# CALSTIS REFERENCE GUIDE



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# **CALSTIS Reference Guide**

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#### 1.0 Introduction

CALSTIS is an IDL procedure that performs the routine reduction of Space Telescope Imaging Spectrograph data. This manual describes how to

- search for data in the data catalog (Section 12)
- read the raw STIS data (Section 2)
- calibrate the data using CALSTIS (Sections 3 through 9)
- read the output data files produced by CALSTIS (Section 10)
- use the wavecals taken with the science data to correct for thermal motion and mode select mechanism non-repeatability (Section 13)

The current version of CALSTIS performs the basic image processing steps (Section 7), and can either perform a point source echelle or first order grating extraction (Section 8) or a first order grating extended source image rectification (Section 9). The current version of CALSTIS does not support echelle extended source extractions or prism spectral extractions.

# 2.0 Raw STIS Data Files

Raw STIS data files are stored in FITS files with names:

```
rootname_raw.fits - uncalibrated images
rootname_spt.fits - engineering data for the raw images
rootname_wav.fits - uncalibrated wavelength calibration lamp images.
rootname_wsp.fits - engineering data for the wavelength calibration images.
rootname_tag.fits - MAMA time-tag data file
rootname_trl.fits - ASCII Table Containing ST ScI Processing History
rootname_lrc.fits - local rate check image
rootname_lsp.fits - engineering data for the local rate check image
```

Further information on the format of these files can be found in the STScI STIS Instrument Handbook and the in ST ScI Data Handbook.

The raw STIS data files are catalogued in the IDL data base STISLOG (Section 12). As received, each observation is assigned an ID number (the entry number in the STISLOG catalog). Data can be referenced when using the STIS IDL software either by the filename or by the ID number.

The routine STIS\_READ is used to read raw STIS data files. Its calling sequence is:

```
STIS_READ, file, header, data, udl, heng where
```

file is the ID number or the raw image filename. header is the output FITS header.

data is the output image array.

udl is binary engineering data and can be typically ignored by the average user.

heng is the FITS header of the engineering (\_spt) file.

For example, to read observation 1000 use:

stis\_read,1000, h, data

or equivalently,

stis\_read,'o3xv03w3m\_raw.fits', h, data

You can then proceed to examine the header and data. For example,

hprint, h tvscl, data ;list header contents ;display data

STIS\_READ has the following optional keyword inputs:

/HIRES to leave MAMA data in the Hi-Res mode. The default is to convert the data to Lo-Res mode.

/NO\_OVERSCAN to disable the CCD overscan bias removal.

READOUT = number of the readout to read. The default is to read the first readout (readout=1)

Multiple readout input data files can be read with STIS\_READ using the keyword parameter *READOUT*. This method of reading the data is fine when only one of the readouts is to be read. If however, you are going to process all the readouts, this method can be very inefficient, particularly if the input raw data is stored in compressed format. When you read each readout, the data is uncompressed into a temporary file, the readout read, the file closed, and the temporary file deleted. A better approach is to uncompress, open the file once, and then read all readouts before closing and deleting the temporary uncompressed file. This can be done using the routines STIS\_OPEN and STIS\_CLOSE. To open the file use:

STIS\_OPEN, '<filename>', fcb, nreads=nreads

where *fcb* (file control block) is an output data structure which will be passed to the other routines. The optional keyword output, *nreads*, gives the number of readouts in the file. Now you can start reading the data:

for i=1,nreads do begin STIS\_READ, fcb, h, d, readout=i

end

when done with the file, close it with:

#### STIS\_CLOSE, fcb

#### 3.0 Using CALSTIS

CALSTIS can be called from the IDL command line or from another IDL procedure with the following calling sequences:

1) Point Source Spectral Extraction

CALSTIS, id, params, h, m, wave, flux, epsf, errf, gross, back, net, pos, image, eps, err

2) Extended Source Spectral Extraction

CALSTIS, id, params, h, m, wave, flux, epsf, errf, sens,x,y, unused, image,eps,err

3) Imaging Mode Spectral Extraction

CALSTIS, id, params, h, unused, unused, image, eps, err

where:

id is the input filename or the ID number in the STISLOG data base (See section 2.0). params is a user specified parameter string used to control the reduction steps (Section 4.0). h is the output FITS header string array.

m is a vector of spectral order numbers for the echelle modes. For the first order grating modes, m has the value of 1.

wave is a vector of wavelengths for first order grating modes. For echelle modes, it contains a two dimensional array giving the wavelengths for each spectral order.

flux is an array of extracted flux values for point source spectra, a rectified image for extended source extractions, or a flat fielded image for imaging mode observations.

epsf is a data quality array for flux. It is the same size as flux and has values as defined in Appendix A.

errf is the propagated statistical error array for flux.

gross is the extracted (point source) gross spectrum in count rates.

back is the extracted (point source) background spectrum in count rates.

net is the net spectrum (gross minus the smoothed background spectrum).

pos is an array of spectral extraction positions for point source extractions. The center of the extraction slit in raw data image coordinates for flux(i,j) is at position x=i, y=pos(i,j).

image is a flat fielded count rate image.

eps is a data quality array for image.

err is the propagated statistical error array for image.

sens is the sensitivity vector for an extended source extraction.

x is the raw image x-positions for the center of each pixel in an extended source rectified image

y is the raw image y-positions for the center of each pixel in an extended source rectified image unused indicates that the output parameter position is not used for this type of extraction.

The only required input/output parameter that must be supplied to CALSTIS is *id*. If CALSTIS is called by:

#### CALSTIS, id

all default processing will be performed and the results will be placed in output FITS files.

CALSTIS also has the following optional IDL keyword input parameters:

**READOUT** specifies which readout is to be processed for files containing multiple readouts. Unless specified, CALSTIS will processing the first readout (READOUT=1). To process the second readout, for example, use:

**QUICK** can be used to get a quick spectral extraction. Specifying QUICK causes the bias image removal, dark rate image subtraction, and flat fielding to be turned off. Its usage is:

CR\_REJECT is used to specify that multiple readouts in a CCD data file are to be combined with cosmic ray removal by the routine STIS\_CR. Use:

AUTOWAVE is used to control the zero point shift in the wavelengths using an associated wavecal observation. See section 13.1 for its usage.

Additional keyword inputs, *HEADER*, *DATA*, *EPSIN*, and *ERRIN*, allow you to explicitly supply the data to be processed instead of having CALSTIS read it from an input file. Their usage is described in section 11.

As CALSTIS executes, it will print a processing history to the screen. This same history is placed into the output FITS header. To turn off the screen printing, set !dump to 0 before executing CALSTIS:

!dump = 0 calstis,id,....

To turn the screen printing back on, set !dump to 1.

# **4.0 CALSTIS Reduction Parameters**

The second parameter in the calling sequence of CALSTIS, PARAMS, is the user specified parameters to CALSTIS. These parameters control how CALSTIS reduces the data, including which reduction steps are performed and which calibration reference files are used. PARAMS can have any of the following forms:

VALUE	
0	- use all default parameters
1	- interactively select the parameter values
'string'	- parameter definition of the form:
	'parname=value,parname=value,'
' <filename>'</filename>	- name of a text file with the contents:
	parname=value
	parname=value

Table 4.1 gives a listing of the parameter names and the default values. Further information on them can be found in the individual reduction sections which follow.

TABLE 4.1 CALSTIS Parameter Selection

Parameter Name	Default Value	Description	Sections
outspec	DEF	Spectral file name (DEF, none, or a filename)	10
outimage	none	Output flat fielded file name (DEF, none, or a filename)	10
outgeom	none	Output geometrically corrected camera mode image file name (DEF, none, or a filename)	7.12
extended	0	Extended Source Extraction (0 - no, 1 - yes)	9
hires	0	MAMA Hires processing (0 - no, 1 - yes)	
minerry	0	Minimum MAMA raw data variance	7.3
nospec	0	Skip Spectral Extraction or extended source image rectification (0 - no, 1 - yes)	7.5
soffset	0	Spatial offset (pixels) from nominal spectral position to center of the search area used to locate a point source spectrum.	8.1
dist1	200 (G140L) 100 (all other modes)	Maximum distance from center of search area (in the spatial direction) to look for a point source spectrum.	8.1

Parameter Name	Default Value	Description	Sections
gwidth	11.0 (MAMA) 7.0 (CCD)	Width of the region defining the point source gross spectrum (i.e. height of the extraction slit)	8.2
bwidth	5.0 (first order) 3.0 (echelle)	Width of the regions defining the point source background spectra (i.e. height of the upper and lower background extraction slits)	8.2
bdist	300	distance to the center of the background extraction slit from the center of the gross spectrum (pixels).	8.2
trace	1	Display tracing of the location of the point source extraction (0 - no, 1-yes)	8.1
woffset	0	Wavelength offset in pixels. Typically determined using STIS_WAVECAL (section 13.0)	8.4
b_median	7 (CCD)	Median filter width used for smoothing the background spectra	8.3
b_mean1	0 (MAMA) 15	Mean filter width used for smoothing the background spectra	8.3
b_mean2	15	Mean filter width used for smoothing the background spectra	8.3
b_order	-1	Order of the polynomial used to fit the background spectra (-1 for no fit).	8.3
rate	1	Perform count rate conversion (0 - no, 1 - yes)	7.7
goomcorr	0	Apply camera mode geometric correction	7.12
geomcorr subback	1	Subtract the background spectrum in point source spectral processing (0 - no, 1-yes)	8.3
blazecorr	1	Echelle blaze correction (0-no, 1-yes)	0.6
helio	1	Heliocentric correction of the wavelengths (0 - no, 1-yes)	8.6
biasfile	DEF	Name of the CCD bias image file	5.0, 7.6
ccdpar	DEF	Name of the CCD parameter table	5.0, 7.3, 7.7
darkfile	DEF	Name of the detector dark image file	5.0, 7.8
hpxtab	DEF	Name of the hot pixel table	5.0, 7.10
bpxtab	DEF	Name of the bad pixel table	5.0, 7.10
hotthresh	0.2	Dark rate threshold for flagging CCD pixel as hot (counts/second)	7.10
warmthresh	0.05	Dark rate threshold for flagging CCD pixel as warm (counts/second)	7.10
hrepair	0	Repair threshold for automatic repair of hot pixels (counts/second)	7.10

Parameter Name	Default Value	Description	Sections
pfltfile	DEF	Pixel to pixel flat field file name	5.0, 7.9
lfltfile	DEF	Low frequency variation flat field file name	5.0, 7.0
dfltfile	DEF	Delta flat field file name	5.0, 7.9
geomfile	DEF	Camera mode distortion file	7.12
exttab	DEF	Spectral extraction position table	5.0, 8.1,
dctab	DEF	Dispersion coefficient table	5.0, 8.4,
iactab	DEF	Incidence angle table for correcting for non-concentric aperture locations	5.0, 8.4
angletab	DEF	Table of longslit angles	8.4
scaletab	DEF	Table of image plate scales	8.4
senstab	DEF	Sensitivity table	5.0, 8.5,
pcttab	DEF	Slit GWIDTH photometric correction table	8.5
apttab	DEF	Aperture throughput table	5.0, 8.5
apdtab	DEF	Aperture description table	5.0, 8.4
esdeltaw	0	Extended source extraction wavelength spacing (0 to use value of esdeltas)	9.1
esdeltas	1	Extended source extraction pixel size in the wavelength direction for the output extended source image (in input pixels); Ignored if esdelaw is non-zero	9.1
esdeltal	1	Extended source extraction pixel size in the spatial direction for the output extended source image (in input pixels)	9.1
esnsout	1,064	Number of samples in the output extended source image	9.1
esnlout	1,064	Number of lines in the output extended source image	9.1
escentw	0	Central wavelength of the output extended source image (default is the wavelength at the center of the input image)	9.1
escentl	511.5	Central line position of the output extended source image	9.1
eslogl	0	Output wavelengths on an equal delta log(wavelength) scale (0-no, 1-yes)	9.1
surfaceb	0	Convert extended source extraction to surface brightness flux units	9.2
esgwidth	600	Extraction region width (extraction slit height) for adjusting extended source flux levels	9,2

Parameter Name	Default Value	Description	Sections
essnwave	0	Wavelength of isolated line for selection of the sensitivity value for extended source extraction flux calibration	9.2
geomrot	0	Camera mode rotation angle	7.12
geomxsiz	0	Number of columns in output geometrically corrected camera mode image (if not specified it is selected by routine depending on rotation and magnification)	7.12
geomysiz	0	Number of rows in output geometrically corrected camera mode image	7.12
geomxc	511.5	X-Pixel location in the input image to be mapped to the center of the output image	7.12
geomyc	511.5	Y-Pixel location in the input image to be mapped to the center of the output image	7.12
geommag	1	Magnification factor for output camera mode image	7.12
geomxmag	0	X- magnification factor for output camera mode image	7.12
geomymag	0	Y- magnification factor for output camera mode image	7.12

CALSTIS adds all of these parameters to the FITS header which is included in the output data files. In the case of reference file names with a value of DEF, CALSTIS replaces the DEF with the actual reference file name selected.

