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THE BLENDED FINITE ELEMENT METHOD FOR MULTI-FLUID PLASMA MODELING

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U.S. AIR FORCE





OUTLINE

1 THE MULTI-FLUID PLASMA MODEL

2 BLENDED FINITE ELEMENT METHOD

- Blended Finite Element Method
- Nodal Continuous Galerkin
- Modal Discontinuous Galerkin
- Model Verification

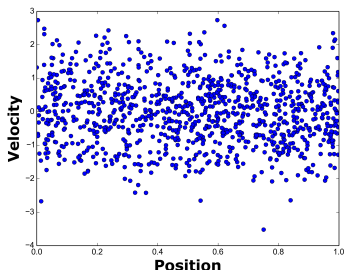
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- 1D Soliton problem
- Electromagnetic Plasma Shock Problem

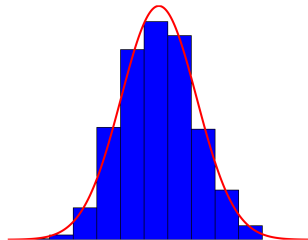
4 APPLICATION



THERE ARE MULTIPLE PLASMA MODELS.



- 3-Dimensions + 3-Velocities
- Evolve the particles position and velocity
- e.g. Particle-In-Cell models



- Ensemble average of particles distribution, $f_s(\mathbf{x}, \mathbf{v}, t)$
- Evolve the distribution function
- e.g. Vlasov-Maxwell models



PHYSICAL DESCRIPTION OF A FLUID.

- Modeling each particle velocity and position is not practical.
- Instead an average is performed to give a statistical description.
- Calculate the number of particles per unit volume having approximately the velocity $\mathbf{v}$ near the position $\mathbf{x}$ and at time t , distribution function $f(\mathbf{v}, \mathbf{x}, t)$

$$\rho_s = m_s \int f_s(\mathbf{v}) d\mathbf{v}$$

$$\rho_s \mathbf{u}_s = m_s \int \mathbf{v} f_s(\mathbf{v}) d\mathbf{v}$$

$$\mathbb{P}_s = \mathbf{P}_s = m_s \int \mathbf{w} \mathbf{w} f_s(\mathbf{v}) d\mathbf{v}, \quad p_s = \frac{1}{3} m_s \int w^2 f_s(\mathbf{v}) d\mathbf{v}$$

$$\mathbf{H}_s = m_s \int \mathbf{w} \mathbf{w} \mathbf{w} f_s(\mathbf{v}) d\mathbf{v}, \quad \mathbf{h}_s = \frac{1}{2} m_s \int w^2 \mathbf{w} f_s(\mathbf{v}) d\mathbf{v}$$

$$\mathbf{w} = \mathbf{v} - \mathbf{u}_s$$



BOLTZMANN EQUATION EVOLVES f_s .

- The Boltzmann eqn:

$$\frac{\partial f_s}{\partial t} + \mathbf{v} \cdot \frac{\partial f_s}{\partial \mathbf{x}} + \frac{q_s}{m_s} (\mathbf{E} + \mathbf{v} \times \mathbf{B}) \cdot \frac{\partial f_s}{\partial \mathbf{v}} = \left. \frac{\partial f_s}{\partial t} \right|_c$$

- Take the 0^{th} , 1^{st} , 2^{nd} moments of the Boltzmann Eqn.

$$m_s \int \mathbf{v}^n \frac{\partial f_s}{\partial t} d\mathbf{v} + m_s \int \mathbf{v}^{n+1} \cdot \frac{\partial f_s}{\partial \mathbf{x}} d\mathbf{v} + q_s \int \mathbf{v}^n (\mathbf{E} + \mathbf{v} \times \mathbf{B}) \cdot \frac{\partial f_s}{\partial \mathbf{v}} d\mathbf{v} = m_s \int \mathbf{v}^n \left. \frac{\partial f_s}{\partial t} \right|_c d\mathbf{v}$$

- Each moment of the Boltzmann eqn gives an equation for the moment variable, and introduces the next higher moment variable
- This process can go on indefinitely



BOLTZMANN EQUATION EVOLVES f_s .

