

stitut de Mathématiques de Toulon



Adaptive mesh refinement method : numerical density of entropy production and automatic thresholding.

Ersoy, M.¹ Pons, K.^{1,2}

¹EA 2134, IMATH, Université de Toulon

²Principia S.A.S.,13705 La Ciotat cedex

2017, 17 February, LAMA - Chambéry

- a. ersoy@univ-tln.fr
- b. pons@univ-tln.fr

MOTIVATIONS

 $\bullet\,$ numerical simulations of real-life large scale fluid flows in dimension d=1,2,3



(a) Monster waves (Nazare)

(b) Wave-breaking (Nazare)



(c) Tsunami (Japan)

(d) tsunami (Brazil)

MOTIVATIONS

- $\bullet\,$ numerical simulations of real-life large scale fluid flows in dimension d=1,2,3
- \bullet Simulating fluids in large-scale \rightarrow high memory and computer requirements.
 - ▶ how to compute fast and accurate : \rightarrow meshing or moving the mesh only in "desired" regions with a suitable mesh refinement criterion

MOTIVATIONS

- $\bullet\,$ numerical simulations of real-life large scale fluid flows in dimension d=1,2,3
- $\bullet\,$ Simulating fluids in large-scale $\rightarrow\,$ high memory and computer requirements.
- Mathematical motivations : introducing new tool for numerical purpose !

OUTLINE OF THE TALK

1 Adaptive Mesh Refinement method

- Generality
- An example

2 Automatic mesh refinement threshold



1 Adaptive Mesh Refinement method

- Generality
- An example

2 Automatic mesh refinement threshold



ADAPTIVE MESH REFINEMENT METHOD Generality

• An example

2 Automatic mesh refinement threshold

HYPERBOLIC EQUATIONS AND ENTROPY INEQUALITY

We focus on general non linear hyperbolic conservation laws

$$\left\{ \begin{array}{l} \frac{\partial \boldsymbol{w}}{\partial t} + \frac{\partial \boldsymbol{f}(\boldsymbol{w})}{\partial x} = 0, \ (t, x) \in \mathbb{R}^+ \times \mathbb{R} \\ \boldsymbol{w}(0, x) = \boldsymbol{w}_0(x), \ x \in \mathbb{R} \end{array} \right.$$

 $oldsymbol{w} \in \mathbb{R}^d$: vector state,

f : flux governing the physical description of the flow.

We focus on general non linear hyperbolic conservation laws

$$\begin{cases} \frac{\partial \boldsymbol{w}}{\partial t} + \frac{\partial \boldsymbol{f}(\boldsymbol{w})}{\partial x} = 0, \ (t, x) \in \mathbb{R}^+ \times \mathbb{R} \\ \boldsymbol{w}(0, x) = \boldsymbol{w}_0(x), \ x \in \mathbb{R} \end{cases}$$

Weak solutions satisfy

$$S = \frac{\partial s(\boldsymbol{w})}{\partial t} + \frac{\partial \psi(\boldsymbol{w})}{\partial x} \begin{cases} = 0 & \text{for smooth solution} \\ = 0 & \text{across rarefaction} \\ < 0 & \text{across shock} \end{cases}$$

where (s,ψ) stands for a convex entropy-entropy flux pair :

$$(\nabla \psi(\boldsymbol{w}))^T = (\nabla s(\boldsymbol{w}))^T \ D_{\boldsymbol{w}} \boldsymbol{f}(\boldsymbol{w})$$

We focus on general non linear hyperbolic conservation laws

$$\begin{cases} \frac{\partial \boldsymbol{w}}{\partial t} + \frac{\partial \boldsymbol{f}(\boldsymbol{w})}{\partial x} = 0, \ (t, x) \in \mathbb{R}^+ \times \mathbb{R} \\ \boldsymbol{w}(0, x) = \boldsymbol{w}_0(x), \ x \in \mathbb{R} \end{cases}$$

Weak solutions satisfy

$$S = \frac{\partial s(\boldsymbol{w})}{\partial t} + \frac{\partial \psi(\boldsymbol{w})}{\partial x} \begin{cases} = 0 & \text{for smooth solution} \\ = 0 & \text{across rarefaction} \\ < 0 & \text{across shock} \end{cases}$$

where (s,ψ) stands for a convex entropy-entropy flux pair :

$$\left(
abla \psi(oldsymbol{w})
ight)^T = \left(
abla s(oldsymbol{w})
ight)^T \ D_{oldsymbol{w}} oldsymbol{f}(oldsymbol{w})$$

Entropy inequality \simeq "smoothness indicator"

Croisille J.-P., Contribution à l'Étude Théorique et à l'Approximation par Éléments Finis du Système Hyperbolique de la Dynamique des Gaz Multidimensionnelle et Multiespèces, PhD thesis, Université de Paris VI, 1991 We focus on general non linear hyperbolic conservation laws

$$\begin{cases} \frac{\partial \boldsymbol{w}}{\partial t} + \mathsf{div}(\boldsymbol{f}(\boldsymbol{w})) = 0, \ (t, x) \in \mathbb{R}^+ \times \mathbb{R}^d\\ \boldsymbol{w}(0, x) = \boldsymbol{w}_0(x), \ x \in \mathbb{R}^d \end{cases}$$

Weak solutions satisfy

$$S = \frac{\partial s(\boldsymbol{w})}{\partial t} + \operatorname{div}(\psi(\boldsymbol{w})) \begin{cases} = 0 & \text{for smooth solution} \\ = 0 & \text{across rarefaction} \\ < 0 & \text{across shock} \end{cases}$$

where (s,ψ) stands for a convex entropy-entropy flux pair :

$$(\nabla \psi_i(\boldsymbol{w}))^T = (\nabla s(\boldsymbol{w}))^T D_{\boldsymbol{w}} \boldsymbol{f}_i(\boldsymbol{w}), \quad i = 1, \dots, d$$

Entropy inequality \simeq "smoothness indicator"

Croisille J.-P., Contribution à l'Étude Théorique et à l'Approximation par Éléments Finis du Système Hyperbolique de la Dynamique des Gaz Multidimensionnelle et Multiespèces, PhD thesis, Université de Paris VI, 1991



FIGURE: a cell C_k

Finite volume approximation :

$$\boldsymbol{w}_{k}^{n+1} = \boldsymbol{w}_{k}^{n} - rac{\delta t_{n}}{h_{k}} \left(\boldsymbol{F}_{k+1/2}^{n} - \boldsymbol{F}_{k-1/2}^{n}
ight)$$

with

$$\boldsymbol{w}_k^n \simeq rac{1}{h_k} \int_{C_k} \boldsymbol{w}\left(t_n, x
ight) \, dx \text{ and } \boldsymbol{F}_{k+1/2}^n pprox rac{1}{\delta t} \int_{C_k} \boldsymbol{f}(t, w(t, x_{k+1/2})) \, dx$$