# **5.0 CALSTIS Reference Files**

The following calibration reference files can be specified in the user parameters:

biasfile - CCD bias image

ccdpar - CCD parameter table

darkfile - detector dark rate image

hpxtab - hot pixel table

bpxtab - bad pixel table

pfltfile - pixel to pixel flat field

IfItfile - low frequency flat field

dfltfile - delta flat field

geomfile - camera mode geometric distortion maps

exttab - extraction position table (spatial distortion table for extended source calibrations)

dctab - dispersion coefficient table

iactab - incidence angle correction table

angletab - table of angles for the long slits
scaletab - table of plate scales for each grating mode
senstab - sensitivity table
pcttab - sensitivity correction versus extraction GWIDTH
apttab - point source aperture throughput table
apdtab - aperture description table

The user can specify a *filename*, *DEF*, or *NONE* for the parameter value giving a reference file. A value of *NONE* indicates the processing step should be skipped. A value of *DEF* indicates that the default reference file should be selected for the observation's date and observing configuration. The default reference files can be found in the text file, *calstis\_ref.txt*, tabulated by reference file type, observing configuration, and Modified Julian Date (giving the use after date). The file *calstis\_ref.txt* can be found in the software library in a directory specified by the environmental variable, SCAL. SCAL also shows the location of where the calibration reference files are stored. On the STIS team computer systems at GSFC, *calstis\_ref.txt* can be found in:

#### /stis0/data/cal/calstis\_ref.txt

This file can be printed or examined with your favorite editor.

# **6.0 CALSTIS Processing Steps**

CALSTIS performs the following image processing steps on all observations to generate a flat fielded count rate image (details can be found in section 7):

- Conversion of MAMA data to Lo-Res
- Overscan bias level subtraction (CCD only)
- Statistical error initialization
- Cosmic ray rejection (CCD only for multiple readout observations when the keyword input /CR\_REJECT is set)
- Generation of the Doppler smoothing function for smoothing the dark rate file, the flat fields, and the bad pixel table files (MAMA observations only with Doppler correction enabled)
- Bias image subtraction (CCD only)
- Conversion to count rates
- Subtraction of a dark rate image
- Division by a flat field image
- Bad/hot pixel processing

For camera mode images, the following additional step is available:

Correction for geometric distortion (Section 7.12)

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For a point source spectral extraction, the following additional steps are performed (details can be found in section 8).

- point source spectral location
- gross and background spectral extraction
- background processing and subtraction
- wavelength assignment
- · conversion to absolute flux units
- heliocentric wavelength correction

For extended source extractions (section 9) as indicated by the user parameter, *EXTENDED=1*, the following steps are performed:

- wavelength assignment
- heliocentric wavelength correction
- 2-D image rectification

# 7.0 Image Processing Steps

Routine image processing generates a flat fielded image in units of surface brightness. For the MAMA detector, the units are counts/second/Lo-res pixel. For the CCD, the units are electrons/second/unbinned CCD pixel. These units can easily be converted to other units using available keyword values in the output header. For example, Let H be the output header from CALSTIS and IMAGE be the flat fielded image. To convert MAMA data to counts/Lo-Res Pixel use:

```
IMAGE = IMAGE * sxpar(H, 'EXPTIME')
```

To convert CCD data to counts/original binned pixel size use:

```
binx = sxpar(H,'BINAXIS1')>1
biny = sxpar(H,'BINAXIS2')>1
gain = sxpar(H,'CCDGAIN')
exptime = sxpar(H,'EXPTIME')
IMAGE = IMAGE * binx * biny * exptime / gain
```

The following sections describe the image processing steps performed on all data.

# 7.1 Conversion of MAMA Data to Lo-Res

MAMA data are taken in Hi-Res mode, but typically reduced after conversion to Lo-Res. The conversion is done by binning the raw 2048 x 2048 Hi-Res array to a 1024 x 1024 array. Each Lo-Res pixel contains the total counts (not the average) of 4 Hi-Res pixels.

#### 7.2 Overscan Subtraction

For all CCD observations, CALSTIS subtracts a bias level computed from the serial overscan. The image is first rotated to detector coordinates so that the trailing overscan is at the right side of the image and the parallel overscan is at the top of the image. A bias level is computed for each row of the image by taking the median data value in the trailing serial overscan. The following table gives the pixel range used for the various binning factors (keyword BINAXIS1) in the serial direction (Note: the first pixel in the row is pixel 0):

Serial Bin Factor	Overscan pixels
0	1043 to1061
1	1043 to 1053
2	522 to 531
4	261 to 270

A serial bin factor of 0 indicates no binning in either direction. The median value of all overscan values (one value per image row) is computed. Values more than twice the median or less than half of the median are ignored. A linear fit versus row number is made to the remaining values and evaluated for each row of the image. The fitted overscan value for each row is then subtracted from all pixels within the row. The overscan regions are trimmed from the image and the image is then rotated back to the original data coordinates.

For full frame images (including binned ones), an overscan plane is created using both the serial and parallel overscan regions. The table below lists the pixel ranges used in the parallel direction:

Parallel Bin Factor	Overscan pixels
0	1024 to 1043
1	1024 to 1033
2	512 to 521
4	256 to 265

As with the serial overscan, a linear fit versus column number is made to the un-rejected overscan pixels. The overscan plane is computed using the serial overscan vector as the intercept for the last column in the plane. The parallel overscan vector is used to define the slope of each row. The overscan plane is subtracted from the image.

#### 7.3 Statistical Error Initialization

CALSTIS assigns an error to each raw data value using the formula

$$Error = \sqrt{\frac{D}{Gain} + (\frac{\sigma_r}{Gain})^2}$$

where,

D is the data value clipped to be greater than or equal to MINERRV.

Gain is the CCD gain factor

 $\sigma_r$  is the CCD readout noise in electrons per second.

MINERRV is the minimum Poisson variance (default = 0)

The values of *Gain* and  $\sigma_r$  are extracted from the CCD parameter table controlled by user parameter CCDPAR. For MAMA observations, Gain = 1 and  $\sigma_r = 0$ .

The statistical initialization step also flags saturated CCD pixels with a data quality (epsilon) of 190. The saturation levels for the various CCD amplifier and gain settings are stored in the CCD parameter table, CCDPAR.

#### 7.4 Cosmic Ray Removal

If the keyword parameter /CR\_REJECT is supplied in the calling sequence, CALSTIS will combine the multiple readouts of a CCD observation with cosmic ray removal using the routines STIS\_CR and CR\_REJECT. STIS\_CR formats the raw images into a data cube and a statistical error cube. These are passed to CR\_REJECT which uses an iterative cosmic ray rejection algorithm.

# 7.5 Doppler Smoothing Function

When on-board Doppler compensation is used, each output data value can be accumulated from data taken at multiple detector pixel locations. In this case, it is necessary to "smooth" the input reference files.

- 1) The observation is divided into 1 second intervals from observation start to observation end.
- 2) The Doppler shift (in Hi-Res pixels) is computed for each 1 second interval by:

$$Shift = Doppmag \times \sin(\frac{2\pi T}{Period})$$

where,

**Doppmag** is the Doppler magnitude in Hi-Res pixels (header keyword DOPPMAG) **Period** is the orbital period (header keyword ORBITPER)

T is the time since the Doppler Zero time (header keyword DOPPZERO)

- 3) A smoothing vector DS of length Doppmag\*2+1 is accumulated by adding 1 to DS(Doppmag+Shift) for the shift computed for each one second interval. This is a histogram of the shifts.
- 4) DS is then normalized by the total number of 1 second intervals.

The Doppler smoothing function, DS, now contains the fractional amount of time that each shift was used and can be convolved with a Hi-Res calibration reference image to generate an appropriate Doppler smoothed reference image.

#### 7.6 Bias Image Subtraction

A "super" bias image is subtracted from CCD observations as specified by the user parameter BIASFILE. The default bias image is selected by the values of the header keywords CCDAMP, CCDGAIN, CCDOFFST, BINAXIS1, and BINAXIS2. If the input image was a combination of readouts (cosmic ray removal), then the bias image is multiplied by the number of images combined (as stored in the header keyword NCOMBINE) before subtraction.

#### 7.7 Conversion to Count Rates

The input image in counts is converted to count rates (counts/second/Lo-Res pixel for the MAMA detectors and electrons/second/pixel for the CCD) by dividing by the integration time, multiplying by the CCD Gain, and dividing by a binning factor. The binning factor has the value of BINAXIS1 times BINAXIS2 for the CCD, 1.0 for MAMA Lo-Res processing, and 0.25 for MAMA Hi-Res processing. The calibrated value of the CCD Gain is read from the CCD parameter table, CCDPAR. Statistical error propagation is done by applying the same operations to the propagated error image.

#### 7.8 Dark Rate Image Subtraction

CALSTIS subtracts a dark rate image selected by the parameter *DARKFILE*. If the dark rate calibration reference image is for a different binning, it is rebinned to the same binning as the observation before subtraction. Normally, the reference files will be unbinned CCD dark rate and Lo-Res MAMA dark rate images. If Doppler compensation is enabled, the MAMA darks are expanded to 2048 by 2048 then smoothed by the Doppler smoothing function (section 7.5). If Lo-Res processing is being performed, the dark is then rebinned to Lo-Res before subtraction.

No changes to the propagated statistical errors are done during dark rate subtraction.

### 7.9 Flat Field Processing

The STIS flat fields consist of three components: a pixel-to-pixel flat field; a low frequency variation flat field; and a delta flat field. These three components are selected based on the values

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of the parameters: **PFLTFILE**, **LFLTFILE**, and **DFLTFILE**. CALSTIS combines the three components, optionally smoothes by the Doppler smoothing function, rebins the flats to the binning of the observation being processed, and divides the rebinned flat into the data and the associated statistical error array.

### 7.10 Bad/Hot Pixel Processing

The parameters **BPXTAB** and **HPXTAB** specify the bad pixel and hot pixel tables used by CALSTIS. **BPXTAB** contains columns **AXIS1**, **AXIS2**, and **FLAG**, where **AXIS1** and **AXIS2** give the column and row of the bad pixel, and **FLAG** gives the data quality or epsilon value for the pixel. For MAMA observations, **AXIS1** and **AXIS2** are given in Lo-Res pixel coordinates. A given data quality flag is therefore propagated to 4 Hi-Res pixels during Hi-Res processing. CALSTIS flags each image pixel with the maximum **FLAG** value associated with the pixel (in the case of binned data, multiple detector pixels contribute a single data value). If the data quality value of the pixel is already greater than any of the flags in the bad pixel table, it is left as is. If Doppler compensation is enabled, bad pixel data quality flags are propagated to any data value in which the detector element contributed.

The hot pixel table contains columns, AXIS1, AXIS2, and RATE. RATE gives the estimated dark rate for the pixel in counts or electrons per second. Two CALSTIS parameters, **HOTTHRESH** and **WARMTHRESH** are used to determine which pixels are flagged as HOT or WARM. Any pixels with a dark rate greater than the thresholds will be flagged with a quality value of 180 if HOT, and 150 if WARM.

CALSTIS also has the option to repair hot pixels using the values of the neighboring pixels. Care should be taken when interpreting results with the hot pixels repaired. The parameter, HREPAIR, is used to specify hot pixel repair. If HREPAIR = 0.0, no hot pixel repair is performed.

