Pipelines for radio interferometric data reduction

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# Outline

- Why pipelines ?
- AIPS based pipelines
- Writing your own AIPS pipeline
- CAPTURE : CASA Pipeline-cum-Toolkit for UGMRT Data Reduction
- Writing tasks in CASA, creating your own pipelines

# Why pipelines ?

- Data reduction is a lengthy process, many parts are similar across datasets: automation !
- Increases reproducibility of the results
- Reduces human errors
- Data sizes are large or going to get larger: easy to port to servers and run remotely.
- Ease of testing one aspect at a time for complex algorithms like "tclean"
- A step towards "open science": https://zenodo.org/record/2631868#.XWiX-JzhVUQ

"Reproducibility and open science in the SKA era" by Rachel Ainsworth

### AIPS



- RUNFILES
- Commands can be put into a text file and the text file can be provided to AIPS
- Quirks: needs to have extension of AIPS userid in e-hex format
- Any programming: e. g. for loops, condition testing etc. in "Parseltongue"

### AIPS



- SPAM: Source Peeling and Atmospheric Modeling (Interna et al 2009)
- http://www.intema.nl/doku.php?id=huibintemas pam

# CASA

- Writing tasks and pipelines: Python
- Writing a task in CASA:
  XML file : sets the input interface taskname.py : actual code
- Use of CASA tasks and toolkit functionalities
- Excellent documentation for learning.

# CASA

- Writing tasks and pipelines: Python
- Pipeline:

- automating the process that you do interactively at the terminal

- automation of decision making is crucial !

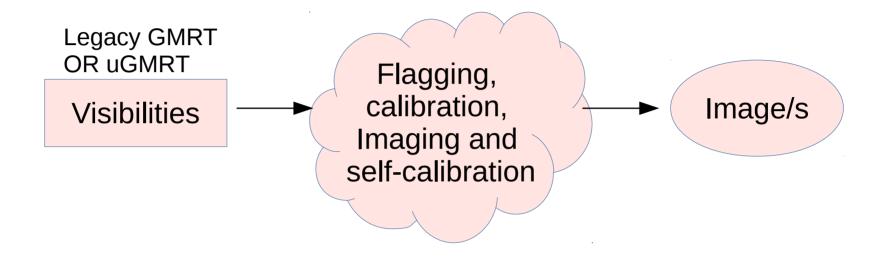
### CAPTURE

### CASA Pipeline-cum-Toolkit for UGMRT Data Reduction

- Automation of routine processes: conversion from Ita to MS, flagging of bad antennas, standard calibration, splitting target source data.
- Efficient elimination of RFI while not overdoing it.
- Automated self-calibration but still giving enough freedom to the to choose the strategy.
- Easily tailored for special needs: for e.g. for online RFI excision system testing: can deal with only calibrator data, half or one fourth of an array of data

### CAPTURE

https://github.com/ruta-k/uGMRT-pipeline



Python 2.7 and Common Astronomy Software Applications

(CASA, McMullin, J. P. et al 2007)

