

M3D-C1 ZOOM Meeting

02/07/2022

Upcoming Meetings

CS Issues

1. LBL Report
2. Perlmutter status
3. Traverse status --
4. Mesh adaptation update
5. NERSC Time
6. Timing tests
7. Changes to github master since last meeting
8. Regression tests

Physics Studies

1. Hybrid simulation with kinetic thermal ions – Chang Liu
2. Update on Soft beta limit study

Note: [meeting minutes posted on m3dc1.pppl.gov](https://m3dc1.pppl.gov)

In attendance

Steve Jardin
Adelle Wright
Dingyun Liu
Hank Strauss
Nate Ferraro
P Sinha
Chang Liu
Anders Kleiner
Brendan Lyons
Jin Chen
Chen Zhao

Mark Shephard
Seegyoung Seol
Morteza Siboni
Usman Riaz

Sam Williams
Sherry Li
Yang Liu

Upcoming Meetings

- Sherwood April 4-6 Santa Rosa, CA (in person)
 - CTTS Sunday April 3
- ITPA: MHD, Disruptions, Control April 4-8
- EPS Conference on Plasma Physics June 27—July 1 (online)
 - Abstract submission by February 25
- IAEA Technical Meeting on Plasma Disruptions and their Mitigation 19-22 July
 - In person at ITER HQ in France
 - Abstract submission by May 31

LBL Report

Perlmutter status

- Perlmutter is available for testing, but is not in a production mode
- Maintenance of the cooling system for the next 6 weeks means that only 500 nodes will be available to users
- I have done some timing tests and find it to be unreliable...code crashes after some timesteps for no apparent reason
- No file `/usr/common/usg/bin/nersc_host`
- Problem with installing PETSc with NVIDIA...Seegyong

TRAVERSE GPU Status (Jan 27 from Chang Liu)

I have pushed the code to do gpu matrix assembling to the master branch. They are named `ludef_t_gpu.f90` and `metricterms_new_gpu.f90`. They should not affect the code compiling on non-GPU machines. You can test it on traverse using

`ARCH=traverse_gpu`, or do the following

```
export M3DC1_CODE_DIR=<your path to M3DC1 directory>
```

```
cd $M3DC1_CODE_DIR
```

```
module use $M3DC1_CODE_DIR/unstructured/modules/traverse
```

```
module load m3dc1/devel-gpu
```

You can then compile the code on traverse and run regression tests. Currently, only part of the code is ported, and it seems that only the RMP test can pass. I will continue working on it and make more tests pass.

Note: to compile a GPU version you need to enable `ACC=1` when compiling.

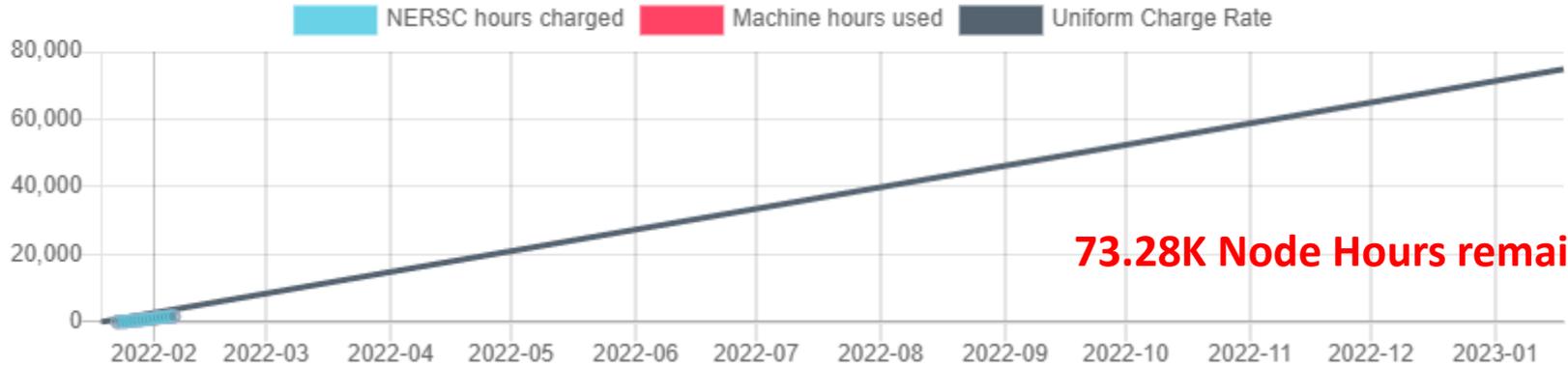
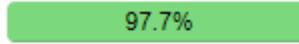
Mesh adaptation update

- Pellet case Brendan
- Interest in Soft-Beta-Limit case ?

NERSC Time

mp288

Remaining %:



- New award period began Jan 19
- We are NESAP Tier 2 for Pearlmuter. . Phase-I w GPUs We have been given a repo m3984 with a small allocation. Presently we are not being charged.
- N9ES-N2 M3D-C1: J. Chen , C. Liu, S. Seol are early users

FY22 allocation (started Jan 19)

Project name: mp288

CPU Node Hours Award: 75,000

GPU Node Hours Award: 7,000

Archive Storage Award (TB): 157

Project CFS Award (TB): 20

75,000 = .34 x 365 x 24 x 25

=> we can use an average of 25 nodes on Haswell continuously

→ For 2022, the Machine Charge Factors are:

→ Perlmutter CPU Nodes: 1.0

→ Cori KNL Nodes: 0.20

→ Cori Haswell Nodes: 0.34

and the charge units are "CPU Node Hours"

Timing Tests

	P	T	N	lundef -s	solve-s	rate/task
Cori Haswell	32	16	48	112	180	$4.45 \times 10^{-6} \text{ s}^{-1}$
	64	32	48	72	95	$3.89 \times 10^{-6} \text{ s}^{-1}$
Perlmutter (CPU)	32	32	24	85	160	$5.37 \times 10^{-6} \text{ s}^{-1}$
Stellar	32	96	8	125	200	$4.04 \times 10^{-6} \text{ s}^{-1}$

P -- # of partitions per plane

T -- # of tasks per node

N --# of nodes

Note: Perlmutter CPU-only
nodes will have twice the
memory of present nodes

$$\text{Rate} = (\text{lundef-s} + \text{solve-s})^{-1}$$

28792 = 170^2 triangles, 24 planes

Changes to github master since 01/10/22

Nate Ferraro:

