
bifrostDocs Documentation

Release 1.0

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CHAPTER 1

helita

Helita documentation can be found here (the installation section is taken from this link):

<http://helita.readthedocs.io/en/latest/index.html>

Helita is a Python library for solar physics focused on interfacing with code and projects from the [Institute of Theoretical Astrophysics](#) (ITA) at the [University of Oslo](#). The name comes from Helios + ITA. Currently, the library is a loose collection of different scripts and classes with varying degrees of portability and usefulness.

1.1 Installation

1.1.1 Pre-requisites

To make use of helita **you need a Fortran compiler** ([GFortran](#) is recommended), because some modules are compiled from Fortran. The packages in the left column should be installed before installing helita. The right column contains recommended packages that allow the user to take advantage of all of the features.

Required	Recommended
<ul style="list-style-type: none">• Python (2.7.x, 3.4.x or later)• Astropy• NumPy• SciPy	<ul style="list-style-type: none">• Matplotlib• h5py• netCDF4• Cython• pandas• BeautifulSoup4

Helita will install without the recommended packages, but functionality will be limited. All of these packages are available through Anaconda, and that is the recommended way of setting up your Python distribution.

1.1.2 Cloning git repository

The easiest way is to use git to clone the repository. To grab the latest version of helita and install it:

```
$ git clone https://github.com/ITA-solar/helita.git  
$ cd helita  
$ python setup.py install
```

Note: The majority of the most updated versions of bifrost.py, ebysus.py, and all things bifrost related are in the <https://github.com/jumasy/helita.git> fork.

1.1.3 Non-root install

If you don't have write permission to your Python packages directory, use the following option with setup.py:

```
$ python setup.py install --user
```

This will install helita under your home directory (typically `~/.local`)

1.1.4 Developer install

If you want to install helita but also actively change the code or contribute to its development, it is recommended that you do a developer install instead:

```
$ python setup.py develop
```

This will set up the package, such as the source files used, from the git repository that you cloned (only a link to it is placed on the Python packages directory). Can also be combined with the `-user` flag for local installs:

```
$ python setup.py develop --user
```

1.1.5 Installing with different C or Fortran compilers

The procedure above will compile the C and Fortran modules using the default gcc/gfortran compilers. It will fail if these are not available in the system. If you want to use a different compiler, please use `setup.py` with the `-compiler=xxx` and/or `-fcompiler=yyy` options, where `xxx`, `yyy` are C and Fortran compiler families (names depend on system). To check which Fortran compilers are available in your system, you can run:

```
$ python setup.py build --help-fcompiler
```

and to check which C compilers are available:

```
$ python setup.py build --help-compiler
```

1.2 bifrost.py

bifrost.py is a set of tools to read and interact with output from Bifrost simulations. LMSAL has also contributed to this.

1.2.1 BifrostData class

bifrost.py includes the BifrostData class (among others). The following allows to create the class for snapshot number 430 from simulation ‘cb10f’ from directory ‘/net/opal/Volumes/Amnesia/mpi3drun/Granflux’ (a network local to LMSAL):

```
[1]: from helita.sim import bifrost as br

[2]: dd = br.BifrostData('cb10f', snap = 430, fdir = '/net/opal/Volumes/Amnesia/mpi3drun/
    ↪Granflux', verbose = False)
```

The snapshot(s) being read can be defined when creating the object, or set/changed anytime later. Snaps can be ints, arrays, or lists.

Getting variables & quantities

`get_var` and `get_varTime` can be used to read variables as well as to calculate quantities (with a call to `_get_quantity`). `iix`, `iiy`, and `iiz` are axis positions, or grid points. They can be specified to slice the return array (they can be ints, arrays, or lists). If `iix`, `iiy`, or `iiz` is not specified, `get_var` will read all the numerical domain along the x, y, or z axis, respectively. If no snapshot is specified, the current snap value will be used (either the initialized one, which is the first in the series, or the most recent snap used in `set_snap` or in a call to `get_var`). To get variable ‘r’ at snapshot 430 (a specific timestep) with only values at `iiy = 200`, `iiz = 5`, and `iiz = 7`:

```
[3]: var1 = dd.get_var('r', snap = 430, iiy = 200, iiz = [5, 7])

WARNING: cstagger use has been turned off, turn it back on with "dd.cstagop = True"
```

`get_varTime` can be used in the same fashion, with the added option of reading a specific variable or quantity from many snapshots at once. Its return arrays have an added dimension for time.