$$\frac{\partial \rho_s}{\partial t} + \nabla \cdot (\rho_s \mathbf{u}_s) = \frac{\partial \rho_s}{\partial t} \Big|_{\Gamma}$$

$$\frac{\partial \rho_s \mathbf{u}_s}{\partial t} + \nabla \cdot (\rho_s \mathbf{u}_s \mathbf{u}_s + p_s \mathbf{I} + \Pi_s) = \frac{\rho_s q_s}{m_s} (\mathbf{E} + \mathbf{u}_s \times \mathbf{B}) - \sum_{s^*} \mathbf{R}_{s,s^*} + \frac{\partial \rho_s \mathbf{u}_s}{\partial t} \Big|_{\Gamma}$$

$$\frac{\partial \varepsilon_s}{\partial t} + \nabla \cdot (((\varepsilon_s + p_s) \mathbf{I} + \Pi_s) \cdot \mathbf{u}_s + \mathbf{h}_s) = \frac{\rho_s q_s}{m_s} \mathbf{u}_s \cdot \mathbf{E} + \sum_{s^*} \mathcal{Q}_{s,s^*} + \frac{\partial \varepsilon_s}{\partial t} \Big|_{\Gamma}$$

- System is truncated by relating higher moment variables to the lower ones
- The fluids are coupled to each other and to the electromagnetic fields through Maxwell's equations and interaction source terms.



ADVANTAGES OF THE MODEL



IDEAL MHD MODEL IS VALID WHEN:

- High collisionality, $\tau_{ii}/\tau \ll 1$
- Small Larmor radius, $r_{Li}/L \ll 1$
- Low Resistivity, $\left(\frac{m_e}{m_i}\right)^{1/2} \left(\frac{r_{Li}}{L}\right)^2 \frac{\tau}{\tau_{ii}} \ll 1$

MULTI-FLUID PLASMA MODEL

- Less computationally expensive than kinetic models
- Multi-fluid effects become relevant at small spacial and temporal scales
- Finite electron mass and speed-of-light effects are included
- There is charge separation is modeled
- Displacement current effects are resolved in the MFPM



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4 APPLICATION



THE MFPM HAS DISPERSIVE SOURCES.

$$\frac{\partial \mathbf{Q}}{\partial t} + \frac{\partial \overleftrightarrow{\mathbf{F}}}{\partial \mathbf{x}} = \mathbf{S}$$

- The source Jacobian $\frac{\partial \mathbf{S}}{\partial \mathbf{Q}}$ has imaginary eigenvalues
- The equation system has dispersive sources
- The dispersion is physical (may be difficult to distinguish from numerical dispersion)
- This dispersion is due to plasma waves that result from ion and electron plasma interactions with electromagnetic fields
- An ideal numerical method for the MFPM should:
 - be high-order accurate
 - couple the flux and the sources
 - capture shocks
 - not impose strict time-step

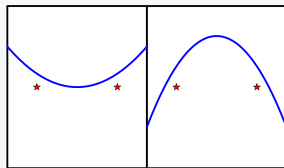
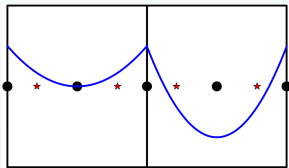


Srinivasan et al, CCP 10 (2011)



BFEM SIMULTANEOUSLY USES CG AND DG.

- Solution to the electron and EM fields is smooth and does not shock



- Continuous Galerkin
- Electron fluid and EM fields

$$\mathbf{Q} = \sum_i \mathbf{q}_i v_i$$

- Discontinuous Galerkin
- Multiple ion and neutral fluids

$$\mathbf{Q} = \sum_i \mathbf{c}_i v_i$$



IMPLICIT CONTINUOUS GALERKIN

- For this implementation the balance law form is cast as

$$\frac{\partial \mathbf{Q}}{\partial t} + \frac{\partial \overleftrightarrow{\mathbf{F}}}{\partial \mathbf{Q}} \cdot \frac{\partial \mathbf{Q}}{\partial \mathbf{x}} = \mathbf{S} + \kappa \nabla^2 \mathbf{Q}_d$$

