

Current Status of BELFEM

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BERKELEY LAB







Outline





- Recap: Thin Shell Approach
- a-formulation vs φ-formulation
- Current BCs using Cohomology cuts
- Implementing jump conditions



Recap: Thin Shells using H-φ



Sirois' Thin Shell Approach: How it Works



copper hastelloy HTS copper



Typical HTS stack

- 20 µm
- 80 µm
 - 2 µm
 - 20 µm





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- normal component of h is assumed to be constant along thickness
- only in-plane component of h is resolved along layers
- Nédélec interpolation of ht in plane
- Lagrange interpolation of **ht** within layers along thickness
- coupled to connected air elements using static condensation









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H-Formulation in tapes

- Directly uses the magnetic field H, no gauging needed! → Important for field-dependent material properties.
- Can resolve individual layers of the tape ➡ No need for homogenization methods.
- In 3D: "Out-of-the-box" support for inter-layer current sharing
- Very simple thermal coupling using "thermal resistor mesh" "Out-of-the-box" support for quenching. (Unpublished!)



Sirois' TS approach : Differences to T-Formulation













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Sirois' TS approach : Differences to T-Formulation











Non-Conducting domain A-Formulation vs. ϕ -Formulation



A vs φ Formulation for non-conducting Domains

A-Formulation

• Uses magnetic vector potential

$$b = \nabla \times a$$

- Based on Faraday's law $\nabla \times e = -\dot{b}$
- Dofs associated with edges (in 3D)
- Pros:
 - Well established
 - Straight forward current BCs
- Cons:
 - More DoFs than H-Formulation





H-Formulation

• Uses magnetic scalar potential

$$h = -\nabla \phi$$

• Based on Ampere's law

 $\nabla \times h = -j$



- Dofs associated with nodes
- Pros:
 - Significantly less DoFs
- Cons:
 - Complicated Current BCs





Recap: Cohomologies



current is applied over Ampere's circuital law:

- homologies represent the loops that can be drawn around the conducting regions that fulfill Ampéré's law
- only integral current "I [A]" needs to be known



- cohomologies are cuts in the domain over which jumps in the magnetic potential ϕ are imposed so that $\Delta \phi = I$.
 - very elegant mathematics!
 - homology definition not user friendly
 - → <u>difficult to implement in commercial codes</u>













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H- ϕ formulation: Jump Visualization









Discontinuity Conditions



3D Tapestack





H- ϕ formulation: Homology Examples

CORC geometry









Lagrange Multipliers

- Well known from contact conditions in structural dynamics
- In COMSOL: "Pointwise Constraint"
- Pros:
 - Very easy to implement
- Cons:
 - Element matrix non-positive definite
 - ➡ More work for solver



Jump Conditions

element topology:



virtual work equation:

 $\Pi = \lambda \left(\hat{\phi}_{+} - \hat{\phi}_{-} - I \right) = 0$

element matrix:

$$\begin{bmatrix} \delta \hat{\phi}_{-} & \delta \hat{\phi}_{+} & \delta \lambda \end{bmatrix} \begin{bmatrix} 0 & 0 & 1 \\ 0 & 0 & -1 \\ 1 & -1 & 0 \end{bmatrix} \begin{bmatrix} \hat{\phi}_{-} \\ \hat{\phi}_{+} \\ \lambda \end{bmatrix} = \begin{bmatrix} \delta \hat{\phi}_{-} & \delta \hat{\phi}_{+} & \delta \lambda \end{bmatrix}$$











• Very promising results, however, the Lagrange multipliers are a reason to be concerted about very large models since the non-positive definite matrices will be problematic.