FIGURE: a cell C_k

Finite volume approximation :

$$\boldsymbol{w}_{k}^{n+1} = \boldsymbol{w}_{k}^{n} - rac{\delta t_{n}}{h_{k}} \left(\boldsymbol{F}_{k+1/2}^{n} - \boldsymbol{F}_{k-1/2}^{n}
ight)$$

with

$$\boldsymbol{w}_k^n \simeq rac{1}{h_k} \int_{C_k} \boldsymbol{w}\left(t_n, x\right) \, dx \text{ and } \boldsymbol{F}_{k+1/2}^n pprox rac{1}{\delta t} \int_{C_k} \boldsymbol{f}(t, w(t, x_{k+1/2})) \, dx$$

The numerical density of entropy production :

$$S_{k}^{n} = \frac{s_{k}^{n+1} - s_{k}^{n}}{\delta t_{n}} + \frac{\psi_{k+1/2}^{n} - \psi_{k-1/2}^{n}}{h_{k}} \lessapprox 0$$



FIGURE: a cell C_k

Finite volume approximation :

$$\boldsymbol{w}_{k}^{n+1} = \boldsymbol{w}_{k}^{n} - \frac{\delta t_{n}}{h_{k}} \left(\sum_{a} \boldsymbol{F}(\boldsymbol{w}_{k}^{n}, \boldsymbol{w}_{a}^{n}; n_{k/a}) \right), \quad h_{k} = \frac{|C_{k}|}{\sum_{a} |\partial C_{k/a}|}$$

with

$$\boldsymbol{w}_k^n \simeq \frac{1}{h_k} \int_{C_k} \boldsymbol{w}(t_n, x) \, dx, \quad \text{and } \boldsymbol{F}(\boldsymbol{w}_k^n, \boldsymbol{w}_a^n; n_{k/a}) \approx \frac{1}{\delta t} \int_{\partial C_k} \boldsymbol{f}(t, w) \cdot n_{k/a} \, ds$$



FIGURE: a cell C_k

Finite volume approximation :

$$\boldsymbol{w}_{k}^{n+1} = \boldsymbol{w}_{k}^{n} - \frac{\delta t_{n}}{h_{k}} \left(\sum_{a} \boldsymbol{F}(\boldsymbol{w}_{k}^{n}, \boldsymbol{w}_{a}^{n}; n_{k/a}) \right), \quad h_{k} = \frac{|C_{k}|}{\sum_{a} |\partial C_{k/a}|}$$

with

$$\boldsymbol{w}_k^n \simeq \frac{1}{h_k} \int_{C_k} \boldsymbol{w}(t_n, x) \, dx, \quad \text{ and } \boldsymbol{F}(\boldsymbol{w}_k^n, \boldsymbol{w}_a^n; n_{k/a}) \approx \frac{1}{\delta t} \int_{\partial C_k} \boldsymbol{f}(t, w) \cdot n_{k/a} \, ds$$

The numerical density of entropy production :

$$S_k^n = \frac{s_k^{n+1} - s_k^n}{\delta t_n} + \frac{\sum_a \psi(\boldsymbol{w}_k^n, \boldsymbol{w}_a^n; n_{k/a})}{h_k} \lessapprox 0$$

• Compute w_k^n

- Compute w_k^n
- Compute $S_k^n : S_k^n \neq 0 \Longrightarrow$ the cell is refined or coarsened

- Compute w_k^n
- Compute $S_k^n:S_k^n \neq 0 \Longrightarrow$ the cell is refined or coarsened
- More precisely : for a given mesh refinement threshold lpha

•
$$S_k^n \ge \alpha \overline{S} \Longrightarrow$$
 the cell is refined with, for instance, $\overline{S} = \frac{1}{|\Omega|} \int_{\Omega} S_k^n$



- Compute w_k^n
- Compute $S_k^n:S_k^n \neq 0 \Longrightarrow$ the cell is refined or coarsened
- More precisely : for a given mesh refinement threshold lpha
 - $S_k^n \ge \alpha \overline{S} \Longrightarrow$ the cell is refined with, for instance, $\overline{S} = \frac{1}{|\Omega|} \int_{\Omega} S_k^n$



- Compute w_k^n
- Compute $S_k^n : S_k^n \neq 0 \Longrightarrow$ the cell is refined or coarsened
- More precisely : for a given mesh refinement threshold lpha
 - $S_k^n \ge \alpha \overline{S} \implies$ the cell is refined with, for instance, $\overline{S} = \frac{1}{|\Omega|} \int_{\Omega} S_k^n$
 - $S_k^n < \alpha \overline{S} \Longrightarrow$ the cell is coarsened
 - Dynamic mesh refinement :
 - ★ Dyadic tree (1D)
 - hierarchical numbering : basis 2



- Compute w_k^n
- Compute $S_k^n: S_k^n \neq 0 \Longrightarrow$ the cell is refined or coarsened
- More precisely : for a given mesh refinement threshold α
 - $S_k^n \ge \alpha \overline{S} \Longrightarrow$ the cell is refined with, for instance, $\overline{S} = \frac{1}{|\Omega|} \int_{\Omega} S_k^n$
 - $S_k^n < \alpha \overline{S} \Longrightarrow$ the cell is coarsened
 - Dynamic mesh refinement :
 - * Non-structured grid : macro-cell
 - * Dyadic tree (1D), Quadtree (2D)
 - hierarchical numbering : basis 2,4



- Compute w_k^n
- Compute $S_k^n: S_k^n \neq 0 \Longrightarrow$ the cell is refined or coarsened
- More precisely : for a given mesh refinement threshold lpha
 - $S_k^n \ge \alpha \overline{S} \Longrightarrow$ the cell is refined with, for instance, $\overline{S} = \frac{1}{|\Omega|} \int_{\Omega} S_k^n$
 - $S_k^n < \alpha \overline{S} \Longrightarrow$ the cell is coarsened
 - Dynamic mesh refinement :
 - * Non-structured grid : macro-cell
 - * Dyadic tree (1D), Quadtree (2D), Octree (3D)
 - hierarchical numbering : basis 2,4,8



- Compute w_k^n
- Compute $S_k^n: S_k^n \neq 0 \Longrightarrow$ the cell is refined or coarsened
- More precisely : for a given mesh refinement threshold lpha
 - $S_k^n \ge \alpha \overline{S} \Longrightarrow$ the cell is refined with, for instance, $\overline{S} = \frac{1}{|\Omega|} \int_{\Omega} S_k^n$
 - $S_k^n < \alpha \overline{S} \Longrightarrow$ the cell is coarsened



- Compute w_k^n
- Compute $S_k^n: S_k^n \neq 0 \Longrightarrow$ the cell is refined or coarsened
- More precisely : for a given mesh refinement threshold lpha
 - $S_k^n \ge \alpha \overline{S} \Longrightarrow$ the cell is refined with, for instance, $\overline{S} = \frac{1}{|\Omega|} \int_{\Omega} S_k^n$
 - $S_k^n < \alpha \overline{S} \Longrightarrow$ the cell is coarsened



- Compute w_k^n
- Compute $S_k^n: S_k^n \neq 0 \Longrightarrow$ the cell is refined or coarsened
- More precisely : for a given mesh refinement threshold lpha
 - $S_k^n \ge \alpha \overline{S} \Longrightarrow$ the cell is refined with, for instance, $\overline{S} = \frac{1}{|\Omega|} \int_{\Omega} S_k^n$
 - $S_k^n < \alpha \overline{S} \Longrightarrow$ the cell is coarsened
 - A simple projection method but the scheme is locally non consistent [S088, TW05]



Shu C. W., Osher S., Efficient implementation of essentially nonoscillatory shock-capturing schemes. J. Comput. Phys., 77(2) :439–471, 1988.

