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# Ale with Mixed Elements

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# Introduction

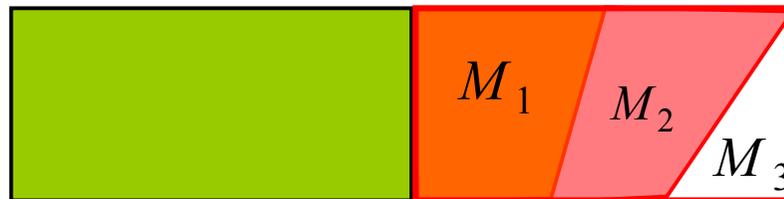
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- Because hydrocodes with lagrangian scheme lack of robustness specially on thin meshes, an ALE formulation with multi-material elements has been implemented to perform simulations with small sizes cells :
  - We introduce ALE blocks with several materials in the same mesh (mixed and pure cells) which can interact with each other or with pure Lagrangian by slide lines.
  - In a same simulation, we can track accurately material boundaries (sliding, void opening and closure) and high material deformation.

$\Omega_1^L$  = Lagrangian

$\Omega_2^A$  = Ale block with M1/M2/M3 materials



## Main features



# Governing equations

<p>Notations</p> <p>Time derivative :</p> <p>Lagrangian :</p> <p>Ale Reference system :</p>	<p><math>u(x, t)</math> : material velocity</p> <p><math>v(x, t)</math> : mesh velocity</p> $\frac{d}{dt} = \frac{\partial}{\partial t} + u \cdot \nabla$ $\frac{D}{Dt} = \frac{\partial}{\partial t} + v \cdot \nabla$	$\left\{ \begin{array}{l} \frac{D\rho}{Dt} + \rho \nabla \cdot u = -(u - v) \cdot \nabla \rho \\ \frac{D\rho u}{Dt} + \nabla p = -(u - v) \cdot \nabla \rho u \\ \frac{D\rho e}{Dt} + p \nabla \cdot u = -(u - v) \cdot \nabla \rho e \end{array} \right.$	$\frac{D}{Dt} = \frac{\partial}{\partial t} - (u - v) \cdot \nabla$
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## In Practice : Splitting method

Lagrangian phase :  $u = v$

$$\left\{ \begin{array}{l} \rho V = Cste \\ \rho \frac{du}{dt} = -\nabla p \\ \frac{de}{dt} + p \frac{dV}{dt} = 0 \\ p = \text{EOS}(e, \rho) \end{array} \right.$$

Advection phase :

- 1 – Mesh smoothing under criteria
- 2 – Remapping : nodal or cell values ( $\rho$ ,  $\rho u$ ,  $\rho e$ ) from the old mesh onto the new mesh

## ● Splitting method

### ■ Lagrangian step : with multi-material cells

- Use the classical Wilkins second order scheme
- Special treatment in multi-material cells :
  - ⇒ Assumption of equal material volumetric strain for all materials
  - ⇒ Interface reconstruction

### ■ Advection step :

- Mesh smoothing under criteria : a new grid is defined.
- Remapping phase : the material quantities are interpolated from the old configuration onto the new one.
  - ⇒ Material by material, after material mesh reconstruction
  - ⇒ Dukowicz method (intersection old mesh / new mesh), second order accuracy

# Lagrangian step with multi-material cells



- **Treatment of the multi-material cell :** Note Volume fraction  $f_\alpha = \frac{V_\alpha}{V}$

- **Assumption : equal volumetric strain for all materials**

Only one velocity for all material :

$$u = u_\alpha$$

→ imply the relation closure for volume fraction in a cell :

$$\frac{dV_\alpha}{dt} = V_\alpha \nabla \cdot u_\alpha = V_\alpha \nabla \cdot u = f_\alpha V \nabla \cdot u = f_\alpha \frac{dV}{dt}$$

→ equal material volumetric strain rate for all materials

$$\frac{df_\alpha}{dt} = 0$$

→ And so with the internal energy balance equation, we have the average pressure definition

$$p = \sum_\alpha p_\alpha f_\alpha$$

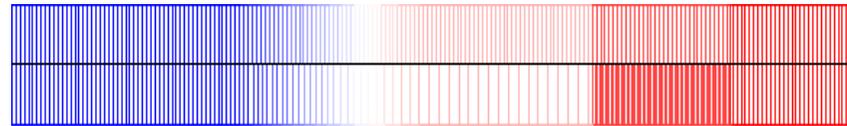
- **More robust multi-material cell treatment : Relaxation pressure**  
we perform an iterative method after Lagrangian step (without volume variation) to balance material pressure in a cell.