#### Input files

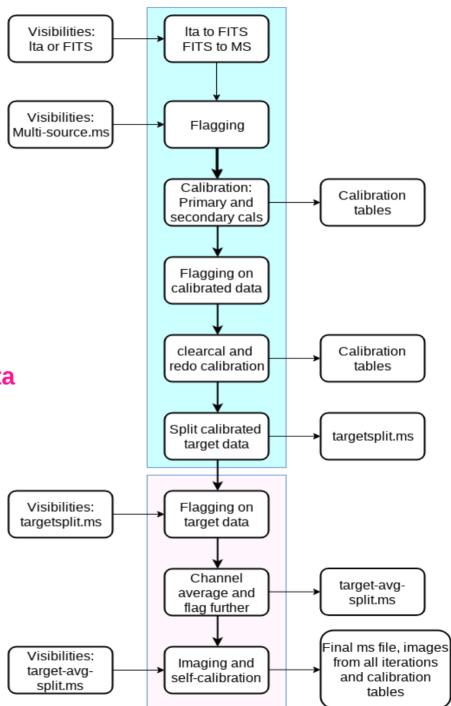
# CAPTURE

#### **Multi-source data**

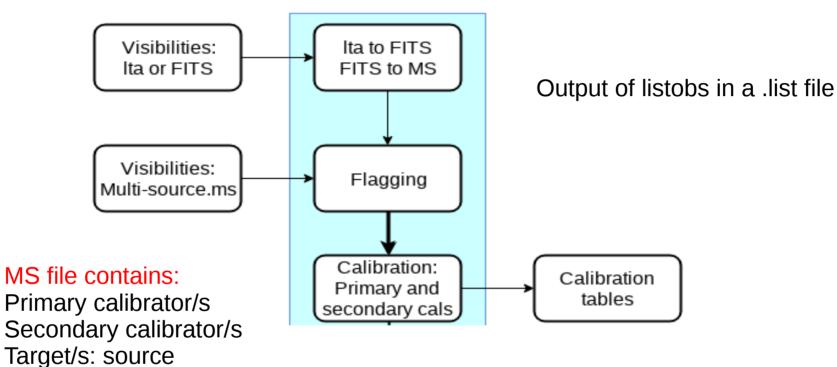
- Initial flagging and calibration
- Further flagging and final calibration