Tang H., Warnecke G., A class of high resolution difference schemes for nonlinear Hamilton-Jacobi equations with varying time and space grids. SIAM J. Sci. Comput., 26(4):1415–1431, 2005.

- Compute w_k^n
- Compute $S_k^n: S_k^n \neq 0 \Longrightarrow$ the cell is refined or coarsened
- More precisely : for a given mesh refinement threshold lpha
 - $S_k^n \ge \alpha \overline{S} \Longrightarrow$ the cell is refined with, for instance, $\overline{S} = \frac{1}{|\Omega|} \int_{\Omega} S_k^n$
 - $S_k^n < \alpha \overline{S} \Longrightarrow$ the cell is coarsened
 - A simple projection method but the scheme is locally non consistent [S088, TW05]
 - * less time time consuming than other methods



Shu C. W., Osher S., Efficient implementation of essentially nonoscillatory shock-capturing schemes. J. Comput. Phys., 77(2) :439–471, 1988.

Tang H., Warnecke G., A class of high resolution difference schemes for nonlinear Hamilton-Jacobi equations with varying time and space grids. SIAM J. Sci. Comput., 26(4):1415–1431, 2005.

- Compute w_k^n
- Compute $S_k^n: S_k^n \neq 0 \Longrightarrow$ the cell is refined or coarsened
- More precisely : for a given mesh refinement threshold lpha
 - $S_k^n \ge \alpha \overline{S} \Longrightarrow$ the cell is refined with, for instance, $\overline{S} = \frac{1}{|\Omega|} \int_{\Omega} S_k^n$
 - $S_k^n < \alpha \overline{S} \Longrightarrow$ the cell is coarsened
 - A simple projection method but the scheme is locally non consistent [S088, TW05]
 - * less time time consuming than other methods
 - * non consistency is almost negligible if the level of adjacent cells is limited by 2



Shu C. W., Osher S., Efficient implementation of essentially nonoscillatory shock-capturing schemes. J. Comput. Phys., 77(2) :439–471, 1988.

Tang H., Warnecke G., A class of high resolution difference schemes for nonlinear Hamilton-Jacobi equations with varying time and space grids. SIAM J. Sci. Comput., 26(4):1415–1431, 2005.



M. Ersoy, F. Golay, L. Yushchenko. Adaptive multi-scale scheme based on numerical entropy production for conservation laws. CEJM, Central European Journal of Mathematics, 11(8), pp 1392–1415, 2013.



F. Golay, M. Ersoy, L. Yuschenko, D. Sous. Block-based adaptive mesh refinement scheme using numerical density of entropy production for three-dimensional two-fluid flows. International Journal of Computational Fluid Dynamics, Taylor & Francis, 2015.

- Compute w_k^n
- Compute $S_k^n: S_k^n \neq 0 \Longrightarrow$ the cell is refined or coarsened
- More precisely : for a given mesh refinement threshold lpha
 - $S_k^n \ge \alpha \overline{S} \Longrightarrow$ the cell is refined with, for instance, $\overline{S} = \frac{1}{|\Omega|} \int_{\Omega} S_k^n$
 - $S_k^n < \alpha \overline{S} \Longrightarrow$ the cell is coarsened
 - A simple projection method but the scheme is locally non consistent [S088, TW05]
 - * less time time consuming than other methods
 - * non consistency is almost negligible if the level of adjacent cells is limited by 2
 - Improvement (cpu-time) : local time stepping [EGY13] · See more details



Tang H., Warnecke G., A class of high resolution difference schemes for nonlinear Hamilton-Jacobi equations with varying time and space grids. SIAM J. Sci. Comput., 26(4) :1415–1431, 2005.





F. Golay, M. Ersoy, L. Yuschenko, D. Sous. Block-based adaptive mesh refinement scheme using numerical density of entropy production for three-dimensional two-fluid flows. International Journal of Computational Fluid Dynamics, Taylor & Francis, 2015.

- Compute w_k^n
- Compute $S_k^n: S_k^n \neq 0 \Longrightarrow$ the cell is refined or coarsened
- ullet More precisely : for a given mesh refinement threshold α
 - $S_k^n \ge \alpha \overline{S} \Longrightarrow$ the cell is refined with, for instance, $\overline{S} = \frac{1}{|\Omega|} \int_{\Gamma} S_k^n$
 - $S_k^n < \alpha \overline{S} \Longrightarrow$ the cell is coarsened
 - A simple projection method but the scheme is locally non consistent [S088, TW05]
 - * less time time consuming than other methods
 - * non consistency is almost negligible if the level of adjacent cells is limited by 2
 - Improvement (cpu-time) : local time stepping [EGY13] See more details
 - 2D-3D computer management : BB-AMR [GEYS] See more details



Tang H., Warnecke G., A class of high resolution difference schemes for nonlinear Hamilton-Jacobi equations with varying time and space grids. SIAM J. Sci. Comput., 26(4) :1415–1431, 2005.





F. Golay, M. Ersoy, L. Yuschenko, D. Sous. Block-based adaptive mesh refinement scheme using numerical density of entropy production for three-dimensional two-fluid flows. International Journal of Computational Fluid Dynamics, Taylor & Francis, 2015.



ADAPTIVE MESH REFINEMENT METHOD Generality

• An example

2 Automatic mesh refinement threshold

AN EXAMPLE : THE ONE-DIMENSIONAL GAS DYNAMICS EQUATIONS FOR IDEAL GAS

$$\begin{split} \frac{\partial\rho}{\partial t} + \frac{\partial\rho u}{\partial x} &= 0\\ \frac{\partial\rho u}{\partial t} + \frac{\partial\left(\rho u^2 + p\right)}{\partial x} &= 0 \quad \text{where} \\ \frac{\partial\rho E}{\partial t} + \frac{\partial\left(\rho E + p\right) u}{\partial x} &= 0\\ p &= (\gamma - 1)\rho\varepsilon \end{split}$$

$$\begin{array}{llll} \rho(t,x) & : & \text{density} \\ u(t,x) & : & \text{velocity} \\ p(t,x) & : & \text{pressure} \\ \gamma := 1.4 & : & \text{ratio of the specific heats} \\ E(\varepsilon,u) & : & \text{total energy} \\ \varepsilon & : & \text{internal specific energy} \\ E & = & \varepsilon + \frac{u^2}{2} \end{array}$$

