



# **reftools Documentation**

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**STScI**

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Set of tools used to support creation of calibration reference files for Hubble Space Telescope.

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## MAKING IMPHTTAB TABLES

Making IMPHTTAB tables requires a connection to CDBS and that the `PYSYN_CDBS` environmental keyword point to CDBS.

Use `reftools.mkimphttab.create_table()` and provide the name of your output file and the base observation mode from which to build a set observation modes for which photometry keywords will be calculated.

Examples of a base observation mode include "acs,hrc", "wfc3,uvis1", or "cos".

### 1.1 Examples

```
>>> from reftools import mkimphttab
>>> mkimphttab.create_table('acs_wfc1', 'acs,wfc1', 'WFC1',
...                        'Mar 01 2002 00:00:00', clobber=True, verbose=False)
>>> mkimphttab.create_table('acs_sbc', 'acs,sbc', 'SBC',
...                        'Mar 01 2002 00:00:00', clobber=True)
```

```
reftools.mkimphttab.create_table(output, basemode, detector, useafter, tmgtab=None, tm-
                                ctab=None, tmttab=None, mode_list=[], nmodes=None,
                                clobber=False, verbose=True)
```

Create an IMPHTTAB reference file for a specified base configuration, basemode.

#### Parameters

**output** : str

Output IMPHTTAB filename. (`_imp.fits` will be appended if only prefix is given.)

**basemode** : str

Base observation mode for which to generate IMPHTTAB (e.g., `acs,hrc`). This is ignored if `mode_list` is given.

**detector** : str

Detector name.

**useafter** : str

Useafter date in the format of `MMM DD YYYY HH:MM:SS`.

**tmgtab, tmctab, tmttab** : str, optional

Graph (TMG), component (TMC), and thermal component (TMT) tables to use. If `None`, the most recent version in CDBS is used.

**mode\_list** : list of str, optional

A list of observation modes which should be used to make an IMPHTTAB reference file. If given, `basemode` is ignored.

**nmodes** : int, optional

Set to limit the number of modes to calculate. This is for testing only, otherwise set to `None`.

**clobber** : bool, optional

Overwrite existing IMPHTTAB?

**verbose** : bool, optional

Display extra information.

`reftools.mkimphttab.create_table_from_table` (*output, useafter, imphttab, \*\*kwargs*)

Use a previously created IMPHTTAB reference file to generate a new IMPHTTAB reference file.

**Parameters**

**output, useafter** : str

See `create_table()`.

**imphttab** : str

File name of `_imp.fits` IMPHTTAB table from which to take observation modes.

**kwargs** : dict

Keywords accepted by `create_table()`, except `mode_list`.

`reftools.mkimphttab.create_nicmos_table` (*output, detector, useafter, pht\_table, \*\*kwargs*)

Use a NICMOS `_pht.fits` table to generate an IMPHTTAB table for observation modes listed in the given table.

**Parameters**

**output, detector, useafter** : str

See `create_table()`.

**pht\_table** : str

File name of `_pht.fits` table from which to take observation modes.

**kwargs** : dict

Keywords accepted by `create_table()`, except `mode_list`.

`reftools.mkimphttab.compute_values` (*obsmode, component\_dict*)

Compute the 3 basic photometric values needed for a given observation mode string using `pysynphot`.

Values calculated:

- PHOTFLAM - Unit response in FLAM.
- PHOTPLAM - Pivot wavelength.
- PHOTBW - Bandwidth.

**Parameters**

**obsmode** : str

Observation mode for which to calculate values.

**component\_dict** : dict

A dictionary in which to cache opened component objects. May be empty.



**Returns****valdict** : dict

Dictionary with photometry keywords as keys.

`reftools.mkimphttab.compute_synphot_values(obsmode)`Calculate the same values as `compute_values()` but using IRAF SYNPHOT.