#### Working with calibrated target source data

- flagging
- averaging in frequency
- imaging and self-calibration



# Initial flagging and calibration



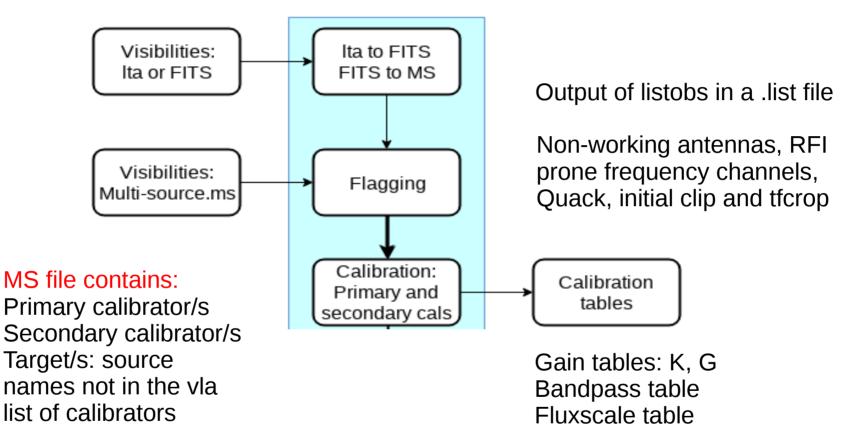
list of calibrators

names not in the vla

**Dependencies**:

If starting from Ita file: listscan, gvfits Calibration part: vla-cals.list

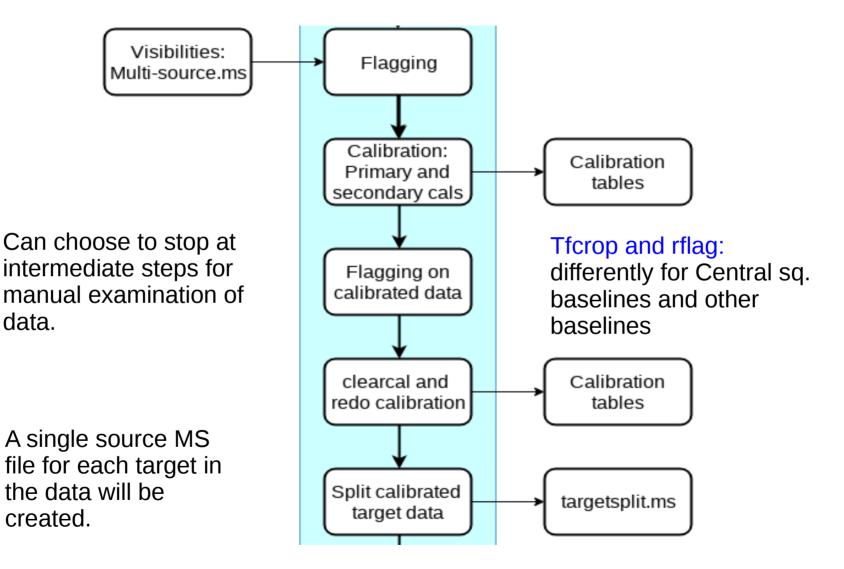
# Initial flagging and calibration



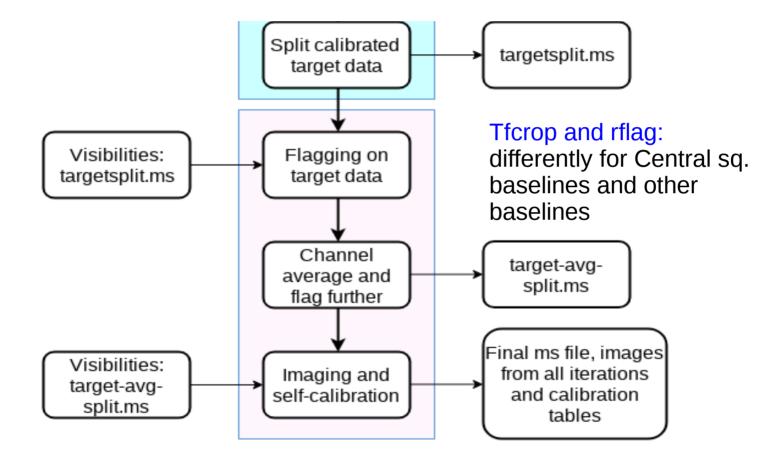
#### **Dependencies**:

If starting from Ita file: listscan, gvfits Calibration part: vla-cals.list

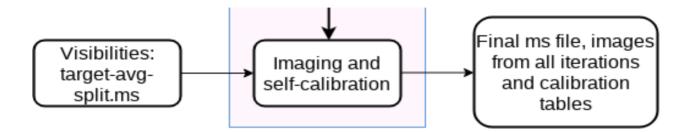
# **Further flagging and calibration**



### Working with calibrated target data



# Working with calibrated target data

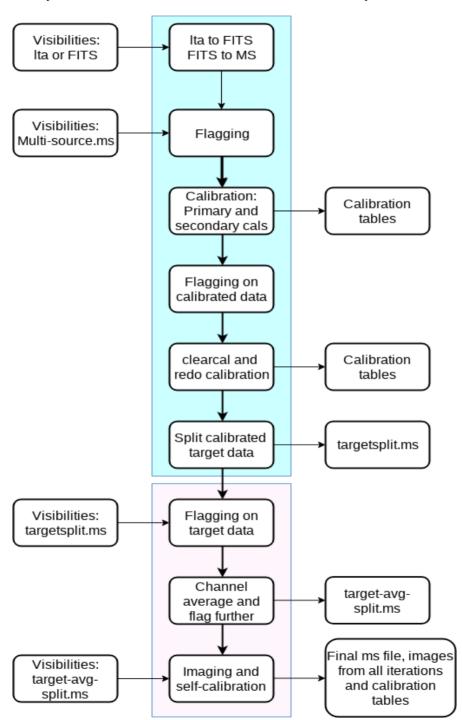


#### Imaging and self-calibration:

- Option to make a dirty image can examine and decide self-calibration strategy.
- Phase-only and amp and phase self-calibration iterations given by the user carried out.
- Flagging on residual data column is carried out in the self-calibration loop.

Input files

Output files



# Structure of the pipeline program

The pipeline is a single python program.

| Initial set-up | These need to be modified according to the type of the data that are to be analysed. |
|----------------|--|
| Inputs         |  |
| Functions      | Modifications only for any special requirements.                                     |
| Applications   |  |

# Initial set-up: initial flagging and calibration

#### 

| fromIta = True                      | # If starting from Ita file set it True.                                     |
|-------------------------------------|--|
| gvbinpath = ['./listscan','./gvfits | ] # set the path to listscan and gvfits if fromIta==True.                    |
| fromraw = True                      | # True if starting from FITS data. Otherwise keep it False.                  |
| fromms = True                       | # True If working with multi-source MS file.                                 |
| findbadants = True                  | # find bad antennas when True  |
| flagbadants= True                   | # find and flag bad antennas when True                                       |
| findbadchans = True                 | # find bad channels within known RFI affected freq ranges when True          |
| flagbadfreq= True                   | # find and flag bad channels within known RFI affected freq ranges when True |
| myflaginit = True                   | # True to flag first channel, quack, initial clips                           |
| doinitcal = True                    | # True to calibrate data   |
| mydoflag = True                     | # True to flag on the calibrated data  |
| redocal = True                      | # True to redo calibration - recommended                                     |
| dosplit = True                      | # True to split calibrated data on target source                             |
| mysplitflag = True                  | # True to flag on the target source  |
| dosplitavg = True                   | # True to average channels   |
| doflagavg = True                    | # True to flag on the channel averaged file                                  |
| makedirty = True                    | # True only if you want to make a dirty image of your target source          |
| doselfcal = True                    | # True if selfcal loop should be run   |
| usetclean = True                    | # True if you want to use tclean (recommended); False will use clean.        |
|                                     |  |