Several of the class’s parameters can be found in the dictionary `params`. This dictionary contains most of the parameters in .idl files. It includes information about time (`t` and `dt`) and the axes (`x`, `y`, `z`, `dx`, `dy`, and `dz`) among other things. When `snap` contains more than one snapshot, many of the parameters in the dictionary contain one entry for each snap. To get time:

```
[4]: time = dd.params['t']
```

To view all available keys:

```
[5]: dd.params.keys()

[5]: dict_keys(['mx', 'my', 'mz', 'mb', 'nstep', 'nstepstart', 'debug', 'periodic_x',
    ↪'periodic_y', 'periodic_z', 'ndim', 'u_l', 'u_t', 'u_r', 'u_p', 'u_u', 'u_kr', 'u_ee',
    ↪', 'u_e', 'u_te', 'u_tg', 'u_b', 'meshfile', 'dx', 'dy', 'dz', 'cdt', 'dt', 't',
    ↪'timestepdebug', 'nul', 'nu2', 'nu3', 'nu_r_xy', 'nu_r_xy_k', 'nu_r', 'nu_r_min',
    ↪', 'nu_r_k', 'nu_ee_xy', 'nu_ee', 'grav', 'eta3', 'ca_max', 'mhddebug', 'do_mhd',
    ↪', 'mhdclean', 'one_file', 'snapname', 'isnap', 'large_memory', 'nsnap', 'nsqr', 'aux',
    ↪', 'dtsnap', 'newaux', 'dtscr', 'tsnap', 'tscr', 'boundarychk', 'max_r', 'smooth_r',
    ↪', 'qmax', 'noneq', 'do_hion', 'gamma', 'tabinputfile', 'do_rad', 'dtrad', 'quadrature
    ↪', 'zrefine', 'maxiter', 'taustream', 'accuracy', 'strictint', 'linear', 'monotonic
    ↪', 'minbin', 'maxbin', 'dualsweep', 'teff', 'timing', 'spitzer', 'debug_spitzer',
    ↪', 'info_spitzer', 'spitzer_amp', 'theta_mg', 'dtgerr', 'ntest_mg', 'tgb0', 'tgb1',
    ↪', 'tau_tg', 'fix_grad_tg', 'niter_mg', 'bmin', 'do_genrad', 'genradfile', 'debug_
    ↪genrad', 'incrard_detail', 'incrard_quad', 'dtincrard', 'debug_incrard', 'bctypelower',
    ↪', 'tau_bcl', 'tau_ee_bcl', 'tau_d2_bcl', 'tau_d5_bcl', 'tau_d6_bcl', 'tau_d7_bcl',
    ↪', 'tau_d8_bcl', 's0', 'e0', 'r0', 'cs0', 'p0', 'nsmooth_bcl', 'naver_bcl', 'nclean_bcl
    ↪', 'nclean_lbl', 'nclean_ubl', 't_bdry', 'rbot', 'ebot', 'bx0', 'by0', 'x0_bcu',
    ↪', 'x1_bcu', 'y0_bcu', 'y1_bcu', 'uz_bcu', 'strtb', 'rtb', 'xotb', 'zotb', 'contines on next page)
    ↪', 'nclean_bcu'])
```

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Table 1: Variables

Simple	r, px, py, pz, e, bx, by, bz, p, tg, i1, i4, qjoule, qspitz
Composite	ux, uy, uz, ee, s

Table 2: Calculated Quantities

Derivatives	dxup, dyup, dzup, dxdn, dydn, dzdn
Centers vector quantity in cells	xc, yc, zc
Module of vector	'mod' + root letter of varname (eg. modb)
Divergence of vector	'div' + root letter of varname (eg. divb)
Squared module	root letter of varname + '2' (eg. u2)
Ratio of two vars	var1 + 'rat' + var2 (eg. rratpx)
Eostab (unit conversion to SI)	ne, tg, pg, kr, eps, opa, temt
Magnitude of vector components // or	root letter of v1 + 'par' or 'per' + root letter of v2 (eg. uparb)
Current/Vorticity	ix, iy, iz, wx, wy, wz
Flux	pfx, pfy, pfz, pfx, pfey, pfez, pfwx, hx, hy, hz, kx, ky, kz
Plasma	beta, s, ke, mn, man, hp
Wave speeds	alf, fast, long, va, cs, vax, vay

