the subcell scale

Definitions

- $\widetilde{\mathbf{F}}_{mp} := \widehat{\mathcal{F}}_{mp}^{\text{FV}} + \Theta_{mp} \underbrace{\left(\widehat{\mathbf{F}}_{mp} - \widehat{\mathcal{F}}_{mp}^{\text{FV}} \right)}_{\Delta\mathbf{F}_{mp}}$ convex blended flux
- $\Theta_{mp} = \begin{pmatrix} \theta_{mp}^\rho & \mathbf{0} & \mathbf{0} \\ \mathbf{0} & \theta_{mp}^q & \mathbf{0} \\ \mathbf{0} & \mathbf{0} & \theta_{mp}^E \end{pmatrix}$
- $\mathcal{F}_{mp}^{\text{FV}} := \mathcal{F} \left(\overline{\mathbf{V}}_m^{c,n}, \overline{\mathbf{V}}_p^{v,n}, \mathbf{n}_{mp} \right)$ Global L-F, Rusanov, HLL(C), ...

Positivity of the density

- $\theta_{mp}^\rho = \theta_{mp}^{(1)} \theta_{mp}^{(2)}$

$$\theta_{mp}^{(1)} \leq \min \left(1, \left| \frac{\gamma_{mp}}{\Delta F_{mp}^\rho} \right| \rho_{mp}^{*, \text{FV}} \right)$$

Positivity of the internal energy

- $A_{mp} = \frac{1}{(\gamma_{mp})^2} \left(\frac{1}{2} \|\Delta \mathbf{F}_{mp}^q\|^2 - \theta_{mp}^{(1)} \Delta F_{mp}^\rho \Delta F_{mp}^E \right)$
- $B_{mp} = \frac{1}{\gamma_{mp}} \left(\mathbf{q}_{mp}^{*, \text{FV}} \cdot \Delta \mathbf{F}_{mp}^q - \rho_{mp}^{*, \text{FV}} \Delta F_{mp}^E - \theta_{mp}^{(1)} E_{mp}^{*, \text{FV}} \Delta F_{mp}^\rho \right)$
- $M_{mp} = \rho_{mp}^{*, \text{FV}} E_{mp}^{*, \text{FV}} - \frac{1}{2} \|\mathbf{q}_{mp}^{*, \text{FV}}\|^2$

$$\theta_{mp}^{(2)} \leq \min \left(1, \frac{M_{mp}}{|B_{mp}| + \max(0, A_{mp})} \right)$$

- $\theta_{mp}^\rho = \theta_{mp}^{\rho(1)} \theta_{mp}^{(2)}$, $\theta_{mp}^{q^x} = \theta_{mp}^{(2)}$, $\theta_{mp}^{q^y} = \theta_{mp}^{(2)}$, $\theta_{mp}^E = \theta_{mp}^{(2)}$



A. RUEDA-RAMÍREZ, B. BOLM, D. KUZMIN AND G. GASSNER, *Monolithic Convex Limiting for Legendre-Gauss-Lobatto Discontinuous Galerkin Spectral Element Methods.* Commun. Appl. Math. Comput., 2024.

LMP

$$\bar{V}_m^{c,n+1} \in [\alpha_m^c, \beta_m^c]$$

- $v \in \{\rho, q^x, q^y, E\}$ conservative variable
- $\alpha_m^c := \min_{S_q^w \in \mathcal{N}(S_m^c)} (\bar{v}_q^{w,n}, v_{mq}^{*,\text{FV}})$ and $\beta_m^c := \max_{S_q^w \in \mathcal{N}(S_m^c)} (\bar{u}_q^{w,n}, v_{mq}^{*,\text{FV}})$

$$\theta_{mp} \leq \min \left(1, \left| \frac{\gamma_{mp}}{\Delta F_{mp}} \right| \begin{cases} \min (\beta_p^v - u_{mp}^{*,\text{FV}}, u_{mp}^{*,\text{FV}} - \alpha_m^c) & \text{if } \Delta F_{mp} > 0 \\ \min (\beta_m^c - u_{mp}^{*,\text{FV}}, u_{mp}^{*,\text{FV}} - \alpha_p^v) & \text{if } \Delta F_{mp} < 0 \end{cases} \right)$$