---

**Note:** This is replaced by `compute_values()` but kept for debugging.

---

**Parameters****obsmode** : str

Observation mode for which to calculate values.

**Returns****valdict** : dict

Dictionary with photometry keywords as keys.



## COMPARE IMPHTTAB TABLES

Tools for comparing two IMPHTTAB tables from the same instrument and detector.

`class reftools.imphtcomp.ImphttabComp` (*tab1*, *tab2*)

Class for comparing two IMPHTTAB tables from the same instrument and detector.

### Parameters

**tab1** : str

Filename of first IMPHTTAB for comparison.

**tab2** : str

Filename of second IMPHTTAB for comparison.

### Attributes

<code>tab1_name</code>	(str) Filename of first IMPHTTAB.
<code>tab2_name</code>	(str) Filename of second IMPHTTAB.
<code>modes</code>	(ndarray) Obsmodes present in both input files.
<code>flams1</code>	(ndarray) PHOTFLAM values from <code>tab1</code> for obsmodes in <code>modes</code> .
<code>plams1</code>	(ndarray) PHOTPLAM values from <code>tab1</code> for obsmodes in <code>modes</code> .
<code>bws1</code>	(ndarray) PHOTBW values from <code>tab1</code> for obsmodes in <code>modes</code> .
<code>flams2</code>	(ndarray) PHOTFLAM values from <code>tab2</code> for obsmodes in <code>modes</code> .
<code>plams2</code>	(ndarray) PHOTPLAM values from <code>tab2</code> for obsmodes in <code>modes</code> .
<code>bws2</code>	(ndarray) PHOTBW values from <code>tab2</code> for obsmodes in <code>modes</code> .
<code>flamdiff</code>	(ndarray) Percent differences between <code>flams1</code> and <code>flams2</code> calculated as $(flams1 - flams2) / flams1$ .
<code>plamdiff</code>	(ndarray) Percent differences between <code>plams1</code> and <code>plams2</code> calculated as $(plams1 - plams2) / plams1$ .
<code>bwdiff</code>	(ndarray) Percent differences between <code>bws1</code> and <code>bws2</code> calculated as $(bws1 - bws2) / bws1$ .

`make_plot` (*outname*='imphttab\_comp.pdf')

Make a plot with histograms of the percent differences between PHOTFLAM, PHOTPLAM, and PHOTBW for the IMPHTTAB tables.

Differences plotted are  $100 * (table1 - table2) / table1$ .

### Parameters

**outname** : str, optional

Filename of output plot, including extension. Defaults to 'imphttab\_comp.pdf'.

`print_diffs` (*orderby*='photflam', *lines*=25)

Print obsmodes and parameters ordered by *orderby* parameter, with the largest absolute differences in that parameter at the top. This is for seeing which obsmodes have the largest difference in the specified parameter. Prints the number of modes given in the *lines* parameter.

Differences shown are calculated as  $100 * (table1 - table2)/table1$ .

**Parameters**

**orderby** : str, optional

The parameter by which to order the printed results, with modes having the largest absolute difference in this parameter printed at the top.

May be one of: 'photflam', 'photplam', or 'photbw'. An `ImphtcompError` is raised if the input does not match one of these.

Defaults to 'photflam'.

**lines** : int, optional

The number of lines to print. Defaults to 25.

**Raises****ImphtcompError**

If `orderby` does not match a valid option.

`reftools.imphtcomp.plot_table_diffs` (*table1*, *table2*, *outname*='imphttab\_comp.pdf')

Make a plot with histograms of the percent differences between PHOTFLAM, PHOTPLAM, and PHOTBW for the IMPHTTAB tables.

Differences plotted are  $100 * (table1 - table2) / table1$ .

**Parameters**

**table1** : str

Filename of first IMPHTTAB for comparison.

**table2** : str

Filename of the second IMPHTTAB for comparison.

**outname** : str, optional

Filename of output plot, including extension. Defaults to 'imphttab\_comp.pdf'.