# Initial set-up: input files

fromIta = True # If starting from Ita file set it True.

gvbinpath = ['./listscan','./gvfits'] # set the paths to listscan, gvfits

fromraw = True

# True if starting from FITS data.

fromms = True

#True if working with multi-source MS

# Initial set-up: initial flagging and calibration

from Ita = True# If starting from Ita file set it True. gvbinpath = ['./listscan','./gvfits'] # set the path to listscan and gvfits if from ta==True. fromraw = True # True if starting from FITS data. Otherwise keep it False. fromms = True # True If working with multi-source MS file. # find had antennas when True findbadants = True flagbadants= True # find and flag bad antennas when True findbadchans = True # find bad channels within known RFI affected freg ranges when True # find and flag bad channels within known RFI affected freg ranges when True flagbadfreg= True myflaginit = True # True to flag first channel, quack, initial clips doinitcal = True# True to calibrate data mydoflag = True # True to flag on the calibrated data redocal = True # True to redo calibration - recommended dosplit = True # True to split calibrated data on target source mysplitflag = True # True to flag on the target source dosplitavg = True # True to average channels doflagavg = True # True to flag on the channel averaged file makedirty = True # True only if you want to make a dirty image of your target source doselfcal = True# True if selfcal loop should be run usetclean = True# True if you want to use tclean (recommended); False will use clean.

# Initial set-up: initial flagging and calibration

findbadants = True # find bad antennas when True flagbadants= True # find and flag bad antennas when True findbadchans = True # find bad channels within known RFI affected freq ranges when True flagbadfreq= True # find and flag bad channels within known RFI affected freq ranges when True myflaginit = True # True to flag first channel, quack, initial clips doinitcal = True # True to calibrate data

# Initial set-up: flagging, calibration and split

#### 

from Ita = True# If starting from Ita file set it True. gvbinpath = ['./listscan','./gvfits'] # set the path to listscan and gvfits if from ta==True. fromraw = True # True if starting from FITS data. Otherwise keep it False. fromms = True # True If working with multi-source MS file. # find bad antennas when True findbadants = True flagbadants= True # find and flag bad antennas when True findbadchans = True # find bad channels within known RFI affected freg ranges when True flagbadfreg= True # find and flag bad channels within known RFI affected freg ranges when True myflaginit = True # True to flag first channel, quack, initial clips doinitcal = True# True to calibrate data mydoflag = True # True to flag on the calibrated data redocal = True # True to redo calibration - recommended dosplit = True # True to split calibrated data on target source mysplitflag = True # True to flag on the target source dosplitavg = True # True to average channels doflagavg = True # True to flag on the channel averaged file makedirty = True # True only if you want to make a dirty image of your target source doselfcal = True# True if selfcal loop should be run usetclean = True# True if you want to use tclean (recommended); False will use clean.

# Initial set-up: flagging, calibration and split

- mydoflag = True # True: flags on the calibrated data
- **redocal = True** # True to redo calibration
- dosplit = True

# True to split calibrated target data

# Initial set-up: frequency avg, flagging

#### 

fromIta = True # If starting from Ita file set it True.

gvbinpath = ['./listscan','./gvfits'] # set the path to listscan and gvfits if fromIta==True.