- Lagrange polynomials are used for basis functions, v_r

$$\int_{\Omega} v_r \frac{\partial \mathbf{Q}}{\partial t} dV = \mathcal{L}_r(\mathbf{Q}) = \int_{\Omega} v_r \mathbf{S} dV - \int_{\Omega} v_r \frac{\partial \overleftrightarrow{\mathbf{F}}}{\partial \mathbf{Q}} \cdot \frac{\partial \mathbf{Q}}{\partial \mathbf{x}} dV + \kappa \int_{\Omega} v_r \nabla^2 \mathbf{Q}_d dV$$

- θ -method time integration

$$\mathcal{R}(\mathbf{Q}^n) = \overleftrightarrow{\mathbf{M}} \frac{\mathbf{Q}^{n+1} - \mathbf{Q}^n}{dt} - \theta \mathcal{L}_r(\mathbf{Q}^{n+1}) - (1 - \theta) \mathcal{L}_r(\mathbf{Q}^n) = 0$$

- $\theta = 0.5$ is used for 2nd order accuracy

$$\overleftrightarrow{\mathbf{J}}(\mathbf{Q}^n) = \frac{\partial \mathcal{R}(\mathbf{Q}^n)}{\partial \mathbf{Q}^n}, \quad \overleftrightarrow{\mathbf{J}}(\mathbf{Q}^n) \Delta \mathbf{Q} = -\mathcal{R}(\mathbf{Q}^n), \quad \mathbf{Q}^{n+1} = \mathbf{Q}^n + \Delta \mathbf{Q}$$



Reddy, An Introduction to the Finite Element Method (2006)



$$\frac{\partial \mathbf{Q}}{\partial t} + \frac{\partial \overleftrightarrow{\mathbf{F}}}{\partial \mathbf{x}} = \mathbf{S}$$

- Legendre polynomials are used for basis functions, v_p
- The hyperbolic equation is multiplied by the basis function,

$$\int_{\Omega} v_p \frac{\partial \mathbf{Q}}{\partial t} dV = \mathcal{L}_p(\mathbf{Q}) = \int_{\Omega} v_p \mathbf{S} dV - \oint_{\partial \Omega} v_p \overleftrightarrow{\mathbf{F}} \cdot d\mathbf{A} + \int_{\Omega} \overleftrightarrow{\mathbf{F}} \cdot \nabla v_p dV$$

- Explicit Runge-Kutta time integration
- $CFL = c\Delta t/\Delta x \leq 1/(2p - 1)$, p is the polynomial order

$$\mathbf{Q}^* = \mathbf{Q}^n + \Delta t \mathcal{L}_p(\mathbf{Q}^n),$$

$$\mathbf{Q}^{n+1} = \frac{1}{2}\mathbf{Q}^* + \frac{1}{2}\mathbf{Q}^n + \frac{1}{2}\Delta t \mathcal{L}_p(\mathbf{Q}^*).$$

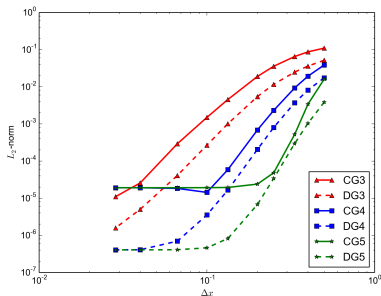




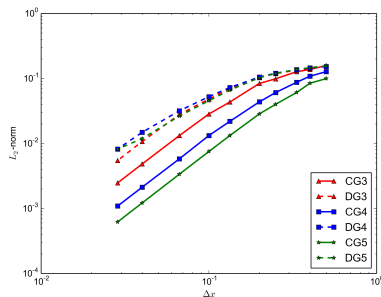
CONVERGENCE OF THE BFEM.

$$\frac{\partial Q}{\partial t} + \frac{\partial Q}{\partial x} = 0, \quad Q(x, 0) = e^{-10(x-8)^2}, \quad \|\Delta Q\|_2 = \sqrt{\frac{1}{n} \sum_{i=1}^n (\hat{Q} - Q_i)^2}$$

Spatial Convergence



Temporal Convergence



• Simulations at fixed time-step

• Simulations at fixed CFL=1



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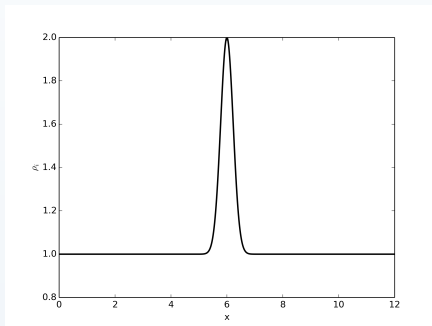
4 APPLICATION