Small Demo

Here, we plot dz versus z, illustrating the non-uniform gaps in the z axis.

```
[6]: from helita.sim import bifrost as br
import matplotlib.pyplot as plt
import numpy as np

rootname='l2d90x40r_it' #this is for the 2D case
fdirectory='/net/opal/Volumes/Amnesia/mpi3drun/2Druns/genohm/rain/l2d90x40r'
dd=br.BifrostData(rootname,fdirectory, verbose = False) #this loads the structure
var=dd.get_var('r',305) #this reads the density (r) for the instant 305

# gets the z coordinates of data points and makes an empty array of the same length
# for values of dz
zarr = dd.z
length = zarr.shape[0]
dzarr = np.empty([length])

# iterates through zarr and sets each entry in dzarr as the difference between the
# next and current values of z
# for the final value of dzarr, sets it to be the same as the previous value of dzarr
for i, val in enumerate(zarr):
    if i < length - 1 :
        dzarr[i] = zarr[i + 1] - val
        i = i + 1

    else :
        dzarr[i] = dzarr[i - 1]

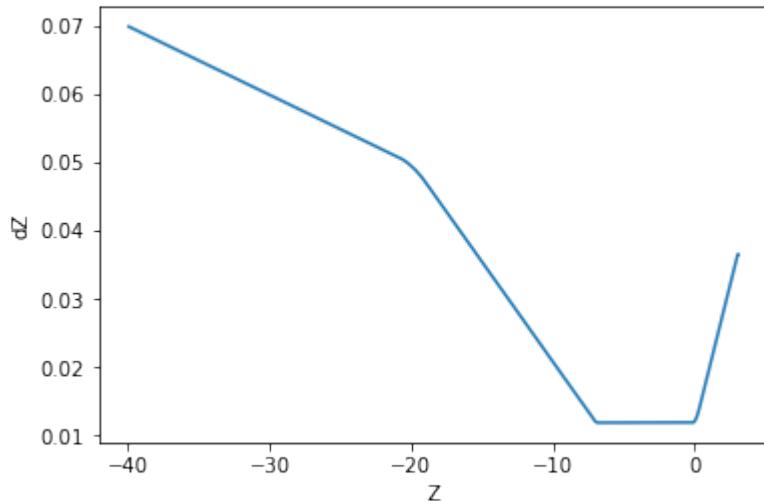
# plots z vs dz and labels the axes
```

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```
plt.plot(zarr, dzarr)
plt.xlabel('z')
plt.ylabel('dz')
plt.show()
```

WARNING: cstagger use has been turned off, turn it back on with "dd.cstagop = True"



In this example, we interpolate temperature data from the simulation to conform to a uniform z axis and plot the data with both the original, and the even, z axes.

```
[7]: import scipy.interpolate as sp
# bifrost already imported and dd initialized after previous example

temp = dd.get_var('tg', 300)

zarr = dd.z
xarr = dd.x

# at instant 305 - x is uniform, y is kept constant, and z is non uniform

# converts the 3d array with width of 1 to 2d array (and transposes because otherwise
# image is sideways)
temp2d = np.transpose(temp[:, 0])

f = sp.interp2d(xarr, zarr, temp2d)

# makes a new uniform z, x is already uniform
zlength = zarr.shape[0]
zstart = zarr[0]
zend = zarr[zlength - 1]
newz = np.linspace(zstart, zend, zlength)

# makes the interpolated data with the previously created interpolation function,
# acting on the new uniform axes
uniform_a = f(xarr, newz)
```

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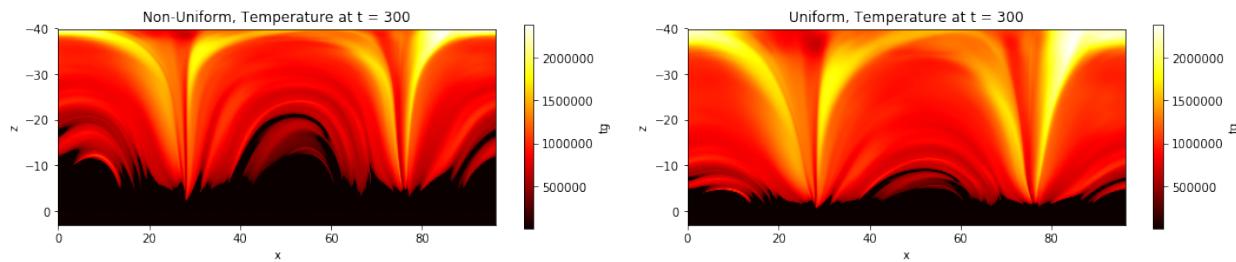
```

fig = plt.figure(figsize = (15, 3))