02/03/22: Fixed bug in plot_surfmn introduced when plot_surfmn was extended to 3D data

02/03/22: Improvements to plot_br for 3D data

02/03/22: Fixed error in field_data.pro parsing kprad_particle_source ionization state

02/04/22: Added ntor parameter to plot_bmn for compatibility with plotting 3D data

02/04/22: Fixed bug plotting linear bmn data

02/04/22: Corrected sign of toroidal field perturbation with irmp=1 in 3D version

02/04/22: Updated field_spectrum.pro to use only positive ntor values

Updated schaffer_plot.pro to correctly normalize 3D data

Jin Chen:

01/24/22 : ST=1 option on Perlmutter

Steve Jardin:

02/04/22: Added kappari_fac; ion parallel thermal conductivity is kappari_fac x electron value

Chang Liu:

01/27/22: Enable matrix assembling using GPU on traverse

01/31/22: Fix a type in lundef_t.f90

Seegyong Seol

02/06/22: minor change in mesh collapse not to mandate a model file

New Variable

kappari_fac

Ion parallel thermal conductivity is $\text{kappari_fac} \times \text{electron value}$
(only for numvar=3 and ipres=1)

Local Systems

- PPPL centos7(02/07/22)
 - 7 jobs **PASSED**
- PPPL greene (02/07/22)
 - 5 jobs **PASSED**
- STELLAR (02/07/22)
 - 7 regression tests **PASSED** on stellar
- TRAVERSE(02/07/22)
 - 7 regression tests **PASSED** on (01/24/22)
 - 7 regression tests **FAILED** on (02/07/22)

Other Systems

- Cori-KNL (02/07/2022)
 - 7 regression tests **PASSED**
- Cori-Haswell (02/07/2022)
 - 7 regression tests **PASSED**
- Perlmutter (01/29/2022)
 - 6 regression tests **PASSED**
 - NCSX timed out
 - “make all” does not work
 - Had to modify the PATH command, replacing ‘pwd’ with actual directory

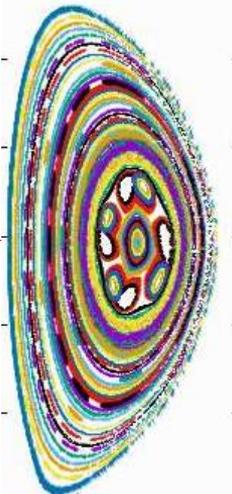
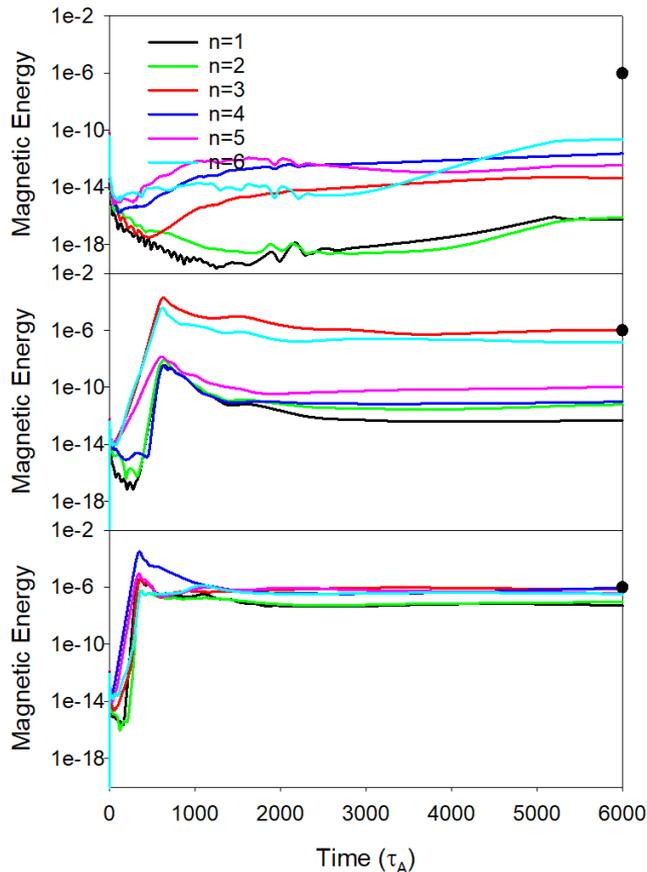
Hybrid simulation with kinetic thermal ions

