status\_0521.md 2025-05-21

# SCRIPTS FOR SOLAR ACTIVE REGIONS MAGNETIC FIELD ANALYSIS



🥯 🦈 🥯 Active Regions Dynamics Analysis Pipeline 🥯 💖



Step 1. Getting the data for the Period

Add your JSOC\_EMAIL to the ENV in bash:

export JSOC\_EMAIL="your@email.com"

Modify parameters like analysis time & download dir in the script

python magnetic/jsoc\_query.py

Then run it in a screen interface. SDO/HMI 720s files will be downloaded in the output\_dir.

When at least a part of files downloaded, run this script again to prepare  $t \times t$  files, which are necessary for IDL.

Step 2. Manually setting parameters of analysis with IDL

Create a project in IDL from mmc folder of this repo and compile it.

All following scripts are related to mmc/vmf subdir.

#### Select a region for analysis

This steps should be done for the First and Last moments in time: for at least two EVENT\_TIME's.

```
prepare_hmi, DATA_DIR, EVENT_TIME, REFERENCE_TIME, SAVE_ALL_DIR
```

#### where

DATA\_DIR -- a folder with downloaded SDO/HMI data;

EVENT\_TIME -- an observation time in format 20170903\_090000\_TAI;

REFERENCE\_TIME -- time of the first observation in the same format;

SAVE\_ALL\_DIR -- output directory, where all data will be stored.



Fix the observed center of the Sun

```
;FIND CENTER HERE
dx0 = 0.5*(956.5 - 949.5)
dy0 = 0.5*(950.0 - 953.0)
```

via plotting in ranges dx, dy and vice-versa  $\sim +/-$  (940 -- 960) and  $\sim$  (-10 -- 10).

```
; DRAW A SUBREGION HERE. USED FOR SOLAR LIMB COORDINATES FINDING
sub_map, mapbz, smapbz, xrange=[-250, 50.0], yrange=[-400, -100.0]
```

After all you also should select a region of analysis here.

status\_0521.md 2025-05-21

```
X Look at the vector fields directions
```

```
plot_vmap,/over,smapbx,smapby,mapbz=smapbz,limit=180,scale=0.012,iskip=15,j
skip=15,$
  v_color=255,axis_color=0,/Nolabels,v_thick=2.0 ;,/No_arrow_head
;,/Noaxis
```

☑ Differential rotation will apply automatically if EVENT\_TIME differs from REFERENCE\_TIME

```
IF (REFERENCE_TIME NE EVENT_TIME) THEN BEGIN
  ; Load smapbz from sav file
  restore, SAVE_ALL_DIR + '/' + REFERENCE_TIME + '/bxyz_submap.sav',/ver
  mapbz = drot_map(mapbz, time=smapbz.time)
```

#### **Building map projection**

In IDL command line run

```
proj, IN_DIR
```

Where IN\_DIR is SAVE\_ALL\_DIR/REFERENCE\_TIME.

If projected map doesn't satisfy, you could go to the first step, to re-select the Region-of-interest. You also should . sav inside IN\_DIR.

Should I use cmd\_lat.sav from the REFERENCE (initial) TIME moment, picked from the beginning of analysis, for other time moments? Because, for some moments in time, a size of data may be different. For example, at 9AM [576, 480] at 10AM with newly evaluated, but same coords range, may become [577, 480]. What could be the reason for such behaviour? Should I recalculate the center of the sun every hour?

#### Choosing a region for analysis with respect to projection effects

In IDL command line:

```
plot_submap_m, IN_DIR
```

IN DIR is picked from previous step.



X ROI selection

```
restore, IN_DIR+'/mapbxyz_m.sav',/ver
xrange = [-244., 44]
yrange = [-395., -155.5]
res = [0.5, 0.5]
```

? In some future moments in time, since REFERENCE (initial) TIME, a size of data may be different. For example, at 9AM [\*576\*, 480] at 10AM with newly evaluated , but same coords range, may become [\*577\*, 480]. What could be the reason for such behaviour? Should I recalculate the center of the sun every hour? Not only rely on the reference time SOLAR\_CENTER estimation? Am I correct that I should

status\_0521.md 2025-05-21

keep the xrange and yrange the same (taking drot in account), for all the period of analysis [3 days]? But I also need to control the SUN\_CENTER on the step1 for every time in the period?



The shape of output data should be a multiple of a good number like 16, 32,...

```
; CHECK HERE THE SHAPE OF DATA. IT SHOULD BE MULTIPLE OF A GOOD NUMBER * 4
; Select the RANGE in the begin of file.
help, smapbz.data
```

#### Preprocessing data for a Force-Free extrapolation

```
In IDL run pro creb_lv3_, IN_DIR
```

This script will create a level3\_data (smoothed over 4x4 kernel) which will be used for a simulation.