Intel(R) Core(TM) i5-2500 CPU @ 3.30GHz

AN EXAMPLE : THE ONE-DIMENSIONAL GAS DYNAMICS EQUATIONS FOR IDEAL GAS

$$\begin{array}{cccc} & \rho(t,x) & : & \text{density} \\ & \frac{\partial\rho}{\partial t} + \frac{\partial\rho u}{\partial x} = 0 & \mu(t,x) & : & \text{velocity} \\ & u(t,x) & : & \text{pressure} \\ & \frac{\partial\rho u}{\partial t} + \frac{\partial\left(\rho u^2 + p\right)}{\partial x} = 0 & \text{where} & \gamma := 1.4 & : & \text{ratio of the specific heats} \\ & E(\varepsilon,u) & : & \text{total energy} \\ & \frac{\partial\rho E}{\partial t} + \frac{\partial\left(\rho E + p\right) u}{\partial x} = 0 & \varepsilon & : & \text{internal specific energy} \\ & p = (\gamma - 1)\rho\varepsilon & E & = & \varepsilon + \frac{u^2}{2} \end{array}$$

• Conservative variables

$$\boldsymbol{w} = \left(\rho, \rho u, \rho E\right)^t$$

convex continuous entropy

$$s(oldsymbol{w}) = -
ho \ln\left(rac{p}{
ho^{\gamma}}
ight) ext{ of flux } \psi(oldsymbol{w}) = u \, s(oldsymbol{w}) \ .$$

Intel(R) Core(TM) i5-2500 CPU @ 3.30GHz

SOD'S SHOCK TUBE PROBLEM

 12
 13
 13
 1- 5000

 14
 14
 14

 15
 14
 1

 16
 13
 1

 17
 14

 18
 14

 19
 14

 10
 14

 10
 14

 11
 14

 12
 14

 13
 14

 14
 14

 15
 14

 16
 14

 17
 14

 18
 14

 19
 14

 10
 14

 10
 14

 11
 14

 12
 14

 13
 14

 14
 14

 15
 14

 16
 14

 17
 14

 18
 14

 19
 14

 10
 14

 10
 14

 11
 14

 12
 14

 14
 14

 15
 14

 14
 14

 15
 14

 16
 14

 17
 14

 18
 14

 19
 14

 10
 14<

ACCURACY



(a) Density and numerical density of en- (b) Mesh refinement level, numerical tropy production. density of entropy production and local error.

FIGURE: Sod's shock tube problem : solution at time t = 0.4 s using the AB1M scheme on a dynamic grid with $L_{\rm max} = 5$ and the AB1 scheme on a uniform fixed grid of 681 cells.

Properties of S : shock criterion type

Theorem ([P04])

Consider a p^{th} convergent scheme. Let S_k^n be the corresponding numerical density of entropy production and $\Delta t = \lambda h$ be a fixed time step where h stands for the meshsize.

Then

$$\lim_{n \to \infty} S_k^n = \begin{cases} O(\Delta t^p) \\ O\left(\frac{1}{\Delta t}\right) \end{cases}$$

if the solution is smooth,

if the solution is discontinuous.

Puppo G., Numerical entropy production for central schemes. SIAM J. Sci. Comput., 25(4) :1382-1415, 2004.

Properties of S : shock criterion type

Theorem ([P04])

Consider a p^{th} convergent scheme. Let S_k^n be the corresponding numerical density of entropy production and $\Delta t = \lambda h$ be a fixed time step where h stands for the meshsize.

Then

$$\lim_{n \to \infty} S_k^n = \begin{cases} O(\Delta t^p) \\ O\left(\frac{1}{\Delta t}\right) \end{cases}$$

if the solution is smooth,

if the solution is discontinuous.

PROPERTIES

Consider a monotone scheme. Then, for almost every k, every $n, S_k^n \leq 0$.
Properties of S : shock criterion type

Theorem ([P04])

Consider a p^{th} convergent scheme. Let S_k^n be the corresponding numerical density of entropy production and $\Delta t = \lambda h$ be a fixed time step where h stands for the meshsize.

Then

$$\lim_{n \to \infty} S_k^n = \begin{cases} O(\Delta t^p) \\ O\left(\frac{1}{\Delta t}\right) \end{cases}$$

if the solution is smooth,

if the solution is discontinuous.

PROPERTIES

Consider a monotone scheme. Then, for almost every k, every n, $S_k^n \leq 0$.

Thus, even if locally S^n_k can take positive value, one has $S^n_k\leqslant C\Delta t^q,\quad q\geqslant p$. $\ensuremath{\bullet}$ See example



1 Adaptive Mesh Refinement method

- Generality
- An example

2 Automatic mesh refinement threshold

3 NUMERICAL SIMULATIONS



.

.

Mesh refinement criterion S^n_k and the mesh refinement threshold $\boldsymbol{\alpha}$

•
$$S_k^n \ge \alpha \overline{S} \implies$$
 the cell is refined with $\overline{S} = \frac{1}{|\Omega|} \int_{\Omega} S_k^n$
• $S_k^n \le \alpha \overline{S} \implies$ the cell is coarsened

Mesh refinement criterion S_k^n and the mesh refinement threshold α

•
$$S_k^n \ge \alpha \overline{S} \implies$$
 the cell is refined with $\overline{S} = \frac{1}{|\Omega|} \int_{\Omega} S_k^n$

• $S^n_k < \alpha \overline{S} \Longrightarrow$ the cell is coarsened

• how to set α suitably?

Mesh refinement criterion S_k^n and the mesh refinement threshold α

•
$$S_k^n \ge \alpha \overline{S} \Longrightarrow$$
 the cell is refined with $\overline{S} = \frac{1}{|\Omega|} \int_{\Omega} S_k^n$

•
$$S_k^n < \alpha \overline{S} \Longrightarrow$$
 the cell is coarsened

- how to set α suitably? ideally
 - automatically
 - accuracy and computational time are "well-balanced"
 - * catch the smallest maxima but not the smallest

Mesh refinement criterion S_k^n and the mesh refinement threshold α

•
$$S_k^n \ge \alpha \overline{S} \Longrightarrow$$
 the cell is refined with $\overline{S} = \frac{1}{|\Omega|} \int_{\Omega} S_k^n$

• $S_k^n < \alpha \overline{S} \Longrightarrow$ the cell is coarsened

- how to set α suitably? ideally
 - automatically
 - accuracy and computational time are "well-balanced"
 - * catch the smallest maxima but not the smallest
 - simple method ("low-cost")

.

• Mean method or deviation

$$S(x,t) > \alpha \frac{1}{|\Omega|} \int_{\Omega} S(x,t) \ dx$$

where α is a tunable dimensionless threshold parameter or deviation method

$$S(x,t) > \alpha \frac{1}{|\Omega|} \int_{\Omega} S(x,t) \ dx + \beta \sigma(x,t)$$

where σ is the standard deviation with β is a tunable dimensionless threshold parameter



Kallinderis, Y.G., Baron, J.R. : Adaptation methods for a new navier-stokes algorithm. AIAA journal 27(1), 37-43 (1989)

Ersoy, M., Golay, F., Yushchenko, L. : Adaptive multiscale scheme based on numerical density of entropy production for conservation laws. Cent. Eur. J. Math. 11(8), 1392–1415 (2013).

