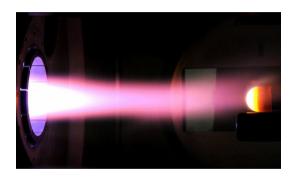
Development of a multi-domain hybridized discontinuous Galerkin solver for inductively coupled plasma



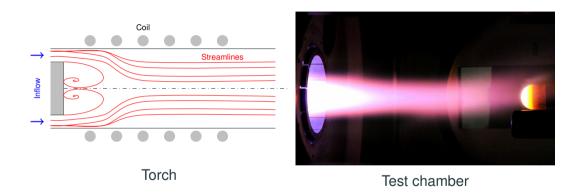
Corthouts Nicolas, Hillewaert Koen, May Georg, Magin Thierry





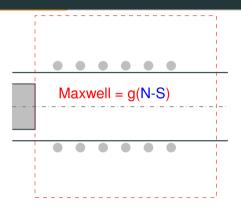


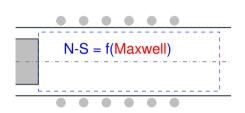
Context of research



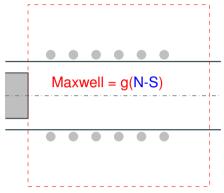
 $\textit{Re} \sim 100 \quad \textit{Ma} \sim 0.001 \quad \rho \simeq \rho(\textit{T}).$

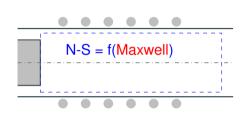
ICP: segregated approach of previous solvers





ICP: segregated approach of previous solvers

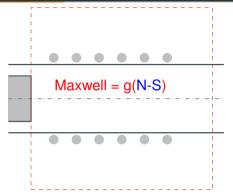


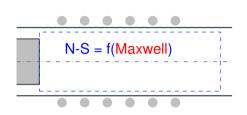


Pros

- It works (see Magin, 2004).
- Allows to freeze the electric field in unsteady simulations.

ICP: segregated approach of previous solvers





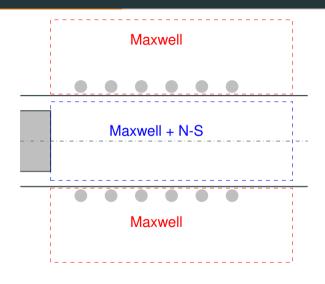
Pros

- It works (see Magin, 2004).
- Allows to freeze the electric field in unsteady simulations.

Cons

 Convergence can be hard to achieve (O(1000) iterations with COOLFluiD).

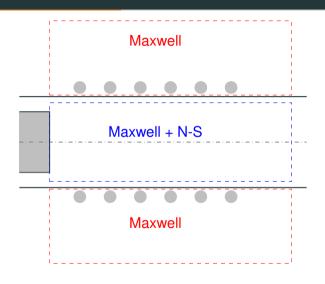
A multi-domain solver



Two approaches

- MONOLITHIC: system solved as a whole.
- COUPLED: two solvers that exchange interface data.

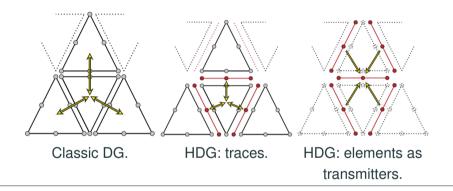
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Two approaches

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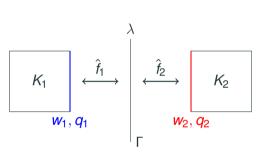
The numerical method: HDG



- 1. **Local systems** of element size solved directly.
- 2. A global system smaller than the global DG system.

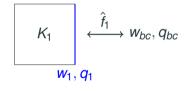
Multi-domain HDG

Conservativity



$$\int_{\Gamma} \left[\hat{f}_1(w_1, q_1, n_1) + \hat{f}_2(w_2, q_2, n_2) \right] \mu dS = 0.$$

Multi-domain HDG



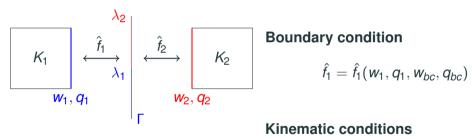
Conservativity

$$\int_{\Gamma} \left[\hat{f}_1(w_1, q_1, n_1) + \hat{f}_2(w_2, q_2, n_2) \right] \mu dS = 0.$$

Boundary condition

$$\hat{f}_1 = \hat{f}_1(w_1, q_1, w_{bc}, q_{bc})$$

Multi-domain HDG



Conservativity

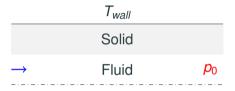
$$\int_{\Gamma} \left[\hat{f}_1(w_1, q_1, n_1) + \hat{f}_2(w_2, q_2, n_2) \right] \mu dS = 0.$$