- Smooth extrema relaxation to preserve accuracy

Sod shock tube test case

10 cells

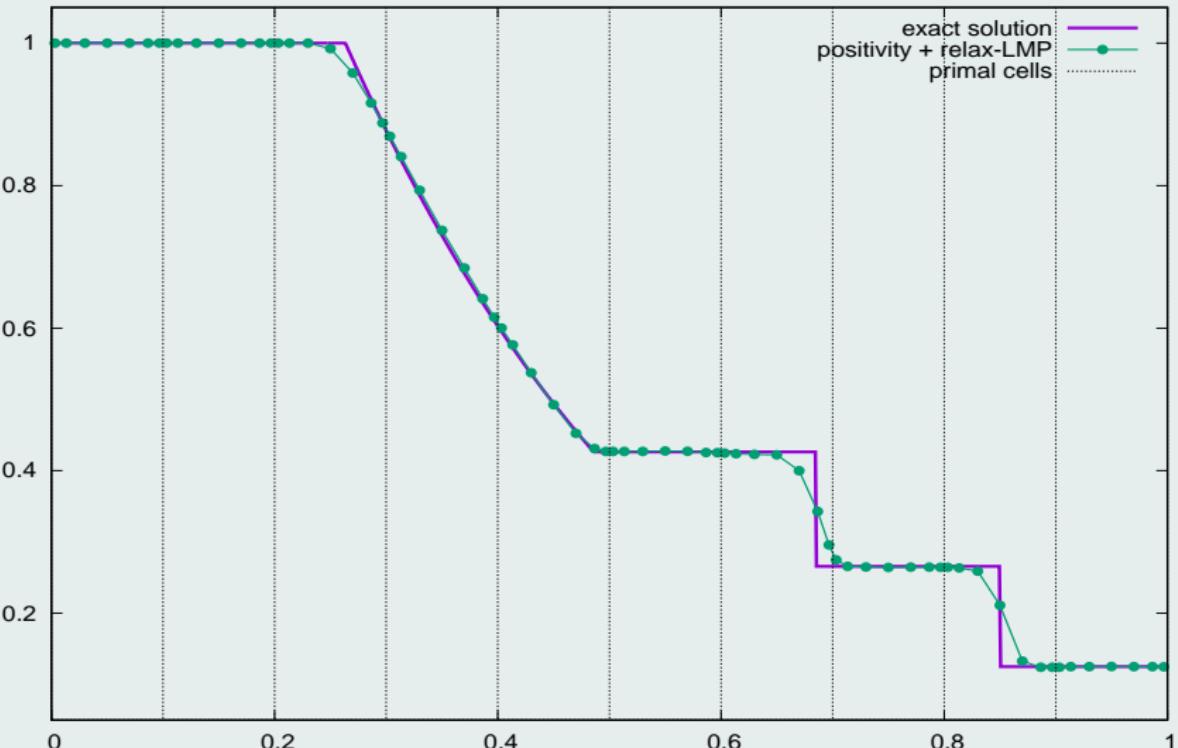
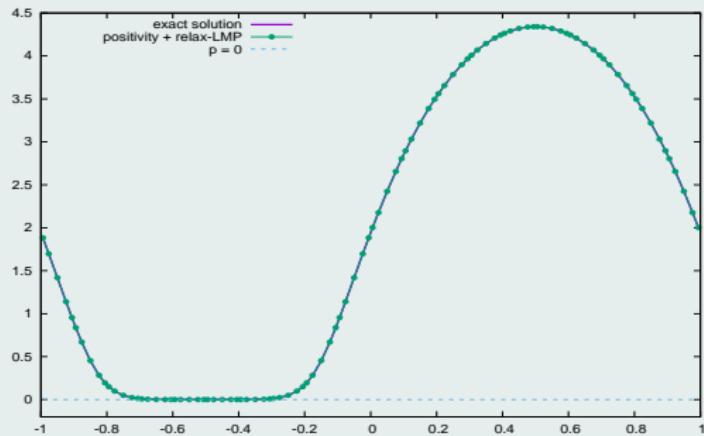