- Mean method or deviation \rightarrow pb : set α or (α, β)
- Filtering : two-steps





Aftosmis, M.: Upwind method for simulation of viscous flow on adaptively refined meshes. AIAA journal 32(2), 268-277 (1994)

Warren, G.P., Anderson, W.K., Thomas, J.L., Krist, S.L. : Grid convergence for adaptive methods. AIAA paper 1592, 1991 (1991)

- Mean method or deviation \rightarrow pb : set α or (α, β)
- Filtering : two-steps \rightarrow pb : two-step ? Is it enough ? still to set α
- Filtering : wavelets



FIGURE: Illustration of the wavelet transformation for a given mesh refinement criterion computed with the Daubechies wavelet with four vanishing moments (warm colors correspond to large coefficients and cold colors to small coefficients)



- Mean method or deviation \rightarrow pb : set α or (α, β)
- Filtering : two-steps ightarrow pb : two-step ? Is it enough ? still to set lpha
- Filtering : wavelets \rightarrow pb : efficient but the cost ! still to set α
- Local maxima \rightarrow pb : efficient but the cost ! still to set α



 FIGURE : Illustration of the local maxima method for a mesh refinement criterion involving multiple scales

• Assumptions and notations

 \blacktriangleright S is smooth and has p local maxima.

•
$$S(0) = S(L) = S'(0) = S'(L) = 0$$

•
$$0 < S_{\infty} = \max_{x \in (0,L)} S(x) < \infty$$



- Assumptions and notations
 - \blacktriangleright S is smooth and has p local maxima.

•
$$S(0) = S(L) = S'(0) = S'(L) = 0$$

•
$$0 < S_{\infty} = \max_{x \in (0,L)} S(x) < \infty$$

• Then

$$Z_{\alpha} = \{x \in (0,L); \varphi_{\alpha}(x) = S(x) - \alpha = 0 \text{ and } S'(x) \neq 0\} \neq \emptyset$$
$$= \{x_0(\alpha) < x_1(\alpha) < \dots < x_{2p_{\alpha}-2}(\alpha) < x_{2p_{\alpha}-1}(\alpha)\}$$
$$\#Z_{\alpha} = 2p_{\alpha}$$



- Assumptions and notations
 - S is smooth and has p local maxima.

•
$$S(0) = S(L) = S'(0) = S'(L) = 0$$

•
$$0 < S_{\infty} = \max_{x \in (0,L)} S(x) < \infty$$

• Then

$$Z_{\alpha} = \{x \in (0,L); \varphi_{\alpha}(x) = S(x) - \alpha = 0 \text{ and } S'(x) \neq 0\} \neq \emptyset$$
$$= \{x_0(\alpha) < x_1(\alpha) < \dots < x_{2p_{\alpha}-2}(\alpha) < x_{2p_{\alpha}-1}(\alpha)\}$$

$$\blacktriangleright \# Z_{\alpha} = 2p_{\alpha}$$

• there exist sequence $(x_k^*)_{1 \le k \le p}$



- Assumptions and notations
 - S is smooth and has p local maxima.

•
$$S(0) = S(L) = S'(0) = S'(L) = 0$$

•
$$0 < S_{\infty} = \max_{x \in (0,L)} S(x) < \infty$$

$$Z_{\alpha} = \{x \in (0, L); \varphi_{\alpha}(x) = S(x) - \alpha = 0 \text{ and } S'(x) \neq 0\} \neq \emptyset$$
$$= \{x_0(\alpha) < x_1(\alpha) < \dots < x_{2p_{\alpha}-2}(\alpha) < x_{2p_{\alpha}-1}(\alpha)\}$$

$$\blacksquare \# Z_{\alpha} = 2p_{\alpha}$$

 \blacktriangleright there exist sequences $(x_k^*)_{1\leqslant k\leqslant p}$ and $(\alpha_k^*)_{1\leqslant k\leqslant p}$ such that

$$\star \quad \forall k = 1, \dots, p$$

*
$$S(x_k^*) = \alpha_k^*$$
 and $S'(x_k^*) = 0, \ S''(x_k^*) < 0$

$$\star \ x_k^*
otin Z_lpha$$
 and $lpha_p^* = S_\infty$



- Assumptions and notations
 - S is smooth and has p local maxima.

•
$$S(0) = S(L) = S'(0) = S'(L) = 0$$

•
$$0 < S_{\infty} = \max_{x \in (0,L)} S(x) < \infty$$

• With these settings, we define the distribution d

$$\alpha \in [0, S_{\infty}] \mapsto d(\alpha) := \begin{cases} 1 & \text{if} \quad \alpha = 0 \ ,\\ \frac{1}{L} \sum_{k=1}^{p_{\alpha}} x_{2k+1}(\alpha) - x_{2k}(\alpha) & \text{if} \quad 0 < \alpha < S_{\infty} \ ,\\ 0 & \text{if} \quad \alpha = S_{\infty} \ . \end{cases}$$



Properties of d and α_{PE} threshold

Properties

$$\begin{aligned} \bullet & \int_0^{S_\infty} d(\alpha) \, d\alpha = \int_0^L S(x) \, dx. \\ \bullet & d \in C^0([0, S_\infty], \mathbb{R}^+) \text{ and } d' \text{ satisfies} \\ \bullet & \forall \alpha \in [0, S_\infty], \, d'(\alpha) < 0 \\ \bullet & \forall k \in [\![0, p]\!], \, \lim_{\alpha \to \alpha_k^*} d'(\alpha) = -\infty \text{ with the convention } \alpha_0^* := 0 \end{aligned}$$

•
$$d \in C^{l}(D_{*}, \mathbb{R}^{+})$$
 on the set $D^{*} := \bigcup_{k=0}^{p-1} (\alpha_{k}^{*}, \alpha_{k+1}^{*})$.