- 1D soliton is a two-fluid plasma problem
- The solution is smooth, therefore artificial dissipation can be small
- The simulation uses 512 second-order elements
- $B_z = 1.0$, $T_e = T_i = 0.01$,
 $\mathbf{u}_i = \mathbf{u}_e = \mathbf{0}$
- $n_e = n_i = 1.0 + e^{-10(x-6)^2}$



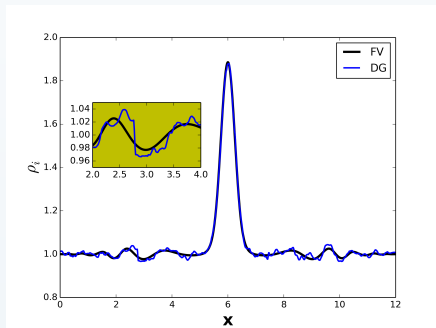
Baboolal, Math. and Comp. Sim. 55 (2001)



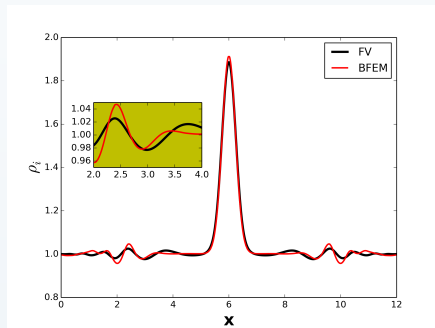


DG AND BFEM COMPARISON WITH A SOLUTION

$$\frac{m_i}{m_e} = 1836, \frac{c}{c_{si}} = 1000\sqrt{2}, \text{FV 5000 cells}$$



- DG Solution is very dispersive



- BFEM is less dissipative than the converged solution



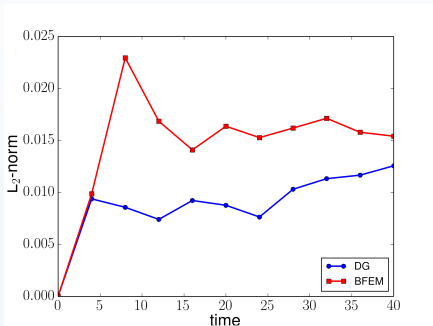
Hakim et al, JCP 219 (2006)



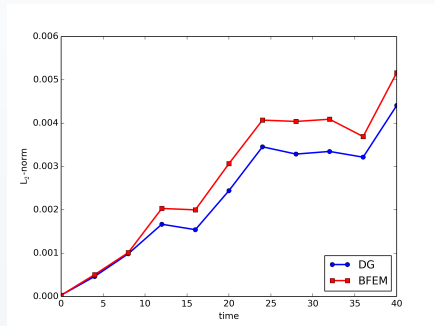
BFEM COMPUTATIONAL COST SAVINGS

case	m_i/m_e	c/c_{si}	DG time(s)	BFEM time (s)	BFEM cost over DG
1	25	$10/\sqrt{2}$	0.32	37.7	+11681%
2	100	$10/\sqrt{2}$	1.28	37.7	+2845%
3	500	$10/\sqrt{2}$	6.82	37.7	+452.8%
4	1000	$10/\sqrt{2}$	12.4	38.2	+208.1%
5	1836	$10/\sqrt{2}$	23.5	40.4	+71.91%
6	3672	$10/\sqrt{2}$	47.2	39.2	-16.95%
7	3672	$100/\sqrt{2}$	520	265	-49.04%
8	3672	$1000/\sqrt{2}$	5274	2735	-48.14%