- fromraw = True # True if starting from FITS data. Otherwise keep it False.
- fromms = True # True If working with multi-source MS file.
- findbadants = True # find bad antennas when True
- flagbadants= True # find and flag bad antennas when True
- findbadchans = True # find bad channels within known RFI affected freq ranges when True
- flagbadfreq= True # find and flag bad channels within known RFI affected freq ranges when True
  - # True to flag first channel, quack, initial clips
- doinitcal = True # True to calibrate data

myflaginit = True

mydoflag = True

redocal = True

dosplit = True

mysplitflag = True

dosplitavg = True

doflagavg = True

makedirty = True

doselfcal = True

usetclean = True

- # True to flag on the calibrated data
- # True to redo calibration recommended
  - # True to split calibrated data on target source
  - # True to flag on the target source
- # True to average channels
  - # True to flag on the channel averaged file
  - # True only if you want to make a dirty image of your target source
  - # True if selfcal loop should be run
    - # True if you want to use tclean (recommended); False will use clean.

# Initial set-up: frequency avg, flagging

mysplitflag = True

# True to flag on the target source

dosplitavg = True

# True to average channels

doflagavg = True

# True to flag on the channel

averaged file

# Initial set-up: imaging and self-calibration

makedirty = True

# True only if you want to make a dirty

image of your target source

doselfcal = True

# True if selfcal loop should be run

usetclean = True

# True if you want to use tclean

(recommended); False will use clean.

# Inputs

| Itafile ="  | # Ita file   |  |
|---|--|--|
| rawfile = "   | # TEST.FITS or provide the name of the FITS file if you already have;                              |  |
| myfile1 ="  | # MS file (REQUIRED if starting from multi-source MS file)   |  |
| mysplitfile ="  | # target source file name (split file)   |  |
| mysplitavgfile = "  | # target source file name after averaging; REQUIRED if starting from this file                     |  |
| # Inputs for flagging and calibration   |  |  |
| myquackinterval = 10.   | .0 # time in s to flag at the beginning of a scan and at the end of the scan.                      |  |
| clipfluxcal =[0.0,60.0]   | # in Jy. typically twice the expected flux; only to remove high points                             |  |
| clipphasecal =[0.0,60.  | 0] # in Jy. typically twice the expected flux; only to remove high points                          |  |
| cliptarget =[0.0,30.0]  | # in Jy. typically four times the expected flux; only to remove high points                        |  |
| clipresid=[0.0,10.0]  | # in Jy. 10 times the rms for single channel and single baseline                                   |  |
| myrefant = 'C00'  | # choose a reference antenna - make sure it is one of the working antennas.                        |  |
| uvracal ="  | # Leave it to "; will apply it to all the calibrators in the current version of the pipeline       |  |
| # Inputs for post split averaging of channels   |  |  |
| mywidth2 = 10   | # number of channels to average - choose aptly to avoid bandwidth smearing.                        |  |
| # Inputs for imaging and self-calibration : You will need to change relevant advanced controls if you change the values here. |  |  |
| scaloops = 8  | # Total number of self-cal loops (including both phase-only and amp-ph)                            |  |
| mypcaloops = 4  | # Number of p-only selfcal loops; should be <= scaloops. The remaning loops will and a&p self-cal. |  |
| mythresholds = $0.1$  | # in mJy. Global flux threshold – starting threshold – will change with iteration.                 |  |
| mycell = ['2.0arcsec']  | # Set the cellsize for imaging.  |  |
| myimsize = [12000]  | # Set the size of the image in pixel units. Should cover the primary beam.                         |  |
|   |  |  |

### **More Inputs**

# Further control on imaging and self-calibration

mynterms = 2 # nterms used in tclean; not tested for nterms >2.

mywproj2 = -1 # Number of wprojection planes- leave it to -1 so that it is determined internally in tclean

# Solint used for self-cal: provide solints for each self-cal iteration : edit according to the number of self-cal loops you

# have chosen. Has to be of the same length as nscaloops

mysolint2 = ['8.0min','4.0min','2.0min','1.0min','8.0min','4.0min','2.0min','1.0min']

uvrascal=" # uvrange cutoff used in self-calibration – will use in the task gaincal.

# Structure of the pipeline program

The pipeline is a single python program.

| Initial set-up<br>Inputs | These need to be modified according to the type of the data that are to be analysed. |
|--------------------------|--|
| Functions                | A list of python function calling CASA tool-kit and tasks                            |
| Applications             | Main processing block: runs the functions with the inputs given by the user.         |

# How to run the pipeline ?

- Works in CASA versions 5.0 and above. Likely also in earlier ones but not tested.
- Copy all the files from github to the directory from which you are going to run the pipeline.
- In the python program file (.py), make the "Initial set-up" set the True or False states of the parameters and then provide the "Inputs".