• Hard to define α using only d



- Hard to define α using only d
- However, one can set $\alpha = \alpha_D$ or
 - either $\alpha = \alpha_D = \min\{\alpha \mid d'(\alpha) = -1\}$





Dannenhoffer, J.F. : Grid adaptation for complex two-dimensional transonic flows. Ph.D. thesis, Massachusetts Institute of Technology (1987)

- Hard to define α using only d
- However, one can set $\alpha = \alpha_D$ or $\alpha = \alpha_P$
 - either $\alpha = \alpha_P = \min\{\alpha \mid d''(\alpha) = 0\}$





Powell, K.G., Murman, E.M. : An embedded mesh procedure for leading-edge vortex flows. In : NASA, Langley Research Center, Transonic Symposium : Theory, Application, and Experiment, vol. 1 (1989)

M. Ersoy (IMATH)

- Hard to define α using only d
- However, one can set $\alpha = \alpha_D$ or $\alpha = \alpha_P$
- In practice, d is defined by $d(\alpha) = \sum_{j=0}^{M-1} d_j \ \mathbb{1}_{(\alpha_j, \alpha_{j+1})}(\alpha)$ with

$$d_j = \#\{k \ ; \ S_k^n > \alpha_j\} \text{ and } (\alpha_j)_{0 \leqslant j \leqslant M} = \left(S_m \left(\frac{j}{M}\right)^{\beta}\right)_{0 \leqslant j \leqslant M} \quad \beta \geqslant 1$$





- Hard to define α using only d
- However, one can set $\alpha = \alpha_D$ or $\alpha = \alpha_P$
- α is usually too small (too sensitive to β)!!! \rightarrow detects small fluctuations

- Hard to define α using only d
- However, one can set $\alpha = \alpha_D$ or $\alpha = \alpha_P$
- α is usually too small (too sensitive to β)!!! \rightarrow detects small fluctuations

 $\alpha = \alpha_{PE} = \min\{\alpha \in [0, \overline{S}] \mid \max\{\alpha d(\alpha)\}\}$

We propose to set



Pons, K., Ersoy, M. Adaptive mesh refinement method. Part 1 : Automatic thresholding based on a distribution function Pons, K., Ersoy, M. Adaptive mesh refinement method. Part 2 : Application to tsunamis propagation

Why $\alpha d(\alpha)$?

PROPERTIES

Assume that S is twice differentiable and has p local maxima. Then, there exists

$$\forall k = 0 \dots p - 1, \ \alpha_{k+1}^{**} \in (\alpha_k^*, \alpha_{k+1}^*)$$
 such that $d''(\alpha_{k+1}^{**}) = 0$

and the function f has p local maxima $\overline{\alpha}_1, \ldots, \overline{\alpha}_p$ such that

$$\forall k = 1 \dots p, \ \overline{\alpha}_k \in (\alpha_k^{**}, \alpha_k^*)$$
.

As a consequence $\alpha_{PE} > \alpha_P > \alpha_D$.



Illustration : α_{PE} threshold for "discontinuous" flows



FIGURE: The function f for the mesh refinement criterion $S(x) = 200 \exp(-1000(x - 1.25)^2) + 1.25 \exp(-5(x - 3.75)^2)$ representing a shock type solution

Illustration : α_{PE} threshold for "smooth" flows



FIGURE: The function f for the mesh refinement criterion $S(x) = 2 \exp(-10(x - 1.25)^2) + 1.25 \exp(-5(x - 3.75)^2)$ representing a smooth flow



1 Adaptive Mesh Refinement method

- Generality
- An example

2 Automatic mesh refinement threshold

3 NUMERICAL SIMULATIONS

NUMERICAL VALIDATION [PE]

A Dam-break problem for the Saint-Venant equations :

	$\left\{ \begin{array}{c} \partial_t \\ \partial_t \end{array} \right.$	$\begin{aligned} h + \partial_x(hu) &= 0, \\ hu) + \partial_x \left(hu^2 + gh^2/2\right) &= -ghZ'(x) \end{aligned}$
where	$ \begin{array}{c} h(t,x) \\ u(t,x) \\ g \\ Z \end{array} $	 density velocity of the water column gravity strength topography

Kévin Pons, Mehmet Ersoy. Adaptive mesh refinement method. Part 1 : Automatic thresholding based on a distribution function

NUMERICAL VALIDATION [PE]



FIGURE: Numerical results for the water height at time t = 2 s

NUMERICAL VALIDATION [PE]



FIGURE: Time evolution of the threshold parameter and the mean value S_m

TEST CASE I : SOLITARY WAVE PROPAGATION OVER A TWO DIMENSIONAL REEF



FIGURE: Experimental settings and wave gauges locations



Kévin Pons, Mehmet Ersoy, Frédéric Golay, Richard Marcer. Adaptive mesh refinement method. Part 2 : Application to tsunamis propagation

Roeber, V., Cheung, K.F. : Boussinesq-type model for energetic breaking waves in fringing reef environments. Coastal Engineering 70, 1-20 (2012)

 M. Ersoy (IMATH)
 AMR
 2017, 17 February, LAMA - Chambéry
 20 / 22

	Adaptive mesh simula- tion	Uniform mesh simula- tion
Simulation time	240	240
Number of cells	200-560	1000
Re-meshing time step	0.05 s	not applicable
Time order integration	1	1
Space order integration	1	1
CFL	0.99	0.99

TABLE: Numerical parameters










Test case I : Solitary wave propagation over a two dimensional reef



FIGURE: Surface profiles of solitary wave propagation over an exposed reef crest. Confrontation of experimental data (blue circles) to numerical data computed on a uniform grid (solid green line) and on an adaptive grid (solid red lines). The solid cyan line represents the mesh level and the black one the bathymetry

Test case I : Solitary wave propagation over a two dimensional reef



M. Ersoy (IMATH)

Test case I : Solitary wave propagation over a two dimensional reef



FIGURE: Time evolution of the mesh refinement threshold and the number of cells : 95 s against 210 s



FIGURE: Experimental settings



Kévin Pons, Mehmet Ersoy, Frédéric Golay, Richard Marcer. Adaptive mesh refinement method. Part 2 : Application to tsunamis propagation.

Lynett, P.J., Swigler, D., Son, S., Bryant, D., Socolofsky, S. : Experimental study of solitary wave evolution over a 3d shallow shelf. Coastal Engineering Proceedings 1(32), 1 (2011).

M. Ersoy (IMATH)

2017, 17 February, LAMA - Chambéry 21 / 22

SHELF

	Adaptive mesh simula- tion	Uniform mesh simula- tion
Simulation time	30 s	30 s
Number of blocks	128	128
Number of cells	7 500-25 000	33 000
Re-meshing time step	0.25 s	not applicable
Time order integration	2	2
Space order integration	2	2
CFL	0.5	0.5

TABLE: Numerical parameters



(a) t=0.5s



(a) t=2.5s



(a) t=5.75s



(a) t=23.75s

SHELF



21 / 22

SHELF



FIGURE: Time evolution of the mesh refinement threshold and the number of cells : speed up the computation by 2.5 time

Test case III : Tsunami runup onto a complex three dimensional

MONAI-WALLEY BEACH



FIGURE: Settings



Kévin Pons, Mehmet Ersoy, Frédéric Golay, Richard Marcer. Adaptive mesh refinement method. Part 2 : Application to tsunamis propagation

TEST CASE III : TSUNAMI RUNUP ONTO A COMPLEX THREE DIMENSIONAL

MONAI-WALLEY BEACH

	Adaptive mesh simula- tion	Uniform mesh simula- tion
Simulation time	30 s	30 s
Number of blocks	240	240
Number of cells	8 000-40 000	62 000
Re-meshing time step	0.25 s	not applicable
Time order integration	2	2
Space order integration	1	1
CFL	0.5	0.5