`reftools.imphtcomp.print_table_diffs` (*table1*, *table2*, *orderby*='photflam', *lines*=25)

Compare two IMPHTTAB tables and print their differences to the terminal. Prints any obsmodes which are in either table but not in both.

Also prints the obsmodes and parameters for the modes which most differ in the parameter given in `orderby`. This is for seeing which obsmodes have the largest percent difference in the specified parameter. Prints the number of modes given in the `lines` parameter.

Differences shown are calculated as  $100 * (table1 - table2)/table1$ .

**Parameters**

**table1** : str

Filename of first IMPHTTAB for comparison.

**table2** : str

Filename of the second IMPHTTAB for comparison.

**orderby** : str, optional

This specifies one of 'photflam', 'photplam', 'photbw', or 'all'. The printed results are ordered according to the absolute difference in the specified parameter, with the mode with the largest absolute difference at the top of the list.

Specifying 'all' will print 3 tables, one ordered by each of the parameters.

Defaults to 'photflam'.

**lines** : int, optional

Number of lines of differences to print. Defaults to 25.

**Raises**

**ImphtcompError**

If `orderby` does not match a valid option.



## COMPARE SYNPHOT AND PYSYNPHOT

Tools for comparing pysynphot and synphot photometry calculations.

**class** `reftools.synpysyncomp.SynPysynComp` (*imphttab*)

Compare synphot and pysynphot values for obsmodes listed in the input IMPHTTAB. Only obsmodes with three or fewer listed components are compared. For example, `acs,wfc1,f555w` would be processed but not `acs,wfc1,f555w,MJD#`.

This can be a long process for large numbers of obsmodes so it is recommended that users use the CSV functionality of this class to save the results to a .csv file and then use the supporting functions `print_synpysyn_diffs` and `plot_synpysyn_diffs` in this module to investigate them.

**Parameters**

**imphttab** : str

Filename of IMPHTTAB from which to take obsmodes for comparison.

**calculate\_diffs** (*verbose=False*)

Calculate diffs for all obsmodes and return.

**Parameters**

**verbose**: bool, optional

If True, print obsmodes as they are processed. Defaults to False.

**Returns**

**res** : dict

Dictionary containing lists of all pysynphot and synphot calculated values and their differences calculated as  $(\text{pysynphot} - \text{synphot})/\text{synphot}$ .

**comp\_synpysyn** (*mode*)

Returns a dictionary of pysynphot and synphot values and their percent differences calculated as  $(\text{pysynphot} - \text{synphot})/\text{synphot}$ .

**Parameters**

**mode** : str

Obsmode string.

**Returns**

**comp** : dict

Dictionary containing calculated values.

**get\_pysyn\_vals** (*mode*)

Get comparison values from pysynphot.

**Parameters**

**mode** : str

Obsmode string.

**Returns**

**ret** : dict

Dictionary containing calculated values.

**get\_syn\_vals** (*mode*)

Get comparison values from synphot.

**Parameters**

**mode** : str

Obsmode string.

**Returns**

**ret** : dict

Dictionary containing calculated values.

**write\_csv** (*outfile*, *verbose=False*)

Write a CSV file containing the pysynphot and synphot values for all obsmodes and their percent differences calculated as  $(\text{pysynphot} - \text{synphot})/\text{synphot}$ .

**Parameters**

**outfile** : str

Name of file to write to.

**verbose** : bool, optional

If True, print obsmodes as they are processed. Defaults to False.

**class** reftools.synpysyncomp.**SynPysynPlot** (*csvfile*)

Make a plot from a CSV file created by `SynPysynComp.write_csv` illustrating differences between synphot and pysynphot calculated products.

The plots show differences between 7 parameters calculated by both synphot and pysynphot. The differences are shown as histograms. The difference plotted is calculated as  $100 * (\text{pysynphot} - \text{synphot}) / \text{synphot}$ .