Figure: \mathbb{P}^6 -DG/FV scheme with GMP and relaxed-LMP: submean values

Smooth isentropic solution

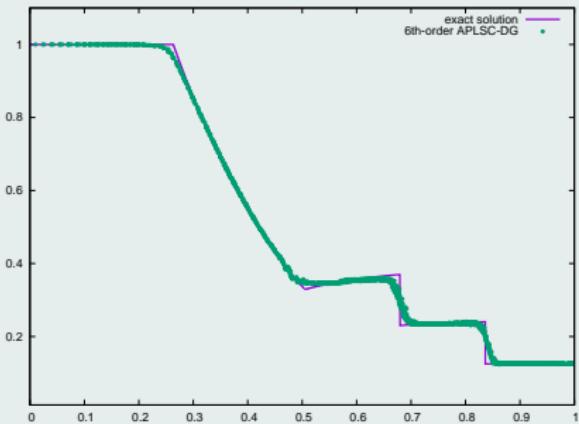
$$\rho_0 = 1 + 0.9999999 \sin(2\pi x)$$



	L_1		L_2		θ_{mp}	
h	$E_{L_1}^h$	$q_{L_1}^h$	$E_{L_2}^h$	$q_{L_2}^h$	min. θ_{mp}	aver. θ_{mp}
$\frac{1}{10}$	9.07E-4	5.86	1.23E-3	5.90	2.00E-1	0.981
$\frac{1}{20}$	1.56E-5	4.03	2.05E-5	3.83	1.92E-1	0.997
$\frac{1}{40}$	9.53E-7	4.89	1.44E-6	4.85	5.65E-4	0.999
$\frac{1}{80}$	3.21E-8	4.80	5.00E-8	4.87	3.48E-5	0.999
$\frac{1}{160}$	1.15E-9	-	1.71E-9	-	1.00	1.00

Table: Convergence rates of the \mathbb{P}^4 -DG/FV scheme with positivity and relaxed-LMP