# optional plotting of the non uniform data to show the difference
ax0 = fig.add_subplot(1, 2, 1)
im0 = ax0.imshow(temp2d, cmap = 'hot', extent = (xarr[0], xarr[xarr.shape[0] - 1], zstart, zend), aspect = 'equal')
ax0.set_title('Non-Uniform, Temperature at t = 300')
ax0.set_xlabel('x')
ax0.set_ylabel('z')
c0 = fig.colorbar(im0, ax = ax0)
c0.set_label('tg')

ax1 = fig.add_subplot(1, 2, 2)
im1 = ax1.imshow(uniform_a, cmap = 'hot', extent = (xarr[0], xarr[xarr.shape[0] - 1], zstart, zend), aspect = 'equal')
ax1.set_title('Uniform, Temperature at t = 300')
ax1.set_xlabel('x')
ax1.set_ylabel('z')
c1 = fig.colorbar(im1, ax = ax1)
c1.set_label('tg')
plt.tight_layout()
plt.show()

```



1.2.2 create_new_br_files class

1.2.3 bifrost_units class

1.2.4 Rhoeetab class

1.2.5 Opatab class

1.2.6 FFTData class

This class can be found within bifrost_fft.py. It performs operations on Bifrost simulation data in its native format. After creating a class for a specific snap root name and directory (much like with BifrostData), one can get a dictionary of the frequency and amplitude of the Fourier Transform for a certain quantity over a range of snapshots.

We have defined 3 variables that allow us to decompose the velocity in Alvenic, fast mode and longitudinal component ('alf', 'fast', and 'long'). Here, we show the transformation of 'alf',

```
[8]: from helita.sim import bifrost_fft as brft
```

```
[9]: dd = brft.FFTData(file_root = 'cb10f', fdir = '/net/opal/Volumes/Amnesia/mpi3drun/
    ↪Granflux')

[10]: transformed = dd.get_fft('alf', snap = [430, 431, 432], iix = 5, iiy = 20)
WARNING: cstagger use has been turned off, turn it back on with "dd.cstagop = True"

[11]: transformed.keys()
[11]: dict_keys(['freq', 'ftCube'])
```

Depending on the number of snaps and the size of the cube, using cuda or python multiprocessing may speed up the calculation.

1. Using cuda

If pycuda is available, the code imports reikna (a python library that contains fft functions using pycuda). In order to make use of the GPU, use the function run_gpu(). The default is to **not** use the GPU, even if there is one available.

```
[12]: dd.run_gpu() # to use GPU
dd.run_gpu(False) # to stop use of GPU
```

When get_fft() is called, the GPU will be used in accordance with the last call to run_gpu(). If the GPU has limited memory, the user can specify numBlocks in the call to get_fft(). This will send the calculation over to the GPU in several blocks as opposed to all at once. To use 5 different blocks, a call would look like this:

```
[13]: usingBlocks = dd.get_fft('bx', snap = [400, 401, 402], numBlocks = 5)
```

2. Using python multiprocessing

This can be used whether or not pycuda is available, as multiprocessing is a library that comes with python. It makes use of threading on the CPU. In order to use a multiprocessing threadpool when calculating the Fourier Transform, specify numThreads with a number greater than 1, when calling get_fft():

```
[14]: usingThreads = dd.get_fft('bx', snap = [400, 401, 402], numThreads = 10)
```

FFT demo #1

This first demo tests the get_fft() method with standard functions: a sine wave, a gaussian curve, and y = 0. It pre-sets dd.preTransform and

```
[15]: import numpy as np
import helita.sim.cstagger
from helita.sim.bifrost import BifrostData, Rhoeetab, read_idl_ascii
from helita.sim.bifrost_fft import FFTData
import matplotlib.pyplot as plt