If *HREPAIR* is greater than 0, all pixels with a dark rate in the hot pixel table greater than *HREPAIR* are replaced by the average of its 8 neighbors with data quality values less than 125. Repaired pixels are flagged with a data quality of 125. If all neighbors have data quality values greater than or equal to 125, the pixel value is set to 0 and the data quality is set to 252.

If HREPAIR is set to a value less than zero, The absolute value of HREPAIR is used as the threshold, but only the left- and right-hand neighbors (instead of all 8 neighbors) are averaged when repairing the value. HREPAIR < 0 is useful for repairing point source spectra.

# 7.11 Image Processing Outputs

The calling sequence of CALSTIS with the outputs of the image processing steps shown in bold is:

CALSTIS, id, params, h, m, wave, flux, epsf, errf, gross, back, net, pos, image, eps, err

+ Darke Schiff.

Where: *image* gives the flat fielded count rate image in units of counts or electrons/second/ unbinned pixel; *eps* is the epsilon or data quality image containing coded values for each image pixel as described in appendix A; and *err* is a propagated statistical error image for *image*.

In the case where no spectral processing is performed (either a camera mode image or an image processed with parameter nospec=1), *flux*, *epsf*, and *errf* will have the same values as *image*, *eps*, and *err*.

The outputs of the image processing steps can also be automatically written to a FITS file (see section 10).

#### 7.12 Camera Mode Geometric Distortion

To turn on geometric distortion correction for camera mode images, set *GEOMCORR=1* (the default is *GEOMCORR=0*):

calstis,id,'geomcorr=1',h,m,w,image,eps,err

The output *image* will now be a geometrically corrected flat fielded count rate image. Note: for camera mode reductions, output parameters m and w are not used. The value of GEOMCORR has no effect on the processing of images not taken in a camera mode.

You can also have CALSTIS write the geometrically corrected image to a file by setting the parameter *OUTGEOM* to a file name or to DEF.

calstis,id, 'geomcorr=1,outgeom=myfile.fits'

or

calstis,id,'geomcorr=1,outgeom=def'

If OUTGEOM is set to DEF, the output fits image file will have the name gim\_<id>.fits.

If you want both the flat fielded image before and after geometric correction, use:

calst is, id, 'geomcorr=1', h, m, w, gimage, epsg, errg, g, b, net, pos, image, eps, errg, g, b, net, pos, image, epsg, errg, errg,

gimage, epsg, and errg will contain the geometrically corrected image and its associated data quality image and propagated statistical error image. image, eps and err will contain the flat fielded image without geometric correction. m, w, g, b, net, and pos are not used for carnera mode images.

Note, that if *GEOMCORR* was not set to 1, *gimage* and *image* would both contain the same image (the flat fielded image without geometric correction).

Camera mode images can also be calibrated independently of CALSTIS by using the routine CALSTIS\_GEOM. Its calling sequence is:

calstis\_geom, h, image, eps, err, image\_out, eps\_out, err\_out

where h is the image header and *image*, *eps* and *err* are the input image, data quality, and errors. *image\_out*, *eps\_out*, and *err\_out* are the geometrically corrected results. For example:

calstis, id, 'GEOMCORR=1', h, m, w, gimage, geps, gerr

is equivalent to:

calstis,id,0,h,m,w,image,eps,err calstis\_geom,h,image,eps,err,gimage,geps,gerr

To geometrically correct a raw image use:

stis\_read,id,h,image calstis\_geom,h,image,0,0,gimage

During the geometric correction of a camera mode image, you can optionally change plate scale, center pixel and rotation angle using the following parameters. This is useful for matching your data to the orientation and platescale of an existing image without having to perform another interpolation on the output from CALSTIS.

camera mode rotation angle (default = 0.0) geomrot Number of rows in output camera mode image geomxsiz default = 1024 \* sqrt(2) \* geomxmag \* max(abs([cos(geomrot),sin(geomrot)])) Number of columns in output camera mode image geomysiz default = 1024 \* sqrt(2) \* geomymag \*max(abs([cos(geomrot),sin(geomrot)])) X Position of the center of the output image in the input image geomxc (default = 511.5)Y Position of the center of the output image in the input image geomyc (default = 511.5)magnification factor for output image (default = 1) geommag x-direction magnification factor for output image (default = geommag) geomxmag y-direction magnification factor for output image (default = geommag) geomymag

#### NOTES:

- 1. The magnification is applied after the rotation.
- 2. The header astrometry is correctly updated.

- 3. **geomxc** and **geomyc** are the center of a geometrically corrected raw image with no rotation or magnification. These parameters are useful for generating a fractional pixel shift.
- 4. The geometric correction, rotation, and magnification are all applied with only one bilinear interpolation of the image.

For Example, to geometrically correct a flat fielded count rate image with a 0.25 pixel shift in the x-direction:

```
calstis,id,'geomcorr=1,geomxc=511.75',h,m,w,image
```

To geometrically correct with a factor of 2 increased sampling:

```
calstis,id, 'geomcorr=1,geommag=2',h,m,w,image
```

To geometrically correct an image with a 10 degree rotation, 1.01 magnification in the y-direction, and a 1024x1024 output image size:

```
par = 'geomcorr=1,geomrot=10.0,geomymag=1.01,geomxsiz=1024,geomysiz=1024' calstis,id,par,h,m,w,image
```

To geometrically correct with cubic interpolation:

```
calstis,id,'geomcorr=1',h,m,w,image,eps,err,sens,x,y,junk,fimage cimage = interpolate(fimage,x,y,/cubic,missing=0.0)
```

fimage contains the flat fielded image, image contains the geometrically corrected image using bilinear interpolation, and cimage contains the geometrically corrected image using cubic interpolation

## 8.0 Point Source Spectral Extraction

### **8.1 Point Source Spectral Location**

The first step in point source spectral extraction is to locate the center of the spectrum in the spatial direction. The reference file *exttab* contains tracings (spatial distortions) of the spectrum at various image lines (positions) for the first order gratings and tracings for each spectral order at the nominal position for echelle gratings. For the first order spectral modes, line position 512 is used for the nominal position. The default *exttab* is selected by the optical element and central wavelength of the observation. The spectral location is found as follows.

The nominal position is modified by an optional user supplied offset, parameter **soffset**. **Soffset** should be supplied in pixels (Lo-Res pixels for the MAMA even if a Hi-Res reduction is being performed). This offset is added to the nominal spectral location to define the center of the region to search for the spectrum.

Figure 8.1 shows the shift of the nominal position (lower dotted line) to the new position (upper dotted line).

The spectral tracing at the center of the search area (upper dotted line in figure 8.1) is used as a spectral template. The region plus or minus the parameter dist1 about the center of the search area is searched by cross correlation with this template. This region is between the two dark solid lines in Figure 8.1. At each line position in this region, the total counts along the tracing shifted in increments of 1 pixel in the spatial direction are summed. The maximum sum, computed to sub-pixels by quadratic refinement, defines the center of the point source spectrum (solid line at the center of the spectrum in Figure

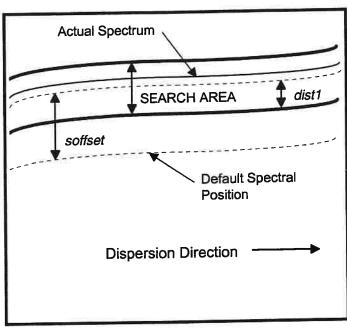


Figure 8.1
Point Source Spectral Location

8.1). For echelle spectra, the correlation is made using all orders simultaneously to obtain a single offset.

Once CALSTIS has completed the spectral location, it will display the image with an overlay so that the user can verify that the spectrum was properly located. This overlay will show the regions to be extracted as the gross and the background spectra. To turn off this display, set the user parameter *trace* to 0. For weak spectra, particularly when residual cosmic rays or hot pixels are present, spectral location may fail, and the trace display is a very useful verification tool.

In the event that the spectral location does fail, the user can recover using a number of strategies.

- 1. Limit the size of the search area by specifying a smaller value for dist1.
- 2. Supply an approximate offset using soffset to allow dist1 to be decreased.
- 3. If all else fails, set dist1 to 0 and specify the final center of the spectrum using soffset.

The following example shows the use of the point source location parameters with a user supplied offset of 75 and a small search area of 10 pixels.

calstis,id,'soffset=75,dist1=10',h,m,wave,flux

# 8.2 Gross and Background Spectral Extraction

Figure 8.2 illustrates the various parameters which control the point source extraction. For each spectral order, three extractions are made illustrated by the regions between the three sets of solid lines in figure 8.2. The gross spectral region has a width in the spatial direction of gwidth and is centered on the spectral order shown by the dotted line determined in section 8.1. The upper background spectral region has a width of bwidth and is centered bdist above the center of the gross spectral region. The lower background region also has a width of bwidth and is centered bdist below the center of the gross spectral region. For echelle data, bdist is not used. Instead, the upper and lower background

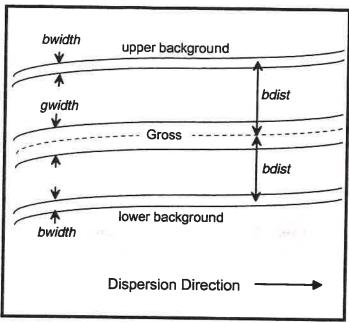


Figure 8.2
Point Source Spectral Extraction Parameters

spectra are centered exactly between the spectral orders. Thus, the upper background for one order is the lower background for the order immediately above it. Gwidth, bwidth, and bdist are

always specified in units of unbinned CCD pixels or Lo-Res MAMA pixels. They are not restricted to be integers.

For each sample position in the dispersion direction, an extraction slit is constructed for the gross and two background spectra. This slit is illustrated in Figure 8.3. The count rates within the slit are summed. The upper and lower ends of the slit may cover only a portion of a pixel. In this case the count rate is weighted by the area of the pixel covered. As the sum of the count rates in the extraction slit is computed, the statistical errors are propagated using the same weights.

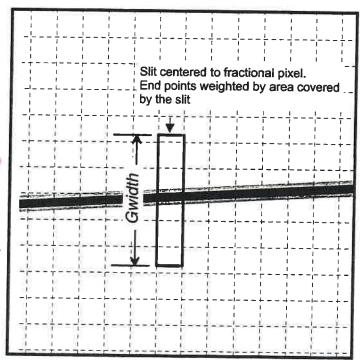


Fig. 8.3 Point Source Extraction Slit

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The output of this extraction is normalized by the binning factor in the spatial direction to give units of counts/second/pixel in the extraction slit.

The data quality value of the extracted data is set to the largest data quality value (the worst condition) of all pixels within the extraction slit. If a portion of the spectral order is outside of the image, the data quality values for that portion of the spectrum are set to 254.

# 8.3 Background Processing and Subtraction

Before the background is subtracted from the gross spectrum, the upper and lower backgrounds are averaged, smoothed, and then normalized to account for differences in the size of gwidth and bwidth. When averaging the upper and lower background, if they do not have the same data quality value, the better of the two values is used.

CALSTIS has a multi-step smoothing process for the background. It performs a median filter, followed by two mean filters, followed by a polynomial fit to the background. Any or all of the smoothing steps can be disabled. The median filter width is specified by the parameter  $b\_median$ . A value of 0 or 1 indicates no median filtering. The widths of the mean filters are controlled by parameters  $b\_mean1$  and  $b\_mean2$ . Again a value of 0 or 1 will disable the filtering. The final optional smoothing is a polynomial fit to the background of each spectral order. This is controlled by the parameter  $b\_order$ . Setting  $b\_order$  to -1 disables the polynomial fitting. If  $b\_order$  is set to 0, the average of all background data points is subtracted from every pixel in the gross spectrum.

The background is normalized to the area of the extraction slit of the gross spectrum by multiplying it by gwidth/bwidth and then subtracting it from the gross spectrum to give the net spectrum.

# 8.4 Wavelength Assignment

CALSTIS assigns wavelengths using the dispersion relation:

$$s = a_0 + a_1 m \lambda + a_2 m^2 \lambda^2 + a_3 m + a_4 \lambda + a_5 m^2 \lambda + a_6 m \lambda^2 + a_7 m^3 \lambda^3 + a_8 m^2 + a_9 \lambda^2$$

where:

s is the pixel location in the dispersion direction, m is the spectral order (1 for the first order gratings),  $\lambda$  is the wavelength, and  $a_0$ ,  $a_1$ ,  $a_2$ , ... are dispersion coefficients

The dispersion coefficients are read from the table specified by user parameter *dctab*. The default dispersion coefficient table is selected based on optical element and central wavelength. Within each table, the dispersion coefficients are tabulated as a function of image position along the slit

(image row or line number). If the desired image line number is not in the table, linear interpolation between the coefficients versus line or slit position is used.