# Lagrangian step with multi-material cells



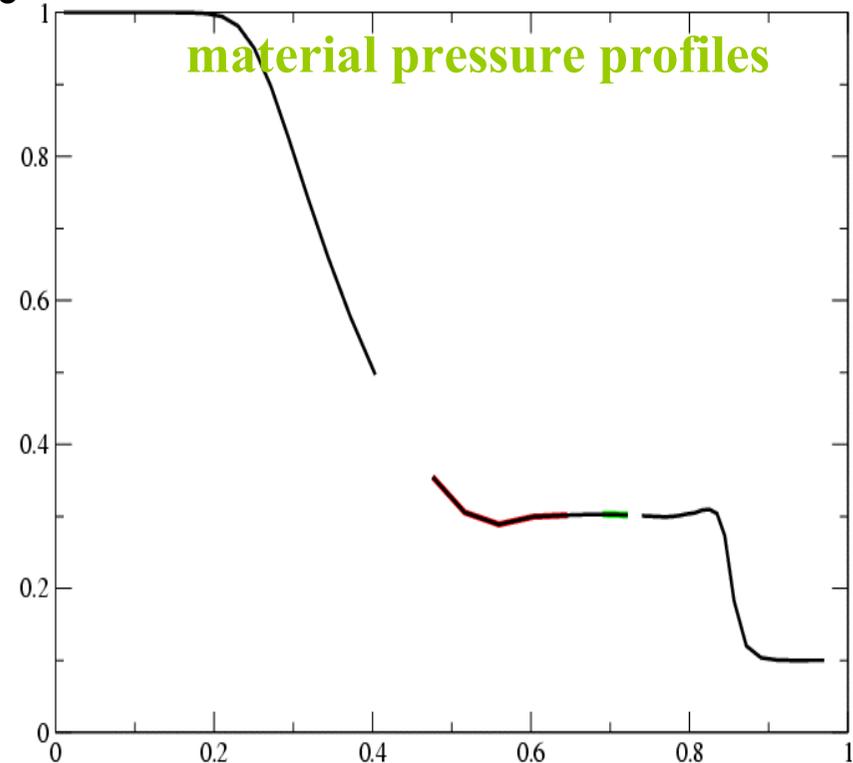
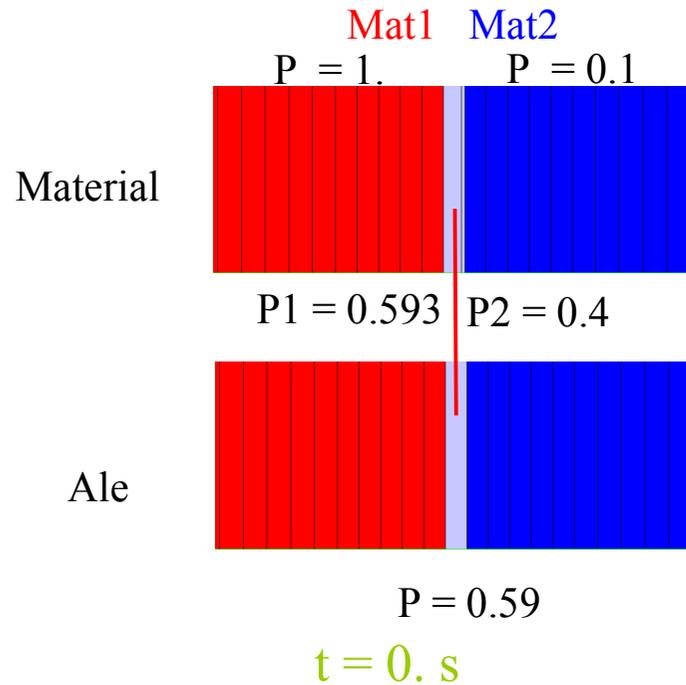
## Sod shock tube problem

ALE



Lagrange

material pressure profiles



t = 0.2 s

# Lagrangian step with multi-material cells



- **Elastic-plastic treatment**

- For each elastic material  $\alpha$ , characterized by mechanical characteristics  $(\mu_\alpha, Y_\alpha)$ , we compute the deviatoric stress tensor:

Same deformation Assumption



$$\left[ \begin{array}{l} \text{Strain rate } \overline{\overline{D}}_\alpha^{n+1/2} = \overline{\overline{D}}^{n+1/2}; \quad D_{ij} = \frac{1}{2} \left( \frac{\partial u_i}{\partial x_j} + \frac{\partial u_j}{\partial x_i} \right) \\ \text{Spin tensor } \overline{\overline{W}}_\alpha^{n+1/2} = \overline{\overline{W}}^{n+1/2}; \quad W_{ij} = \frac{1}{2} \left( \frac{\partial u_i}{\partial x_j} - \frac{\partial u_j}{\partial x_i} \right) \end{array} \right.$$

$$\left\{ \begin{array}{l} \overline{\overline{S}}_{el\alpha}^{n+1} = \overline{\overline{S}}_\alpha^n + \overline{\overline{R}}_\alpha^n + 2\mu_\alpha \left[ \overline{\overline{D}}^{n+1/2} - \frac{1}{3} \text{tr}(\overline{\overline{D}}^{n+1/2}) \cdot \overline{\overline{I}} \right] \\ \overline{\overline{S}}_\alpha^{n+1} = f(Y_\alpha) \overline{\overline{S}}_{el\alpha}^{n+1} \end{array} \right.$$

- On the Ale Block : For the momentum conservative equation

$$\left\{ \begin{array}{l} \text{Deviatoric stress tensor } \overline{\overline{S}}^{n+1} = \sum_\alpha f_\alpha \cdot \overline{\overline{S}}_\alpha^{n+1} \\ \text{Stress tensor } \overline{\overline{\sigma}}^{n+1} = -P \cdot \overline{\overline{I}} + \overline{\overline{S}}^{n+1} \end{array} \right. \quad \text{Average quantities}$$

The average stress is the average of the individual material stresses weighted by the volume fraction.