Sod shock tube problem in cylindrical geometry



(b) Density profile

(a) Density map

Figure: \mathbb{P}^5 -DG/FV with positivity and relaxed-LMP on a 110 cells mesh

Sedov point blast problem in cylindrical geometry

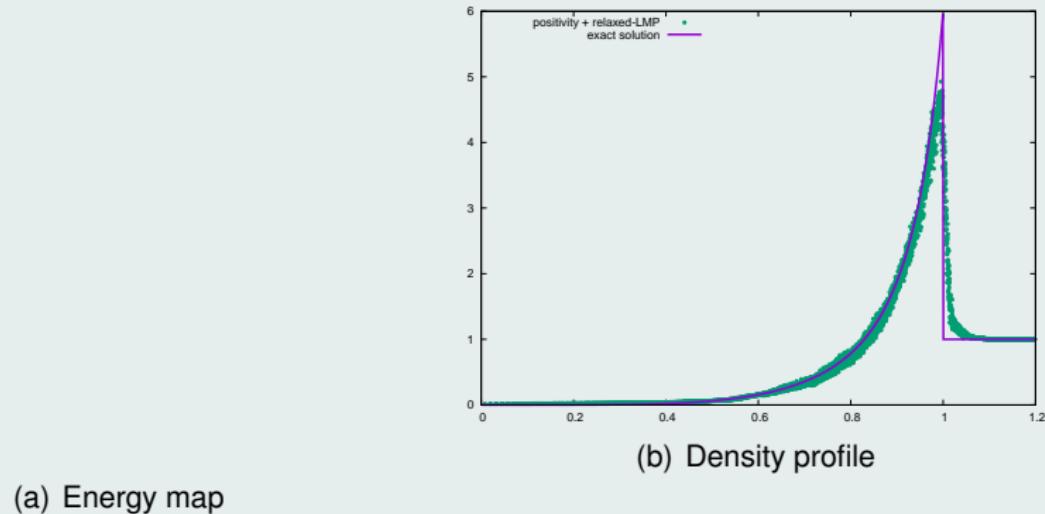


Figure: \mathbb{P}^5 -DG/FV with positivity and relaxed-LMP on a 271 cells mesh

Mach 3 forward-facing step

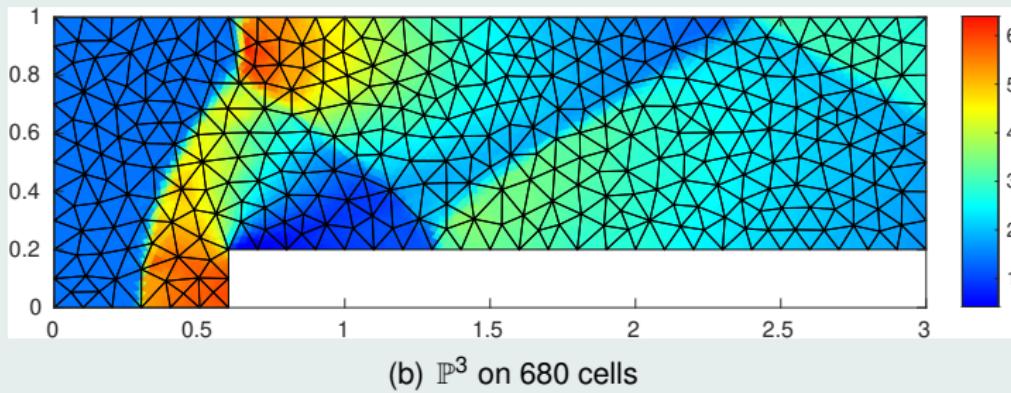
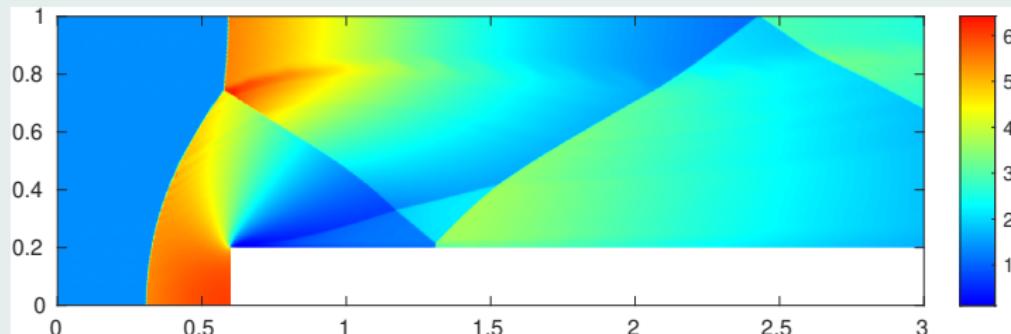


Figure: Monolithic subcell DG/FV scheme with positivity and relaxed-LMP

Mach 20 hypersonic flow over half cylinder

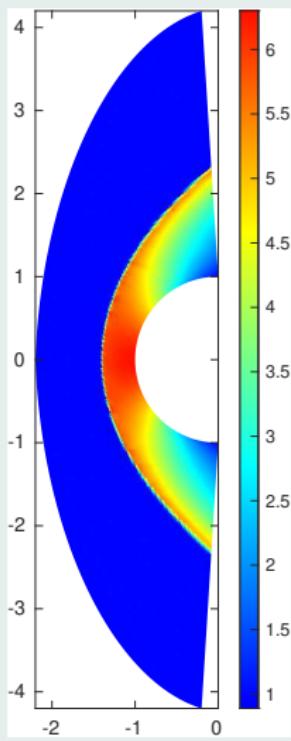
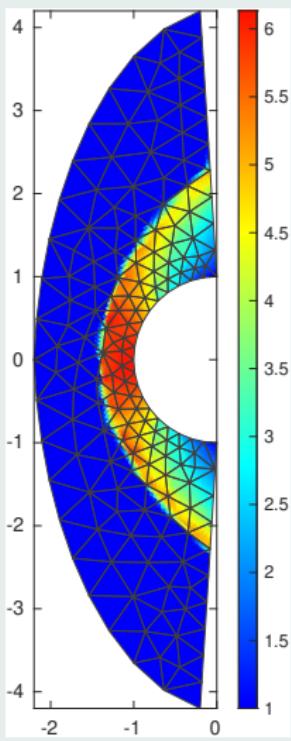
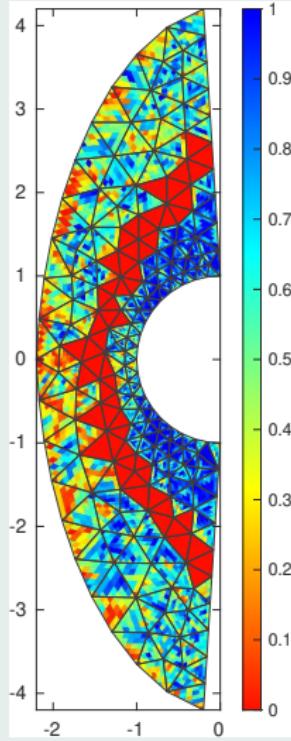
(a) \mathbb{P}^0 on 25266 cells(b) \mathbb{P}^5 on 292 cells(c) \mathbb{P}^5 on 292 cells