TABLE: Numerical parameters



(a) t = 11.25 s



(a) t = 13.25 s





(a) t = 17.5 s



 $\label{eq:FIGURE: Free surface results at different positions: experimental data versus numerical simulation with and without mesh adaptivity$

Test case III : Tsunami runup onto a complex three dimensional

MONAI-WALLEY BEACH



FIGURE: Time evolution of the mesh refinement threshold and the number of cells : speed up the computation by 3 time

Thank you

for your

VOU

attention

Let us consider the transport equation :

$$\begin{cases} w_t + w_x &= 0\\ w(0, x) &= w_0(x) \end{cases}$$

Back to slide

Let us consider the transport equation :

$$\begin{array}{rcl} w_t + w_x &=& 0\\ w(0,x) &=& w_0(x) \end{array}$$

and the Godunov scheme :

$$w_k^{n+1} = w_k^n - \frac{\delta t}{\delta x} \left(w_k^n - w_{k-1}^n \right)$$

Let us consider the transport equation :

$$\begin{array}{rcl} w_t + w_x &=& 0\\ w(0,x) &=& w_0(x) \end{array}$$

and the Godunov scheme :

$$\begin{cases} w_k^{n+1} &= w_k^n - \frac{\delta t}{\delta x} \left(w_k^n - w_{k-1}^n \right) \\ S_k^{n+1} &= \frac{s(w_k^{n+1}) - s(w_k^n)}{\delta t} + \frac{\psi(s(w_k^n)) - \psi(s(w_{k-1}^n))}{\delta x} \end{cases}$$

with $s(w) = w^2$ and $\psi(w) = w^2$.

Let us consider the transport equation :

$$\begin{array}{rcl} w_t + w_x &=& 0\\ w(0,x) &=& w_0(x) \end{array}$$

and the Godunov scheme :

$$\begin{cases} w_k^{n+1} &= w_k^n - \frac{\delta t}{\delta x} \left(w_k^n - w_{k-1}^n \right) \\ S_k^{n+1} &= \frac{s(w_k^{n+1}) - s(w_k^n)}{\delta t} + \frac{\psi(s(w_k^n)) - \psi(s(w_{k-1}^n))}{\delta x} \end{cases}$$

with $s(w) = w^2$ and $\psi(w) = w^2$. Substituting w_k^{n+1} into S_k^{n+1} , we get

$$S_k^{n+1} = -\varepsilon \left(\frac{w_k^n - w_{k-1}^n}{\delta x}\right)^2 \leqslant \mathbf{0} \text{ with } \varepsilon = \delta x \left(1 - \frac{\delta t}{\delta x}\right) > 0.$$

Back to slide

• Explicit adaptive schemes : time consuming due to the restriction

$$\|w\| \frac{\delta t}{h} \leqslant 1, \quad h = \min_k h_k$$

TIME RESTRICTION, LOCAL TIME STEPPING APPROACH

• Explicit adaptive schemes : time consuming due to the restriction

$$\|w\|rac{\delta t}{h}\leqslant 1, \quad h=\min_k h_k$$

- Local time stepping algorithm : save the cpu-time
 - Sort cells in groups w.r.t. to their level



Muller S., Stiriba Y., Fully adaptive multiscale schemes for conservation laws employing locally varying time stepping. SIAM J. Sci. Comput., 30(3) :493 7531, 2007.

TIME RESTRICTION, LOCAL TIME STEPPING APPROACH & AIMS

• Explicit adaptive schemes : time consuming due to the restriction

$$\|w\| rac{\delta t}{h} \leqslant 1, \quad h = \min_k h_k$$

- Local time stepping algorithm : save the cpu-time
 - Sort cells in groups w.r.t. to their level
 - Update the cells following the local time stepping algorithm.

Muller S., Stiriba Y., Fully adaptive multiscale schemes for conservation laws employing locally varying time stepping. SIAM J. Sci. Comput., 30(3) :493 7531, 2007.





FIGURE:
$$t_{n_1} = t_n + \delta t_r$$

with

$$\delta F_{k-1,k,k+1}^n := \left(F_{k+1/2}^n(w_k, w_{k+1}) - F_{k-1/2}^n(w_{k-1}, w_k) \right)$$

$$\boldsymbol{w}_{k_{000}}^{n_{2}} = \boldsymbol{w}_{k_{000}}^{n_{1}} - \frac{\delta t_{n}}{h_{k_{000}}} \delta \boldsymbol{F}_{k_{000},k_{000},k_{000}}^{n_{1}}$$
$$\boldsymbol{w}_{k_{00}}^{n_{2}} = \boldsymbol{w}_{k_{00}}^{n_{1}} - \frac{\delta t_{n}}{h_{k_{000}}} \delta \boldsymbol{F}_{k_{000},k_{000},k_{000}}^{n_{1}}$$
$$\boldsymbol{w}_{k_{000}}^{n_{2}} = \boldsymbol{w}_{k_{000}}^{n_{1}} - \frac{\delta t_{n}}{h_{k_{000}}} \delta \boldsymbol{F}_{k_{000},k_{000},k_{000}}^{n_{1}}$$
$$\boldsymbol{w}_{k_{000}}^{n_{2}} = \boldsymbol{w}_{k_{000}}^{n_{1}} - \frac{\delta t_{n}}{h_{k_{000}}} \delta \boldsymbol{F}_{k_{000},k_{000},k_{001}}^{n_{1}}$$

FIGURE:
$$t_{n_2} = t_n + 2\delta t_n$$

with $\delta F_{k-1,k,k+1}^n := \left(m{F}_{k+1/2}^n(m{w}_k,m{w}_{k+1}) - m{F}_{k-1/2}^n(m{w}_{k-1},m{w}_k)
ight)$



FIGURE:
$$t_{n_3} = t_n + 3\delta t_n$$

with

$$\delta F_{k-1,k,k+1}^n := \left(F_{k+1/2}^n(w_k, w_{k+1}) - F_{k-1/2}^n(w_{k-1}, w_k) \right)$$



LOCAL TIME STEPPING ALGORITHM

foreach $i \in \{1, 2^N\}$ do Let j be the biggest integer such that 2^{j} divides i foreach interface $x_{k+1/2}$ such that $\mathcal{L}_{k+1/2} \ge N - j$ do • compute the integral of $F_{k+1/2}(t)$ on the time interval $2^{N-\mathcal{L}_{k+1/2}}\delta t_n$, **2** distribute $F_{k+1/2}(t_n)$ to the two adjacent cells, • update only the cells of level greater than N - j. end end

Shu and Osher test case

Mesh coarsening parameter α Mesh refinement parameter \bar{S}

CFL

0.001

Ω

0.219,



EFFICIENCY OF THE LOCAL TIME STEPPING METHOD

	\mathcal{P}	$\ \rho - \rho_{ref}\ _{l^1_x}$	cpu-time	$N_{L_{\max}}$	maximum number of cells
AB1	0.288	4.7410^{-2}	181	1574	2308
AB1M	0.288	4.8010^{-2}	120	1572	2314

