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On the coupling of Spectral Element Method with Discontinuous Galerkin approximation for elasto-acoustic problems

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¹ Team project Magique.3D, INRIA.UPPA.CNRS, Pau, France.

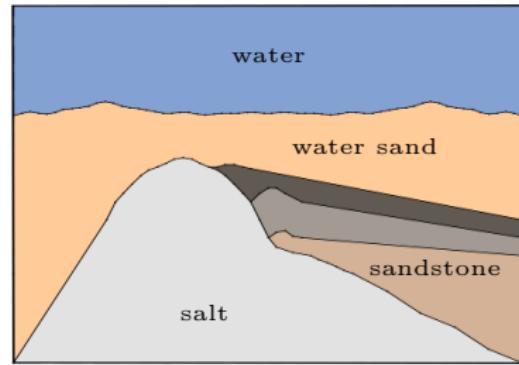
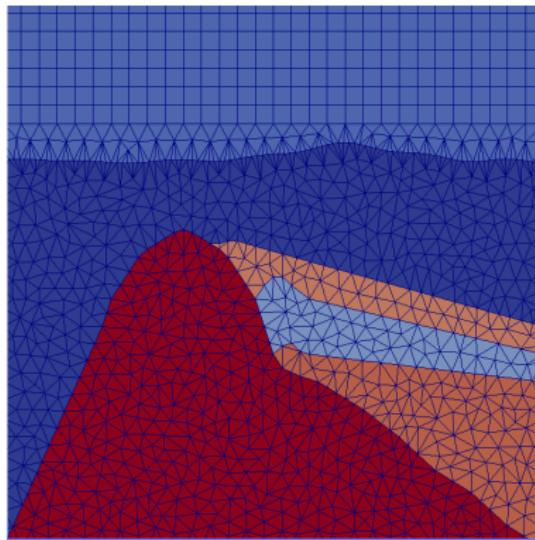
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Why using hybrid meshes?



- Useful when the use of unstructured grid is non-sense (e.g. medium with a layer of water)
- Well suited for the coupling of numerical methods in order to reduce the computational cost and improve the accuracy

Elastodynamic system

$$\begin{cases} \rho(x) \frac{\partial \textcolor{blue}{v}}{\partial t}(x, t) = \nabla \cdot \underline{\underline{\sigma}}(x, t) \\ \frac{\partial \underline{\underline{\sigma}}}{\partial t}(x, t) = \underline{\underline{C}}(x) \underline{\underline{\epsilon}}(\textcolor{blue}{v}(x, t)) \end{cases}$$

With :

- $\rho(x)$ the density
- $C(x)$ the elasticity tensor
- $\epsilon(x, t)$ the deformation tensor
- $v(x, t)$, the wavespeed
- $\underline{\underline{\sigma}}(x, t)$ the strain tensor

Elasticus software

Written in **Fortran** for wave propagation simulation in the **time domain**

Features

- Using various types of meshes (**unstructured triangle and tetrahedra**)
- Modelling of various physics (**acoustic, elastic and elasto-acoustic**)

- **Discontinuous Galerkin Method (DG)** based on **unstructured triangle and unstructured tetrahedra**
- with various time-schemes : **Runge-Kutta (2 or 4), Leap-Frog**
- with **p-adaptivity, multi-order computation...**

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- 5 3D extension**

1 Numerical Methods

- Discontinuous Galerkin Method (DG)
- Spectral Element Method (SEM)
- Advantages of each method

Discontinuous Galerkin Method

Use discontinuous functions :



mesh



continuous

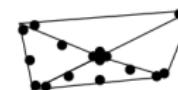


discontinuous

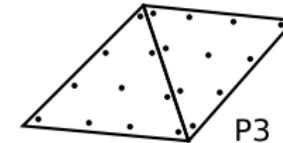
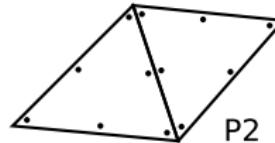
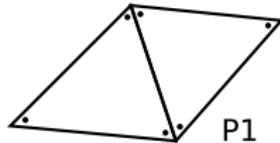
h adaptivity :



p adaptivity :