Prior to use, the dispersion coefficients must be corrected for any offset due to the slit non-concentricity. A table of relative slit positions, apdtab, and a table of slit angles for the long slits, angletab, are used to determine the offset,  $\delta$ , between the wavelength calibration dctab slit and the observation's slit in arcsecs.

$$\delta = \delta_0$$
 short slits

$$\delta = \delta_0 + P(y - y_c)(\tan(\theta_{obs}) - \tan(\theta_{dc})) \qquad long slits$$

Where  $\delta_0$  is the slit offset in arcsec at the center of the format from table *apdtab*, P is the platescale in the dispersion direction (arcsec/pixel) from the table *scaletab*, y is the line position in the observation,  $y_c$  is the y-position for the center of the format adjusted for any monthly MSM offsets,  $\theta_{obs}$  is the slit angle for the observations slit, and  $\theta_{dc}$  is the slit angle for the wavecal observation's slit used to construct the dispersion coefficient file.

The table, *iactab*, contains the coefficients  $C_i$  and  $D_i$  which are used to adjust the dispersion coefficients by the relations:

$$a_i = a_i + C_i \delta$$
 for each  $a_i$ 

$$a_1 = a_1 + D_1 \delta + D_2 \delta^2$$

Additionally  $a_0$  is adjusted for MSM non-repeatability and thermal motion by the user supplied offset, parameter **woffset**, given in pixels:

$$a_0 = a_0 + woffset$$

The value of woffset can be determined from a wavelength calibration observations taken with the science data using the routine STIS\_WAVECAL. (Section 13).

#### 8.5 Conversion to Absolute Flux Units

Point source spectra are converted to absolute flux units using the sensitivity table selected by the parameter *senstab*. To turn off conversion to absolute flux units use:

If the target aperture is different than the aperture of the sensitivity table, senstab, CALSTIS corrects the sensitivity for the relative point source aperture throughput between the senstab aperture and your target aperture. The relative point source throughputs are stored in a table

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selected by *apttab*. It also adjusts the sensitivity for your specified GWIDTH using the adjustment tables in the *pcttab* table.

CALSTIS uses linear interpolation between wavelengths in both the sensitivity table and the relative aperture throughput table. It does not extrapolate either table. The flux values of data points outside of the wavelength range of the tables are set to 0 and their data quality values are set to 251.

# 8.6 Heliocentric Wavelength Correction

After the sensitivity correction is completed, CALSTIS will apply a heliocentric correction to the wavelengths provided that the user parameter helio is not set to 0. The Earth's velocity toward the target, v, is computed (in km/sec) and added to the output FITS header in keyword EARTHVEL. Wavelengths are then multiplied by the factor 1+v/c where c is the speed of light. If helio is set to 0, CALSTIS will still populate the value of the keyword EARTHVEL but will leave the wavelengths uncorrected.

When computing sensitivity functions from observations of standard stars, *helio* should be set to zero:

#### calstis,id,'helio=0,senstab=none',h,m,w,f

The wavelengths, ws, of the standard star flux table should then be converted to the velocity frame of the observation by:

$$ws = ws/(1+sxpar(h, 'earthvel')/2.99792458D5)$$

CALSTIS does not convert wavelengths to an AIR reference frame. If you need air wavelengths, run the routine *vactoair* on the output wavelengths of CALSTIS:

vactoair,w

# 8.7 Point Source Spectral Extraction Outputs

The calling sequence of CALSTIS with the outputs of the point source extraction processing shown in bold is:

CALSTIS, id, params, h, m, wave, flux, epsf, errf, gross, back, net, pos, image, eps, err

h is the output FITS header and includes all the keywords present in the input data set, a set of keywords giving the values of all parameters specified by params updated to reflect the selected reference file names, history records describing the processing, and a few additional keyword values added by the processing steps of CALSTIS. This header can be examined with the IDL routine hprint:

#### hprint,h

Individual keyword values extracted for use in post-reduction analysis can be examined with *sxpar*:

value = sxpar(h,'<keyword name>')

or

print,sxpar(h,'<keyword name>')

m is the spectral order numbers. For a first order grating extraction, m will have the scalar value of 1. For echelle observations, m will be a vector of spectral order numbers for the spectral orders extracted.

wave contains the wavelengths. For first order grating mode point source extractions, wave will be a 1024 point vector (2048 point vector for Hi-Res extractions). For echelle mode extractions, wave will be a 2-dimensional array of size 1024 or 2048 by N, where N is the number of spectral orders extracted. wave(\*,i) will have the wavelengths for spectral order m(i). wave is a vector for first order grating extended source extractions. Assuming all processing steps are executed, wave will be heliocentric wavelengths in Angstroms.

flux is the calibrated flux vectors or calibrated image. For first order grating point source extractions, flux will be a vector with the same length as wave. For echelle point source extractions flux will be an array of the same size as wave where flux(\*,i) contains the calibrated spectrum for order m(i). Assuming all processing steps are executed, the output units of flux for point source extractions is ergs/sec/cm²/Angstrom. If conversion to absolute flux is not performed, flux will be identical to net.

epsf is the data quality or epsilon value array for flux. It has the same size as flux and is coded with the values given in Appendix A.

**errf** is the propagated statistical error array for *flux* and has the same size and units as *flux*. It contains the statistical errors (as defined in section 7.3) propagated through the calibration process. It does not include statistical errors introduced from the statistical or systematic errors in calibration reference files.

gross gives the extracted gross spectrum as described in section 8.2. It has the same size and dimension as wave and flux and is in units of counts/second/pixel in the dispersion direction.

**back** is an ns by N+1 array with the upper and lower background vectors, where ns is the number of elements in the spectra and N is the number of spectral orders. back(\*,i) and back(\*,i+1) give the lower and upper background spectra for gross(\*,i). back is in the same units as gross but has **not** been normalized to the slit size of gross (section 8.2).

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**net** is the **gross** minus the smoothed average of the upper and lower background spectra. It has the same size and units as **gross**. **net** is not used for extended source reductions. CALSTIS does not return the smoothed background. It can easily be obtained after running CALSTIS by using:

#### smoothed\_background = gross - net

**pos** is an array of the positions of the center of the extraction slit for the gross spectrum. It has the same size as **gross** and is defined such that the center of the extraction slit for gross(i,j) is at raw data pixel i, pos(i,j). For example, if you want to overplot the positions of the spectral orders used for extraction, use:

tvscl,alog10(image>0.1)
for i=0,N-1 do plots,findgen(1024),pos(\*,i),/dev

# 8.8 Echelle Scattered Light Correction

CALSTIS uses the interorder flux to estimate the dark rate and scattered light in a spectral order. Although this is a good first approximation, it can result in either under or over subtraction of the background depending on the flux distribution of the source. A routine, ECHELLE\_SCAT (which can be run after CALSTIS), constructs a scattered light model, removes it from the image, and re-extracts the spectrum. To run the routine, first run CALSTIS with parameter *outimage* set to DEF, followed by ECHELLE\_SCAT.

calstis, id, 'outimage=def,...', h, m, w, f,... echelle\_scat, id

outimage=def tells CALSTIS to write the flat fielded count rate image to a file with the name im\_<id>.fits. ECHELLE\_SCAT uses this file and the output extracted spectral file to remove compute a model image and a scattered light model image. The scattered light model is subtracted from the flat fielded count rate image (file im\_<id>.fits). A new spectrum is extracted from the difference image using the same extraction parameters specified in the original call to CALSTIS. The extracted spectral output is then written to a FITS binary table named newspec\_<id>.fits. This file has the same format as the original spec<id>.fits written by CALSTIS (see section 10.3).

#### **IMPORTANT NOTE:**

- Echelle\_scat requires a lot memory. You should probably have at least 512 MB of memory on the machine to run it. You can use less but it may significantly slow the process down.
- Execution time is on the order of 10 to 20 minutes.
- ❖ E230H is not currently supported

Briefly, the method used by ECHELLE\_SCAT is as follows:

- 1) Using the original net flux produced by CALSTIS, create an estimated source flux distribution by correcting each spectral order for the echelle blaze and merging the orders. Use this as a first guess of the flux distribution.
- 2) Perform the following iteration on the flux distribution (3 iterations are currently used)
  - Using the current flux distribution, construct a model light image using reference models for the echelle scatter function, telescope PSF, detector halo and grating isotropic scattering, the cross disperser scatter function, and detector ghosts.
  - Extract a net spectrum from the model using the same parameters specified in the original call to CALSTIS. Use this difference to adjust the current flux distribution and repeat the iteration.
  - If it is the last iteration, also construct an object model image (without scattered light) by using the core of the echelle scattered light function (light within 5 pixels of the center) and the 13 x13 pixel core of the combined telescope PSF, detector PSF, isotropic grating scatter, and cross-disperser scatter. Subtract this from the model image to give a model scattered light image.
- 3) Subtract the model scattered light image from the original flat fielded count rate image. Extract the spectrum from this difference and flux calibrate using the same parameters specified in the original call to CALSTIS.

If, in addition to constructing a scattered light corrected spectrum, you would like to see the final model images, use the keyword parameter/model\_file in the call to ECHELLE\_SCAT.

echelle\_scat,id,/model\_file

ECHELLE\_SCAT will then write a FITS file, model\_<id>.fits containing two extensions with names MODEL and SCAT. To read these images from the file use:

fits\_read, 'model\_<id>.fits', model, header, extname='MODEL' fits\_read, 'model\_<id>.fits',scat, header, extname='SCAT'

Where **<id>** is replaced with the observation id number or name.

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Source Extr.

9.0 Extended Source Spectral Extraction

CALSTIS extended source spectral extractions are done by including extended=1 in the parameters:

calstis,id,'extended=1...', h, m, wave, flux, epsf, errf, sens,x,y, unused, image,eps,err

Extended source spectral extraction is used to create a two dimensional map with wavelength in one dimension and position along the slit in the other dimension. Extended source extraction is appropriate for both extended sources observed in a long slit or for a slitless spectra. Fig. 9.1

illustrates a spectrum taken in the long slit mode. Optical/detector distortions have been exaggerated. Pixels of constant wavelength and pixels of constant spatial position do not lie in a straight line due to optical and detector distortions.

#### 9.1 Extended Source Image Rectification.

CALSTIS will generate a rectified image with wavelength running from left to right (horizontally) and spatial position (in the direction along the slit) running vertically. The rectification uses the dispersion coefficients (which vary with position along the slit) and the spectral tracings table, EXTTAB (which gives the spatial distortion at various positions along the slit), to compute the position of each pixel in the output rectified image in raw data coordinates (output parameters x, y). The value for each output pixel is then

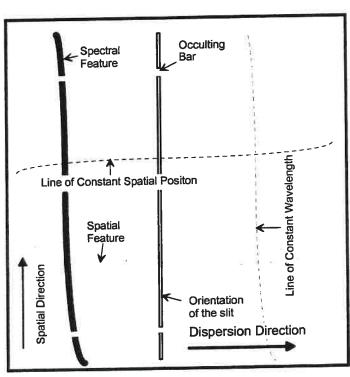


Figure 9.1 Long Slit Extended Source Spectrum

computed using bilinear interpolation in the flat fielded input image. Wavelengths are computed using the same dispersion relation as described in section 8.4 and are converted to the heliocentric system as described in section 8.6.

The default size of the rectified image is 1064 x 1064. This has been increased from 1024 x 1024 so as not to lose any information at the edges of the image during rectification. In the default rectification, the output wavelengths are equally spaced using the average dispersion. The default spacing in the cross-dispersion direction is equal to one pixel in raw image coordinates.

A number of parameters allow the user to control the size of the output rectified image, the spacing between output pixels, and the center position of the output image in wavelength and spatial position along the slit.

The size of the output image is controlled by parameters esnsout and esnlout. For example, to make the output image 1024 x 1024, use:

calstis,id,'extended=1,esnsout=1024,esnlout=1024',...