# Advection step : **Mesh smoothing under criteria**

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- Object : improve the mesh quality at the end of the lagrangian step, by finding some ideal position for the nodes, ideal in the sense of minimizing element distortions.
- We use zones definitions which allow user to specify and to adjust locally the mesh smoothing.
  - We define nodal remapping criteria to identify the nodes which need to be relaxed :
    - For boundary nodes : distance criterion
    - For internal nodes : volume criterion, angular criterion ...
  - The boundary nodes are moved by specific algorithms (specially for the sliding material boundaries) ;
    - To apply an equidistant node distribution along the boundary,
    - To preserve the shape of the boundary.
  - The new position of internal nodes is calculated from classical iterative mesh smoothing algorithms, with a possibility of remapping constraint :
    - Winslow's equipotential relaxation method
    - Simple average method

# Multi-material cell advection step



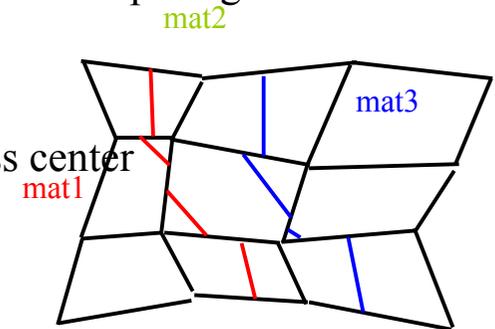
## ● Remapping phase :

### ■ Interface Material Reconstruction in mixed cells

- Multi-material cells are characterized by their volume fractions.
- To track the material interface, we use the Young's method : the slope is given by the gradient of the volume fractions field :

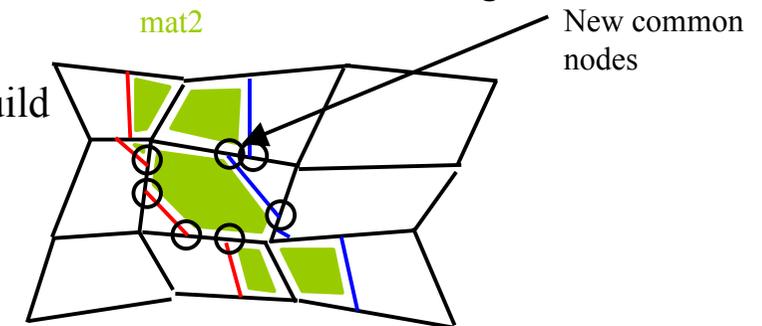
$$\vec{N} = -\vec{\nabla}f_{\beta} \text{ where } \beta \text{ realizes } \max_{\alpha=1,\dots,\text{nbmat}} \|\vec{\nabla}f_{\alpha}\|$$

- Materials are ordered in a cell by locating each material mass center along the normal direction.
- Only **one** normal for all interfaces : onion skin method



### ■ Material mesh reconstruction

- From the old grid, we build nodal-cell connectivity for each material :
  - ⇒ We create the nodes between the material interface and the mesh edge.
  - ⇒ The common nodes are merged.
  - ⇒ The new material lagrangian mesh is build with non structured connectivity

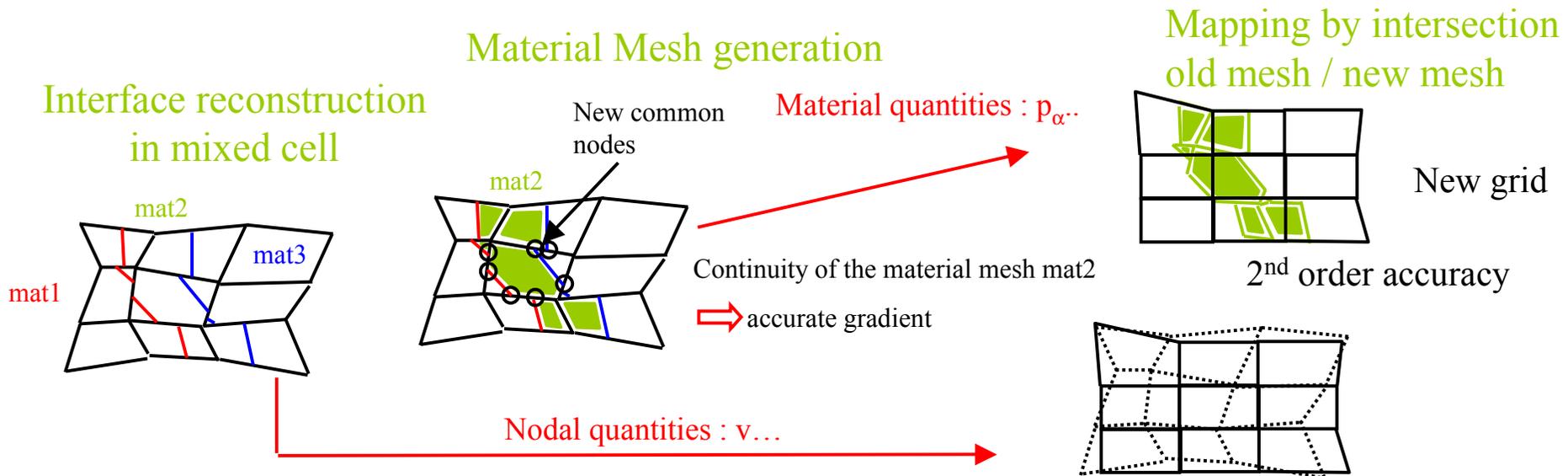


# Multi-material cell advection step



- **Remapping step** : we perform the second order Dukowicz method (intersection old mesh / new mesh) :

- Interfaces Reconstruction
- For each material, we
  - generate an old material mesh,
  - map material cell quantities from this old material mesh onto the new grid.
- Nodal quantities velocities are mapped from the old dual grid (global mesh) onto the new dual one.