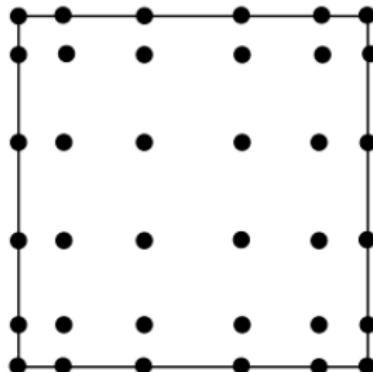
Degrees of freedom on each cell :



Spectral Element Method

General principle

- Finite Element Method (FEM) discretization + Gauss-Lobatto quadrature
- Gauss-Lobatto points as degrees of freedom (exponential convergence on L^2 -norm)



- $\int f(x)dx \approx \sum_{j=1}^{N+1} \omega_j f(\xi_j)$
- $\varphi_i(\xi_j) = \delta_{ij}$

Spectral Element Method

Main change with DG

- DG discontinuous, SEM continuous
- Need of defining local to global numbering
- Global matrices required by SEM
- Basis functions computed differently

Advantages of each method

DG

- Element per element computation (hp -adaptivity)
- Time discretization quasi explicit (block diagonal mass matrix)
- Simple to parallelize

SEM

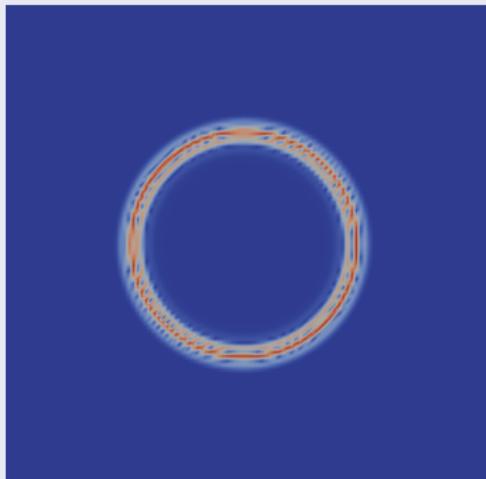
- Couples the flexibility of FEM with the accuracy of the pseudo-spectral method
- Reduces the computational cost when using structured meshes in comparison with DG

2 Comparison DG/SEM on structured quadrangle mesh

- DG/SEM comparison on quadrangle mesh
- Description of the test cases
- Comparative tables

Description of the test cases

Physical parameters



P wavespeed	1000 m.s^{-1}
Density	1 kg.m^{-3}

Second order Ricker Source in P wave
($f_{peak} = 10\text{Hz}$)

General context

- Acoustic homogeneous medium
- Four different meshes : **10000 cells, 22500 cells, 90000 cells, 250000 cells**
- CFL computed using **power iteration** method
- Leap-Frog time scheme
- **Four threads parallel execution with OpenMP**

Comparative tables

- Error computed as the difference between an analytical and a numerical solution for each method

Quadrangle mesh 10000 elements:

	CFL	L2-error	CPU-time	Nb of time steps
DG	1.99e-3	2.5e-2	19.30	500
SEM	4.9e-3	1.3e-1	0.36	204

Quadrangle mesh 22500 elements:

	CFL	L2-error	CPU-time	Nb of time steps
DG	1.33e-3	1.8e-2	100.48	750
SEM	3.26e-3	7e-2	1.19	306

Comparative tables

- Error computed as the difference between an analytical and a numerical solution for each method

Quadrangle mesh 10000 elements:

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SEM	4.9e-3	1.3e-1	0.36	204
SEM(DG CFL)	1.99e-3	4.8e-2	0.73	502

Quadrangle mesh 22500 elements:

	CFL	L2-error	CPU-time	Nb of time steps
DG	1.33e-3	1.8e-2	100.48	750
SEM	3.26e-3	7e-2	1.19	306
SEM(DG CFL)	1.33e-3	1.2e-2	2.82	751

SEM fifty time faster than DG on a mesh with 22500 cells

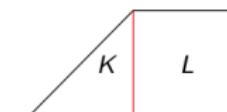
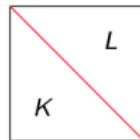
3 DG/SEM coupling

- Hybrid meshes structures
- Variational formulation
- Space discretization

Hybrid meshes structures

- Aim at coupling P_k and Q_k structures.
- Need to extend or split some of the structures (e.g. neighbour indexes)
- Define new face matrices