Chang Liu to present

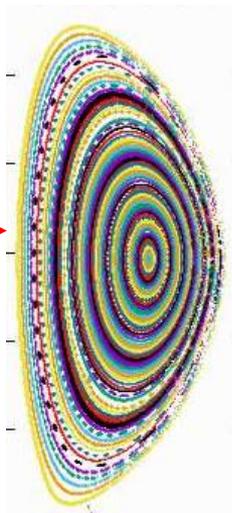
Update on Soft Beta Limit Study

Comparison of the 3 scaled equilibria

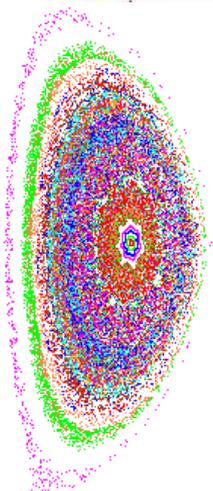
BS=1.1
 $\beta=5.8\%$



Good
-surfaces
except $n=3$
-in center



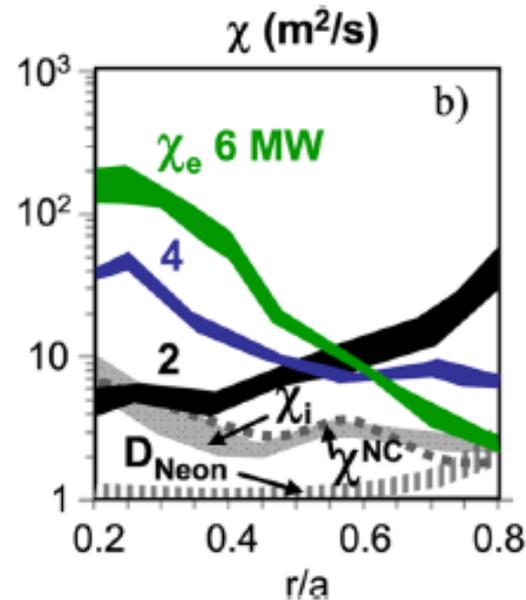
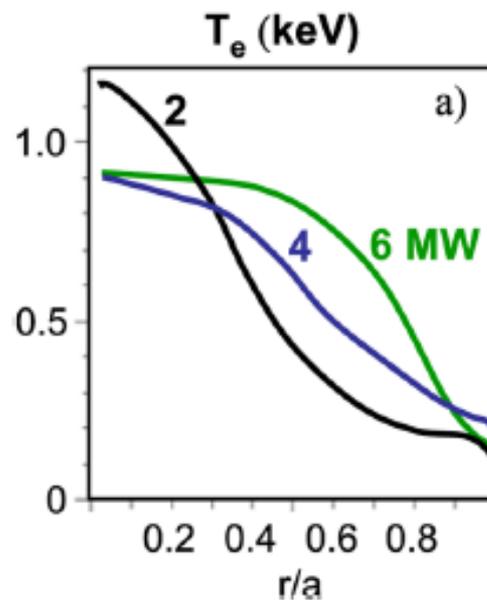
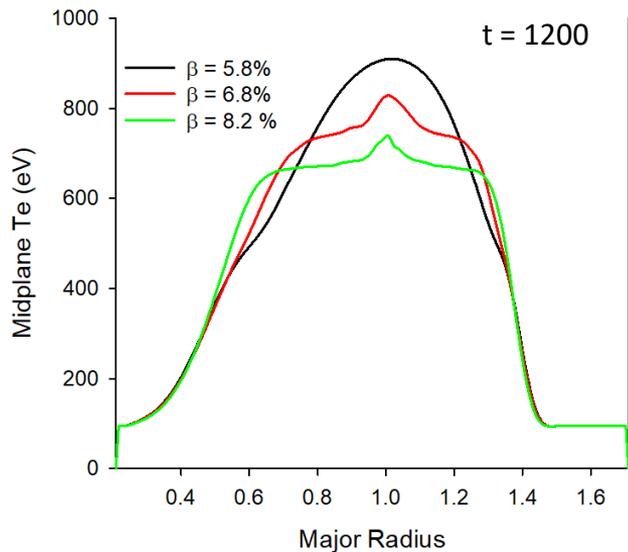
Good
surfaces
everywhere



Poor
surfaces
with
multiple n -
modes

Trend is similar to experiments on NSTX

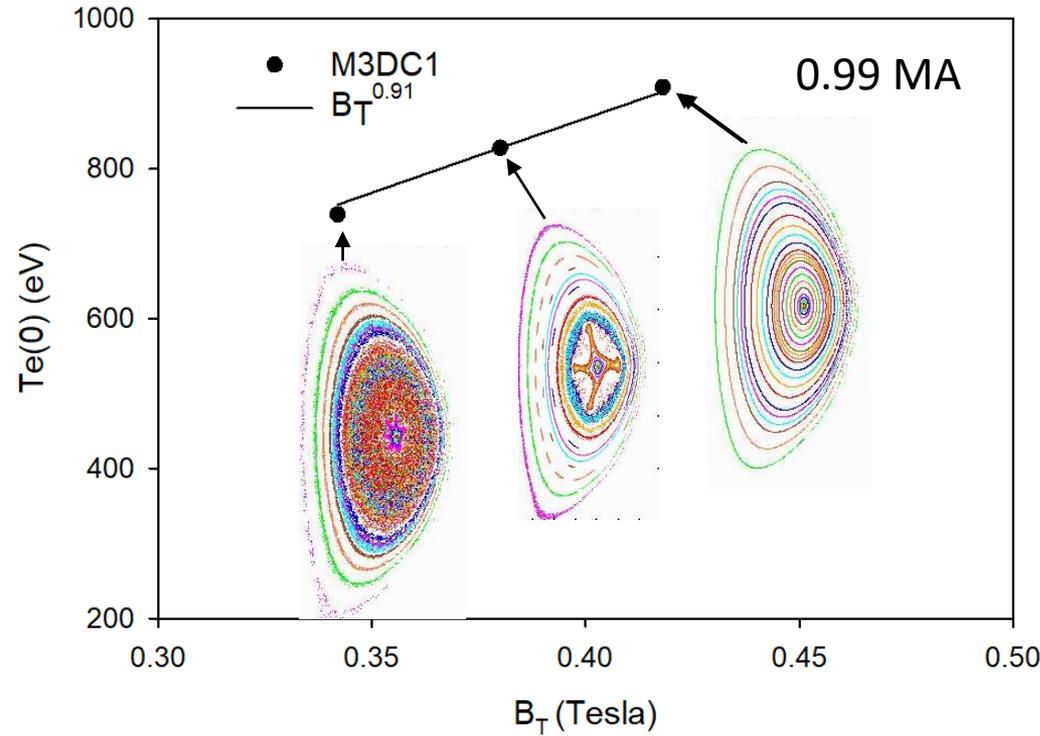
3 Bateman scaled NL runs



- M3D-C1: Central temperature decreases with β
- Exp data: Central transport increases with β

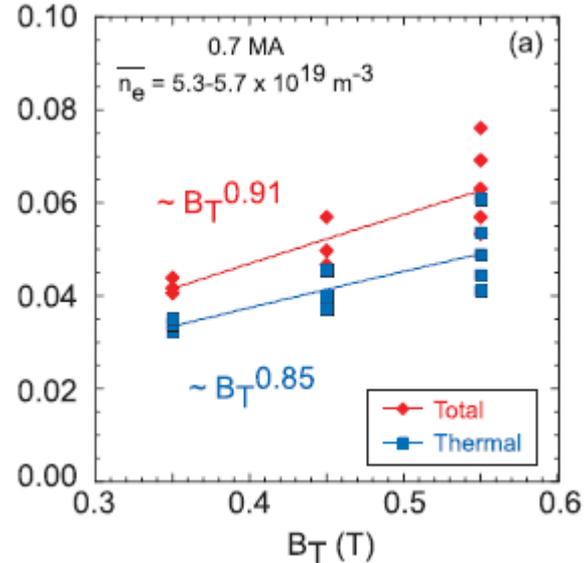
Stutman, et al. PRL (2009)

