M3D-C1 ZOOM Meeting

6/22/2020

Agenda

- 1. Announcements
- 2. CS Issues
 - 1. Local installations
 - 2. Test of pskip and nskip
 - 3. NERSC Time
 - 4. Changes to GIT master since last meeting
- 3. Physics Studies
 - 1. New Toroidal Current plot at different toroidal planes
 - 2. Lyons DIII-D case led to a bug fix
 - 3. ITER VDE with new structure
 - 4. Chang Liu –New Structure Preserving Algorithm based on Slow Manifold
 - 5. Runaways with sources
 - 6. M3D-C1 coupling to KORC
 - 7. Other

Announcements

- Laboratory closed except for Advance Team Site Preparations
- No more partition=greene at PPPL
 - Instead use #SLURM –partition=m3dc1 (job will run on greene, 8 GB/core)
- APS-DPP Meeting is All Virtual: November 9-13, 2020
 - Deadline for Contributed papers is June 29 (Week from today)
- IAEA Technical Meeting on Disruptions and their mitigation
 - Will be held remotely 20-23 July
 - Talks are to be pre-recorded by July 13: (4:3) or (16:9) invited 25 min
- SciDAC PI meeting scheduled for July 28-30 July 2020 is *cancelled*
 - Replaced by a half-day remote panel session on July 29
- ITPA MHD Meeting at IO October 14-16 2020
 - Open to Remote Participation
- IAEA Fusion Energy Conference postponed to May 2021

Local Installations

- CENT_OS-7 on greene
 - Prentice requested we move to OS-7 on greene
 - All regressions tests passed
 - I tested a longer complex case and it passed
 - Any objections?
- Dawson (or general)
 - Still a hdf5 problem with regression tests
 - They also fail on greene
 - Should we send in a ticket to the pppl help?

Test of pskip and nskip

I did some tests with nskip. Here is an example you gave me before. I put the simulation paths in eddy:

nskip=5:/scratch/gpfs/liuchang/STing2 nskip=1:/scratch/gpfs/liuchang/STing2_test We can see that the two simulations agrees pretty well in the linear phase. The nskip=5 case runs much faster.

Note that here I set eqsubtract=1, but for eqsubtract=0 it also works. However, if I set up a current or density control scheme, then the simulation with nskip>1 will get a diverge in the solving process very quickly.

I think we need to do some more work on the control procedure. Anyway, I think we should not remove it from the work.

Chang Liu email June 18

NERSC



M3163



mp288. now consistent with linear usage rate M3163 to expire in 8 days: 2M HRS remaining (clauser, kleiner, lyons, strauss)

Changes to github master since last week

- Chen: ported to cent OS7
 - Added regressions tests and readme file
- Clauser:
 - Some IDL field improvements
 - Bug fix for E_par in 3D
- Seegyoung
 - Updated makefile and readme for SDUMONT
- Jardin
 - KPRAD change to be discussed later

Documented changes in NEWDOC-latest: m3dc1.pppl.gov

New Toroidal Current plot



Cesar Clauser: Toroidal current at different toroidal angles for 3D VDE benchmark problem

Lyons: DIII-D Case C1_28868

(crashes in KPRAD)



- A. dt=1, hyperi=hyperv=1.e-9
- B. Same as A, but start from earlier time
- C. Same as B, but with code fix A. (next slide)



В

3000

Code fix for KPRAD bug

```
! Line Radiation (0 if not advancing KPRAD at that point)
temp79b = merge(dw_rad(:,kprad_z), 0., advance_kprad) / dti
where(temp79b.ne.temp79b) temp79b = 0.
!Check for and delete spurius values (scj 6/21/20)
do i=1,MAX_PTS
    if(abs(temp79b(i)) .gt. 1.e0) then
    temp79b(i) = 0.
    endif
    enddo
!
    dofs = intx2(mu79(:,:,0P_1),temp79b)
    call vector_insert_block(kprad_rad%vec, itri,1,dofs,VEC_ADD)
```

Without this fix, there were isolated very large values of kprad_rad, causing spikes in pressure and magnetic field, causing the solves not to converge. Spikes could be 10^10 or 10^20 when the normal values were 10^-3 to 10^-8.

ITER VDE with new structure



New Structure Preserving Algorithm based on Slow Manifold

Chang Liu to present

Runaways with Sources



Chen: June 9, 2020

Nucl. Fusion 57 (2017) 066038

Reduced fluid simulation of runaway electron generation in the presence of resistive kink modes

A. Matsuyama^a, N. Aiba and M. Yagi



 $\frac{\partial n_{\rm RE}}{\partial t} + (v_{\rm RE}\mathbf{b} + \mathbf{u}) \cdot \nabla n_{\rm RE} = S_{\rm Dreicer} + S_{\rm avalanche},$

 $E = \eta (J - en_{\rm RE}c),$

M3D-C1 coupling to RE code KORC

• Plan to target DIII-D shot 177053 after Chen has a full simulation with fluid runaway electrons



- KORC can now run using fields, densities, and temperatures from M3D-C1 hdf5 files using Nate's Fusion-IO routines
- ZOOM call held on Thursday June 18
 - Matt Beidler was able to login to portal and transfer Cesar's NSTX-U files to NERSC
 - However, these C-pellet runs do not generate large electric fields and so there is no point in running them with KORC
 - Cesar will do some 2D runs with varying background Carbon to see what impurity level is needed to produce strong thermal and current quench

That's All I have

Anything Else ?