Figure: Monolithic subcell DG/FV scheme: density and blending coefficients

1 Introduction

2 DG as a subcell FV

3 Monolithic subcell DG/FV scheme

4 Entropy stabilities

5 Maximum principles

6 Conclusion

Monolithic local subcell DG/FV scheme

- Reformulate DG schemes as subgrid FV-like schemes:
 - regardless the type of mesh used
 - regardless the space dimension (*in theory...*)
 - regardless the cell subdivision ($N_s \geq N_k$)
- Combine high-order reconstructed fluxes and 1st-order FV fluxes
 - ensuring a maximum or positivity preserving principle at the subcell scale
 - ensuring different entropy stability inequalities
 - reducing significantly the apparition of spurious oscillations
 - preserving the very accurate subcell resolution of DG schemes

Questions

- Is an entropy inequality for one entropy enough?

 **Generally, no**

- Is entropy inequality absolutely needed?

 **Maybe not** \implies **GMP + relaxed LMP**

Articles on this topic

-  **F.V**, *A Posteriori Correction of High-Order DG Scheme through Subcell Finite Volume Formulation and Flux Reconstruction*. JCP, 2018.
-  **A. HAIDAR, F. MARCHE AND F.V**, *A posteriori Finite-Volume local subcell correction of high-order discontinuous Galerkin schemes for the nonlinear shallow-water equations*. JCP, 2022.
-  **F.V AND R. ABGRALL**, *A posteriori local subcell correction of DG schemes through Finite Volume reformulation on unstructured grids*. SIAM SISC, 2023.
-  **A. HAIDAR, F. MARCHE AND F.V**, *Free-boundary problems for wave-structure interactions in shallow-water: DG-ALE description and local subcell correction*. JSC, 2023.
-  **A. HAIDAR, F. MARCHE AND F.V**, *A robust DG-ALE formulation for nonlinear shallow-water interactions with a floating object*. JSC, 2024.
-  **F.V**, *Monolithic local subcell DG/FV convex property preserving scheme on unstructured grids and entropy consideration*. JCP, 2024.

Articles on this topic

-  **F.V**, *A Posteriori Correction of High-Order DG Scheme through Subcell Finite Volume Formulation and Flux Reconstruction*. JCP, 2018.
-  **A. HAIDAR, F. MARCHE AND F.V**, *A posteriori Finite-Volume local subcell correction of high-order discontinuous Galerkin schemes for the nonlinear shallow-water equations*. JCP, 2022.
-  **F.V AND R. ABGRALL**, *A posteriori local subcell correction of DG schemes through Finite Volume reformulation on unstructured grids*. SIAM SISC, 2023.
-  **A. HAIDAR, F. MARCHE AND F.V**, *Free-boundary problems for wave-structure interactions in shallow-water: DG-ALE description and local subcell correction*. JSC, 2023.
-  **A. HAIDAR, F. MARCHE AND F.V**, *A robust DG-ALE formulation for nonlinear shallow-water interactions with a floating object*. JSC, 2024.
-  **F.V**, *Monolithic local subcell DG/FV convex property preserving scheme on unstructured grids and entropy consideration*. JCP, 2024.