M3DC1 shows similar scaling with B_T as experiment



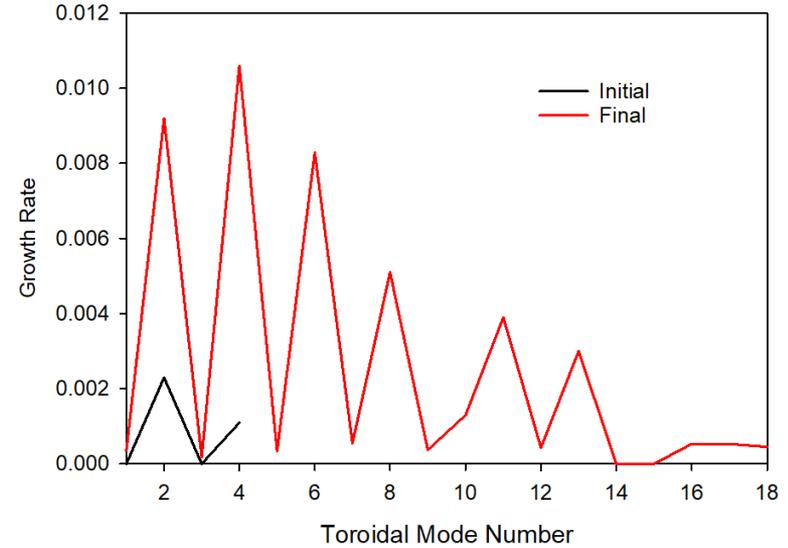
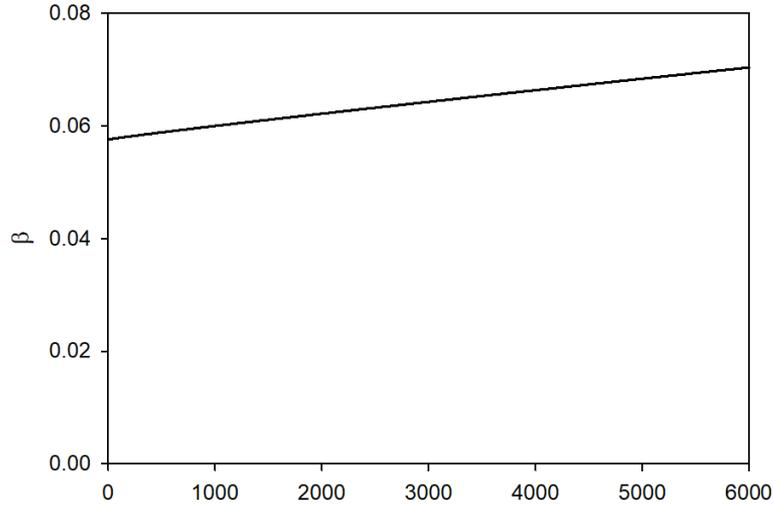
Note: Plot on left is $T_e(0)$. On the right is τ_E

Kaye, et al, PRL (2007)

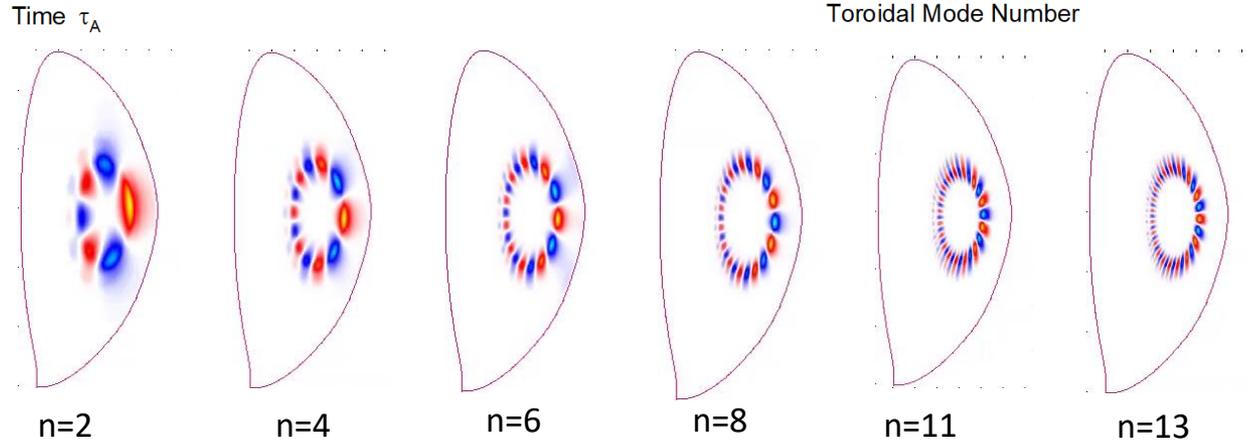


me of the discharges in this study exhibit both low amplitude low-n MHD activity as well as the fast ion driven Alfvén eigenmode (AE) activity,”