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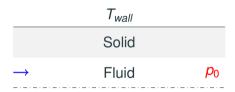
Kinematic conditions

$$\int_{\Gamma}\mathcal{F}(\lambda_{\mathsf{1}},\lambda_{\mathsf{2}})\mu d\mathcal{S}=0$$

Application: Conjugate heat transfer



Application: Conjugate heat transfer

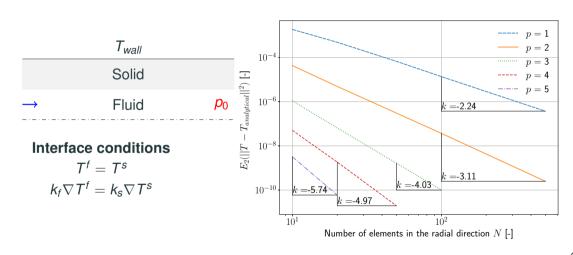


Interface conditions

$$T^f = T^s$$

$$k_f \nabla T^f = k_s \nabla T^s$$

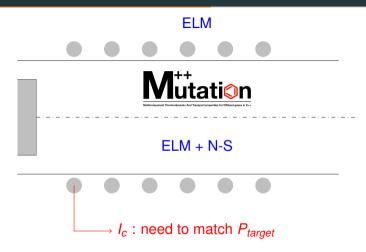
Application: Conjugate heat transfer

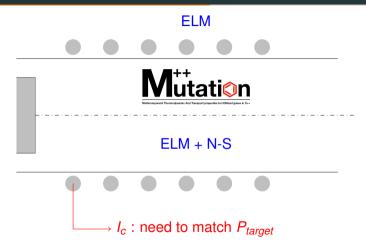










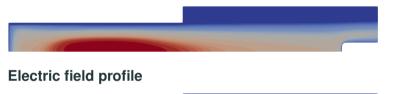


AUSM numerical + low-mach preconditioning (Magin 2004) and Damped Newton-Raphson method.

Temperature profile

$$T_{min} = 350 \text{ K}$$
 $T_{max} = 11000 \text{ K}$

Temperature profile



$$T_{min} = 350 \text{ K}$$

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$$E_{min} = 0 \text{ V}$$

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Temperature profile



$$T_{min} = 350 \text{ K}$$
 $T_{max} = 11000 \text{ K}$

Electric field profile



$$E_{min} = 0 \text{ V}$$

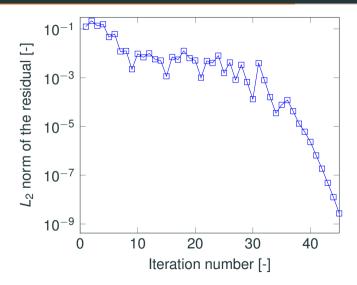
 $E_{max} = 3650 \text{ V}$

Power dissipated in the facility



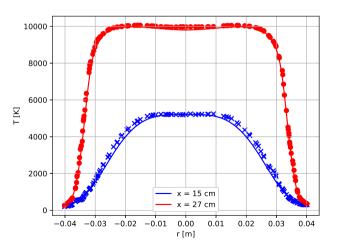
$$P_{min} = 0 \text{ W/m}^3$$
$$P_{max} = 10^{11} \text{ W/m}^3$$

Convergence history



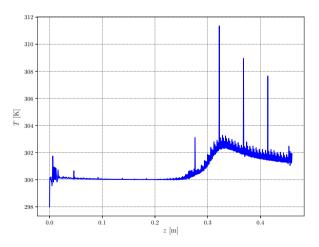
Application to ICP: quantitative results for the mini-torch

Comparison with results of previous ICP code (AUSM flux, p = 2, swirl = 45°).



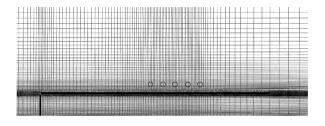
Application to ICP: oscillations near the wall

Temperature oscillations in the near wall region.

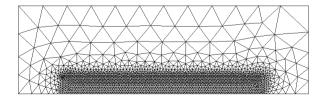


ICP: mesh comparison

FV mesh



ICP mesh



Conclusions and future work

- A versatile tool has been implemented in the HDG code.
- High order ICP flows are now possible.
- Possibility of extending to various physical situation.
- High order methods are prone to oscillations. We are working on them.
- Overall, the code is not very robust, and a slight change in mesh configuration or simulation parameters can lead to unstable simulations.