# note: this calls bifrost_fft from user, not /sanhome

dd = FFTData(file_root='cb10f',
             fdir='/net/opal/Volumes/Amnesia/mpi3drun/Granflux')
```

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```

# test 1: ft of y = sin(8x)
x = np.linspace(-np.pi, np.pi, 201)
dd.preTransform = np.sin(8 * x)
dd.freq = np.fft.fftshift(np.fft.fftfreq(np.size(x)))
dd.run_gpu(False)
# preTransform is already set
tester = dd.get_fft('not a real var', snap='test')
fig = plt.figure(figsize=(15,10))

numC = 3
numR = 2

# plotting original sin signal
ax0 = fig.add_subplot(numC, numR, 1)
ax0.plot(x, dd.preTransform)
ax0.set_title('original signal' + '\n\nsine wave')

# plotting transformation sin signal
ax1 = fig.add_subplot(numC, numR, 2)
ax1.plot(tester['freq'], tester['ftCube'])
ax1.set_title('bifrost_fft get_fft() of signal' + '\n\n ft of sine wave')
ax1.set_xlim(-.2, .2)

# test 2: ft of gaussian curve
n = 30000 # Number of data points
dx = .01 # Sampling period (in meters)
x = dx*np.linspace(-n/2, n/2, n) # x coordinates

stanD = 2 # standard deviation
dd.preTransform = np.exp(-0.5 * (x/stanD)**2)

# plotting original gaussian signal
ax2 = fig.add_subplot(numC, numR, 3)
ax2.plot(x, dd.preTransform)
ax2.set_xlim(-25, 25)
ax2.set_title('gaussian curve')

# plotting transformation of gaussian signal
dd.freq = np.fft.fftshift(np.fft.fftfreq(np.size(x)))
ft = dd.get_fft('not a real var', snap='test') # preTransform is already set
ax3 = fig.add_subplot(numC, numR, 4)
ax3.plot(ft['freq'], ft['ftCube'])
ax3.set_xlim(-.03, .03)
ax3.set_title('ft of gaussian curve')

# test 3: ft of y = 0
# plotting original horizontal line
x = np.linspace(-20, 20, 50)
dd.preTransform = [0] * 50
ax4 = fig.add_subplot(numC, numR, 5)
ax4.plot(x, dd.preTransform)
ax4.set_title('y = 0')

# plotting transformed signal
dd.freq = np.fft.fftshift(np.fft.fftfreq(np.size(x)))
ft = dd.get_fft('not a real var', snap='test') # preTransform is already set
ax5 = fig.add_subplot(numC, numR, 6)

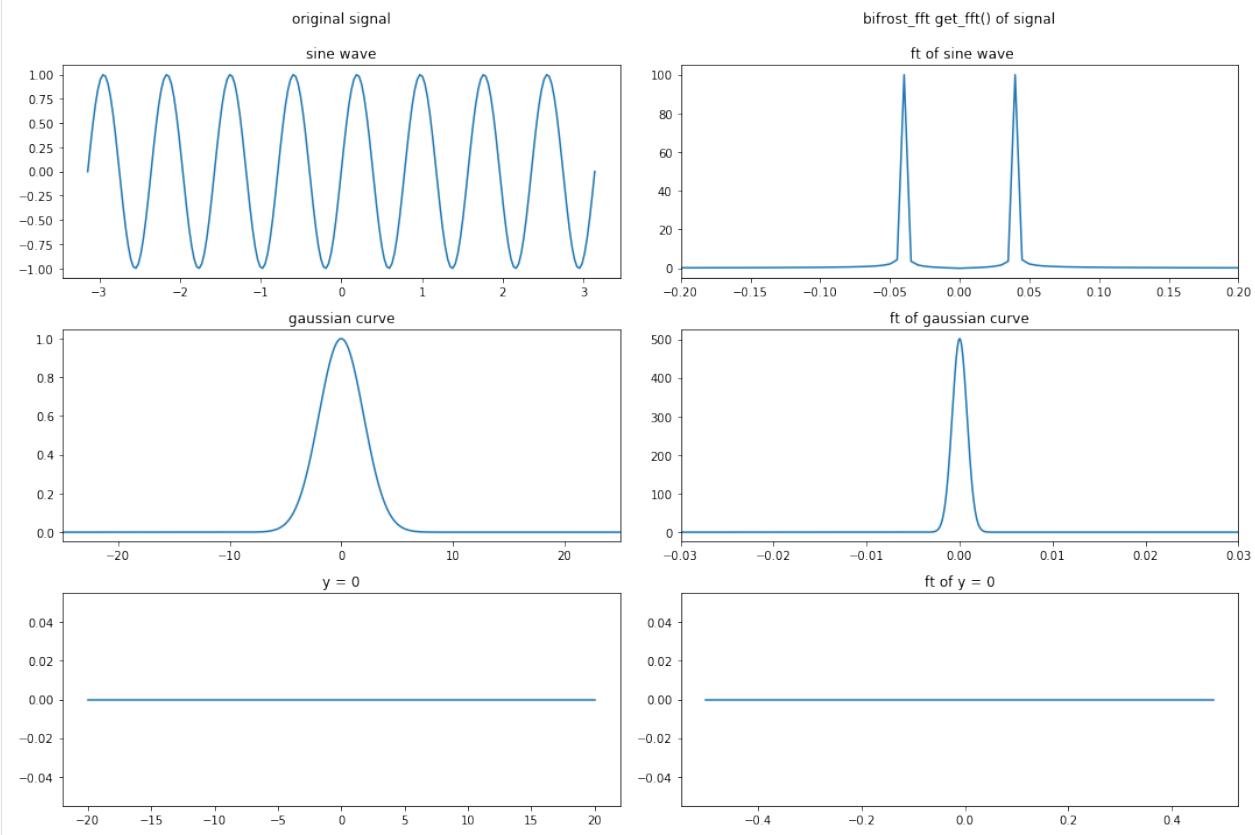
```