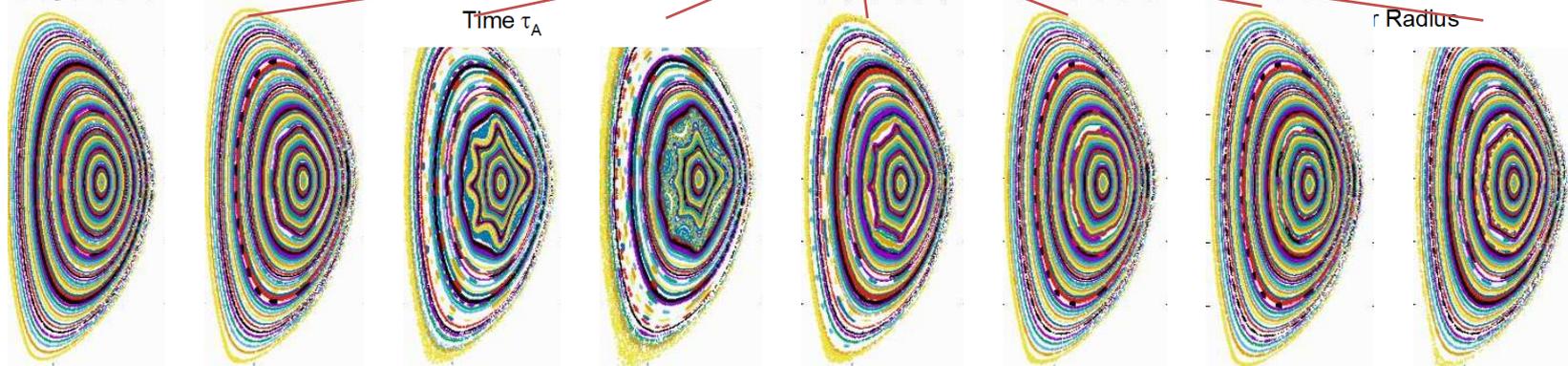
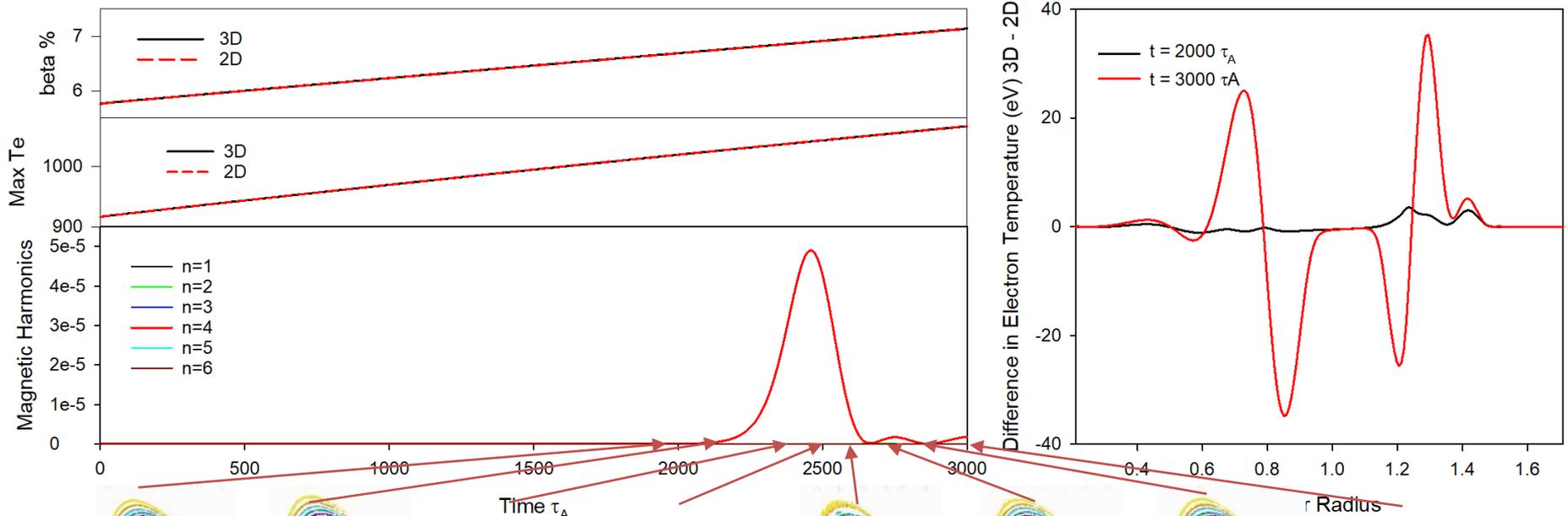
More realistic: Start with stable equilibrium and apply heating power: First in 2D



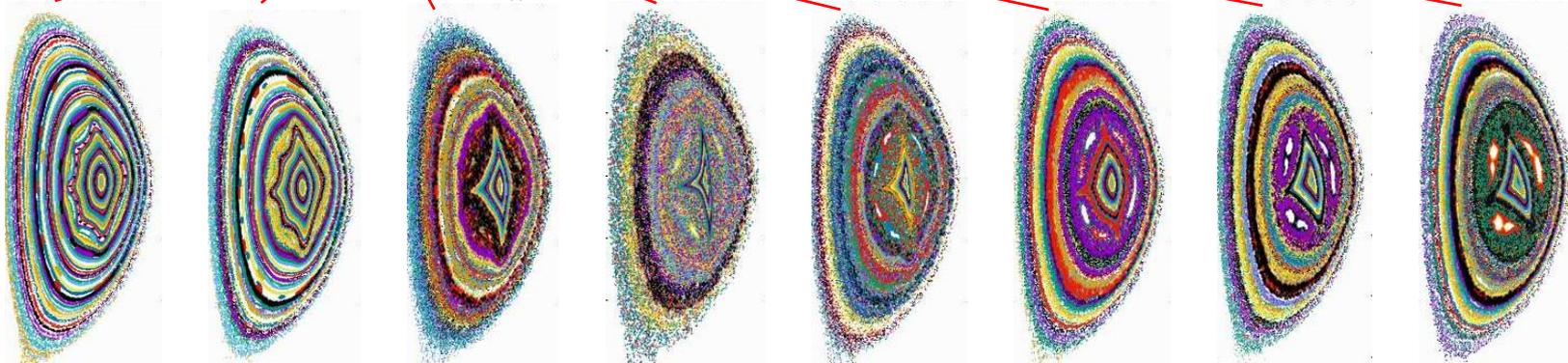
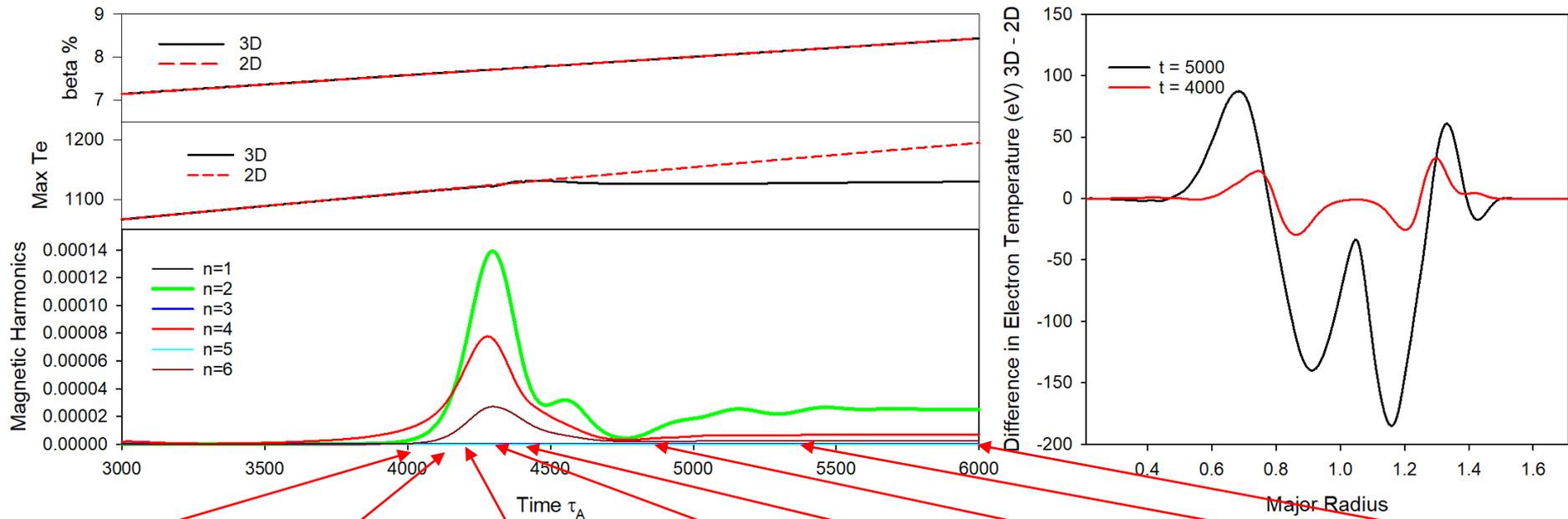
- Start with stable Bateman scaled equilibrium with $\beta = 5.8\%$
- Run in 2D with heating source, increasing β to 7%
- Linear analysis shows final equilibrium unstable to many modes (shown on right)
- Now repeat with 3D run. Do these saturate nonlinearly?