2 Still preserves the raw scale - justifiens

The spacing between pixels can be controlled by the parameters: esdeltas, esdeltal, esdeltaw, and eslogl. esdeltal and esdeltas give the sample spacing of the output image in the cross-dispersion direction and the dispersion direction, respectively. They are specified in units of raw data pixels. They can be used to change the spacing of the rectified image such that the arcsec/output pixel is the same in both directions or to simply increase or decrease the sampling interval. esdeltaw (specified in Angstroms) can be used to specify the sampling spacing in the output image in the dispersion direction. If esdeltaw is specified, then esdeltas is ignored.

The center of the rectified image can be controlled by the parameters: **escentl** (spatial center in raw image coordinates); **escents** (center in the dispersion direction in raw data coordinates); and **escentw** (central wavelength). If **escentw** is specified, then **escents** is ignored. These parameters, along with the output image size parameters, can be used to extract multiple emission lines from the same or different observations on the same output image scale to allow easy comparison of the different spectral lines.

The output wavelengths of the extended source extraction are on a linear scale (constant delta wavelength between pixels). In some cases, it may be advantageous to place the wavelengths on a constant radial velocity scale (constant log(lamda)). This is done by setting the parameter eslogl to 1. When eslogl is set to 1, the spacing between pixels is such that the delta log<sub>10</sub> of the wavelength is equal to the average difference of the log<sub>10</sub> of the maximum wavelength and log<sub>10</sub> of the minimum wavelength divided by 1024 times esdeltas. Alternatively, you can specify the spacing when eslogl is set to 1 by setting esdeltaw to the log<sub>10</sub> of your desired spacing.

#### 9.2 Extended Source Flux Calibration

If senstab is not set to NONE and a default sensitivity table is available for the optical element, CALSTIS will convert extended source extractions to absolute flux units in the surface brightness units of ergs/cm²/sec/Ang/cross-dispersion detector pixel, where a detector pixel is Lo-Res pixels for the MAMA detector. CALSTIS converts the rectified count rate image (counts/second/pixel) to flux units by dividing by the calibrated point source sensitivity function adjusted to an "infinite" extraction slit height (i.e. GWIDTH = 600) using table apttab. To convert the CALSTIS flux calibrated rectified image to flux per unit surface area for a diffuse continuum source (referred to as specific intensity) divide the output of CALSTIS by the cross-dispersion plate scale (arcsec/pixel) and divide by the width of the entrance slit.

To extract a point source-like spectrum from the rectified flux image (with units of ergs/cm²/sec/Ang), first find the image lines (rows) in the rectified image containing the object. Call them L1 and L2. Integrate the lines:

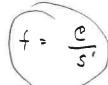
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CALSTIS,id,'extended=1,senstab=def,...', h, m, w, f, eps, err  $ext_flux = total(f(*,L1:L2),2)$ 



To adjust sensitivity calibration for the width of the extraction region compute:

gwidth = (L2-L1+1)/sxpar(h,'edeltal')

edeltal will be 1.0 for the default extraction. It contains the cross-dispersion spacing between rectified image pixels in raw detector pixel units.

Get the sensitivity function for an infinite slit and for your extraction slit:

CALSTIS\_SENS, h, hsens, wsens1, sens1, gwidth=gwidth ;interpolate to observation linterp, wsens1, sens1, w, int\_sens1 CALSTIS\_SENS, h, hsens, wsens2, sens2, gwidth=600 linterp, wsens2, sens2, w, int\_sens2

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Adjust the extracted flux to your extracted slit size:

ext\_flux = ext\_flux\*int\_sens2/int\_sens1 ok as F = net

The units are now ergs/sec/cm<sup>2</sup>/Angstrom.

Alternatively, if you know what extraction slit size you will be using for extracting sources from an extended source rectified image, you can have CALSTIS adjust the flux values ahead of time using the input keyword esgwidth. For example:

CALSTIS,id,'extended=1,senstab=def,esgwidth=15',h,m,w,f  $ext_flux = total(f(*,523:537),2)$ plot,w,ext\_flux

This will give the same flux levels as the point source extraction of the same object:

CALSTIS,id, 'senstab=def',h,m,w,f

An extended source calibration keyword, essnwave (containing a wavelength for sensitivity calibration), can be used to calibrate the entire extended source rectified image with the same sensitivity value (the value at essnwave). Suppose the rectified image is a slitless spectrum of an emission line source. If portions of the source are offset spatially in the dispersion direction, they will appear in the rectified image at different wavelengths when in fact they are approximately the same wavelength. In this case, the sensitivity correction should be made using the actual wavelength of the source. For example, an emission line source at 2500 Angstroms could be calibrated by:

### CALSTIS,id,'extended=1,senstab=def,essnwave=2500.0',h,m,w,f

If you have multiple emission line sources in the same image, it may be necessary to run CALSTIS multiple times, once for each wavelength.

The default output extended source flux units are ergs/sec/cm<sup>2</sup>/Angstrom. Optionally you can select the output to be in surface brightness units by setting parameter surfaceb=1.

CALSTIS,id,'surfaceb=1',h,m,w,f

The output units of f will now be in surface brightness (specific intensity) units (ergs/sec/cm²/Ang/Arcsec²) to the continuum, which depends a second

9.3 Extended Source Extraction Outputs

The calling sequence for an extended source extraction is different than the point source extraction:

calstis,id,'extended=1..., h, m, wave, flux, epsf, errf, sens,x,y, unused, image,eps,err

The output parameters that were used for **gross**, **background**, and **net** spectra are now used for **sens**, **x**, and **y**.

For first order grating extended source extractions, flux will be a 2-D rectified image of size esnsout x esnlout where esnsout and esnlout are parameters with a default value of 1064. Wavelength runs from left to right in this image and spatial position (along the slit) runs from top to bottom. For an extended source extraction, wave will have esnsout values, such that wave(i) is the wavelength for the image column, flux(i,\*). epsf is a data quality image with the same size as flux. The coded values in epsf can be found in Appendix A. errf is the propagated statistical error image. It is important to note that since the values in flux were constructed using bilinear interpolation, the propagated statistical errors in errf can be correlated between neighboring pixels (i.e. have a non-zero covariance). sens contains a sensitivity vector of length esnsout. sens(i) contains the sensitivity used to calibrate column flux(i,\*). x and y are images with the same size as flux. x(i,j) and y(i,j) give the location in the raw image (to a fraction of a pixel) of where the value of flux(i,j) was derived. unused is an unused output parameter. image, eps, and err have the same results as for camera mode images and point source extractions. They contain the flat fielded count rate image along with its associated data quality and error image.

outspec - name to write the rectified flax image.

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# 10.0 CALSTIS Output Data Files

In addition to returning the results in the calling sequence, CALSTIS can also write them into FITS data files. Two parameters control the FITS file output outimage and outspec. outimage tells CALSTIS whether or not to write the flat fielded count rate image to a file. It currently has a default value of NONE (i.e. do not write this file). outspec controls whether or not CALSTIS will write the output of the point source spectral extraction or the extended source rectified image file. The default for outspec is DEF (i.e. write the file using the default file name). outspec and outimage can be set to a file name, NONE, or DEF. If set to a file name (e.g. calstis,id, outimage=myfile.fits',...), CALSTIS will write the results to the specified file, if set to NONE, CALSTIS will not write the file, and if set to DEF, CALSTIS will write the file using the default name. The default filenames are:

# 1068 A

#### **Filetype**

Flat fielded count rate image file Extended source extraction file Point source spectral extraction file

#### **Filename**

im\_<id>.fits espec\_<id>.fits spec\_<id>.fits

where <id> is the observation identification (first input to CALSTIS).

# 10.1 Output Flat Fielded Count Rate Image File

The output flat fielded count rate image is placed into a FITS file with three extensions. The first contains the flat fielded image, the second contains the data quality image with coded values described in Appendix A, and the third contains the propagated statistical error image. This file can be read by IDL using:

```
fits_open,'<filename>',fcb
fits_read,fcb,image,h
fits_read,fcb,eps,h
fits_read,fcb,err,h
fits_close,fcb
```

Or alternatively, the extensions can be randomly accessed by:

```
fits_read,fcb,err,h,exten=3
or fits_read,'<filename>',err,h,exten=3
```

When the file name is supplied to fits\_read, the routine will open and close the file for you.

# 10.2 Output Extended Source Extraction File

The output rectified image from the extended source extraction has the same format for the first three extensions as the flat fielded count rate image except that the images are the rectified

spectral images optionally converted to absolute flux. The file has a fourth extension containing the 1-D wavelength vector.

## 10.3 Extracted Point Source Spectral File

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The point source extractions are written into a FITS binary table. The table has one row per spectral order and has the following columns:

id - the observation ID number or input filename mode\_id - the mode ID (1.1, 1.2, 1.3, ...) cenwave - the commanded central wavelength order - the spectral order number for this row m1 - the first spectral order extracted m2 - the last spectral order extracted line - the central line position (cross-dispersion direction) at the center of the spectrum slitnum - the number of the slit used wavelength - the wavelength vector flux - the flux vector epsf - data quality vector for flux errf - propagated statistical error vector for flux gross - the gross spectrum blower - the lower background spectrum bupper - the upper background spectrum net - the net spectrum (gross - smoothed background) position - the position vector of the center of the extraction slit. position(i) gives the position of the extraction slit for flux(i). coef - the dispersion coefficients used for computing the wavelengths

Some of the columns have been included to allow an extracted point source spectral file to be used as an input dispersion coefficient table (**DCTAB**) or extraction trace table (**EXTTAB**) for subsequent calls to CALSTIS for this observation or for another. For example, to calibrate another observation using the exact same spectral position and dispersion coefficients used to extract observation 9999:

calstis,id,'exttab=spec\_9999.fits',....

The easiest way to read the FITS binary table into IDL is to use the function MRDFITS.

a = mrdfits(filename, 1, header)

This will read the table stored in the supplied filename into an IDL structure, a. For an echelle mode, a will be a structure array. The optional output parameter, header, contains the FITS header. The following examples show how to extract information from the structure. Additional information can be found in the IDL user manuals.

Example 1: Plot the background normalized to the extraction slit of the gross and overplot the smoothed background for a first order grating point source extraction.

```
back = (a.blower + a.bupper)/2*sxpar(header, 'gwidth')/sxpar(header, 'bwidth') \circ sback = a.gross - a.net ;smoothed background plot, a.wavelength, back, psym=1 oplot, a.wavelength, sback
```

Example 2: Put the results of an echelle extraction into 2-D arrays as they would have been returned on the calling sequence to CALSTIS.

```
m = a.order
wave = a.wavelength
flux = a.flux
epsf = a.epfs
errf = a.errf
```

Example 3: Plot the extracted spectrum for echelle spectral order 470.

```
index = where(a.order eq 470)
plot, a(index).wavelength, a(index).flux
```

# 11.0 Supplying Input Data to CALSTIS

In some cases, you may want to perform some pre-processing to the data before calling CALSTIS. To do this, CALSTIS has four optional keyword input variables, which allow you to specify the data to be calibrated. These variables are *data*, *header*, *errin*, and *epsin*. To use them, you are REQUIRED to specify both *data* and a valid *header*. *errin* and *epsin* are optional.

For example, read an observation, do some pre-processing and then call CALSTIS:

```
stis_read,id,hraw,raw ;read raw data
```

; Do some pre-processing (for example, fix some cosmic rays)

calstis, id, 0, h, m, w, f, epsf, errf, data = raw, header = hraw

Even though you are passing the data to CALSTIS, an ID number or input file name must still be supplied. It is used only for identification purposes and for constructing output file names. The input observation is not read by CALSTIS.