**Parameters**

**csvfile** : str

Filename of input CSV file containing comparison results. Should be a file created by `SynPysynComp.write_csv`.

**Attributes**

**fig** : `matplotlib.figure.Figure`

Useful for setting/manipulating figure properties such as the title, which otherwise defaults to `csvfile`.

**save\_plot** (*outname='synpysyn\_comp.pdf'*)

Save plots to a file.

**Parameters**

**outname** : str, optional

Name of file to save to, including extension. Defaults to 'synpysyn\_comp.pdf'.

`reftools.synpysyncomp.plot_synpysyn_diffs` (*csvfile*, *outname='synpysyn\_comp.pdf'*)

Make and save a plot illustrating differences between parameters calculated by both synphot and pysynphot. Data are taken from a CSV file made by `SynPysynComp.write_csv`.



Differences are shown as histograms. The difference plotted is calculated as  $100 * (\text{pysynphot} - \text{synphot}) / \text{synphot}$ .

#### Parameters

**csvfile** : str

Filename of input CSV file containing comparison results. Should be a file created by `SynPysynComp.write_csv`.

**outname** : str, optional

Name of file to save to, including extension. Defaults to 'synpysyn\_comp.pdf'.

`reftools.synpysyncomp.print_synpysyn_diffs(csvfile, orderby='photflam', lines=25)`

Print synphot/pysynphot comparison results from a CSV file produced by `SynPysynComp.write_csv` to the terminal.

Prints the obsmodes and parameters for the modes which most differ in the parameter given in `orderby`. This is for seeing which obsmodes have the largest percent difference in the specified parameter. Prints the number of modes given in the `lines` parameter.

Only prints data for the parameters PHOTFLAM, PHOTPLAM (PIVOT WV), and PHOTBW (RMSWIDTH). Differences are given as  $100 * (\text{pysynphot} - \text{synphot}) / \text{synphot}$ .

#### Parameters

**csvfile** : str

Filename of input CSV file containing comparison results. Should be a file created by `SynPysynComp.write_csv`.

**orderby** : str, optional

This specifies one of 'photflam', 'photplam', 'photbw', or 'all'. The printed results are ordered according to the absolute difference in the specified parameter, with the mode with the largest absolute difference at the top of the list.

Specifying 'all' will print 3 tables, one ordered by each of the parameters.

Defaults to 'photflam'.

**lines** : int, optional

Number of lines of differences to print. Defaults to 25.

#### Raises

##### **SynPysynCompError**

If `orderby` does not match a valid option.

`reftools.synpysyncomp.read_synpysyn(csvfile)`

Read a CSV file created by `SynPysynComp.write_csv` into numpy arrays.

#### Parameters

**csvfile** : str

Filename of CSV to read. Should be a file created by `SynPysynComp.write_csv`.

#### Returns

**res** : dict

Contains one field for each column of the CSV file with a numpy array in that field.



## CALCULATE PHOTOMETRY KEYWORDS FROM IMPHTTAB

getphotpars includes utilities for calculating the photometry keywords PHOTZPT, PHOTFLAM, PHOTPLAM, and PHOTBW for a given obsmode and IMPHTTAB. The calculations are performed in the same way here as they are in hstcal pipelines.

To calculate a single set of keywords use the function `get_phot_pars`.

If you are calculating for several obsmodes from a single IMPHTTAB file it's best to use the `GetPhotPars` class. For example:

```
get_phot = GetPhotPars(imphttab)
for obs in obsmodes:
    photzpt, photflam, photplam, photbw = get_phot.get_phot_pars(obs)
    ...
get_phot.close()
```

4/2014 MLS Made this more general to accept IMPHTTAB files which have any number of extensions, and where the NAME of the phot keyword is the name of the extension (as already is the practice). This is done to accommodate new WFC3 tables with 5 extensions

Right now, the only values that pysynphot computes are PHOTFLAM, PHOTPLAM and PHOTBW

**exception** `reftools.getphotpars.ImphttabError`

Class for errors associated with the imphttab file.