- Per time-step explicit DG is faster than BFEM, but it requires many more time-steps
- BFEM is more efficient only when time-step are considerably larger than explicit DG



$$m_i/m_e = 1836$$



$$m_i/m_e = 1$$

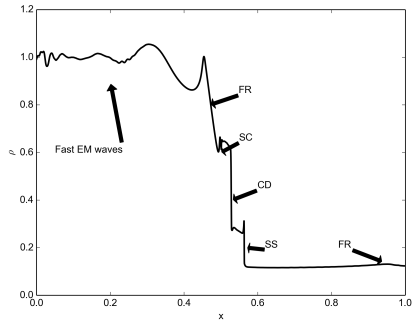
- The BFEM seems to be less accurate than the DG implementation ($\sim 50\%$)
- When the mass ratio is one, the two methods have the same level of accuracy
- The discrepancy is due to the fact that the semi-implicit BFEM does not resolve the plasma frequency in this problem



- Fast rarefaction wave (FR), a slow compound wave (SC), a contact discontinuity (CD), a slow shock (SS), and another fast rarefaction wave (FR)
- The problem exhibits limits of MHD and multi-fluid behavior by changing the Larmor radius, r_L
 - MHD: $r_L \rightarrow 0$
 - Multi-fluid: $r_L \sim L$



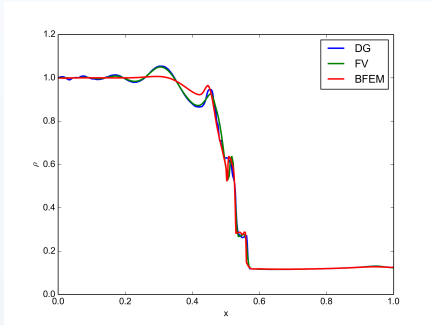
Brio and Wu, JCP 75 (1988)



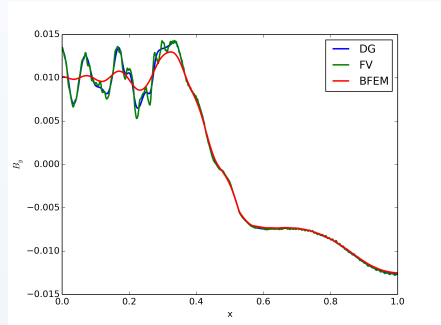


SHOCK IN DENSITY BUT SMOOTH FIELDS.

- $t=0.05/\omega_{ci}$, $c/c_{si}=110$, $m_i/m_e=1836$



mass density



Magnetic field (y-comp.)

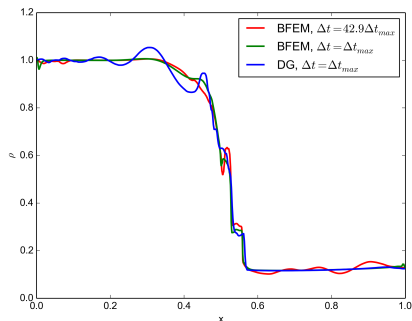
- The main features of the problem are captured by all three methods
- BFEM does not properly resolve the fast electromagnetic waves which require accurately resolving the electron dynamics



MAXIMUM BFEM TIME-STEP

$$\Delta t_{max} = \min \left(\frac{\Delta x}{c_{se}}, \frac{\Delta x}{c_{si}}, \frac{\Delta x}{c}, \frac{0.1}{\omega_{ce}}, \frac{0.1}{\omega_{ci}}, \frac{0.1}{\omega_{pe}}, \frac{0.1}{\omega_{pi}} \right)$$

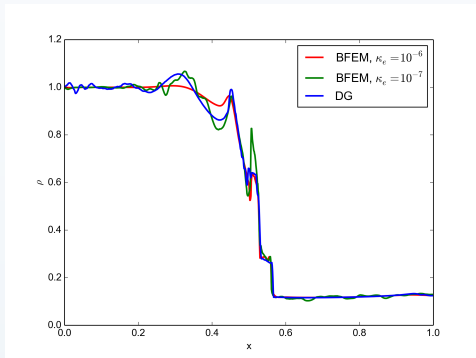
- Δt_{max} corresponds to the maximum value allowed for explicit methods based on the CFL condition
- $\Delta t = 42.9 \Delta t_{max}$ is the maximum time step allowed by the BFEM due to ion dynamics