Say you want to combine all readouts in a list of multiple CCD observations (stored in vector *list*) with cosmic ray removal. You can use:

```
stis_cr,list,header,data,err,eps
calstis,'aeiou', 0, data=data, header=header, epsin=eps, errin=err
```

Notice that a nonsense observation ID was used. Any ID could have been used to identify the data. Also note that since the cosmic ray rejection algorithm can affect the statistical errors calculation in data points where cosmic rays were rejected, an *errin* is supplied to CALSTIS. *epsin* is supplied because it has information on the number of cosmic ray values that were rejected for each pixel. The previous example could also have been done using:

```
stis_cr,list,outfile='myfile.fits' calstis,'myfile.fits'
```

#### 12.0 STIS Data Catalog, STISLOG

All STIS data are stored in an IDL data base catalog, STISLOG. To search, sort, or display portions of this catalog, a set of IDL routines (beginning with characters DB) can be used. These routines include:

```
dbopen - open a catalog.
dbhelp - print help information.
dbfind - search the catalog.
dbsort - sort the catalog or selected entries from the catalog.
dbprint - print the catalog or selected portions of the catalog.
dbext - extract values from the catalog.
```

Additional documentation on these routines can be found in their help files or in the GHRS software users guide found at the GHRS IDT website:

http://hrssun.gsfc.nasa.gov/swg-v2\_1/online\_swg/node29.html

The following examples will get you started:

 $, \\dbprint, sort\_list, 'entry, targname, opt\_elem, cenwave, texptime, nreads, sclamp', textout = 3$ 

Note that the integer values in *list* and *sort\_list* are the same as the value of item *entry*. This is the value that can be used in the call to CALSTIS as the first input parameter instead of supplying a filename. For example, the following two lines will calibrate the same observation.

```
calstis, '042603060_raw.fits'
and
calstis, 2600
```

Some additional search examples are:

1) Find all G140L wavelength calibration observations

```
list = dbfind(`sclamp=[hitm, line], opt\_elem=G140L')
```

2) Find all G750M observations with a central wavelength between (and including) 7500 and 10000:

```
list = dbfind(`opt\_elem=G750L,7500 < cenwave < 10000')
```

3) Average every FUV-MAMA dark:

```
list = dbfind('detector=fuv_mama,targname=dark')
dark = fltarr(1024,1024)
exptime = 0
.run
- for i=0,n_elements(list)-1 do begin
- stis_read,list(i),h,d
- dark = dark+d
- exptime = exptime + sxpar(h,'exptime')
- end
- end
dark = dark/exptime ;convert to count/sec/pixel
tvscl,alog10(dark>0.0000001)
```

#### 13.0 Wavelength Calibration

The dispersion relations used by CALSTIS were derived from engineering wavecals taken both before and after launch. These dispersion relations must be adjusted to account for thermal motion and MSM non-repeatability. A typical STIS observing sequence will have wavecals associated with the science observations. These can be used to generate a zero point adjustment to the library dispersion coefficients using the AUTOWAVE option to CALSTIS or using the routine STIS\_WOFFSET. If the wavecals are deep enough, they can be used to generate new

dispersion coefficients using the routine STIS\_WAVECAL. Most automatic wavecals are not deep enough for generating new dispersion coefficients and should only be used for zero point adjustments.

#### 13.1 CALSTIS Zero Point Wavelength Adjustment

Use the AUTOWAVE keyword parameter to control zero point wavelength adjustments due to MSM non-repeatibility and thermal/temporal image motion. The following five examples of using the AUTOWAVE keyword will give identical results.

Observation 11555 will be processed using the associated wavecal 11556.

- 1) stis\_woffset,11556,results calstis,11555,0,h,m,wave,flux,autowave=results
- 2) calstis,11555,0,h,m,wave,flux,autowave=11556
- 3) calstis,11555,0,h,m,wave,flux,/autowave
- 4) calstis,11555,0,h,m,wave,flux,autowave=-0.74
- 5) calstis,11555,'woffset=-0.74',h,m,wave,flux

In the first example, the routine STIS\_WOFFSET is used to compute the offset. The results are returned in a structure (results) where

```
    results.mjd - Modified Julian Date(s) at the midpoint of the wavecal observation(s)
    results.offsets - zero-point offset for each wavecal results.history - string history array
```

In the second example, CALSTIS will automatically call STIS\_WOFFSET using wavecal observation 11556.

In the third example, CALSTIS will automatically process the wavecal observation associated with the observation. The correct associated wavecal observation filename is found in the header keyword WAVECAL, populated by in the raw data file by the ST ScI. In this case observation 11556 was the associated WAVECAL.

In the fourth and fifth example, the actual zeropoint offset is passed to CALSTIS where the offset was noted from a prior call to CALSTIS. NOTE: if you use the *autowave* keyword parameter, the value of *WOFFSET* parameter is ignored.

If multiple wavecal observations are found in the WAVECAL file, they are all processed. The zero point shift for the science observation is then computed using linear interpolation between

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the wavecal shifts using the time of the observation midpoint. If the science observation midpoint is outside of the range of midpoints for the wavecal observations, the offset of the closest wavecal is used (i.e. no extrapolation).

Calling STIS\_WOFFSET prior to CALSTIS has an advantage over letting CALSTIS make the call. If you need to vary the reduction parameters to CALSTIS, you do not need to waste time reprocessing the wavecal. Also if STIS\_WOFFSET fails to compute a correct offset, you do not waste time reducing the science observation with an incorrect offset.

In almost all cases, STIS\_WOFFSET will give the correct offset. It is however possible for the cross-correlation of the spectral features in the WAVECAL with the template spectrum to fail and find the maximum correlation at the wrong location. This can happen if the spectrum is offset by more than the maximum search distance or if echelle spectral orders are identified with incorrect spectral order numbsers. To assist in identifying the rare cases where the automatic processing fails, STIS\_WOFFSET generates a plot of the results and also optionally returns the calibrated WAVECAL and template spectrum to the user.

In the first order grating modes, STIS\_WOFFSET will plot six regions of the spectrum centered on the brightest spectral lines with the template spectrum overplotted. The observed spectrum is plotted with a solid line connecting diamonds. The template (smoothed to approximately the resolution of the observed WAVECAL) is plotted with a dashed line connecting asterisks. If the plots not not match, STIS\_WOFFSET may have failed. To further investigate if the correlation has failed, you can examine the other optional outputs of STIS\_WOFFSET. The calling sequence is:

#### stis\_woffset,id,results,h,m,wave,flux,wt,t,image

where *h* is the image FITS header, *m* is a vector of spectral order numbers, *wave* is the calibrated wavelength array, *flux* is the observed spectrum, *wt*, *t* are the wavelengths and flux values of the template and *image* is the rectified cosmic ray filtered image in the first order grating modes for the last readout processed. If the routine was successful, *wt* and *t* should overplot on *wave* and *flux* without any offset. Note, however, line strengths may vary from the template and the observed spectrum. They may have different lamps and different current settings. Keyword parameter *MAXOFF* can be used to vary the maximum search distance for cross-correlating the spectra.

In echelle observations, an offset is computed for each spectral order. The returned offset is the median of these offsets. The offsets for each order are printed and also plotted. If most of the offsets are within a pixel of the median offset, the routine was successful. If not, you might try increasing the keyword parameter, MAXOFF, or verifying that the orders were correctly numbered with the proper spectral order numbers. Spectral format plots are available on the team web page:

http://hires.gsfc.nasa.gov/stis/stispage.html

under "Description" to assist you in determining the order numbers of a spectral calibration lamp observation. If the spectral orders are misidentified, STIS\_WOFFSET keyword parameters soffset and dist1 can be used to correct the situation.

#### 13.2 Generating New Dispersion Coefficient Files

New dispersion coefficient files can be determined using the IDL routine STIS\_WAVECAL. The calling sequence for STIS\_WAVECAL is:

stis\_wavecal, id, params, h, m, wave, flux, wtemplate, ftemplate, toffset=toffset

#### where:

id is the observation ID or filename.

params is a parameter string that allows the user to control various aspects of the reduction.

h is an output FITS header string array.

m is a vector of spectral order numbers for echelle observations or a scalar (value=1) for first order grating observations.

wave is the calibrated wavelength vector (or array in echelle mode observations).

flux is the extracted spectral calibration lamp spectrum.

wtemplate is the wavelengths of the template spectrum used to compute a zero-point offset.

ftemplate is the template spectrum normalized to the same total as flux.

toffset is an optional output keyword parameter giving the offset in pixels computed using cross-correlation with a template spectrum.

The input *params* is specified in the same way as the parameter specification of CALSTIS (section 4.0). Table 13.1 gives a list of available parameters and their current defaults.

Table 13.1 STIS\_WAVECAL Control Parameters

Parameter Name	Default Value	Description			
AIR	0	Data taken in air (0 = no, 1=yes) - for use with prelaunch data only			
	Parame	ters for call to CALSTIS			
DCTAB	DEF	Dispersion coefficient table for initial guess			
EXTTAB	DEF	Extraction trace table			
GWIDTH	15	Width of gross spectral region (extraction slit height)			
SOFFSET 0		Spatial offset in cross dispersion direction (pixels) to shift initial guess of the spectral center			

Parameter Name	Default Value	Description			
DIST1	40	Search distance for spectrum in the cross-dispersion			
2.0		direction (pixels)			
TRACE	1	Display spectral tracing on the screen			
WOFFSET	0.0	Initial offset to the wavelengths (in pixels)			
	Template s	pectrum search parameters			
TEMPLATE	DEF	Name of the template spectrum file			
MAXTOFF	40	Maximum offset in dispersion direction to shift the template.			
MINORDER	200	Minimum counts in the spectral order needed to accept a computed template offset.			
	Spectral	line location parameters			
DCFIT	1	Flag: 1 to fit disp. coef., 0 to not fit			
LINLIB	DEF	Name of the reference spectral line library			
MINLINE	30	Minumum peak counts required in a line.			
MININTE	0	Minumum line library intensity required in a line			
SWIDTH	7	Search width (in pixels)			
NT	3	Cross correlation template width			
MAXDEV	3.0	Maximum allowed difference in predicted and found positions			
	Disper	rsion fitting parameters			
NORDER	2	Order of polynomial fit			
NTRY	3	Number of fit tries (with rejection between tries)			
NSIG	2.5	Rejection threshold			
ISEL1	1	Select mλ term for echelle (0=no, 1=yes)			
ISEL2	1	Select m <sup>2</sup> λ <sup>2</sup> term			
ISEL3	1	Select m term			
ISEL4	1	Select λ term			
ISEL5	1	Select m <sup>2</sup> λ term			
ISEL6	1	Select m $\lambda^2$ term			
ISEL7	0	Select m <sup>3</sup> λ <sup>3</sup> term			
ISEL8	0	Select m <sup>2</sup> term			
ISEL9	0	Select λ <sup>2</sup> term			
IDPLOT	1	Generate line ID plot postscript file (0=no, 1=yes)			

#### The basic steps of STIS\_WAVECAL are:

- 1. Calibrate the input spectrum using CALSTIS with the library or user-specified dispersion relation.
- Integrate a spectral template spectrum to the observations wavelength scale and cross-correlate to determine the spectral offset (MSM non-repeatability and thermal motion)
- 3. Find positions of known spectral lines stored in a line library

Fit (by least squares) the found positions of the spectral lines to create a new dispersion

eps 3 and 4 are optional and can be turned off using the parameter DCFIT=0.

# 3.2.1 Initial Wavelength Solution and Spectral Extraction Using CALSTIS

The first step in STIS\_WAVECAL is to call CALSTIS to extract the spectral calibration lamp The first step in \$115\_WAVECRETS to out the wavelengths using a library dispersion coefficient spectrum and to create an initial guess for the wavelengths using a library dispersion coefficient spectrum and to create an initial guess for the wavelengths using a library dispersion coefficient spectrum and to create an initial guess for the wavelengths using a library dispersion coefficient spectrum and to create an initial guess for the spectrum and to create an initial guess for the spectrum and to create an initial guess for the spectrum and to create an initial guess for the spectrum and to create an initial guess for the spectrum and to create an initial guess for the spectrum and to create an initial guess for the spectrum and to create an initial guess for the spectrum and to create an initial guess for the spectrum and to create an initial guess for the spectrum and to create an initial guess for the spectrum and to create an initial guess for the spectrum and to create an initial guess for the spectrum and to create an initial guess for the spectrum and to create an initial guess for the spectrum and to create an initial guess for the spectrum and to create an initial guess for the spectrum and to create an initial guess for the spectrum and to create an initial guess for the spectrum and table. Input parameters actab, extract, 5 man, 5 ma reduction done by CALSTIS. These parameters are subtraction turned off. Turning off the with the conversion to count rates and the background subtracted as if the subtracted as with the conversion to count rates and the state of the conversion allows long slit wavecal is at the center of the decision for a long slit wavecal is at the decision for a long slit wavecal is at the decision for a long slit wavecal is at the decision for a long slit wavecal is at the decision for a long slit wavecal is at the decision for a long slit wavecal is at the decision for a long slit wavecal is at the decision for a long slit wavecal is at the decision for a long slit wavecal is at the decision for a long slit wavecal is at the decision for a lon background subtraction anows long slit wavecal is at the center of the detector in the The default extraction position can be changed if The default extraction position for a long site was the changed if necessary using the soffset parameter, cross-dispersion direction. This position can be changed if necessary using the soffset parameter. cross-dispersion direction. This position can be officed an output spectral file (spec\_<id>.fits). The call to CALSTIS by STIS\_WAVECAL will generate an output spectral file (spec\_<id>.fits). The call to CALSTIS by STIS\_WAY BOTTOM and the specural me (spec\_<id>.fit PLEASE NOTE: if you use this file, the wavelengths are the initial guess to the wavelengths.

