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Mixed Finite Element Formulations for Systems with Superconductors and Ferromagnetic Materials

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Introduction

We model eddy current problems for high-temperature superconductors (HTS) and ferromagnetic materials (FM).

Coated HTS tape



[Solovyov, Supercond. Sci. Technol., 2013]

Magnetic levitation



[Huang, Supercond. Sci. Technol., 2015]

Trapped-field magnet



Magnetic cloak



[Capobianco-Hogan, Nucl. Instrum. Methods Phys. Res., 2018]

[Philipe, Physica C Superconductivity, 2014]

Context

Coupled formulations offer many advantages for HTS-FM modeling:

- improved efficiency for nonlinear system resolution,
- reduced number of DOFs,
- increased flexibility...

However, they enter the framework of mixed formulations, thus requiring to be extremely careful regarding function spaces.

Otherwise, non-physical results must be expected:



Strong form

Magnetodynamic (quasistatic) equations

div
$$\boldsymbol{b} = 0$$
, curl $\boldsymbol{h} = \boldsymbol{j}$, curl $\boldsymbol{e} = -\partial_t \boldsymbol{b}$.

Constitutive relationships

High-temperature superconductors (HTS):

$$\boldsymbol{e} =
ho(\|\boldsymbol{j}\|)\boldsymbol{j}$$
 and $\boldsymbol{b} = \mu_0 \boldsymbol{h},$

with the power law $\rho(\|\boldsymbol{j}\|) = \frac{e_c}{j_c} \left(\frac{\|\boldsymbol{j}\|}{j_c}\right)^{n-1}$.

Ferromagnetic material (FM):

$$\boldsymbol{b} = \mu(\boldsymbol{b}) \boldsymbol{h}$$
 and $\boldsymbol{j} = \boldsymbol{0}$.



Dual formulations

Two classes of formulations with the finite element method:

- h-conform, e.g. h-formulation ,
 - enforces the continuity of the tangential component of h,
 - involves $\boldsymbol{e} = \rho \boldsymbol{j}$ and $\boldsymbol{b} = \mu \boldsymbol{h}$,
 - with curl h = 0 in non-conducting domain ("h- ϕ "+cuts),

$$\left(\partial_t(\mu \boldsymbol{h})\;, \boldsymbol{h}'\right)_\Omega + \left(\rho \operatorname{\mathbf{curl}} \boldsymbol{h}\;, \operatorname{\mathbf{curl}} \boldsymbol{h}'\right)_{\Omega_{\mathrm{c}}} - \left\langle \boldsymbol{e} \times \boldsymbol{n}\;, \boldsymbol{h}'\right\rangle_{\Gamma_e} = 0.$$

b-conform, e.g. *a*-formulation ,

enforces the continuity of the normal component of b,

• involves $\mathbf{j} = \sigma \mathbf{e}$ and $\mathbf{h} = \nu \mathbf{b}$, $(\sigma = \rho^{-1}, \nu = \mu^{-1})$

$$\left(\nu \operatorname{\mathbf{curl}} \boldsymbol{a} \;, \operatorname{\mathbf{curl}} \boldsymbol{a}'
ight)_{\Omega} + \left(\sigma \; \partial_t \boldsymbol{a} \;, \boldsymbol{a}'
ight)_{\Omega_{\mathrm{c}}} - \left\langle \boldsymbol{h} imes \boldsymbol{n} \;, \boldsymbol{a}'
ight
angle_{\Gamma_h} = 0.$$

Nonlinear constitutive laws involved in opposite ways \Rightarrow very different numerical behaviors are expected... and observed.

Best choice for HTS only

Cycles in iterations:



In the *a*-formulation, the diverging slope associated with $\mathbf{j} = \sigma \mathbf{e}$ for $\mathbf{e} \rightarrow 0$ is really difficult to handle.

 \Rightarrow Among the two simple formulations, the <u>*h*-formulation</u> is much more efficient for systems with HTS:

- with an adaptive time-stepping algorithm,
- solved with a Newton-Raphson method.

Dular, J., et al. (2020) TAS 30 8200113.

Ferromagnetic materials

The nonlinearity is in the magnetic constitutive law.

• *h*-formulation the involved law is $b = \mu h$.

$$\mu$$
 \approx σ

 \Rightarrow Often enters cycles with Newton-Raphson.

OK with fixed point, or N-R with relaxation factors but slow.

 $| a - formulation | the involved law is <math>h = \nu b.$

 \Rightarrow Efficiently solved with Newton-Raphson.

The <u>a-formulation</u> is more appropriate for dealing with the nonlinearity, whereas for HTS, the <u>h-formulation</u> is best.