**class** `reftools.getphotpars.GetPhotPars(imphttab)`

This object can be used to get photometry parameters from a given IMPHTTAB reference file. Initialize with the name of an IMPHTTAB reference file and then call this class or the `get_phot_pars` method with a complete obsmode to get the photometry parameters.

**Example obsmodes are:**

'acs,wfc1,f625w,f660n' 'acs,wfc1,f625w,f814w,MJD#55000.0' 'acs,wfc1,f625w,fr505n#5000.0,MJD#55000.0'

### Parameters

**imphttab** : str

Filename and path of IMPHTTAB reference file.

### Attributes

<code>imphttab_name</code>	(str) Filename and path of IMPHTTAB reference file. Same as input <code>imphttab</code> .
<code>imphttab_fits</code>	( <code>pyfits.HDUList</code> ) Open <code>pyfits.HDUList</code> object from <code>imphttab</code> .

**close()**

Close `imphttab_fits` attribute.

**get\_phot\_pars( obsmode )**

Return the required keywords for the specified obsmode

**Parameters**

**obsmode** : str

obsmode string

**Returns**

results\_dict: dict

dictionary of phot keyword results each key corresponds to one keyword and it's value

`reftools.getphotpars.get_phot_pars(obsmode, imphttab)`

Return PHOTZPT, PHOTFLAM, PHOTPLAM, and PHOTBW and any other keywords outlined in the table for specified obsmode and imphttab.

**Parameters**

**obsmode** : str

Complete obsmode string including any parameterized values.

**Example obsmodes are:**

'acs,wfc1,f625w,f660n'

'acs,wfc1,f625w,f814w,MJD#55000.0'

'acs,wfc1,f625w,fr505n#5000.0,MJD#55000.0'

**imphttab** : str

Path and filename of IMPHTTAB reference file.

**Returns**

results\_dict: dict

dictionary containing the photometric keyword results

## MAKE PCTETAB REFERENCE FILE

The PCTETAB reference file contains data in both its primary header and in several table extensions. The parameters that go into the primary header are single numbers that must be specified in the call to *MakePCTETab*.

The data that goes into the table extensions is kept in text files. The names of these text files are given to *MakePCTETab*, which reads them to populate the table extensions. See the documentation for *MakePCTETab* for more detailed descriptions, argument names, and default values.

### 5.1 Primary Header Parameters

- Number of times the readout is simulated to arrive at the corrected image.
- Number of times the pixels are shifted per readout simulation.
- The read noise, in electrons, of the image. This is technically different for each amp but here we pick a single representative value.
- The default value selecting a model for how read noise is handled before CTE correction.
- Threshold for re-correcting over-subtracted pixels. Sometimes the CTE correction removes too much flux from a trail leaving a large divot.

### 5.2 Table Extensions

Functions for ACS PCTETAB reference file.

#### Authors

Pey Lian Lim, Matt Davis

#### Organization

Space Telescope Science Institute

#### History

- 2010-08-31 PLL created this module.
- 2010-11-09 PLL added RN2\_NIT keyword and updated documentation.
- 2011-04-25 MRD updated for new CTE algorithm parameters
- 2011-07-18 MRD updated to handle time dependence
- 2011-11-29 MRD updated with column-by-column CTE scaling
- 2013-08-13 PLL removed deprecated PyFITS calls and cleaned up codes.