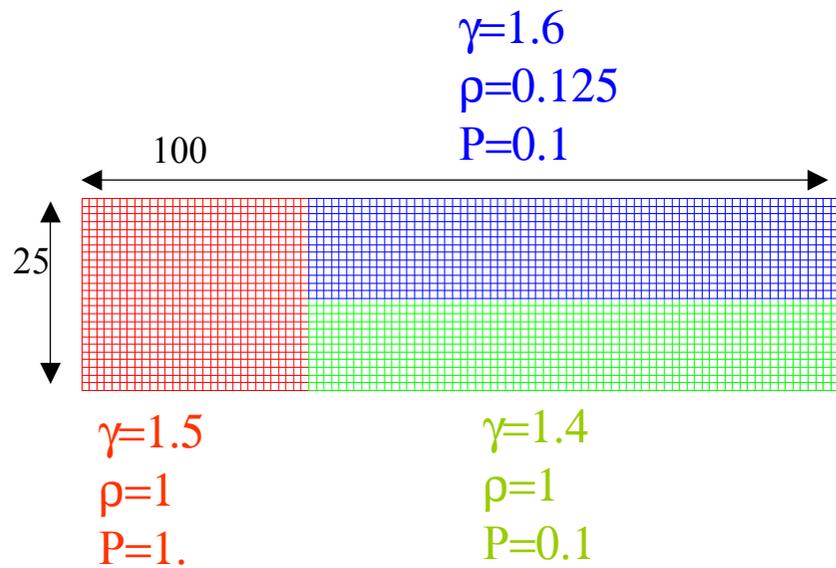
## Some Results

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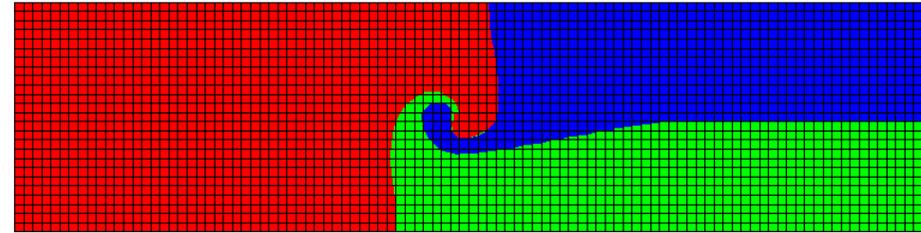


- Instability with 3 materials :
- Taylor bar impact in both Lagrangian and ALE formulation
- Calculation of a dynamic friction experiment