Save and run the file using:

casa -c capture-pipeline-V0.py

Or at the CASA prompt using:

execfile("capture-pipeline-V0.py")

### Images obtained using the pipeline

Pipeline works for:

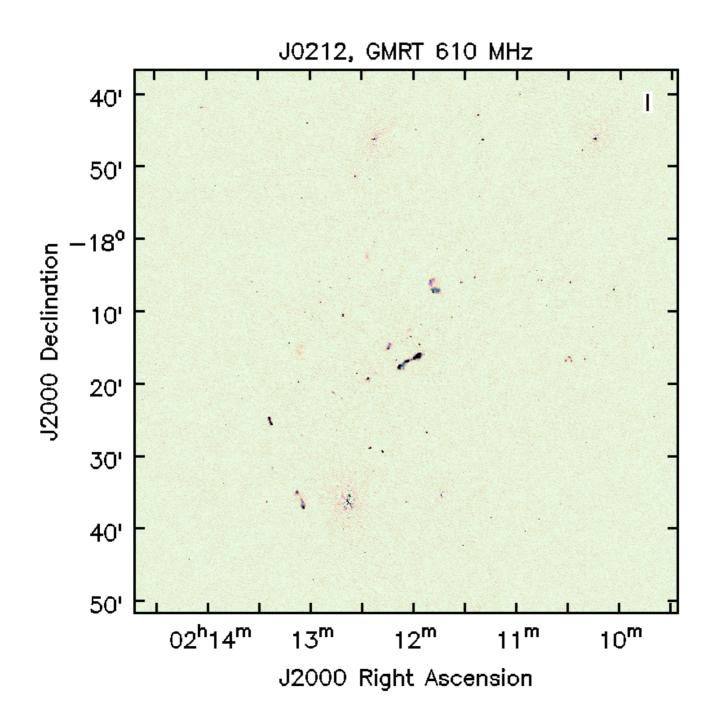
Legacy GMRT : All bands except data taken in dual frequency mode.

Upgraded GMRT: Bands 3, 4 and 5

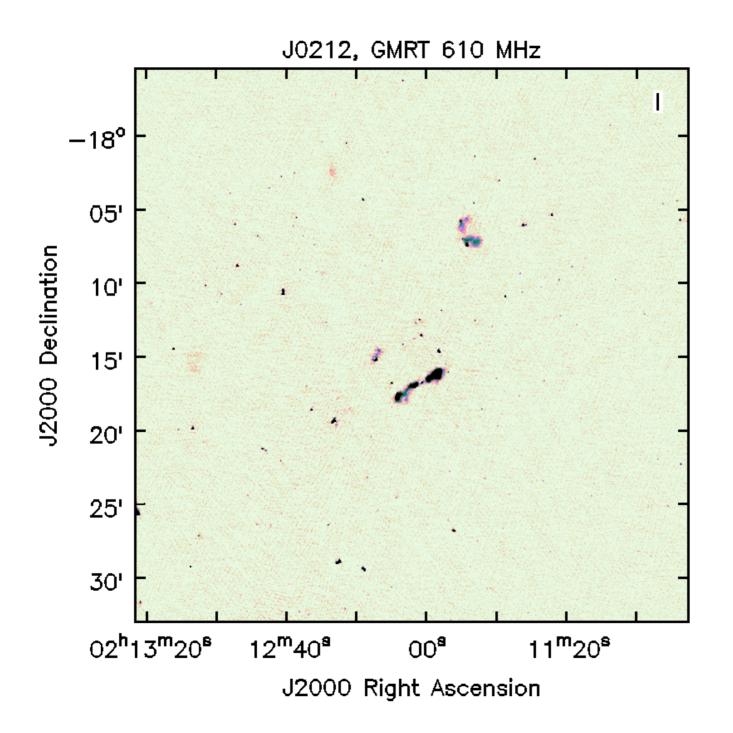
For Band 2: it works after editing the choice of channels to accommodate the notch filter.

Works on sub-array data as well.