## 5.2.1 Examples

```
>>> from reftools import pctetab
>>> pctetab.MakePCTETab(
...     'pctetab_pcte.fits', 'pctetab_dtde.txt',
...     ['pctetab_chgleak-1.txt', 'pctetab_chgleak-2.txt'],
...     'pctetab_levels.txt', 'pctetab_scaling.txt',
...     'pctetab_column_scaling.txt', history_file='pctetab_history.txt')
```

### exception reftools.pctetab.PCTEFileError

Generic exception for errors in this module.

```
reftools.pctetab.MakePCTETab(out_name, dtde_file, chg_leak_file, levels_file, scale_file, col-
                               umn_file, sim_nit=7, shft_nit=7, read_noise=5.0, noise_model=1,
                               oversub_thresh=-10, useafter='Mar 01 2002 00:00:00', pedi-
                               gree='INFLIGHT 01/03/2002 22/07/2010', creatorName='ACS
                               Team', history_file='', detector='WFC')
```

Make the CTE parameters reference file.

#### Parameters

**out\_name** : str

Name of pte fits file being created. May include path.

**dtde\_file** : str

Path to text file containing dtde data.

The file should have 2 columns with the following format:

```
DTDE  Q
float int
... ..
```

Lines beginning with # are ignored.

**chg\_leak\_file** : str or list of str

Path to text file(s) containing charge leak data. If passed as a string the string may contain wild cards so that multiple files are specified.

The input file should contain 5 columns with following format:

```
NODE LOG_Q_1 LOG_Q_2 LOG_Q_3 LOG_Q_4
int float float float float
... .. ... ..
```

Lines beginning with # are ignored.

**levels\_file** : str

Text file containing charge levels at which to do CTE evaluation.

The input file should have a single column with the following format:

```
LEVELS
int
...
```

Lines beginning with # are ignored.

**scale\_file** : str

Text file containing CTE scaling parameters

The input file should have two columns with the following format:

```
MJD      SCALE
float    float
...      ...
```

Lines beginning with # are ignored.

**column\_file** : str

Text file containing CTE column-by-column scaling.

The input file should have 5 columns with the following format:

```
COLUMN  AMPA    AMPB    AMPC    AMPD
int      float  float   float   float
...      ...    ...     ...     ...
```

Lines beginning with # are ignored.

**sim\_nit** : int, optional

Number of iterations of readout simulation per column.

**shft\_nit** : int, optional

Number of shifts each readout simulation is broken up into. A large number means pixels are shifted a smaller number of rows before the CTE is evaluated again.

**read\_noise** : float

Value for RN\_CLIP keyword in PCTEFILE EXT 0. This is the maximum amplitude of read noise used in the read noise mitigation. Unit is in electrons.

**noise\_model** : {0, 1, 2}

Select the method to be used for readnoise removal:

```
0: no read noise smoothing
1: standard smoothing
2: strong smoothing
```

**oversub\_thresh** : float

Value for SUBTHRSH keyword in PCTEFILE header. CTE corrected pixels taken below this value are re-corrected. Unit is in electrons.

**useafter** : str, optional

Value for USEAFTER keyword. Defaults to 'Mar 01 2002 00:00:00'

**pedigree** : str, optional

Value for PEDIGREE keyword. Defaults to 'INFLIGHT 01/03/2002 22/07/2010'

**creatorName** : str, optional

Name of the person generating this fitsFile. Defaults to 'ACS Team'

**historyFile** : str, optional

ASCII file containing HISTORY lines for EXT 0. Include path. Each row will produce one HISTORY line. Defaults to ''

**detector** : str, optional

Supported detector. Defaults to 'WFC'

### Examples

Saving file `pctetab_pcte.fits` with the command:

```
>>> MakePCTETab(  
...     'pctetab_pcte.fits', 'pctetab_dtdel.txt',  
...     ['pctetab_chgleak-1.txt', 'pctetab_chgleak-2.txt'],  
...     'pctetab_levels.txt', 'pctetab_scaling.txt',  
...     'pctetab_column_scaling.txt', history_file='pctetab_history.txt')
```



## INTERPRET DQ FLAGS

The `reftools.interpretdq` module contains functions for interpreting DQ flags.