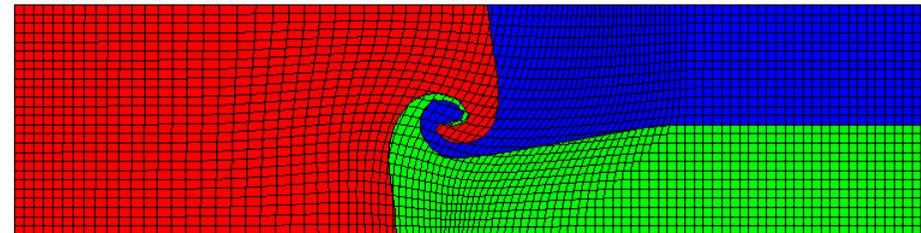
# Instability with 3 Materials



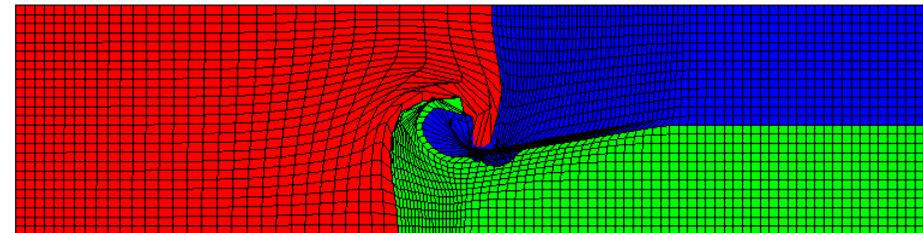
**Eulerian formulation**



**ALE formulation**



**Lagrangian formulation**



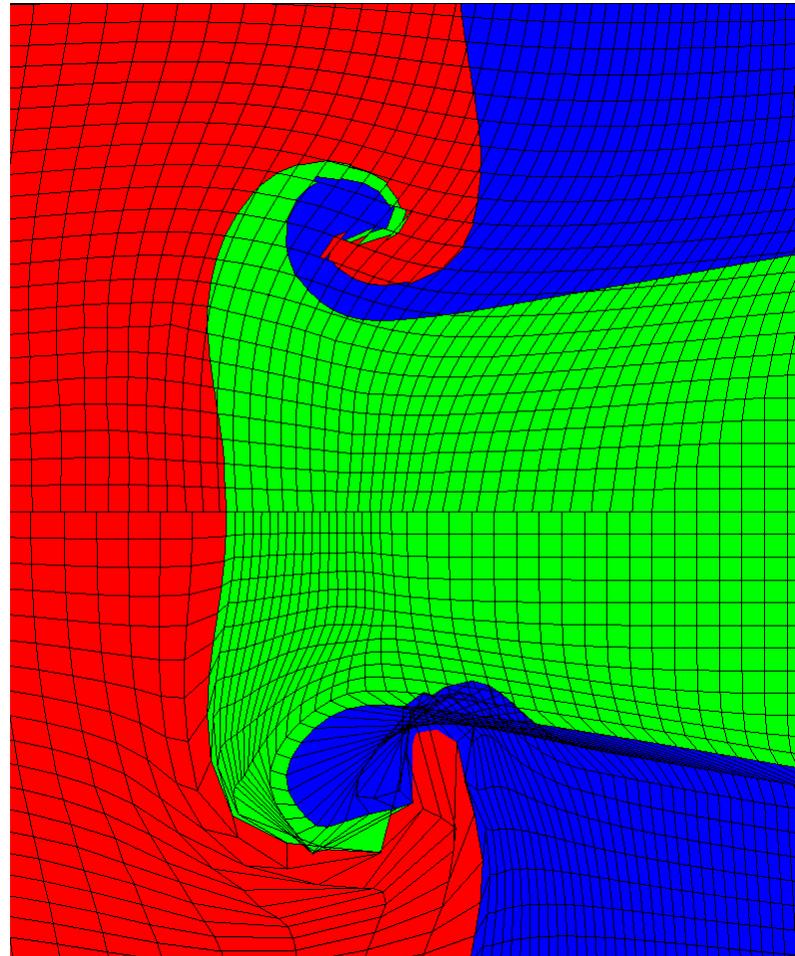
# Instability with 3 materials

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Ale

Lagrangian



Lagrangian Interface  
very distorted.  
The calculation stops.