TABLE: Shu and Osher test case : comparison of numerical schemes of order 1

	\mathcal{P}	$\ \rho-\rho_{ref}\ _{l^1_x}$	cpu-time	$N_{L_{\max}}$	maximum number of cells
AB1	0.288	4.7410^{-2}	181	1574	2308
AB1M	0.288	4.8010^{-2}	120	1572	2314
AB2	0.287	2.7510^{-2}	170	1391	2023
		10 4	5	"si	They are
	2	1	-		A FEE TO

TABLE: Shu and Osher test case : comparison of numerical schemes of order 1 and 2
	\mathcal{P}	$\ \rho - \rho_{ref}\ _{l^1_x}$	cpu-time	$N_{L_{\max}}$	maximum number of cells	
AB1	0.288	4.7410^{-2}	181	1574	2308	
AB1M	0.288	4.8010^{-2}	120	1572	2314	
AB2	0.287	2.7510^{-2}	170	1391	2023	
AB2M	0.286	2.7410^{-2}	108	1357	1994	

TABLE: Shu and Osher test case : comparison of numerical schemes of order 1 and 2

	\mathcal{P}	$\ \rho-\rho_{ref}\ _{l^1_x}$	cpu-time	$N_{L_{\max}}$	maximum number of cells
AB1	0.288	4.7410^{-2}	181	1574	2308
AB1M	0.288	4.8010^{-2}	120	1572	2314
AB2	0.287	2.7510^{-2}	170	1391	2023
AB2M	0.286	2.7410^{-2}	108	1357	1994
RK2	0.285	2.0810^{-2}	299	1375	2005

TABLE: Shu and Osher test case : comparison of numerical schemes of order 1 and 2

M. Ersoy, F. Golay, L. Yushchenko. Adaptive multi-scale scheme based on numerical entropy production for conservation laws. CEJM, Central European Journal of Mathematics, 11(8), pp 1392-1415, 2013.

	\mathcal{P}	$\ \rho - \rho_{ref}\ _{l^1_x}$	cpu-time	$N_{L_{\max}}$	maximum number of cells
AB1	0.288	4.7410^{-2}	181	1574	2308
AB1M	0.288	4.8010^{-2}	120	1572	2314
AB2	0.287	2.7510^{-2}	170	1391	2023
AB2M	0.286	2.7410^{-2}	108	1357	1994
RK2	0.285	2.0810^{-2}	- 299	1375	2005

TABLE: Shu and Osher test case : comparison of numerical schemes of order 1 and 2

M. Ersoy, F. Golay, L. Yushchenko. Adaptive multi-scale scheme based on numerical entropy production for conservation laws. CEJM, Central European Journal of Mathematics, 11(8), pp 1392-1415, 2013.

Reference solution&Numerical results



(a) Density and numerical density of entropy production.

(b) Zoom on oscillating region.

FIGURE: Shu and Osher test case.

Back to slide

- Main difficulty : mesh and data structure. For fast computation, the following are required
 - parallel treatment
 - hierarchical grids

- Main difficulty : mesh and data structure. Some interesting issues :
 - 2D quad-tree [1],
 - 3D octree [2],
 - 2/3D anisotropic AMR [3].

Zhang, M., and W.M. Wu. 2011. A two dimensional hydrodynamic and sediment transport model for dam break based on finite volume method with quadtree grid. Applied Ocean Research 33 (4): 297 – 308.

Losasso, F., F. Gibou, and R. Fedkiw. 2004. Simulating Water and Smoke with an Octree Data Structure. ACM Trans. Graph. 23 (3) : 457–462, 2004.

Hachem, E., S. Feghali, R. Codina, and T. Coupez. Immersed stress method for fluid structure interaction using anisotropic mesh adaptation. International Journal for Numerical Methods in Engineering 94 (9): 805–825, 2013.

- Main difficulty : mesh and data structure.
- The strategy adopted :

- Main difficulty : mesh and data structure.
- The strategy adopted :
 - I domain= 1 block=1 cpu : "failure" → synchronization depends on the finest domain

- Main difficulty : mesh and data structure.
- The strategy adopted :
 - 1 domain= 1 block=1 cpu : "failure" → synchronization depends on the finest domain
 - 1 domain= n × blocks = 1cpu : "good compromise" → each domain has almost the same number number of cells → "better" synchronization

- Main difficulty : mesh and data structure.
- The strategy adopted :
 - I domain= 1 block=1 cpu : "failure" → synchronization depends on the finest domain
 - I domain= n × blocks = 1cpu : "good compromise" → each domain has almost the same number number of cells → "better" synchronization
 - It certainly exists better strategy

- Main difficulty : mesh and data structure.
- The strategy adopted :
- Management of domain's interfaces, projection step,

How it works?

• each domain has almost the same number of cells

- each domain has almost the same number of cells
- domain are defined using Cuthill-McKee numbering



- each domain has almost the same number of cells
- domain are defined using Cuthill-McKee numbering



- each domain has almost the same number of cells
- domain are defined using Cuthill-McKee numbering



- each domain has almost the same number of cells
- domain are defined using Cuthill-McKee numbering



- each domain has almost the same number of cells
- domain are defined using Cuthill-McKee numbering



- each domain has almost the same number of cells
- domain are defined using Cuthill-McKee numbering



- each domain has almost the same number of cells
- domain are defined using Cuthill-McKee numbering



- each domain has almost the same number of cells
- domain are defined using Cuthill-McKee numbering



- each domain has almost the same number of cells
- domain are defined using Cuthill-McKee numbering
- more sophisticated numbering exists

- each domain has almost the same number of cells
- domain are defined using Cuthill-McKee numbering
- more sophisticated numbering exists
- re-numbering and re-meshing being expensive
 - the mesh should kept constant on a time interval

- each domain has almost the same number of cells
- domain are defined using Cuthill-McKee numbering
- more sophisticated numbering exists
- re-numbering and re-meshing being expensive
 - the mesh should kept constant on a time interval
 - AMR time-step computed through the smallest block and not the smallest cell $T_{n+1} T_n = \Delta T_{AMR}$ is given by the CFL

$$\Delta T_{\text{AMR}} \leq \beta \frac{\min_{k} h_{\text{block}_{k}}}{\max_{k} \| \boldsymbol{u}_{\text{block}_{k}} \|}, \ 0 < \beta \leq 1.$$

$$T_{0} - T_{1} - T_{1} - T_{0} - T_{1} - T_{0} - T_$$

- each domain has almost the same number of cells
- domain are defined using Cuthill-McKee numbering
- more sophisticated numbering exists
- re-numbering and re-meshing being expensive
 - the mesh should kept constant on a time interval
 - AMR time-step computed through the smallest block and not the smallest cell
 - Gain is important and numerical stability is conserved !

EXAMPLES :

• A two dimensional example of BB-AMR with 3 domains and 9 blocks.



Back to slide

EXAMPLES :

• A two dimensional example of BB-AMR with 3 domains and 9 blocks.



• A three dimensional example of BB-AMR with 3 domains and 27 blocks.



Back to slide