### 6.1 Examples

```
>>> from reftools.interpretdq import ImageDQ, DQParser
```

Create a class instance for HST/ACS image using pre-defined DQ definitions. Then, display the associated metadata and translation table:

```
>>> acsdq = ImageDQ.from_fits('j12345678qflt.fits', ext=('DQ', 2))
>>> acsdq.parser.metadata
<Table length=1>
  key      val
  str10    str3
-----
INSTRUMENT ACS
>>> acsdq.parser.tab
<Table length=17>
DQFLAG ...          LONG_DESCRIPTION
uint16 ...          str74
-----
  0 ...                      Good pixel
  1 ...                      Lost during compression
  2 ...                      Replaced by fill value
  4 ... Bad detector pixel or beyond aperture or HRC upper-right defect
  8 ...                      Masked by aperture feature or HRC occulting finger
 16 ...                      Hot pixel
... ..                      ...
 512 ... Bad pixel in reference file (FLAT, polarizer, or dust mote)
1024 ...                      Charge trap
2048 ...                      A-to-D saturated pixel
4096 ... Cosmic ray and detector artifact (AstroDrizzle, CR-SPLIT)
8192 ...                      Cosmic ray (ACSREJ)
16384 ... Manually flagged by user
32768 ...                      Not used
```

Interpret DQ value for a single pixel at IRAF-style coordinate X=1457 and Y=170. Also display the original pixel value for comparison:

```
>>> acsdq.interpret_pixel(1457, 170)
<Table length=3>
DQFLAG SHORT_DESCRIPTION LONG_DESCRIPTION
uint16      str9          str74
```

```

-----
  16          HOT          Hot pixel
  32          CTE          CTE tail
 1024         TRAP         Charge trap
>>> acsdq.data[169, 1456]
1072

```

Interpret DQ values for all pixels. Then, extract mask for interpreted DQ value of 16 (hot pixel) and display pixel values for that mask:

```

>>> acsdq.interpret_all()
Parsing DQ flag(s)...
Done!
Run time: 2.822 s
N_FLAGGED: 550656/8388608 (6.564%)
FLAG=1      : 0 (0.000%)
FLAG=2      : 0 (0.000%)
FLAG=4      : 0 (0.000%)
FLAG=8      : 0 (0.000%)
FLAG=16     : 78901 (0.941%)
FLAG=32     : 78915 (0.941%)
FLAG=64     : 343917 (4.100%)
FLAG=128    : 20373 (0.243%)
FLAG=256    : 15202 (0.181%)
FLAG=512    : 26623 (0.317%)
FLAG=1024   : 5128 (0.061%)
FLAG=2048   : 15117 (0.180%)
FLAG=4096   : 0 (0.000%)
FLAG=8192   : 0 (0.000%)
FLAG=16384  : 0 (0.000%)
FLAG=32768  : 0 (0.000%)
>>> hotmask = acsdq.dq_mask(16)
>>> acsdq.data[hotmask]
array([528, 528, 528, ..., 560, 560, 560], dtype=int16)

```

Create a class instance for HST/WFC3 DQ parser (without image). Then, interpret a given DQ value:

```

>>> wfc3dq = DQParser.from_instrument('WFC3')
>>> wfc3dq.interpret_dqval(16658)
<Table length=4>
DQFLAG SHORT_DESCRIPTION          LONG_DESCRIPTION
uint16  str9                      str69
-----
   2          FILLED                Replaced by fill value
  16          HOT                    Hot pixel
 256        SATURATED                Full-well or A-to-D saturated pixel
16384          CRMAX Pixel has more than max CRs, ghost, or crosstalk

```

**class** reftools.interpretdq.**DQParser** (*definition\_file*)

Class to handle parsing of DQ flags.

#### Definition Table

A “definition table” is an ASCII table that defines each DQ flag and its short and long descriptions. It can have optional comment line(s) for metadata, e.g.:

```
# INSTRUMENT = HSTGENERIC
```

It must have three columns:

- 1.DQFLAG contains the flag value (uint16).