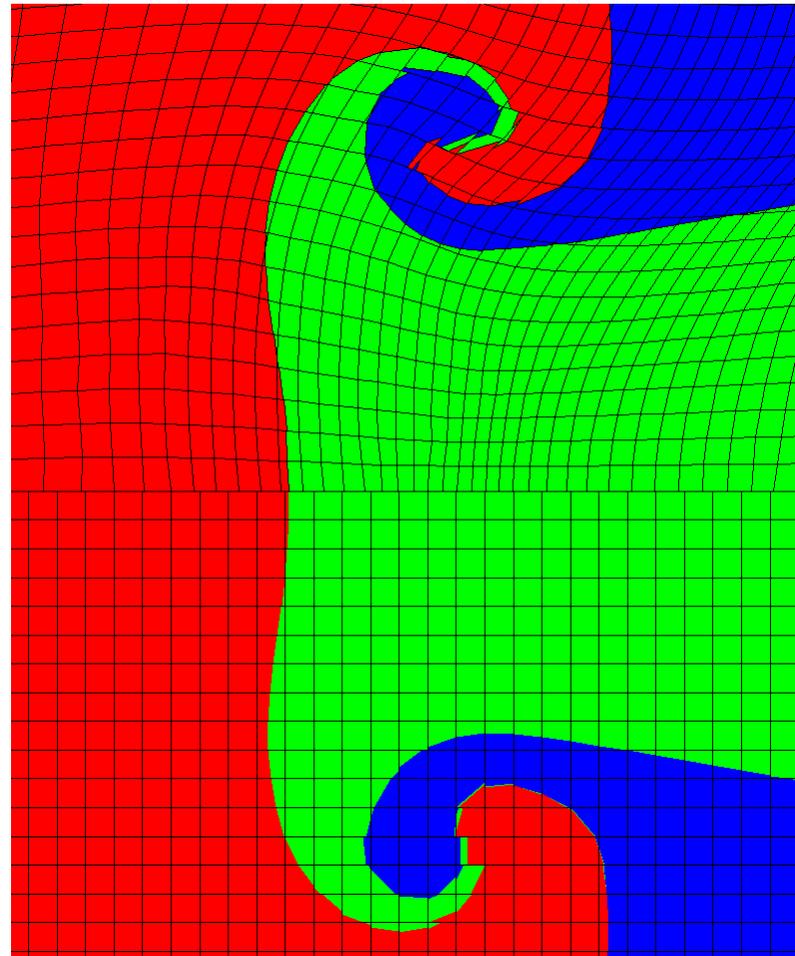
# Instability with 3 materials

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cea

Ale

Eulerian

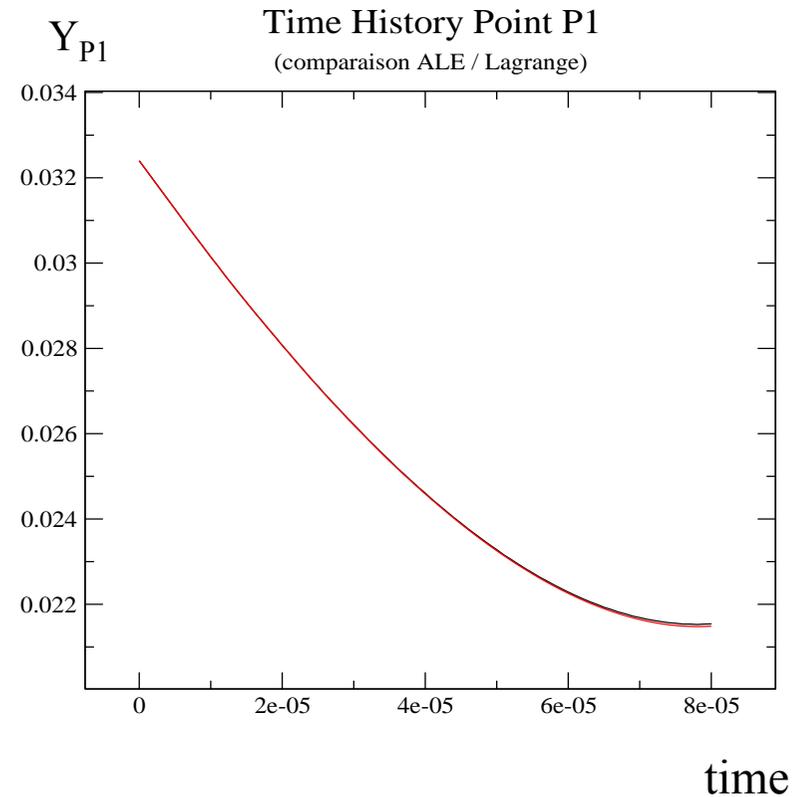
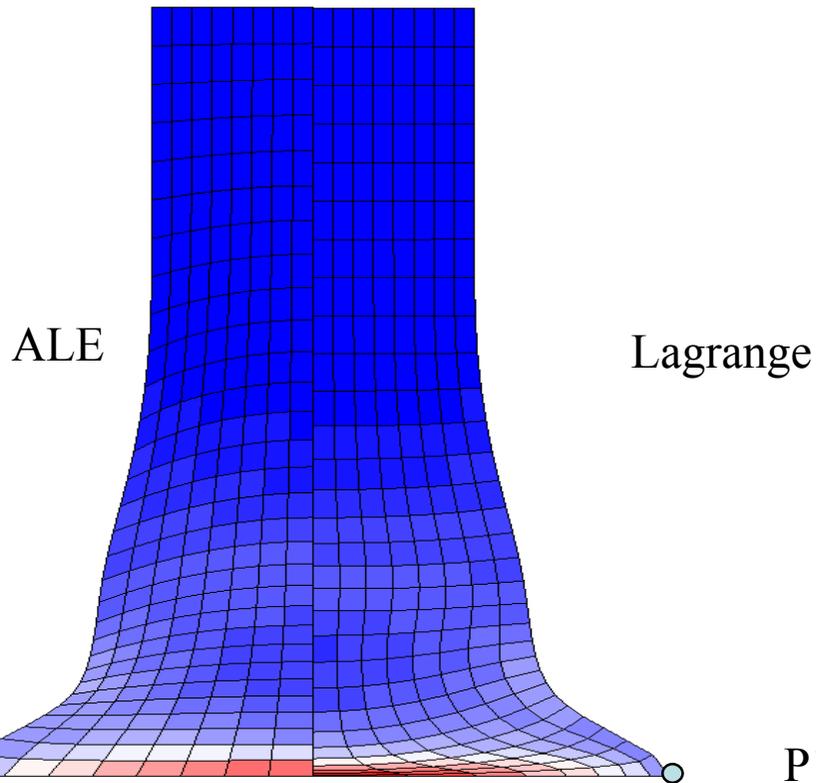


**Same robustness of  
eulerian approach**

# ALE Taylor bar impact



## Plastic strain



Good deformation

Elements severely distorted  
(unphysical mesh distortion  
and hourglassing)

→ Loss of accuracy and reducing the  
time step.

# Dynamic friction experiment

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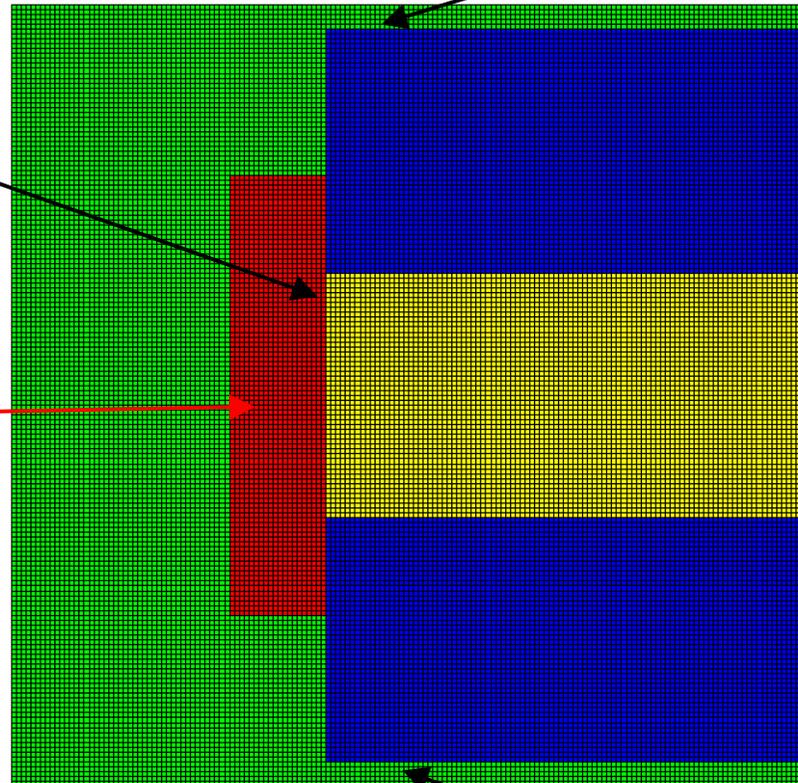