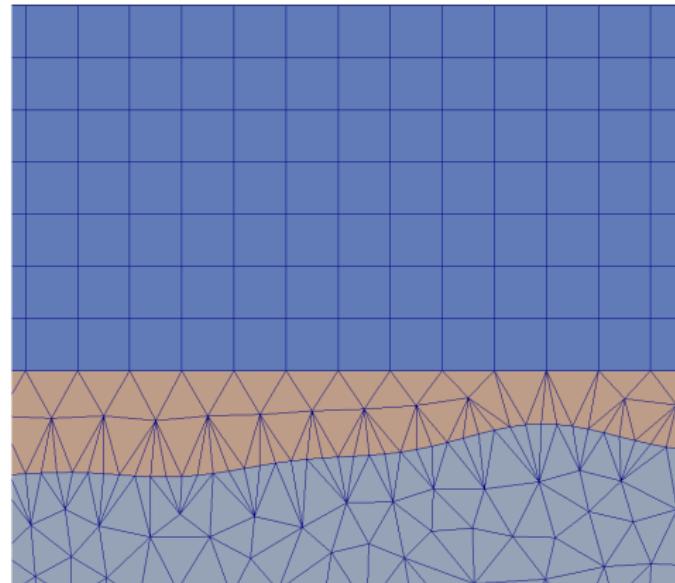
$$M_{ij}^{K,L} = \int_{K \cap L} \phi_i^K \phi_j^L, \quad M_{ij}^{K,L} = \int_{K \cap L} \psi_i^K \psi_j^L, \quad M_{ij}^{K,L} = \int_{K \cap L} \phi_i^K \psi_j^L$$



Variational formulation

Global context

- Domain in two parts : $\Omega_{h,1}$ (**structured quadrangle + SEM**), $\Omega_{h,2}$ (**unstructured triangle + DG**)



Variational formulation

SEM variational formulation :

$$\begin{cases} \int_{\Omega_{h,1}} \rho \partial_t v_1 \cdot w_1 = - \int_{\Omega_{h,1}} \sigma_1 \cdot \nabla w_1 + \int_{\Gamma_{out,1}} (\sigma_1 \mathbf{n}_1) \cdot w_1 \\ \int_{\Omega_{h,1}} \partial_t \sigma_1 : \xi_1 = - \int_{\Omega_{h,1}} (\nabla(C\xi_1)) \cdot v_1 + \int_{\Gamma_{out,1}} (C\xi_1 \mathbf{n}_1) \cdot v_1 \end{cases}$$

DG variational formulation :

$$\begin{cases} \int_{\Omega_{h,2}} \rho \partial_t v_2 \cdot w_2 = - \int_{\Omega_{h,2}} \sigma_2 \cdot \nabla w_2 + \int_{\Gamma_{out,2}} (\sigma_2 \mathbf{n}_2) \cdot w_2 + \int_{\Gamma_{int}} \{\{\sigma_2\}\} [[w_2]] \cdot \mathbf{n}_2 \\ \int_{\Omega_{h,2}} \partial_t \sigma_2 : \xi_2 = - \int_{\Omega_{h,2}} (\nabla(C\xi_2)) \cdot v_2 + \int_{\Gamma_{out,2}} (C\xi_2 \mathbf{n}_2) \cdot v_2 + \int_{\Gamma_{int}} \{\{v_2\}\} [[C\xi_2]] \cdot \mathbf{n}_2 \end{cases}$$

Variational formulation

Add the average of the solution of each part at the interface + put $\sigma_* \mathbf{n}_* = 0$

$$\begin{cases} \int_{\Omega_{h,1}} \rho \partial_t v_1 \cdot w_1 = - \int_{\Omega_{h,1}} \sigma_1 \cdot \nabla w_1 + \frac{1}{2} \int_{\Gamma_{1/2}} (\sigma_1 + \sigma_2) \mathbf{n}_1 \cdot w_1 \\ \int_{\Omega_{h,1}} \partial_t \sigma_1 : \xi_1 = - \int_{\Omega_{h,1}} (\nabla(C\xi_1)) \cdot v_1 + \frac{1}{2} \int_{\Gamma_{1/2}} (C\xi_1 \mathbf{n}_1) \cdot (v_1 + v_2) \end{cases}$$