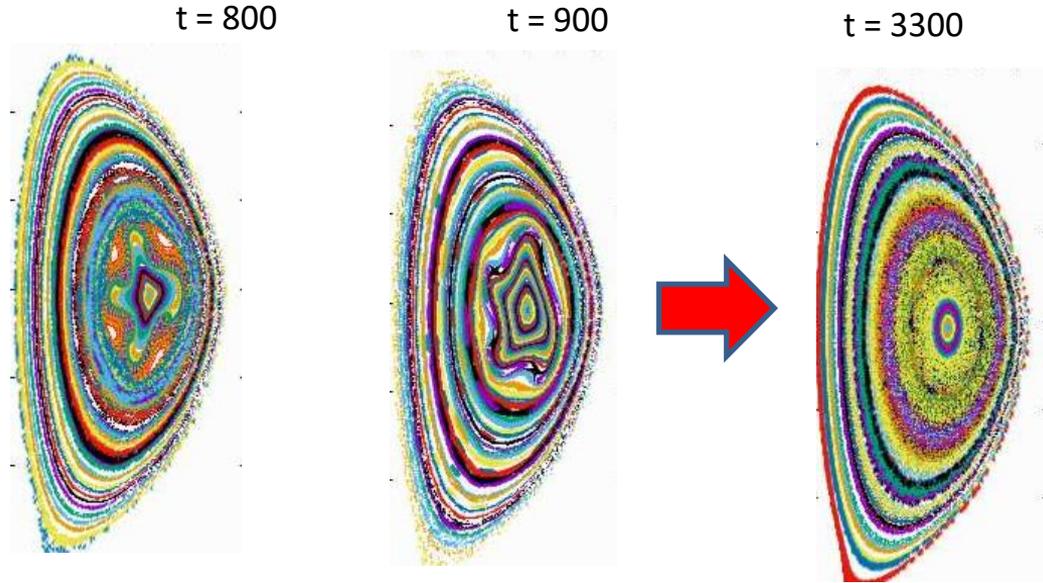
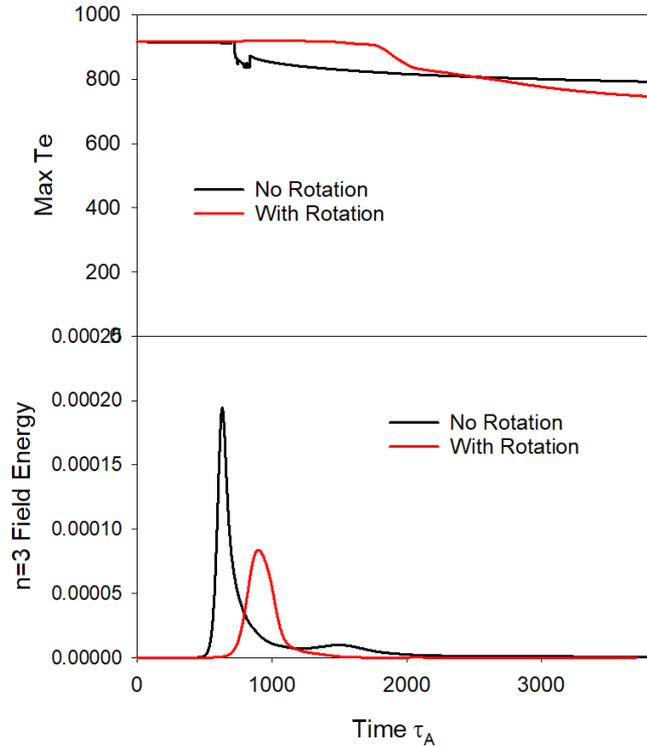
Summary of $0 < t < 3000 \tau_A$



3000 < t < 6000

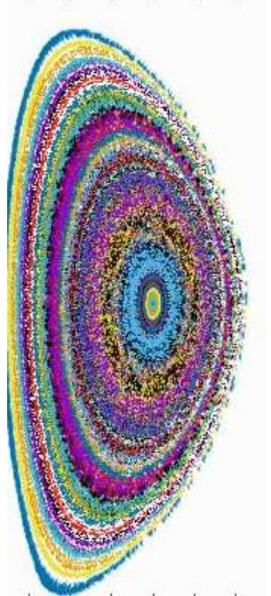
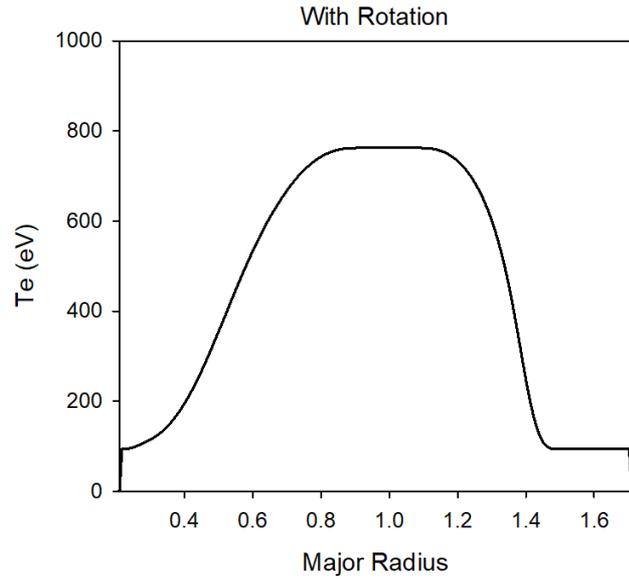
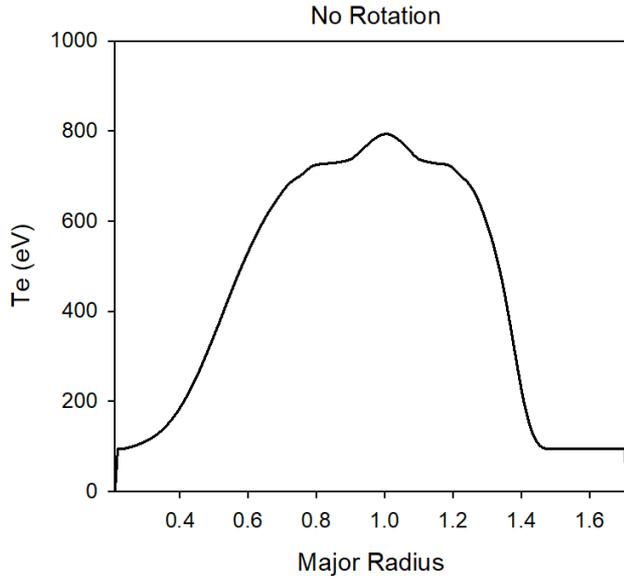
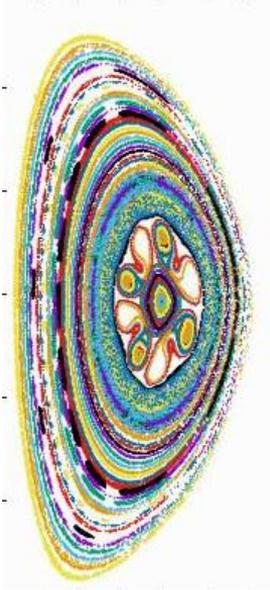