Ale and Lagrangian mesh motion in the same framework

**Sliding with opening**

**Sliding**

**Explosive**



Steel

Aluminium

**Sliding with friction**

ALE = High Explosive / CONF

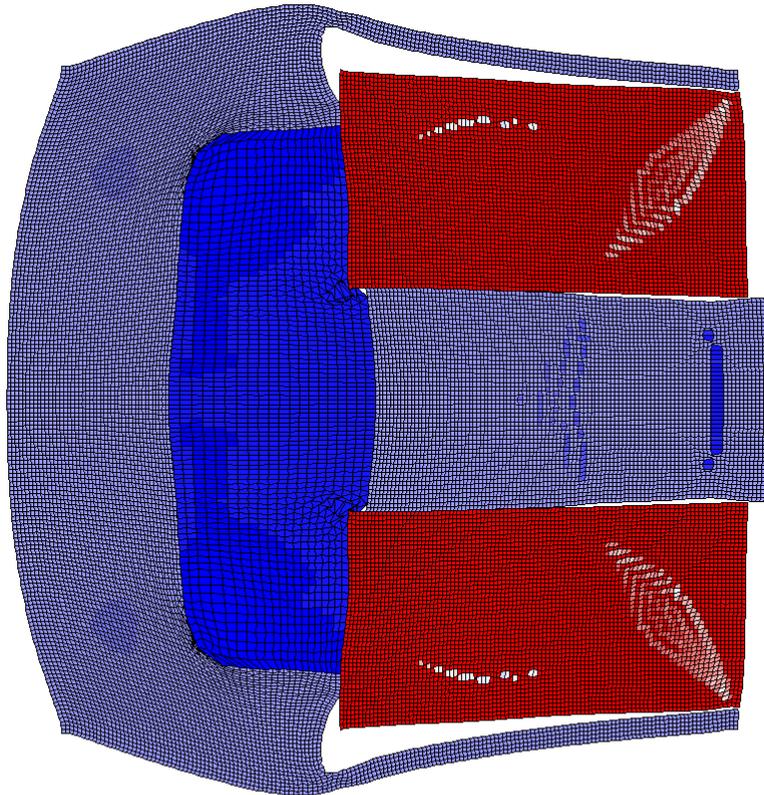
**Sliding with opening**

# Dynamic friction experiment

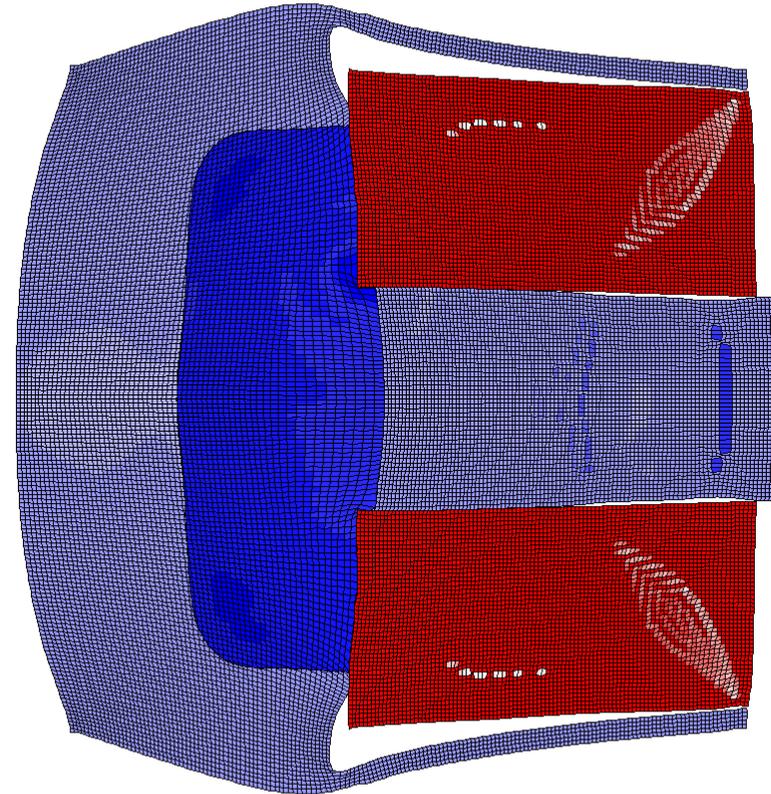
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## Comparison at 30 $\mu$ s