$$\begin{cases} \int_{\Omega_{h,2}} \rho \partial_t v_2 \cdot w_2 = - \int_{\Omega_{h,2}} \sigma_2 \cdot \nabla w_2 + \int_{\Gamma_{int}} \{\{\sigma_2\}\} [[w_2]] \cdot \mathbf{n}_2 - \frac{1}{2} \int_{\Gamma_{1/2}} (\sigma_1 + \sigma_2) \mathbf{n}_1 \cdot w_2 \\ \int_{\Omega_{h,2}} \partial_t \sigma_2 : \xi_2 = - \int_{\Omega_{h,2}} (\nabla(C\xi_2)) \cdot v_2 + \int_{\Gamma_{int}} \{\{v_2\}\} [[C\xi_2]] \cdot \mathbf{n}_2 \\ - \frac{1}{2} \int_{\Gamma_{1/2}} (C\xi_2 \mathbf{n}_1) \cdot (v_1 + v_2) \end{cases}$$

Energy study

Goal : Show that our coupling preserves the energy

- We set $\xi_1 = \sigma_1, \xi_2 = \sigma_2, w_1 = v_1, w_2 = v_2$
- We add the equations of the two parts variational formulation

$$\frac{d}{dt} \mathbb{E} = 0$$

Space discretization : SEM part

- φ_i : SEM basis functions
- ψ_i : DG basis functions

$$\begin{cases} M_{\boldsymbol{\nu}_1} \partial_t \boldsymbol{\nu}_{h,1} + R_{\boldsymbol{\sigma}_1} \boldsymbol{\sigma}_{h,1} + \textcolor{blue}{R_{\boldsymbol{\sigma}_2}^{2,1} \boldsymbol{\sigma}_{h,2}} = 0 \\ M_{\boldsymbol{\sigma}_1} \partial_t \boldsymbol{\sigma}_{h,1} + R_{\boldsymbol{\nu}_1} \boldsymbol{\nu}_{h,1} + \textcolor{blue}{R_{\boldsymbol{\nu}_2}^{2,1} \boldsymbol{\nu}_{h,2}} = 0 \end{cases}$$

- $M_{ij} = \int_{\Omega} \varphi_i \varphi_j \approx \sum_{e \in \text{supp}(\varphi_i) \cap \text{supp}(\varphi_j)} \sum_{k=1}^{(r+1)^d} \omega_k \varphi_i(\xi_k) \varphi_j(\xi_k) = \sum_{e \in \text{supp}(\varphi_i) \cap \text{supp}(\varphi_j)} \omega_i \delta_{i,j}$ the mass matrix
- $R_{p_{ij}} = \int_{\Omega} \varphi_i \frac{\partial \varphi_j}{\partial p}$ stiffness matrix

Matrix of DG/SEM coupling :

$$R_{\boldsymbol{\sigma}_2, ij}^{2,1} = \frac{1}{2} \int_{\partial\Omega_1 \cap \partial\Omega_2} \psi_i \varphi_j$$

Space discretization : DG part

$$\begin{cases} \rho M_{\mathbf{v}_2} \partial_t \mathbf{v}_{h,2} + R_{\sigma_2} \boldsymbol{\sigma}_{h,2} - R_{\sigma_1}^{1,2} \boldsymbol{\sigma}_{h,1} = 0 \\ M_{\sigma_2} \partial_t \boldsymbol{\sigma}_{h,2} + R_{\mathbf{v}_2} \mathbf{v}_{h,2} - R_{\mathbf{v}_1}^{1,2} \mathbf{v}_{h,1} = 0 \end{cases}$$

- $M_{ij}^K = \int_K \psi_i^K \psi_j^K$ mass matrix,
- $R_{p_{ij}}^K = \int_K \psi_i^K \frac{\partial \psi_j^K}{\partial p}$ stiffness matrix,
- $R_{p_{ij}}^{K,L} = \int_{\partial K \cap \partial L} \psi_i^K \psi_j^L n_K \cdot e_p$ the mass-face matrix

Two new matrices which come from the DG/SEM coupling $R_\star^{1,2}$. Block composed :

$$R_{\mathbf{v}_1}^{1,2} = R_{\sigma_1}^{1,2} = -\frac{1}{2} \int_{\partial \Omega_2 \cap \partial K_1} \psi_j^{K_2} \varphi_i \quad (1)$$

4 Comparison between DG/SEM and DG on hybrid meshes

- Experimentation context
- Comparative tables

Context

- Acoustic homogeneous medium
- 54000 triangles
- 21000 quadrangles
- Using Leap-Frog time scheme
- Parallel computation using OpenMP
- Done with different orders of discretization