Effect of Sheared Rotation ~ 20 kHz in Center



Results are similar, but instability growth rates are less and tend to symmetrize final configuration

Comparison of midplane temperatures at $t = 3200 \tau_A$ (1.90 ms)



Effect of 3D on τ_E

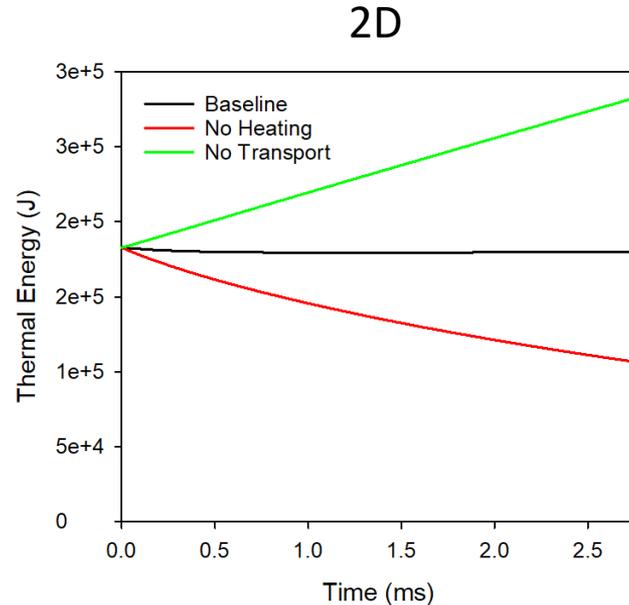
NSTX experimentalists want to know the effect of the 3D instabilities on τ_E

To attempt to answer this, we are running a case with Beams and Transport, first in 2D and then in 3D to see the difference

$S = 36.3 \text{ MW}$, $\kappa_{\perp} = 21.8 \text{ m}^2/\text{s}$, $\tau_E \sim 7.76 \text{ ms}$

Separate ion and electron temperatures

3D in progress



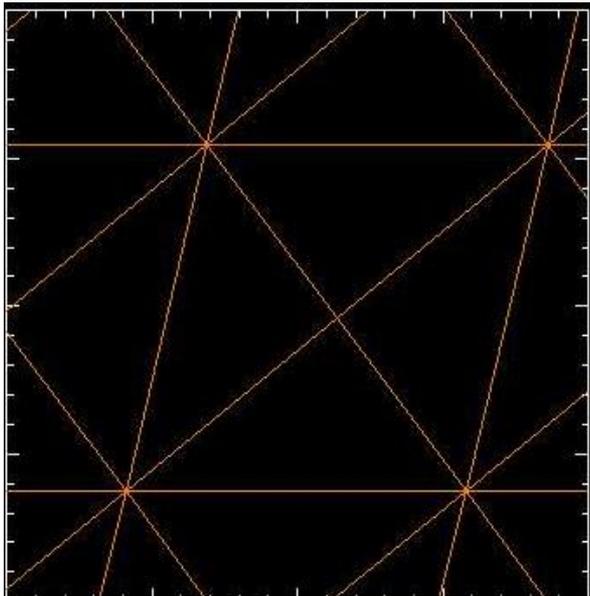
That's All I have

Anything Else ?