Coupled materials - *h-a*-formulation

Use the best formulation in each material

Decompose the domain $\Omega,$ for example into:

- $\Omega^h = \{\text{HTS, Air}\}$
- $\Omega^a = \{\text{Ferromagnet}\}$

and couple via $\Gamma_{\text{m}}=\partial(\text{FM})$:



$$egin{aligned} &ig(\partial_t(\mum{h})\;,m{h}'ig)_{\Omega^h} + ig(
ho\; \mathbf{curl}\;m{h}\;,\mathbf{curl}\;m{h}'ig)_{\Omega^h_{\mathbf{c}}} + ig\langle\partial_t a imes m{n}_{\Omega^h}\;,m{h}'ig
angle_{\Gamma_{\mathsf{m}}} = 0, \ &ig\langlem{h} imes m{n}_{\Omega^a}\;,m{a}'ig
angle_{\Gamma_{\mathsf{m}}} - ig(
u\;\mathbf{curl}\;m{a}\;,\mathbf{curl}\;m{a}'ig)_{\Omega^a} = 0. \end{aligned}$$

Dular, J., et al. (2020) TAS 30 8200113. See also: Brambila R. et al, (2018) TAS 28, 5207511.

Perturbed saddle point problem

$$\begin{split} \left(\partial_t(\mu \boldsymbol{h}) \ , \boldsymbol{h}'\right)_{\Omega^h} + \left(\rho \operatorname{\mathbf{curl}} \boldsymbol{h} \ , \operatorname{\mathbf{curl}} \boldsymbol{h}'\right)_{\Omega^h_c} + \left\langle\partial_t \boldsymbol{a} \times \boldsymbol{n}_{\Omega^h} \ , \boldsymbol{h}'\right\rangle_{\Gamma_m} &= 0, \quad \forall \boldsymbol{h}' \in \mathcal{H}, \\ \left\langle \boldsymbol{h} \times \boldsymbol{n}_{\Omega^a} \ , \boldsymbol{a}'\right\rangle_{\Gamma_m} - \left(\nu \operatorname{\mathbf{curl}} \boldsymbol{a} \ , \operatorname{\mathbf{curl}} \boldsymbol{a}'\right)_{\Omega^a} &= 0, \quad \forall \boldsymbol{a}' \in \mathcal{A}. \end{split}$$

It is a perturbed saddle point problem:

$$\begin{cases} a(u,v) + b(v,p) = \langle f, v \rangle, & \forall v \in V, \\ b(u,q) - c(p,q) = \langle g, q \rangle, & \forall q \in Q, \end{cases} \quad \text{or} \quad \begin{pmatrix} \mathbf{A} & \mathbf{B}^{\mathsf{T}} \\ \mathbf{B} & -\mathbf{C} \end{pmatrix} \begin{pmatrix} \boldsymbol{u} \\ \boldsymbol{p} \end{pmatrix} = \begin{pmatrix} \boldsymbol{f} \\ \boldsymbol{g} \end{pmatrix}.$$

 \Rightarrow Compatibility conditions for numerical stability, otherwise...

First-order functions for *h* and *a*:



D. Boffi, F. Brezzi, et al., Mixed FE methods and applications, Springer, 2013.

Compatibility conditions

$$\begin{pmatrix} \mathbf{A} & \mathbf{B}^{\mathsf{T}} \\ \mathbf{B} & -\mathbf{C} \end{pmatrix} \begin{pmatrix} \boldsymbol{u} \\ \boldsymbol{p} \end{pmatrix} = \begin{pmatrix} \boldsymbol{f} \\ \boldsymbol{g} \end{pmatrix}$$

The solution is stable, i.e., $\|\boldsymbol{u}\|_{V} + \|\boldsymbol{p}\|_{Q} \leq C(\|\boldsymbol{f}\|_{V'} + \|\boldsymbol{g}\|_{Q'})$, if $\exists \alpha, \beta, \gamma > 0$ (strictly) such that

$$\begin{split} \mathbf{v}^{\mathsf{T}} \mathbf{A} \mathbf{v} &\geq \alpha \|\mathbf{v}\|_{V}^{2}, \ \forall \mathbf{v} \in \ker(\mathbf{B}) & (\text{coercivity of } \mathbf{A}), \\ \mathbf{q}^{\mathsf{T}} \mathbf{C} \mathbf{q} &\geq \gamma \|\mathbf{q}\|_{Q}^{2}, \ \forall \mathbf{q} \in \ker(\mathbf{B}^{\mathsf{T}}) & (\text{coercivity of } \mathbf{C}), \end{split}$$

$$\inf_{\boldsymbol{q} \in (\ker(\mathbf{B}^{\mathsf{T}}))^{\perp}} \sup_{\boldsymbol{\nu} \in V} \frac{\boldsymbol{q}^{\mathsf{T}} \mathbf{B} \boldsymbol{\nu}}{\|\boldsymbol{q}\|_{\mathcal{Q}} \|\boldsymbol{\nu}\|_{V}} \ge \beta > 0 \qquad (\text{inf-sup condition}).$$

In our case, the inf-sup condition is the most restrictive.

D. Boffi, F. Brezzi, et al., Mixed FE methods and applications, Springer, 2013.

Inf-sup test

The inf-sup condition is not easy to check analytically.