2D Benchmark from 2023





| | | constant time step | | adaptive tim step | |
|------------------|-----------|-----------------------|--------------|----------------------|-------------|
| code and library | | coarse mesh | fine mesh | coarse mesh | fine mes |
| GetDP | MUMPS | 5:13 | 10:57 | 2:36 | 7: |
| BELFEM | MUMPS | 1:56 | 7:15 | 0:23 | 1: |
| BELFEM | STRUMPACK | 0:55 | 2:54 | 0:11 | 0: |









Better Approach: Eliminate Implicit Dof:





• Node 4 is "hanging" \rightarrow dof 4 is expressed as linear combination of dof 3 and 5

$$\delta \hat{\phi}_{4} \quad \delta \hat{\phi}_{5} \quad \delta \hat{\phi}_{8} \quad \delta \hat{\phi}_{7}] \mathbf{K} \begin{bmatrix} \hat{\phi}_{4} \\ \hat{\phi}_{5} \\ \hat{\phi}_{8} \\ \hat{\phi}_{7} \end{bmatrix} = \begin{bmatrix} \delta \hat{\phi}_{3} & \delta \hat{\phi}_{5} & \delta \hat{\phi}_{8} & \delta \hat{\phi}_{7} \end{bmatrix} \mathbf{T}^{\mathbf{T}} \mathbf{K} \mathbf{T} \begin{bmatrix} \hat{\phi}_{3} \\ \hat{\phi}_{5} \\ \hat{\phi}_{8} \\ \hat{\phi}_{7} \end{bmatrix}$$

• Change of basis us performed using T-Matrix

 $\mathbf{T} = \begin{bmatrix} \frac{1}{2} & \frac{1}{2} & 0 & 0 \\ 0 & 1 & 0 & 0 \\ 0 & 0 & 1 & 0 \\ 0 & 0 & 1 & 0 \end{bmatrix}$

• Method can be used for both domain interfaces and domain cuts

• But: Not a standard method supported in FE solvers!











Hanging Dofs with Abstract Nodes (Unpublished!)











The concept of "Hanging Nodes" is well known in FE theory. Using this technique in for Jump conditions, however, is new, And according to my best knowledge unique to BELFEM!





Hanging Dofs with Abstract Nodes

- The "local dofs" of the yellow element are $[\phi_4, \phi_5, \phi_6]$
- We compute the element Jacobian AI using these dofs.
 - The DOF transformation matrix reads

$$\begin{bmatrix} \hat{\phi}_{4} \\ \hat{\phi}_{5} \\ \hat{\phi}_{6} \end{bmatrix} = \begin{bmatrix} 0 & 0 & 1 & 0 \\ 0 & 1 & 0 & 1 \\ 1 & 0 & 0 & 1 \end{bmatrix} \cdot \begin{bmatrix} \hat{\phi}_{2} \\ \hat{\phi}_{3} \\ \hat{\phi}_{4} \\ I_{0} \end{bmatrix}$$

• The Jacobian with respect to real dofs then reads

$$\mathbf{A} = \mathbf{T}^{\mathrm{T}} \mathbf{A}_{l} \mathbf{T}$$

• Ultimately, I is imposed as Dirichlet condition











Summary and Timeline



Intermediate Summary:



- most aspects.
- the a-formulation.
- Current conditions:



• Thin Shell h- ϕ is still in development, but will probably outperform t-a in

• In 3D, the ϕ -formulation needs significantly less degrees of freedom than

• Current boundary conditions need to be applied over cohomologies Cohomoligies for 3D problems not intuitive, need clever algorithms Developed in cooperation with Polytechnique Montreal

Were implemented with Lagrange Multipliers in the past

Switched to new "Hanging DoF" concept to achieve better conditioning.









Goal:

- model a thin shell tapes tack in 3D after Alves et Al, 2022
- extend model to encompass solder and thermal model
- be able to do the coupled EM-Thermal quenching model by end of the year

Roadmap:

- overhaul data structure for simplified programming of weak governing equations
- improve degree of freedom management system (almost complete)
- first benchmark with 3D tapestack (Hopefully before Christmas!)
- implement solder and thermal model
- benchmark involving quench
- address contact sharing (Spring 2025)



Current Efforts



[Alves et al, 10.1109/TASC.2022.3143076]