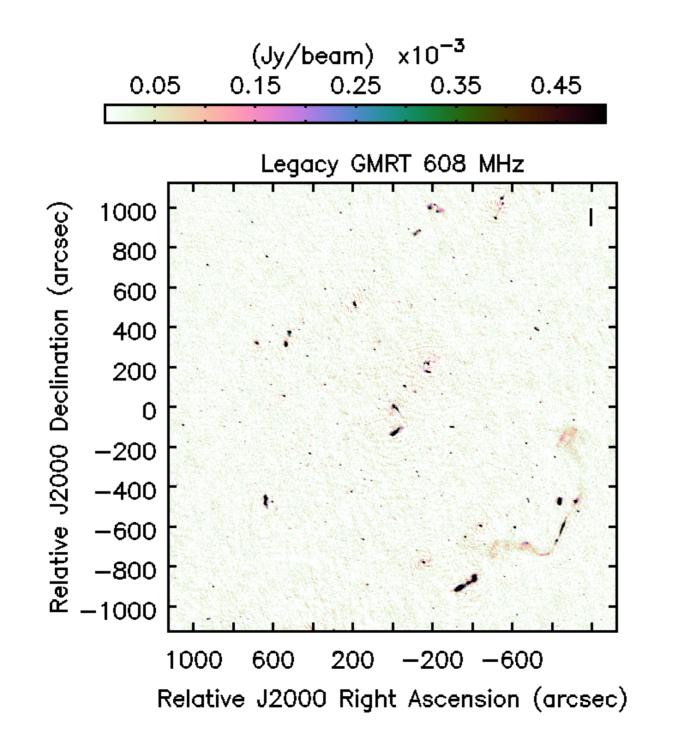
Works when flux and phase calibrator are the same.