 \Rightarrow We perform a numerical inf-sup test.

On progressively refined meshes, for given function spaces:

- 1. Define suitable norms.
- 2. Extract matrices \mathbf{B} , \mathbf{N}_V , and \mathbf{N}_Q , from the FE assembly, with

$$\|\boldsymbol{v}\|_{V}^{2} = \boldsymbol{v}^{\mathsf{T}} \mathbf{N}_{V} \boldsymbol{v},$$
$$\|\boldsymbol{q}\|_{Q}^{2} = \boldsymbol{q}^{\mathsf{T}} \mathbf{N}_{Q} \boldsymbol{q}.$$

3. Solve the eigenvalue problem

$$\left(\mathbf{B}\mathbf{N}_{V}^{-1}\mathbf{B}^{\mathsf{T}}\right)\boldsymbol{q} = \lambda\mathbf{N}_{Q}\boldsymbol{q}.$$

Lowest non-zero eigenvalue = square of the inf-sup value β^{δ} . \Rightarrow How does β^{δ} behave when the mesh is refined?

- It tends to zero \Rightarrow unstable,
- It is bounded from below \Rightarrow stable.

D. Chapelle, K.-J. Bathe, The inf-sup test, C&S 47, 1993.

h-a-formulation Unstable choices

Linear or quadratic elements for both h and $a \Rightarrow$ Unstable.



h-a-formulation Stable choices

One way to stabilize the problem:

 \Rightarrow Increase the discretization order of one field (*h* or *a*).



Increasing the order on the coupling interface only is sufficient.



h-a-formulation Stabilization

First-order functions for *h* and *a* (inf-sup KO):



Second-order for *a*, first-order for *h* (inf-sup OK):



Application 1: HTS bulk magnetization model (3D)

HTS bulk magnetization with a coil, on top of a FM pellet.





	# DOFs	# iterations	Time/it.	Total time
h-formulation	12,172	3,937	1.4s	1h33
a-formulation	26,964	3,147	2.1s	1h48
h-a-formulation	15,776	1,108	2.1s	0h39
h-b-formulation	20,821	1,104	3.2s	0h58

Application 2: magnetic shield model (2D and 3D) Magnetic shield made up of a stack of tape annuli.



Inner radius: 13 mm. Outer radius: 22.5 mm. Height: 14.9 mm.



- Number of tapes: N = 183. One tape: HTS layer + FM substrate.
- Filling factor of the FM: f = 0.92.
- Temperature: 77K.

S. Hahn, 2011. A. Patel, 2016.

Shielding configurations



Magnetic shielding application

h-b-formulation

h-φ in Ω and auxiliary *b* field in the FM domain Ω_m.
 Volume coupling in Ω_m:

$$\begin{split} \left(\mu_0 \partial_t \boldsymbol{h} \ , \boldsymbol{h}'\right)_{\Omega_{\mathrm{m}}^{\mathrm{C}}} &+ \left(\rho \operatorname{\mathbf{curl}} \boldsymbol{h} \ , \operatorname{\mathbf{curl}} \boldsymbol{h}'\right)_{\Omega_{\mathrm{c}}} + \left(\partial_t \boldsymbol{b} \ , \boldsymbol{h}'\right)_{\Omega_{\mathrm{m}}} = 0 \\ & \left(\boldsymbol{h} \ , \boldsymbol{b}'\right)_{\Omega_{\mathrm{m}}} - \left(\nu \boldsymbol{b} \ , \boldsymbol{b}'\right)_{\Omega_{\mathrm{m}}} = 0 \end{split}$$

- If Ω_m is non-conducting, inf-sup condition satisfied with piecewise constant elements for b.
- Much more robust than <u>h-formulation</u>.
- More efficient than <u>h-a-formulation</u> because of large coupling surface:



Homogeneous model: anisotropy

Replace the detailed stack by one homogeneous material.



- Introduce the average h and j fields.
- Introduce anisotropic $\tilde{\rho}(j)$ and $\tilde{\mu}(h)$ tensors.
- Modified *h*-formulation :

$$\left(\partial_t(\tilde{\boldsymbol{\mu}}\,\boldsymbol{h})\;,\boldsymbol{h}'
ight)_\Omega+\left(\tilde{\boldsymbol{
ho}}\,\mathrm{curl}\;\boldsymbol{h}\;,\mathrm{curl}\;\boldsymbol{h}'
ight)_{\Omega_{\mathrm{c}}}=0$$

Not optimal: how to apply the <u>*h-b*-formulation</u> with anisotropy and conducting Ω_m domain?

Conclusion

Coupled formulations help to model HTS and FM efficiently

- Surface coupling \Rightarrow *h*-*a*-formulation
- Volume coupling \Rightarrow *h-b*-formulation
- Thin HTS tapes \Rightarrow *t-a-formulation* (not presented here).

These formulations are mixed \Rightarrow Inf-sup condition for stability.



References

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