This step may fail if data\_size cannot be divided by 4.

#### Preparing boundary conditions for extrapolation in MPI-AMRVAC

In IDL run read\_boundary, IN\_DIR, amrvac\_file, amrvac\_file\_nlfff This script prepares binaries for lfff and nlfff for reading in MPI-AMRVAC.

Carefully set the parameters in the beginning of file:

```
; Parameter needed to be set for your own case
nx = 144
                 ; x-size of the preprocessed magnetic field
ny = 120
                 ; y-size
arcsec2cm=7.29809e7; 2825e7 ; 1 arcsec in cm
xc=-100.23546*arcsec2cm; x-coordinate of the center of the field of view
yc=-275.35374*arcsec2cm ; y-coordinate of the center
dx=4.0*0.50d*arcsec2cm ; x spatial resolution of the preprocessed magnetic
field
                 ; y spatial resolution
dy=dx
```

xc, yc could be found from the output of help, smapbz from previous step. arcsec2cm could be calculated from the known solar radius 6.9634e10 cm & a radius in arcsec from the step 1:  $(|x_l|+|x_r|)/2$ . nx, ny is also the output shape of help, smapbz divided by 4.

This script will update the amrvac files, stored in mmc/vmf/potential, mmc/vmf/nlfff with x, y, z ranges. You should manually update nblocks for parallelization in . par files.

Step 3. Data preparation pipeline for the whole timerange

Automated boundary conditions creation with selected parameters from Step 2.

status\_0521.md 2025-05-21

#### Run in IDL

mmc/vmf/magnetic\_preparation\_pipeline.pro

**X** Setup folders

DATA\_DIR is picked from python script from the Step 1.

X Pipeline fails at 9:48, 4th timestep. Because DATA becomes of shape [577, 480]. This issue is related to previous questions

#### DAVE4VM velocity calculation for each timestep

Run in IDL mmc/idl\_dave4vm/velocity\_preparation.pro

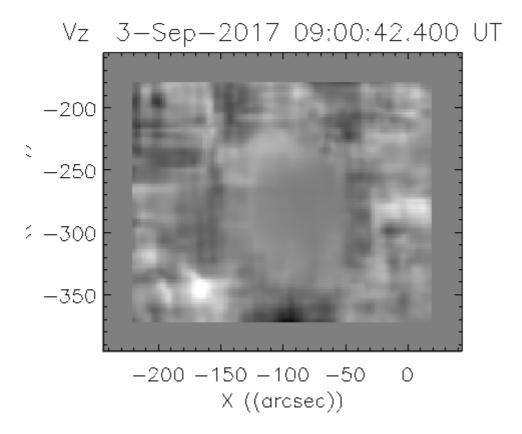
**X** Setup directories

The plasma velocity will be stored in

OUT\_DIR/<EVENT\_TIME>/level3\_data/velocity\_boundary.dat

- ? Am I correct that I use a magnetic field from /level3\_data/allboundaries.dat files for velocity estimation with a DAVE4VM algorithm?
- ? I observe NaNs near the borders of calculated region, which are proportional to the windowsize. Should I increase the size of Region-of-Interest? Currently I substitute NaNs with 0s. See image below.

status\_0521.md 2025-05-21



## ? Should I convert velocity\_boundary. dat to a binary file for MPI-AMRVAC?

Result of data preparation pipeline could be found here: assets/data.

### Step 4. Potential & NLFFF Extrapolation in MPI-AMRVAC

In every Potential subdir like 20170903\_090000\_TAI/level3\_data/extrapolation/potential setup MPI-AMRVAC:

```
setup.pl -d=3 -arch=default
make -j 16
mpirun -np 192 ./amrvac -i amrvac.par
```

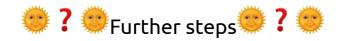
Then in every subdir for NLFFF with Magnetofriction method, run the same sequence.

If having issues run

make clean

If nblocks numbers are not good, update them as well.

status\_0521.md 2025-05-21



Generally, I've completed preparation and downloading of data. I can reconstruct the B-field for every moment in time with MPI-AMRVAC (potential field requires seconds, nlfff requires ~40 minutes). I also can obtain a velocity of plasma on the bottom boundary with an adaptation of DAVE4VM. Some issues for a long time range analysis remain (see above). I believe that they are related a careful RoI choice & Sun center corrections.

? Which next steps should I do to reproduce your paper Data-driven Modeling of a Coronal Magnetic Flux Rope: From Birth to Death?

shainalexander@yandex.ru

Best Regards, Aleksandr (Alex) Shain