2.SHORT\_DESCRIPTION (string).

3.LONG\_DESCRIPTION (string).

Example file contents:

```
# INSTRUMENT = HSTGENERIC
DQFLAG SHORT_DESCRIPTION LONG_DESCRIPTION
0      "OK"                "Good pixel"
1      "LOST"              "Lost during compression"
...    ...                ...
```

The table format must be readable by `astropy.io.ascii`.

#### Parameters

**definition\_file** : str

ASCII table that defines the DQ flags (see above).

#### Attributes

tab	( <code>astropy.table.Table</code> ) Table object from given definition file.
metadata	( <code>astropy.table.Table</code> ) Table object from file metadata.

**classmethod** `from_instrument` (*instrument*)

Use pre-defined DQ flags for the given instrument (HST only for now). The data files are distributed with this module.

#### Parameters

**instrument** : {'ACS', 'WFC3'}

Currently not all instruments are supported. If the instrument is not found, a generic definition is used.

#### Returns

**cls** : *DQParser*

An instance of this class.

**interpret\_array** (*data*, *verbose=True*)

Interpret DQ values for an array.

**Warning:** If the array is large and has a lot of flagged elements, this can be resource intensive.

#### Parameters

**data** : ndarray

DQ values.

**verbose** : bool

Print info to screen.

#### Returns

**dqs\_by\_flag** : dict

Dictionary mapping each interpreted DQ value to indices of affected array elements.

**interpret\_dqval** (*dqval*)

Interpret DQ values for a single pixel.

#### Parameters

**dqval** : int

DQ value.

**Returns****dqs** : `astropy.table.Table`

Table object containing a list of interpreted DQ values and their meanings.

**class** `reftools.interpretdq.ImageDQ` (*data*, *dqparser=None*)

Class to handle DQ flags in an image.

**Parameters****data** : `ndarray`

DQ data array to interpret.

**dqparser** : `DQParser` or `None`

DQ parser for interpretation. If not given, default is used.

**Raises****ValueError**

Invalid image dimension.

**Attributes**

<code>data</code>	( <code>ndarray</code> ) Same as input.
<code>parser</code>	( <code>DQParser</code> ) DQ parser for interpretation.

**dq\_mask** (*dqval*)

Generate mask for the given interpreted DQ value.

**Parameters****dqval** : `int`

Interpreted DQ value.

**Returns****mask** : `ndarray`Boolean mask where the affected pixels are marked `True`.**Raises****ValueError**

Missing interpreter result or invalid DQ value.

**classmethod** `from_fits` (*filename*, *ext=(u'DQ', 1)*, *inskey=u'INSTRUME'*)

Use data from given FITS image and create DQ parser based on header value.

**Parameters****filename** : `str`

Image filename.

**ext** : `str` or `int` or `tuple`Extension identifier, as accepted by `astropy.io.fits`.**inskey** : `str`Keyword value in `PRIMARY` header that defines the instrument.**Returns****cls** : `ImageDQ`

An instance of this class.

**interpret\_all** (*verbose=True*)

Interpret DQ values for all flagged pixels. Results are stored in `self._dqs_by_flag` internal cache.

**Warning:** If the image is large and has a lot of flagged pixels, this can be resource intensive.

**Parameters**

**verbose** : bool

Print info to screen.

**interpret\_pixel** (*x, y*)

Construct a user-friendly table object with interpreted DQ values for a given pixel location (in IRAF format).

**Parameters**

**x, y** : int

IRAF X and Y pixel coordinates.

**Returns**

**dqs** : `astropy.table.Table`

Table object containing a list of interpreted DQ values and their meanings.

**pixlist** ()

Convert cached results from Python indices to list of IRAF-style (*X*, *Y*) pixel coordinates.

**Returns**

**pixlist\_by\_flag** : dict

Dictionary mapping each interpreted DQ value to a list of IRAF coordinates.



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