EFFECTS OF ARTIFICIAL DISSIPATION

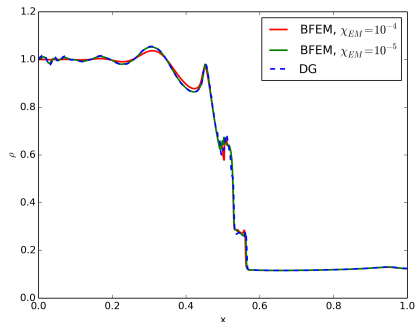
- Varying the artificial dissipation on the electron fluid, κ_e
- Wave-like behavior of the problem is better resolved
- Amplitude of the compound wave increases
- Right fast rarefaction wave is not visible





EFFECTS OF ARTIFICIAL DISSIPATION

- Varying the artificial dissipation on the EM-field, κ_{EM}
- There is better agreement with the DG solution
- This reinforces the point that the wave-like behavior arises from the interaction of the electron fluid with the electromagnetic fields





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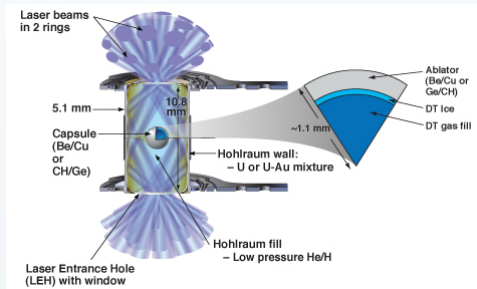
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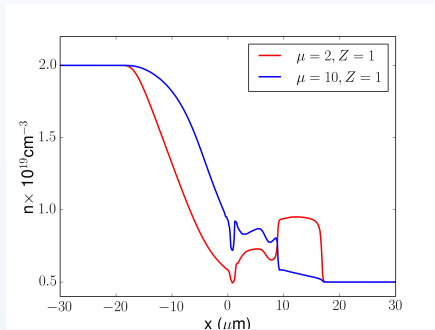
ICF SPECIES SEPARATION

- Deuterium and tritium are heated and compressed to fusion conditions
- The compression is laser-driven
- Deuterium can accelerate faster than the tritium
- Low neutron yield measurements point towards separation
- The phenomenon is not captured by single-fluid plasma models

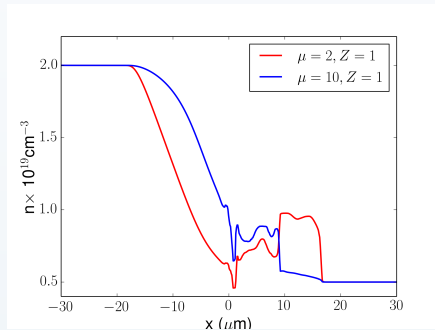




$t=150\text{ps}$



Electron CFL=1



Electron CFL=20

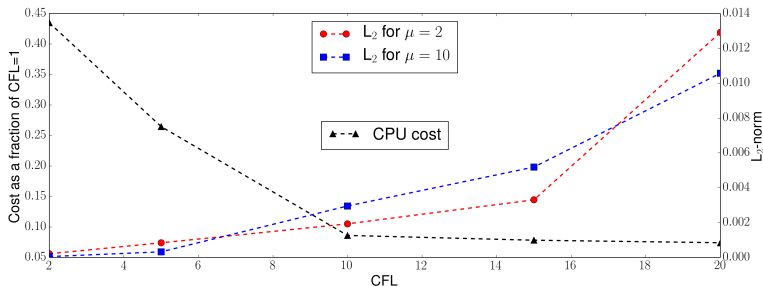
- The ion species separation in both cases is the same although the solution behind the shock fronts differ



Bellei et al., PoP 20 (2013) SOUSA (AFRL)



COST VS. ACCURACY



- The solution error increases as the computational time decreases.



- The blended finite element method (BFEM) is presented
 - DG spatial discretization with explicit Runge-Kutta time integration Ions and neutrals
 - CG spatial discretization with implicit Crank-Nicolson time integration for the electrons and EM fields
 - DG captures shocks and discontinuities
 - CG is efficient and robust for smooth solutions
- Physics-based decomposition of the algorithm yields numerical solutions that resolve the desired timescales
- DG method takes less computational time to advance the solution by one time-step, however Δt is much smaller than that of the BFEM
- Computational cost savings using the BFEM will only occur for relatively large implicit time-steps compared to explicit time-steps

Thank You.