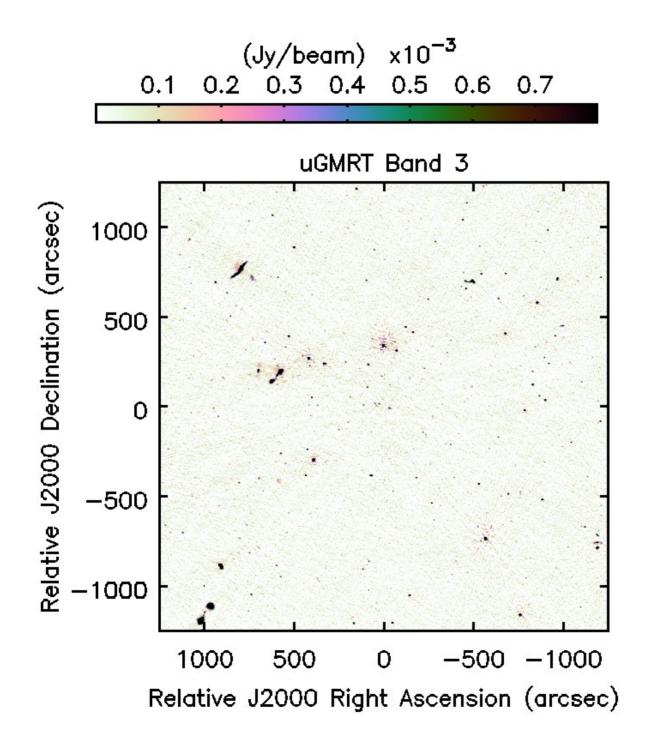
Peak 0.17 Jy/b Rms 0.048 mJy/b



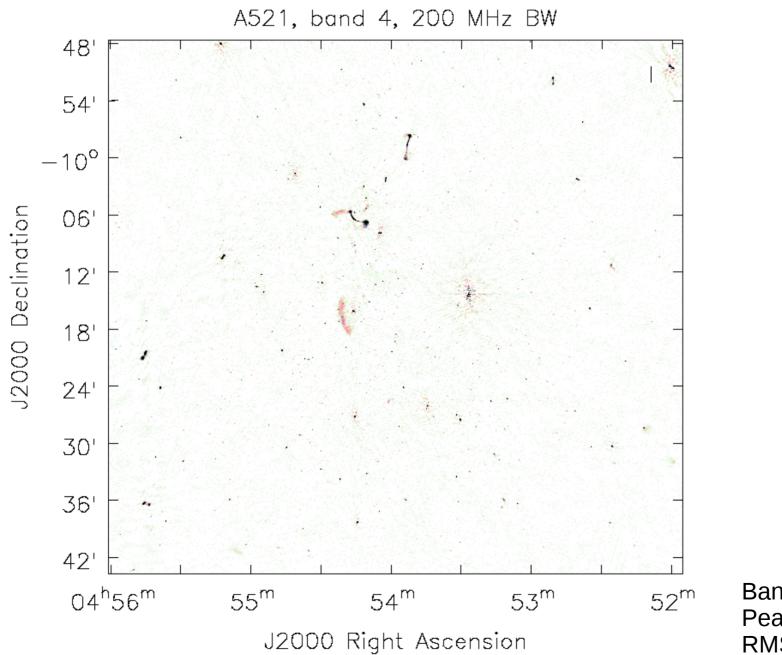
Peak 0.17 Jy/b Rms 0.048 mJy/b

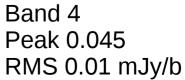


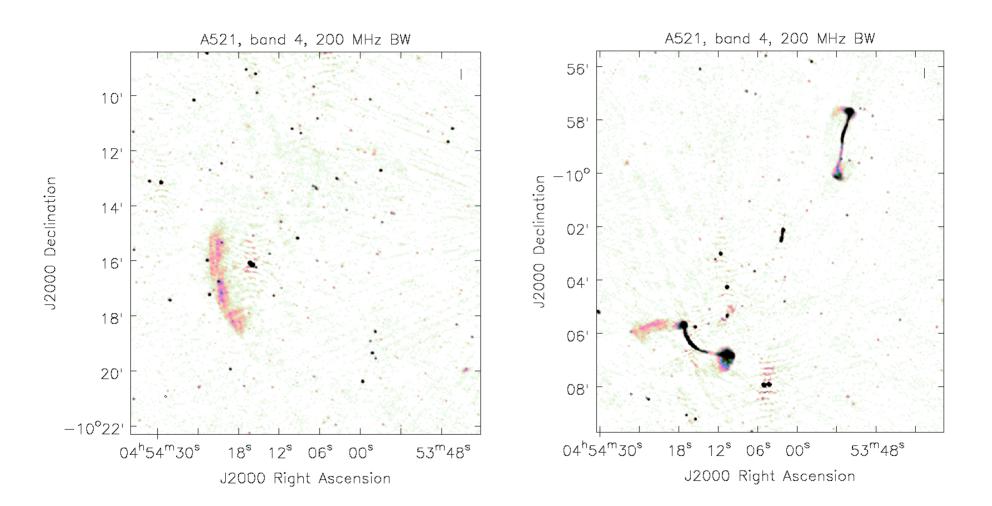
Peak 0.06 Jy/b RMS 0.028 mJy/b



Peak 0.08 Jy/b RMS 0.048 mJy/b







Band 4 Peak 0.045 Jy/b RMS 0.01 mJy/b

# **Pipeline run duration**

Sample numbers:

uGMRT P-band data with on source time of 2.5 hours, cellsize = 1 arcsec, Imsize = 10000 pixels.

Time taken from split file to final self-calibrated image: 32GB RAM, 8 processors, 3.4 GHz : 3.6 days

For a legacy dataset  $\sim$  6 hr duration: End-to-end  $\sim$ 3 days.

For uGMRT dataset ~ 9 hr duration, P-band: End-to-end ~6 days on 128 GB, 24 cores (not an exclusive run)

Memory issues on 32GB machine if imagesize is large ~12000 or so.

### **CAPTURE:** caveats

- Multiple targets: Two step run needed. First step to create calibrated split files for each source and then a separate imaging run for each target is needed.
- If a self-cal run is interrupted, the full imaging run needs to be carried out again.
- For a different choice of parameters in tclean, the python function "mytclean" can be edited appropriately and used (with caution!).
- The pipeline is not tested for nterms > 2.
- Full Stokes data reduction is possible but polarization calibration is not part of it as yet.
- Primary beam correction is a separate task to be run outside the pipeline. https://github.com/ruta-k/uGMRTprimarybeam

# Summary

- Pipelines are essential: reproducibility, automation, data sizes, working on remote servers in general makes life easier !
- Need to be used with caution as at low frequencies each field needs a tailored strategy to obtain the best possible image for the intended science.
- CAPTURE available for uGMRT continuum data reduction can be easily tailored for special needs.
- Writing new CASA tasks, pipelines made easy by Python
- A move towards "open science".