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```
ax5.plot(ft['freq'], ft['ftCube'])
ax5.set_title('ft of y = 0')

plt.tight_layout()
plt.show()
```



FFT demo #2

Here, we use `get_fft()` to find the transformation result for `bx` at each `z` position (from a local network containing 2d simulations).

```
[22]: import numpy as np
import helita.sim.cstagger
from helita.sim.bifrost import BifrostData, Rhoeetab, read_idl_ascii
from helita.sim.bifrost_fft import FFTData
import matplotlib.pyplot as plt

snaps = np.arange(280, 360)
v = 'bx'

dd = FFTData(file_root='12d90x40r_it',
              fdir='/net/opal/Volumes/Amnesia/mpi3drun/2Druns/genohm/rain/12d90x40r/')

# getting ft
transformed = dd.get_fft(v, snaps)
ft = transformed['ftCube']
```

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```

freq = transformed['freq']
zaxis = dd.z

# making empty array to later contain the averages for each z position
zstack = np.empty([np.size(freq), np.shape(ft)[1]])
# filling zstack with average ft for each (x,y) in each z level
for k in range(0, np.shape(ft)[1]):
    avg = np.average(ft[:, k, :], axis=(0))
    zstack[:, k] = avg

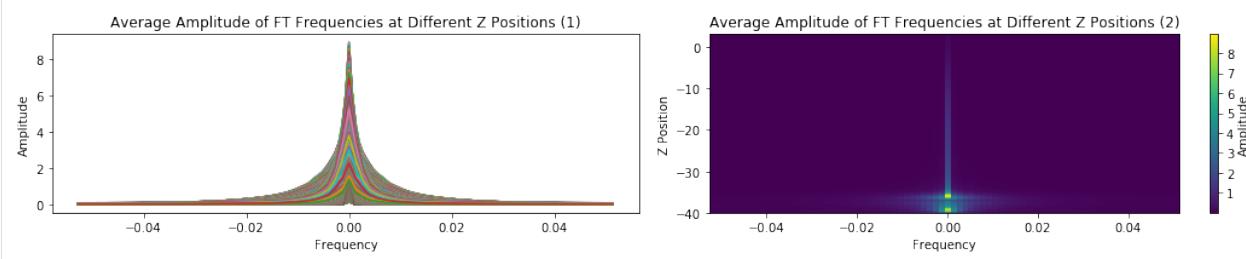
# preparing plots
fig = plt.figure(figsize = (15, 3))
numC = 1
numR = 2

# plotting freq vs amp with multiple lines (1 for each z position)
ax0 = fig.add_subplot(numC, numR, 1)
ax0.plot(freq, zstack)
ax0.set_xlabel('Frequency')
ax0.set_ylabel('Amplitude')
ax0.set_title(
    'Average Amplitude of FT Frequencies at Different Z Positions (1)')
ax0.set_aspect('auto')

# plotting amp at different freq & z with image
ax1 = fig.add_subplot(numC, numR, 2)
im1 = ax1.imshow(zstack.transpose(), extent=[freq[0], freq[-1], zaxis[0], zaxis[-1]])
ax1.set_xlabel('Frequency')
ax1.set_ylabel('Z Position')
ax1.set_title(
    'Average Amplitude of FT Frequencies at Different Z Positions (2)')
ax1.set_aspect('auto')
c1 = fig.colorbar(im1, ax = ax1)
c1.set_label('Amplitude')
plt.tight_layout()
plt.show()

```

WARNING: cstagger use has been turned off, turn it back on with "dd.cstagop = True"



CHAPTER 2

bq_t5_look tool

This tool can be used to visualize and manipulate data sets produced by Bifrost (and Ebysus) simulations. Previous tools were written in IDL, but this tool is written in Python, increasing accessibility. It also has a range of added features including the creation of movies and animations.

