Running the Stacking Code: caustic_stack

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1. Introduction:

Here is a quick tutorial on how to get and run the Stacking code. It is dependent on Dan Gifford's causticpy, as well as other personalized scripts. Having a basic understanding of causticpy will greatly help in understanding caustic_stack.

Within caustic_stack is __init__.py, which is a script that allows you to take the phase spaces of galaxy clusters (the cluster-centric radius vs. line-of-sight velocity of galaxies or DM particles) and overlay them on top of each other (aka "stacking"), creating a joined phase space called an ensemble phase space, or ensemble cluster. The code then utilizes causticpy to run the caustic technique over the ensemble to estimate a mass (Mcrit200), along with other dynamics of the ensemble cluster.

The basic premise of the code, is that the users feeds it an array of phase spaces, and the code then stacks them and runs the caustic technique. Within the caustic_stack directory there will be an example/ folder containing an example of how this is implemented in practice.

2. Dependencies:

See causticpy for relevant dependencies. Along with the same software causticpy is dependent upon, this code is also dependent on DictEZ.py and AttrDict.py, which are included with caustic_stack.

3. Getting the Code:

The code can be downloaded from Github: caustic_stack can be found here, and causticpy can be found here. There are directions to setup causticpy in its README file. To install caustic_stack, you only need to **git clone** the URL, and then add the caustic_stack/ directory to your PATH or PYTHONPATH. In a bash terminal this would be a command in your ~/.bash_profile file that looks something like this:

export PYTHONPATH=/where_ever_this_folder_lives/:\$PYTHONPATH

4. Overview of the Code:

All of the necessary code is in one script called __init__.py. In order to use it, simply add this to the front of a python script:

from caustic_stack import *

This will not only load in the classes defined in __init__.py, but will also load in all of the modules listed at the top of __init__.py.

There are three classes defined in __init__.py:

Data

This class is a class that defines an "empty container" so-to-speak, where one can attach data to it and it will stack or append that data for you. It has a few methods attached to it that allows one to do this. For example, if I have 3 arrays, a b and c, that I want to "stack" together into an array called 'final' (in a sense this is merely concatenating them) then I can do this inside a python script:

```
import numpy as np
a = np.array([1,2,3,4,5])
b = np.array([6,7,8,9,10])
c = np.array([11,12,13,14,15])
D = Data()
D.extend({'final':a})
D.extend({'final':b})
D.extend({'final':c})
print D.final
np.array([1,2,3,4,5,6,7,8,9,10,11,12,13,14,15])
```

Stack

This is the class that contains the function **caustic_stack()**, which is the main function you need in order to run the stacking code.

Universal

This contains functions needed by Stack.caustic_stack() in order to complete its stacking routine. Other functions that may be of interest, but not necessary, are there, such as line_of_sight(), which is a way to make a mock projection of galaxies within a simulation, and print_varibs(), which is a nice function to call at the beginning of any script to print out the values of flags and variables defined in the run in a clean way.

Creating a caustic params.py file:

In order to use this caustic_stack(), you need to define a few flags outside of the function. This is most easily done by creating a parameter file, which can be called caustic_params.py. You should keep this file in the same directory that you run your personal python script that imports caustic_stack. Inside of the parameter file, you should define the following flags:

```
gal_num: integer
```

Number of galaxies to take per individual cluster phase space (Ngal)

```
line_num: integer
```

Number of individual phase spaces to stack into one ensemble (Nclus)

```
scale data: True, False
```

Scale radial data by R200 before stacking?

run_los: True, False

Run caustic technique over individual phase spaces (lines-of-sight)?

avg_meth: 'median', 'mean'

Uses either median or mean to estimate the properties of a Bin.

mirror: True. False

Used in causticpy's estimation of the caustic surface: mirror the v data of the phase space before solving for the caustic surface?

c: 2.9979e5

Speed of light in km/s

h: 1.0

Hubble constant / 100.0

H0: 100*h

Hubble constant

In order to properly load the caustic_params.py file, you should incorporate the following code block with your IMPORT statements:

```
import os, sys
sys.path.insert(0,os.getcwd())
__import__('caustic_params')
from caustic_params import *
print "Loaded caustic_params from",sys.modules['caustic_params']
```

This will take the file called "caustic_params.py" in your working directory and load any variables defined in it into the working session. To check this is the "correct" caustic_params.py, look at the print statement, it should print out the caustic_params.py file in your working directory. An example caustic_params.py file is located in the example/ directory in caustic_stack/

5. Worked Example

Within a custom python script, called *script.py* for example, we can load caustic_stack and work with the function Stack.caustic_stack(*args,**kwargs) to run the stacking code. First we define a basic *caustic_params.py* file in the same directory as *script.py* in order to define certain variables that are needed.

```
----- caustic_params.py -----run_los = True # Run caustic over each individual cluster as well?
```

```
scale data = False
                            # Scale radial data by cluster R200 before stacking?
init_clean = False
                            # Run shiftgapper on individual cluster before stacking?
mirror = True
                            # Mirror phase space over velocity axis before surface estimation?
ens num = 1
                            # Number of ensembles to solve for
gal num = 50
                            # Number of galaxies to stack per line of sight (Ngal)
line num = 5
                            # Number of clusters to stack into one ensemble (Nclus)
method num = 0
                            # Method to choose galaxies from line of sight, 0=Top brightest
avg meth = 'median'
                            # Method to create bin properties, 'mean' or 'median' (fed as string)
c = 2.9979e5
                            # Speed of light in km/s
h = 1.0
                            # hubble constant / 100
H0 = 100*h
                            # hubble constant
----- end caustic params.py -----
```