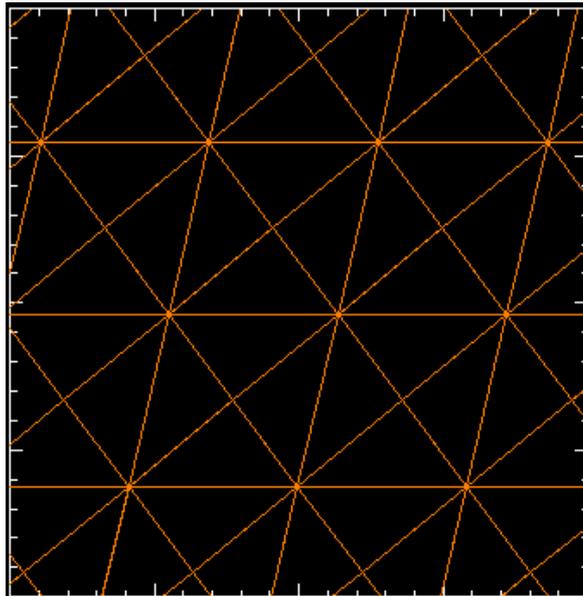
Next Meeting in 3 weeks: Feb 28

Update on Soft-Beta-Limit Study

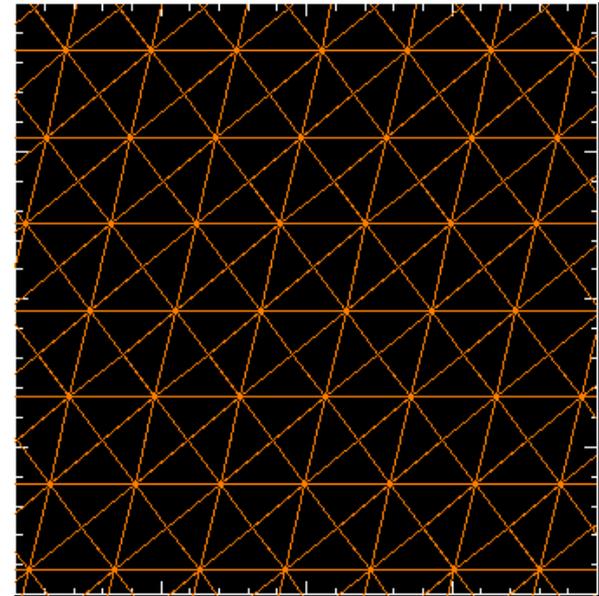
0.8 cm



0.4 cm

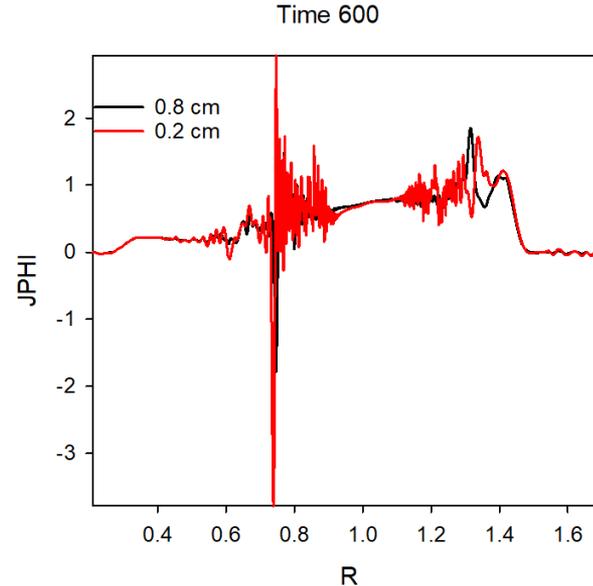
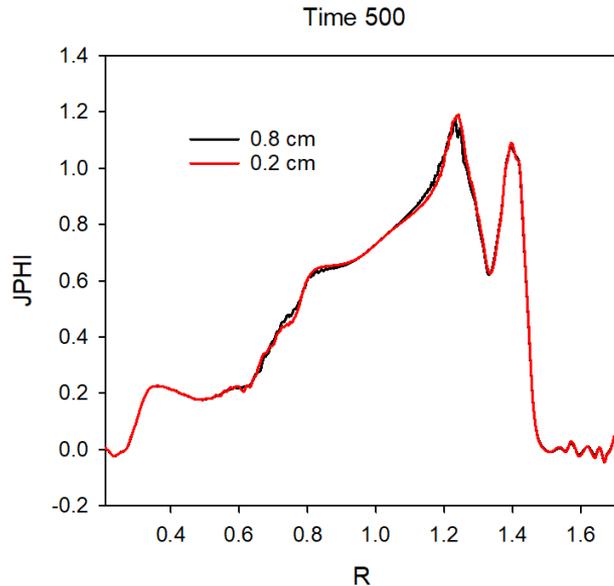


0.2 cm



These are close-ups in center of grid (near magnetic axis)

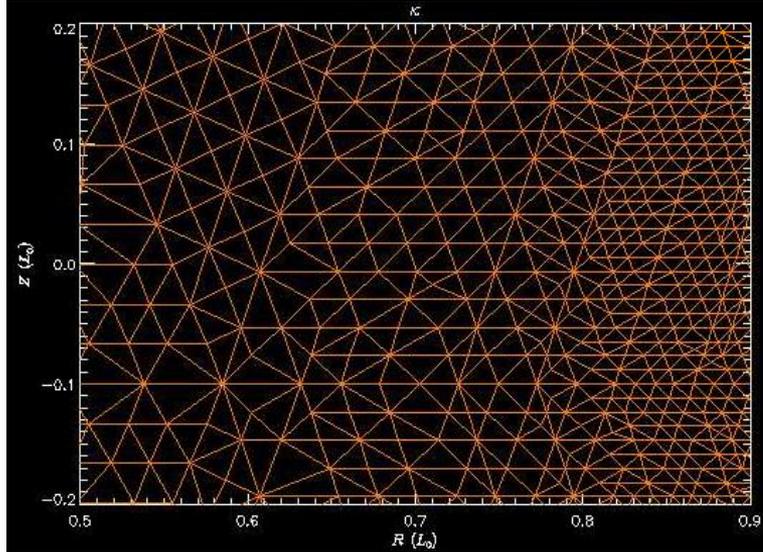
Result of Convergence Study



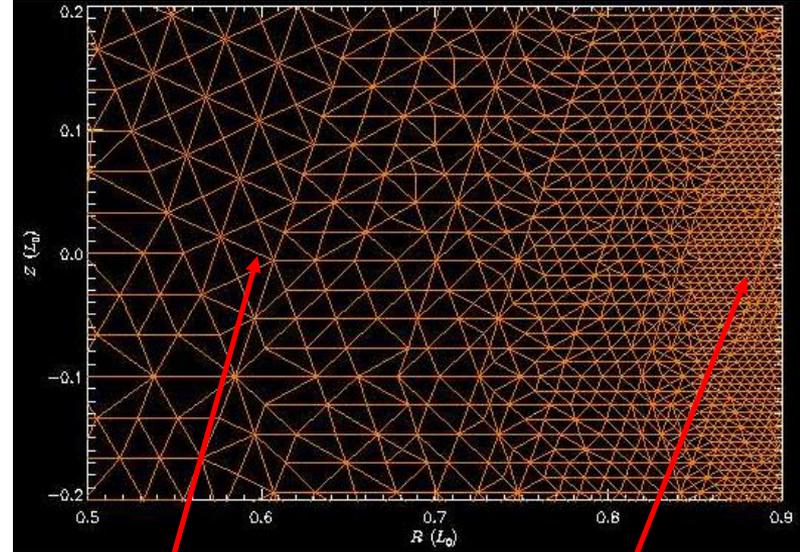
Solution for jphi still very noisy in region $0.5 < R < 0.9$, even for the finest grid with 0.2 cm in center

Grid was not refined where J gets jagged

0.8 cm



0.2 cm



Resolution increased here

But not here

Now producing better grids that are refined where the current gets jagged