2.1 Setup

Note: The Bifrost code, which includes this tool, is **not** publicly available yet.

This tool can be found in Bifrost (which is a separate folder from helita), and requires that helita be installed in order to function. Once you have created a path for the Bifrost folder, add that path to your .cshrc, eg:

```
setenv BIFROSTPATH yourpath
```

2.2 Launch

To launch the tool, call:

```
python path_to_file -i path_to_snapshot
```

Possible extensions include:

-i, --input	required, points to snapshot file
-h, --help	returns help message
-s, --slice	jumps directly to data[:, :, slice], defaults to 0
-z, --depth	finds a slice at the corresponding depth value provided in real [mM] coordinates
-e, --ebysus	code based on Bifrost

Visu

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Abstract

The complexity of spectropolarimetric images from observations of the solar atmosphere require models to interpret them. The example in Figure 1 shows the interface region between the surface of the Sun and the million degree corona, i.e., the chromosphere, **NASA's Interface Region Imaging Spectrograph** (De Pontieu et al 2014). To interpret these observations we use **radiative magnetohydrodynamic numerical simulations**. These numerical simulations generate **data sets** comprised of space and time (Chapter 2 bq_15 look tool) of many variables (e.g. velocity, or magnetic field) and scalar (e.g. density) variables. These data sets can be used in

2.3 Description

open corresponding
settings windows
(1)-(5)

input custom data for
image (4a)

slide through the 3rd
dimension

2.3.1 Initial window

- Slider: moves through third dimension
- Variable choice: select variable to be plotted from dropdown
- Customizable quantity: input custom data for image
- Plane view: select plane view from dropdown
- Control of overlays: quickly add or remove overlays specified in respective window

2.3.2 Overlays & additional settings

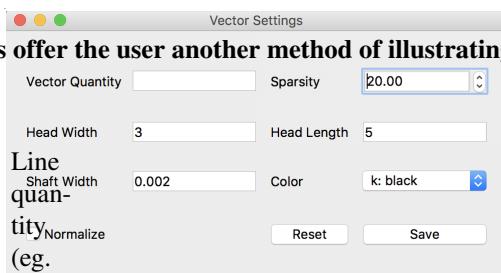
Each of the overlays (vectors, lines, and contour) has its own corresponding pop-up settings window that can be opened from the initial page of the tool. As previously mentioned, the initial window also has the option to quickly hide/reveal a previously loaded overlay. By default, all overlays are shown once a quantity is specified.

1. Vectors

This feature allows the user to visualize non-scalar data. The user can manipulate:

- Vector quantity (eg. p or u)
 - Arrow head width/length
 - Arrow shaft width
 - Sparsity of vectors shown
-

2. Lines



Lines offer the user another method of illustrating non-scalar data besides vector fields. The user can specify:

- Head Width: 3 Head Length: 5
- Line quantity: 0.002 Color: k: black
- Reset Save

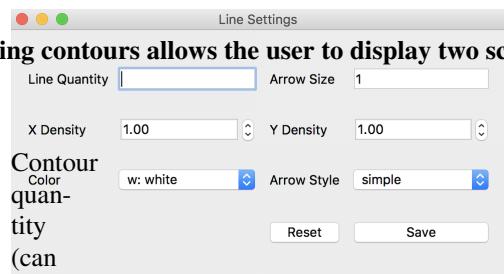
- Fig. 2: Vector Settings Window

Line density in the X and Y directions

2.3. Description

- Line color
 - Arrow style and size
-

3. Contour



Plotting contours allows the user to display two scalar quantities simultaneously. The user can determine:

- X Density 1.00 Y Density 1.00
- Contour Color quantity (can select from dropdown or specify custom quantity)

- Scale (eg. absolute or log)

- Units (CGS)