In this brief example, we can assume that we already have our phase spaces created: in simulation data this means taking the cluster centric galaxy coordinates, making a mock projection away from the cluster center and putting the galaxies in a radius vs. line-of-sight velocity diagram, which is called the phase space. Lets assume that we have 5 individual clusters we want to stack (line_num = 5) and we want to stack 50 galaxies from each cluster (gal_num = 50), and we only want to do this for 1 ensemble (ens_num = 1). We will assume that this data is provided in the form of a 2 dimensional numpy array called Rdata and Vdata, for example:

The same goes for Vdata. Notice that the first index of Rdata--as in Rdata[0] or Rdata[1] etc.--differentiates between unique individual clusters, meaning that the second index--as in Rdata[0][0] or Rdata[0][1]--differentiates between unique galaxies within a unique individual cluster. Note that caustic_stack() will take care of limiting the phase space and building the ensemble according to either **method 0 - take top Ngal brightest galaxies** or **method 1- take a random Ngal galaxies**, but if you want to just stack the Rdata and Vdata fed arrays, you can feed caustic_stack() the keyword stack_raw = True. Therefore, we are not constrained to have every array within Rdata or Vdata be the same length; although we want a final gal_num of 50 galaxies to be stacked per individual cluster, we can feed Rdata as an array that has 300 elements in Rdata[0], and 257 elements in Rdata[1] and 791 elements in Rdata[2], caustic_stack() will handle limiting the sample down to gal_num galaxies within R200 so long as gal_reduce is kept set to True as it is by default. In light of this, it is good to note that numpy does not like having a 2 dimensional array with its 1 dimensional components having different

lengths. To work around this try appending the desired 1 dimensional arrays to a list (as arrays), and then change the final list to an array, as such:

After having the basics, like our phase spaces in Rdata and Vdata arrays, we need the HaloIDs and HaloData at the very minimu.

HaloID is a 1 dimensional array that has the same length as the first axis of Rdata or Vdata, this means that it has the same length as the number of clusters we want to stack into an ensemble cluster (line_num), which in our case is 50. It is a unique identifier, usually a long integer, for each cluster.

HaloData should be fed as a 2 dimensional array, but only has three columns and line_num rows. The three columns should be the M200, R200 and HVD of each cluster, which are defined in simulations. If you don't have this data, i.e. real universe, you still need to feed the HaloData argument, but you should feed it as None.

Lastly, you need to feed the 'stack_num' argument, which is the same as 'line_num' flag, it is the number of cluster that you plan on stacking into one ensemble cluster.

Additionally, if you plan on using method 0 to build an ensemble, you should also use the keyword feed_mags = True. If this is the case you need to feed all three keywords magnitude arrays, G_Mags, R_Mags and I_Mags. If you only have on magnitude array, then **be sure to feed it to R_Mags** and create [None] type arrays of the same shape and feed those to G_Mags and I_Mags.

Finally I'll give a brief example of what a script.py file looks like, how to call the Stack() class and use the caustic_stack() function.

```
## Load Modules
from caustic_stack import *

## Load caustic_params.py file
sys.path.insert(0,os.getcwd())
__import__('caustic_params')
from caustic_params import *
print "Loaded caustic_params from",sys.modules['caustic_params']
```

Make variables defined in caustic_params.py file into a dictionary using DictEZ.py loaded from caustic_stack as ez

'stack_data' is returned as a dictionary, containing variables within it that pertain to the caustic run over the ensemble and individual phase spaces if run_los = True, such as:

ens_caumass: "Ensemble Caustic Mass", float

This is the caustic mass estimate of the ensemble at hand.

ens_caumass_est: "Ensemble Caustic Mass using R200 estimated by code", float

This is the caustic mass estimate using caustipy's estimate of R200, rather than the fed
R200.

ens_edgemass: "Ensemble Edge-detector Mass", float

This is the mass estimated by causticpy's edge detector function.

ens_edgemass_est: "Ensemble Edge-detector Mass using R200 estimate", float

This is the edge detector mass, but using causticpy's estimate of R200, rather than fed.

ens_causurf: "Ensemble Caustic Surface", array

This is an array containing the caustic surface (A(r) in Diaferio 1999).

ens_nfwsurf: "Ensemble NFW Fit Surface", array

This is the NFW that is fit to the caustic surface, mostly outdated and not necessary.

ens_r: "Ensemble R Data", array

This is an array that contains rdata for ensemble phase space.

ens_v: "Ensemble V Data", array

This is an array that contains vdata for ensemble phase space.

ens_gmags,ens_rmags,ens_imags: "Ensemble SDSS (G/R/I) Magnitudes", array
These are arrays containing the SDSS G/R/I magnitudes of the ensemble galaxies.

ens_gal_id: "Ensemble Galaxies ID", array

This is an array containing a unique identifier (index) for each galaxy within the ensemble, with respect to the natural cumulative ordering of galaxies in the fed Rdata and Vdata arrays.

ens_clus_id: "Ensemble Galaxies' cluster ID", array

This is an array containing the galaxies' initial parent cluster HaloID for each galaxy in the ensemble.

Same arrays exist for individual phase spaces, but with "ens" replaced with "ind".

x_range: "Range of x axis for Gaussian Kernel Density Estimation", array

Steps in radius used for creation of density map. Matches up with ens_causurf and

ind_causurf arrays, and used as x-axis in order to plot those arrays.