**Pure Lagrangian**



**Lagrangian and ALE**

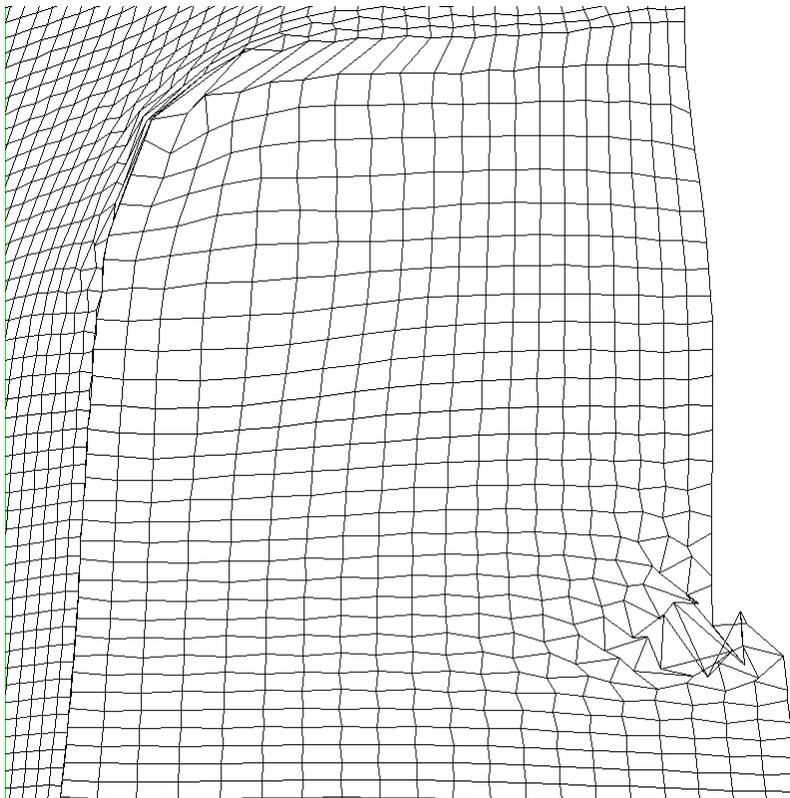
**→ Good agreement with the Lagrangian deformation**

# Dynamic friction experiment

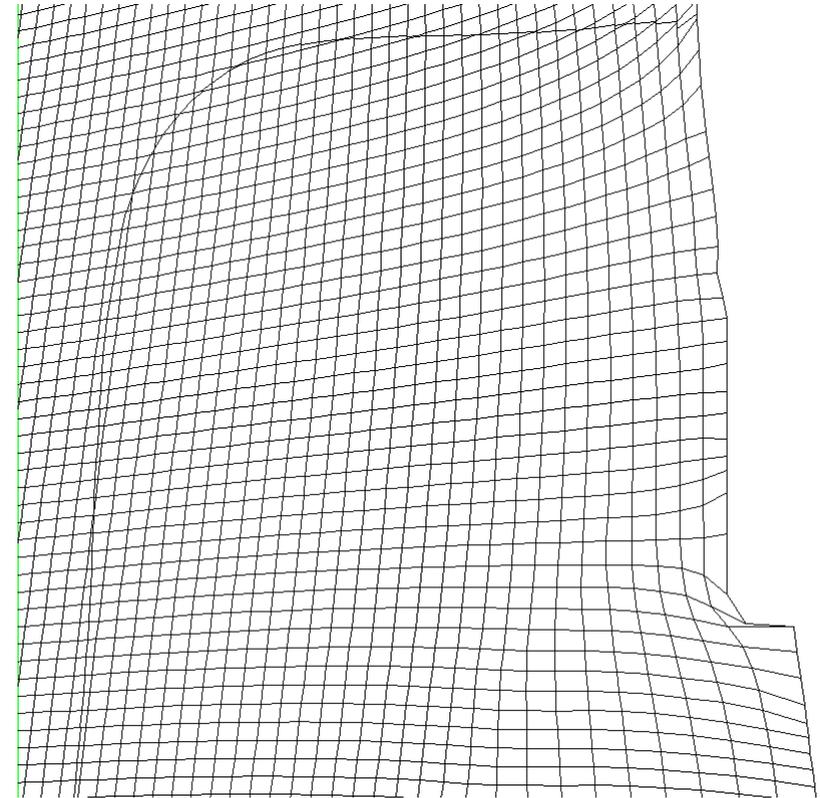
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**Zoom**



Lagrangian materials



ALE block

**→ ALE reduces mesh distortion in the mixed cells but the Ale boundary remains sharp.**

## Conclusion

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- **Ale method with mixed cells :**

- Lagrangian step with treatment of the multi-material cells ( equal volumetric strain assumption + relaxation pressure)
- Full 2D advection step with interface reconstruction in multi-material cells

- **2 different treatments for materials interfaces :**

- Lagrangian interfaces with slidelines
- Multi materials cells.

⇒ Provides the accuracy of Lagrangian formulation and the robustness of the Eulerian Multi materials cells

- **Future works :**

- Extension to the 3D