The file does not have the final wavelengths. PLEASE NOTE: it you use this the, the wavelengths does not have the final wavelengths determined using the library dispersion coefficients. The file does not have the final wavelengths determined using the notary dispersion controlled the passed back in the calling sequence. by the routine. The final wavelengths are passed back in the calling sequence.

# 13.2.2 Wavelength Offset Computed Using a Template Spectrum.

The next step in wavelength calibration is to determine MSM non-repeatability and thermal

The template spectrum (selected using parameter template) is read and resampled to the The template spectrum (selected using parameter to the wavelengths of the spectrum as supplied by CALSTIS using trapezoidal integration and then motion offsets. wavelengths of the spectrum as supplied of the observed wavecal. The resampled smoothed using a mean filter to match the slit width of the observed wavecal. The resampled smoothed using a mean filter to material the sit the extracted spectral calibration lamp spectrum template spectrum is then cross-correlated with the extracted spectral calibration lamp spectrum.

This zero point offset is printed to the screen and optional. template spectrum is then cross-contract that are spectrum as the cross-contract that the spectrum is then cross-contract that are spectrum as the cross-contract that the spectrum is then cross-contract that the spectrum is the cross-contract that the cros to determine a zero point offset. This zero point, the server and optionally returned in the wavecal keyword output variable, toffset. The resampled template spectrum is returned in the wavecar keyword output trained, the resampled template spectrum returned in output parameters wtemplate and ftemplate. If the total counts in the extracted returned in output parameters wtemplate and ftemplate returned in output parameters with a value of the parameter with a state of the returned in output parameters wienipute and femiples. It the total counts in the extracted wavecal observation is less than the value of the parameter mincounts, STIS\_WAVECAL will wavecal observation is less than the value of the parameter mincounts, again by increasing midtle. wavecal observation is less than the value of the point, you can try again by increasing gwidth to get abort (insufficient spectral counts). At this point, you can try again by increasing gwidth to get abort (insurficient spectral counts). At this point, job and a gain by increasing gwidth more counts or by decreasing the value of mincounts. However, for very weak wavecal exposures, the accuracy of the computed offset may be questionable.

In the case of echelle observations, a separate offset is computed for each spectral order with at In the case of echelle observations, a separate of the computed as the median value of the offsets least mincounts total counts. The final offsets for each order can be printed by setting. least mincounts total counts. The individual offsets for each order can be printed by setting for all valid spectral orders. The individual offsets for the for all valid spectral orders. The marviage of the individual orders is !dump=2 before executing STIS\_WAVECAL. Examining the offsets for the individual orders is !dump=2 before executing structure have correctly identified the spectral orders with the correct !dump=2 before executing 5115\_WAY and the spectral orders with the correct order a good way to verify that you have correctly identified the spectral order numbers the order numbers the spectral extraction. If you have correctly identified the order numbers and numbers the order numbers and numbers the order numbers and numbers are numbers and numbers are numbers and numbers and numbers are numbers and numbers are numbers and numbers are numbers and numbers are numbers and nu a good way to verify that you have correctly identified the order numbers, the numbers during spectral extraction. If you have correctly identified If the spectral orders will be nearly identical. If the spectral orders will be nearly identical. numbers during spectral extraction. It job have oblighted the order numbers, the computed offsets for the spectral orders will be nearly identical. If the spectral orders were 39

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mis-numbered, the offsets will vary significantly. If the spectral orders were misidentified, adjusted the parameters soffset and dist1 to correct the problem.

The outputs of STIS\_WAVECAL can be used to verify that the cross correlation with the

```
stis_wavecal,1712,'dcfit=0',h,m,w,f,wtemplate,ftemplate,toffset=off
 print, off
plot,w,f & oplot,wtemplate,ftemplate
plot,w,f,xrange=[2700,2800] & oplot,wtemplate,ftemplate,thick=2
```

## 13.2.3 Wavelength Calibration Spectral Line Location

If parameter defit is set to 0, STIS\_WAVECAL will stop after finding the template spectrum offset. Otherwise, the routine will continue and generate a new dispersion fit. Although you should normally use the offset generated by the cross-correlation with the spectral template, the dispersion fit can be used to evaluate that the correct offset was generated. If the template search failed, few spectral lines will be found and the final fit will have a poor RMS.

The first step in generating a new dispersion fit is to locate individual spectral lines with known wavelengths. These known wavelengths are stored in a spectral line library table controlled by user parameter linlib. The position of each library line within the observed spectrum is computed using the wavelengths adjusted by cross-correlation with the template spectrum. Call this position SAPP. The measured line position is then computed by cross-correlation with a square template of width controlled by the user parameter nt. nt should be set to approximately the projected width of the slit in pixels (i.e. the FWHM of the spectral lines in the wavecal). The parameter swidth (in pixels) controls the size of the area to search for the line. The position of maximum correlation (computed to a sub-pixel level with quadratic refinement) is determined. Call this position SFOUND. The line is rejected if any of the following conditions occur:

- 1. The maximum correlation is found at the edge of the search area. (In this case, quadratic
- 2. The absolute value of SFOUND minus SAPP is greater than the user selectable parameter
- 3. The peak counts in the line is less than the user selectable parameter minline.
- 4. The intensity of the line (as stored in the line library table) is less than the user selectable

If a significant number of lines are rejected for reasons 1 or 2, it may be an indication that the If a significant many near indication that cross correlation with the template spectrum failed to determine a valid wavelength offset.

#### 13.2.4 Fitting a New Dispersion Relation

A table of found line positions (as determined in section 13.2.3) is used to determine a new dispersion relation. The coefficients for the dispersion relation are described in section 8.4. The coefficients are determined by a least squares fit of the found line locations and their known spectral order and wavelength. The fitting process is controlled by input parameters *norder*, *ntry*, *nsig*, *isel1*, *isel2*, ..., and *isel9*. The least squares fit is attempted up to *ntry* times with bad data rejection between each try. If the absolute value of the observed minus the fitted position is more than *nsig* times the RMS of the fit, lines are rejected and the fit retried. For first order gratings, *norder* controls the degree of the polynomial to fit. For echelle observations, the value of *norder* is ignored and the values of *isel1* through *isel9* control which terms to fit.

The success of the fit can be evaluated by the final RMS of the fit, the number of lines fit, and a postscript lineID plot of the final fit. STIS\_WAVECAL generates the following files with the results of the fit.

dc\_<id>.fits - dispersion coefficient table (FITS binary table)

wc\_<id>.fits - FITS binary table giving the results for each line fit including the residual of the fit and the reason any line was rejected.

wc\_<id>.prt - a text version of wc\_<id>.fits which can be examined with a text editor or printed.

wc\_<id>.ps - a line ID plot with the library lines identified on the spectrum plotted with the final fitted wavelength scale. This file can be examined with a postscript previewer or printed to a postscript printer. For echelle modes, a separate plot is done for each spectral order (do not print unless it is really necessary)

#### 14. Introduction to Time-Tag Data Processing

The following routines are available for analysis of STIS time-tag data.

TTAG\_INFO - routine to extract information about a time-tag data set

TIME\_TAG - routine to read time-tag data sets

TTAG\_IMAGE - routine to construct an image from the time-tag data

TTAG\_BIN - routine to generate the count rate versus time for time interval

TTAG\_CENTROID - routine to compute centroid of a feature versus time

TTAG\_MOVIE - routine to construct a data cube (movie) for a selected image region of a time-tag dataset

To get a summary of a time-tag dataset use:

#### TTAG\_INFO,id,out

Where *id* is the observation ID (entry number in STISLOG) or the file name of the time\_tag data set (e.g. 'o3xl08bnq\_tag.fits'). The routine will print a summary and also place the information in the structure *OUT*. Refer to the helpfile for TTAG\_INFO for the format of *OUT*.

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WARNING - Do not try to read a very large time-tag data set entirely into memory at once. It may cause your computer to choke. If TTAG\_INFO reports that the data set has more then 10 to 20 million events, refer to "Working with very large time-tag data sets" found in section 14.1.

To read the time-tag data, use the routine:

#### TIME\_TAG,id,h,time,x,y,xdop

where *id* is the observation ID (entry number in STISLOG) or the filename (post-launch time-tag data files have names <rootname>\_tag.fits). *h* is the output FITS header, *time* is a vector of times for each event (in units of seconds since the exposure start), *x* is the vector of x-locations of the events (in Hi-Res coordinates from 0 to 2047), *y* is the y-locations of the events, and *xdop* is the Doppler corrected x positions for echelle spectral observations.

To construct an image from the time-tag data, use:

#### TTAG\_IMAGE,h,x,y,image

where *h*,*x* and *y* are input header, x and y locations (as output from TIME\_TAG). *image* is the output Hi-Res image. You can tell TTAG\_IMAGE to use only events occurring within a specified time interval using optional keyword parameters (*time*, *tmin*, and *tmax*). For example to generate an image from the events from 5 to 6 seconds after the exposure start, use:

#### TTAG\_IMAGE,h,xdop,y, time=time, tmin=5.0, tmax=6.0

where time is the output of TIME\_TAG.

To generate a plot of count rates versus time, use the routine:

#### TTAG\_BIN,time,tout,rate

where *time* is the time for each event (from TIME\_TAG) and *tout* and *rate* are vectors giving count rate (*rate*) versus time (*tout*). The default is to use events anywhere within the image and to integrate rate on 0.01 second intervals. Keyword parameter *interval* can be used to set a different time interval. Keyword *range*=[xmin,xmax,ymin,ymax] can be used to examine events in a subset of the image, and keyword /counts can be used to get output counts/time interval instead of count rates. Keywords *tmin* and *tmax* can be used to limit the time period processed, For example:

### TTAG\_BIN,time, t, c, range=[1500,1510,800,810], interval=0.001, \$ tmin = 50.0, tmax = 52.0, /counts

To find the centroid of a feature versus time, use TTAG\_CENTROID.

TTAG\_CENTROID, time, x, y, tout, xcent, ycent, xrange=[xmin,xmax], \$
yrange = [ymin, ymax]

where *time*, x, and y are the input times and positions of the events, and *tout*, xcent, and ycent are the output times, x-centroid, and y-centroid on one second intervals. xrange and yrange are keyword inputs giving the range of x/y coordinates to centroid. To change the time interval of the centroids use keyword *interval*. To only centroid a portion of the data (in time) use keyword parameters tmin and tmax.

Much of the archived data is compressed. It will significantly slow you down if every time you read an archived time-tag file, it gets uncompressed. To avoid this problem, use the /KEEP keyword parameter in your call to TTAG\_INFO or TIME\_TAG. If the file was compressed, an uncompressed version will be left in your current directory. This version will be used for subsequent calls to TIME\_TAG or TTAG\_INFO.