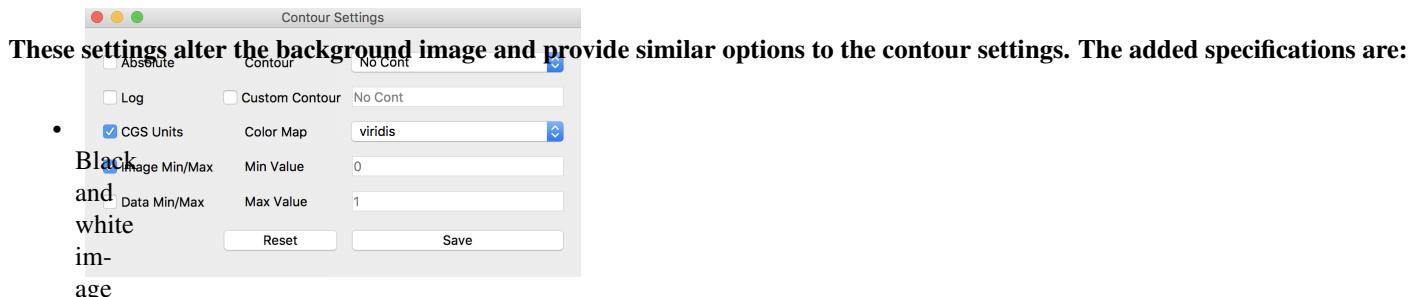
- Minimum and maximum used (based on initial distribution)

Fig. 3: Line Settings Window

vid-
ual
slice, whole data cube, or custom values)

- Color map
-

4. Additional display features

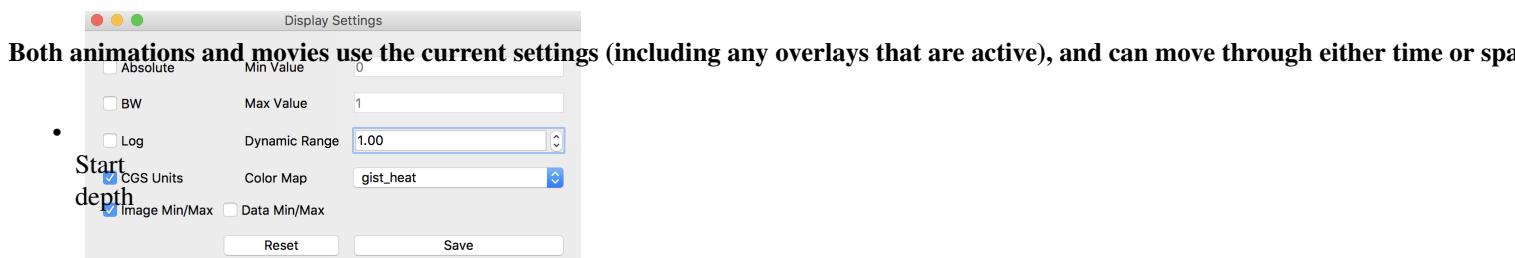


- Fig. 4: Contour Settings Window
- Dynamic range

The repeated features are:

- Scale
 - Units
 - Minimum and maximum used
 - Color Map
-

5. Movies and animation



2.3. Description

Fig. 5: Display Settings Window

and
end
depth

- OR
start
time
and
end
time
(if
“Through
Time”
box
is
checked)

Animation features:

- Make
an-
i-
ma-
tion
be-
gins
the
an-
i-
ma-
tion
on the display window
- User
can
pause/play
cur-
rent
an-
i-
ma-
tion
at
any
point
- Animation
will
loop
un-
til

paused

Movie features:

- File
des-
ti-
na-
tion
(de-
fault
is
home
di-
rec-
tory)
- Frames
per
sec-
ond
- Whether
pic-
tures
should
be
saved
or
re-
moved
(the
movie
is
created from saved pictures)

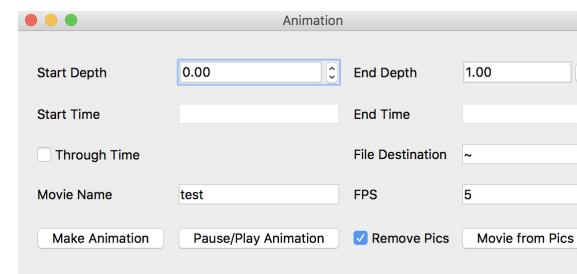


Fig. 6: Animation Settings Window