Comparative tables

$P_1 - Q_3$ computation :

	CFL	L_2 -error	CPU-time
DG	1e-5	0.03	7343.92
DG/SEM	1e-5	0.03	823.22

$P_3 - Q_1$ computation :

	CFL	L_2 -error	CPU-time
DG	3e-5	0.009	3078.15
DG/SEM	3e-5	0.01	2951

$P_2 - Q_3$ computation :

	CFL	L_2 -error	CPU-time
DG	1e-5	0.002	9452.73
DG/SEM	1e-5	0.003	1393.80

$P_3 - Q_2$ computation :

	CFL	L_2 -error	CPU-time
DG	1e-5	5.4e-4	9951.60
DG/SEM	1e-5	0.007	3122

General settings

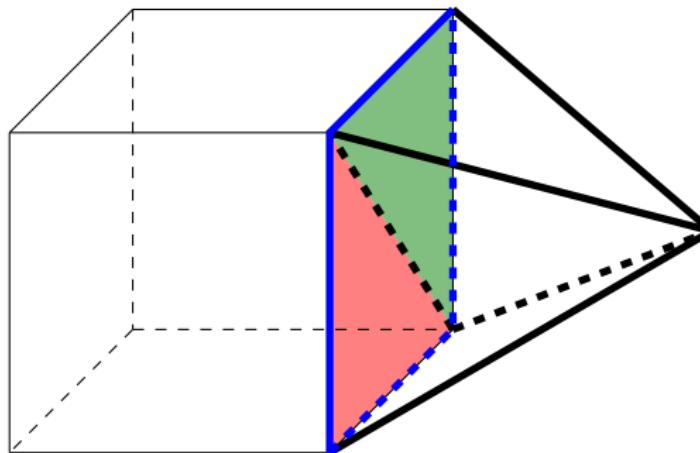
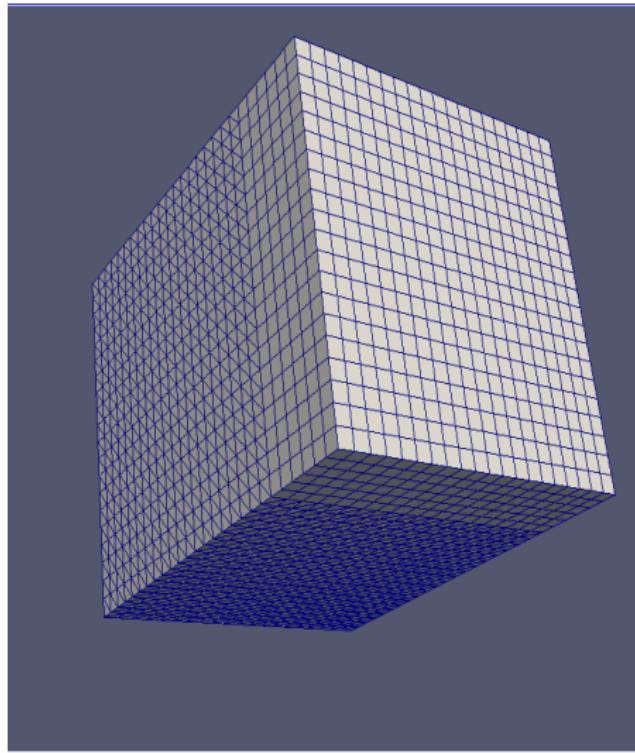


Figure: Hexa/Tet boundary configuration

- Only deal with a simple case of 3D hybrid meshes : one hexahedra has only two tetrahedra as neighbour
- Extend SEM in 3D (basis functions...)
- Require introducing a new matrix which handles the rotation cases between two elements



Conclusion and perspectives

Conclusion

- ① Build a variational formulation for DG/SEM coupling and find a CFL condition that ensures stability
- ② As expected, SEM is more efficient on structured quadrangle mesh than DG
- ③ Show the utility of using hybrid meshes and method coupling (reduce computational cost,...)

Perspectives

- Implement DG/SEM coupling on the code (2D) ✓
- Develop DG/SEM coupling in 3D ✓
- Add a local time-stepping scheme
- Develop PML in the hexahedral part

Thank you for your attention !

Questions?