**EXAMPLE SESSION 1**: Make a movie of a Lyman Alpha for a G140L observation of target JUP-AUR-OVALN and perform an extended source flux calibration for a selected time period of the image.

```
; get information about the observation and keep the uncompressed version ; if the file was compressed ;

id = 16164
    ttag_info,id,out,/keep ;

; read the events ;

time_tag, id, h, time, x, y ;

; construct a Hi-Res image from the events ;

ttag_image, h, x, y, image ;

; use STISLOOK to find a region of interest ;

stislook, image region = [92,491,1020,1419] ; [xmin,xmax,ymin,ymax]
```

```
; plot the total count rate versus time for the region averaged in 10
; second intervals
  ttag_bin, time, tout, rate, x=x, y=y, range=region, interval=10
  plot, tout, rate
; create a data cube (movie) of the region by integrating in 20 second
; intervals with data binned by 4(2 \times 2 \text{ lores pixels})
  ttag_movie, time, x, y, cube, interval=20.0, range=region, binsize=4
; load cube into IDL movie widget which will also let you save it as an
; MPEG file
  s = size(cube) \& nx = s(1) \& ny = s(2) \& nframes = s(3)
  xinteranimate, set=[nx,ny,nframes], /showload
  for i=0,nframes-1 do xinteranimate, frame=i, $
               image = bytscl(cube(*,*,i),min=0,max=5,top=!d.n\_colors)
  xinteranimate,/keep_pixmaps
; create an image and calibrate it for the period of 100 to 300 seconds since
; the observation start.
  ttag_image, h, x, y, image, time=time, tmin=100, tmax = 300
   sxaddpar,h,'exptime',200
   sxaddpar,h,'integ',200
   calstis,id,'extended=1',h,m,wave,flux,eps,err, header=h, data=image
```

#### **EXAMPLE SESSION 2:** Examine Band 2 MAMA dark microbursts.

```
; read time tag data set
;

time_tag, 2657, h, time, x, y
;
; create an image from the time tag data and display
;

ttag_image, h, x, y, image
tvscl, rebin(float(image),512,512) < 1
```

```
; determine count rates in 0.01 second time bins and plot rate versus time
        ttag_bin, time, tout, rate, interval=0.01
        plot, tout, rate
; find the time bin with the peak rate and generate an expanded plot of
; counts versus fine time interval (125 microseconds) for a 0.04 second
; interval containing the peak
       print, max(rate,position)
       t = tout(position)
                                      ;time of maximum count rate
       ttag\_bin, time, t1, c1, tmin = t-0.02, tmax = t+0.02, \$
                       /counts, interval=0.000125
       plot, t1, c1, psym=-4
; It looks like a micro-burst; plot a scatter plot of the x and y coordinates
; of the events for the time period of the microburst.
; (i.e. where we got 4 counts/fine time interval)
       burst = where(c1 ge 4)
       start\_burst = min(t1(burst)) - 0.000125
       stop\_burst = max(t1(burst)) + 0.000125
       burst = where((time ge start_burst) and (time le stop_burst))
       plot, x(burst), y(burst), psym=4
```

#### 14.1 Working with Very Large Time-Tag Data Sets.

Most of the time-tag data sets in the archive have only a few million or fewer number of events. Some, however, will have many millions of events. In these cases, unless your computer has a very large amount of memory, it will be very inefficient to read the entire time-tag data set into IDL. Use TTAG\_INFO to determine if this is the case.

To process the very large data sets, you will need to read the image in parts. The routine TIME\_TAG has two methods to allow you to read parts of the data set. The time-tag data in the raw data file is partitioned into separate groups or file extensions where each extension contains an on-board memory dump. Each group is a good size to read without getting into memory problems.

To read a single group use the *GROUP* keyword parameter when calling TIME\_TAG (*GROUP* runs from 1 to the number of groups). TTAG\_INFO will print the number of groups in the file and also return it in an output data structure. For example, to create an image from a large E140M time-tag data set, use:

```
ttag_info,2222,out,/keep
image = lonarr(2048,2048)
```

Note that in the example, the Doppler corrected x position was used for the echelle time-tag observation.

To create a movie of a selected region of a large time tag observation use:

To create a vector of count rate versus time, use:

An alternate way of selecting which events to use is to specify the number of the first event and the number of events to read. For example:

time\_tag,2222,h,time,x,y,xdop,first=4000000,number=2000000

#### Appendix A **Data Quality Flag Values**

A value of 0 indicates no anomalous condition. Flags are coded using the same basic principal used by the FOS and GHRS ST ScI calibration pipelines. Values from 1 to 99 are minor conditions which are reflected in the propagated statistical errors. Values from 100 to 199 are more serious conditions which are not properly reflected in the propagated statistical error array. Values from 200 to 255 represent data that is totally useless. Each data value is flagged with its 07 Aug 19 - 575,J

ACS	ACC	CONVE	VTION CONVERSIO
Flag	Condition	Notes	
255	Fill data (telemetry dropout)		05 Pet 11
254	Data point outside boundary of the image		em usefue released to
252	Unrepairable hot pixel		1
251	For absolute flux vector (data outside tabulated sensitivity curve)		C 6 D
239	All pixels rejected during cosmic ray removal/ image combination		2 - fill 176,5
240	Dead detector element of bad CCD column flat < 0.01 oz Aug	/ bpx ·	2 - filler
190	Saturated pixel		2.000
195	Cosmic ray detected but not removed	NYI	Sal Transis union
185	Data masked by occulting bar	NYI	176- py yarr
180	Hot pixel (CCD dark rate > 0.2 counts/second) MARA > ~10 ? ? ]		Vericely by
175	Severe flat field blemish	bpx	- alust,
174	Data corrupted by the band 1 repellor wire	bpx	
170	More than 20% uncertainty in non-linearity correction	NYI -	- de Gross 213
160	Slitless spectra contamination by nearby source	NYI	SD +13 0
155	Data near edge of detector		Carrent
150	Warm pixel (CCD dark rate > 0.05 counts/second)		The state of the s
130	More than 5% uncertainty in non-linearity correction	NYI	large
125	Hot pixel repaired by interpolation		Carr.
100	Reed-Solomon Telemetry decoding error	NYI	C + P. C
65	5 or more pixels rejected during CR-SPLIT image combination		
64	4 pixels rejected during CR-SPLIT image combination		-
63	3 pixels rejected during CR-SPLIT image combination		
62	2 pixels rejected during CR-SPLIT image combination		
61	1 pixel rejected during CR-SPLIT image combination		
0	No anomalous condition noted		

NYI - not yet implemented

bpx - flagged if included in the bad pixel table

| STS T - add: Masks >> 8

| See Fastie Fingert Coron. Ars Translationto STSCI ACS ISR 99-08 by Hack 47 & Should Not hart in FF Stuped 1:X

## Appendix B Sample Sessions

#### **B.1 Finding and Reading STIS Data**

; List available catalogs

#### dbhelp

3cr	abell	ans	bruzual
cycle7_abs	cycle7_exp	donas	faust
stisgto	stislog	td1	td1_spec
td1_x	uit_bright	uvgalaxy	virgo
wdwarf	wdwarf_iue	yale_bs	yale_bs_rmks

; open stis catalog and list items stored for each observation

#### dbopen,'stislog' dbhelp,1

ITEM ITEM ENTRY ROOTNAME FILENAME FILETYPE NREADS TARGNAME RA_TARG DEC_TARG	TYPE int*4 char*10 char*20 char*3 int*4 char*18 real*4	DESCRIPTION Entry or Record Number rootname of the observation set  file type RAW, TAB, or WAV number of image readouts in the file proposer's target name right ascension of the target (deg) (J2000) declination of the target (deg) (J2000)
DEC_TARG	real*4	declination of the target (deg) (J2000)
PROPOSID	int*4	PEP proposal identifier

; Find observations from proposal 7123

#### list = dbfind('proposid=7123')

12 entries found in STISLOG

stislook,d

```
; print selected information about the observations
```

## dbprint, list, \$ 'entry,filename,nreads,targname,mode\_id,cenwave,aperture,sclamp'

STIS observation log

ENTRY	FILENAME	n read	TARGNAME	mode _id	cen wave	APERTURE	sc lamp
933	 o3x107a3m_raw.fits	 3	LMC-SN1987A-OFF	3 6		 F28X50LP	NONE
	o3x107010_raw.fits		LMC-SN1987A	4.2	_	52X2	NONE
935	o3x107010_wav.fits	1	WAVELINE	4.2	6581	52X0.1	LINE
	o3x107a8m_raw.fits	1	WAVEHITM	4.2	6581	52X0.1	HITM:
	o3x107020_raw.fits	2	LMC-SN1987A	3.2	4961	52X2	NONE
938	o3x107020_wav.fits	1	WAVELINE	3.2	4961	52X0.1	LINE
939	o3x107acm_raw.fits	1	WAVEHITM	3.2	4961	52X0.1	HITM:

```
; read one of them
; stis_read, 934, h, d
; or
stis_read, 'o3xl07010_raw.fits', h, d
; does the same thing
hprint,h
; display header
window, xs=1024, ys=1024
tvscl, alog10(d>0.1)
; display image
stis_read, 934, h, d, readout=2
; get second readout
stis_cr, 934, h, d
; cosmic ray rejection using
; both readouts
```

; interactively examine it

#### **B.2 First Order Grating Point Source Spectral Extraction**

```
; calibrate observation 730 with cosmic ray rejection for the cr-split science observation
; using the associated wavecal, id=731, to adjust the wavelength zero point.

calstis, 730, 0, h, m, wave, flux, eps, err,/cr_reject,autowave=731

plot, wave, flux

; read calstis extracted spectrum file

a = mrdfits('spec_730.fits',1,h)

; scale upper and lower background to gross extraction slit

bupper = a.bupper / sxpar(h,'bwidth')*sxpar(h,'gwidth')

blower = a.blower / sxpar(h,'bwidth')*sxpar(h,'gwidth')

plot, a.wavelength, bupper, psym=2

oplot, a.wavelength, blower, psym=4

; overplot the smoothed background used by CALSTIS

oplot,a.wavelength, a.gross - a.net, thick=3
```

#### **B.3 G750M Extended Source Extraction of a Long Slit Observation**.

```
; Perform extended source extraction on G750M observation 934 using wavecal 935 to adjust ; the wavelength zero point. Use STIS_CR to combine the cosmic ray splits with the results ; written into file myimage.fits.
```

stis\_cr, 934, outfile='myimage.fits'
stis\_woffset,935,results
calstis, 'myimage.fits', 'extended=1', h, m, wave, flux, \$
autowave=results

; extract a single spectral line into a 40 x 100 image. Adjust the x-direction plate scale in the ; output image to match the y-direction plate scale. The plate scales in th raw images are ; x = 0.54 arcsec/pixel and y = 0.5 arcsec/pixel. Use an x sample spacing of 0.926 (ratio of y ; plate scale over the x plate scale) to match the y-sample spacing.

## **B.4 Long Slit Extended Source Extraction (M84) with a Constant Delta Radial Velocity**

; Data Preparation, Cosmic Ray and Hot Pixel removal

stis\_cr, 863, h, data,err,eps imgclean, data,h

; Extended source extraction on constant delta Log(Wavelength) Scale so that radial velocity ; shifts are equal at all wavelengths. we will provide the input data to CALSTIS as IDL ; variables instead of using the data in the raw data file

## calstis, 863, 'extended=1,eslogl=1', h, m, wave, flux, \$ data=data, header=h, epsin=eps, errin=err

; Create a template spectrum for cross correlation using the rest wavelengths of the spectral ; lines present in the observation. The line at rest wavelength 6583.4 is approximately twice ; as bright as the others.

sig = 7.0/2.3548 ; seven pixel FWHM gauss, wave, 6548.1, sig, 1.0, l1 gauss, wave, 6562.8, sig, 1.0, l2 gauss, wave, 6583.4, sig, 1.0, l3 gauss, wave, 6716.4, sig, 1.0, l4 gauss, wave, 6730.8, sig, 1.0, l5 template = l1 + l2 + 2\*l3 + l4 + l5

; cross correlate each spatial position with template

; convert to slit position and radial velocities and plot

```
slitpos = (findgen(1000)-534.0) * 0.05
rv = -offsets * (wave(500)-wave(499))/wave(500) * 2.9979e5
plot, slitpos, rv, psym=-4, xrange=[-2.0,2.0], yrange=[700,1600], $
    min_value=700, max_value=1600, $
    ytitle = 'Radial Velocity (km/sec)', $
    xtitle = 'Position Along Slit (arcsec)'
```

#### **B.5 G140L Spectra of Q0302 With Geocoronal Emission Subtracted**

```
; run calstis with a modified background subtraction. We do not want
; to smooth the geocoronal emission before subtraction. To improve
; background S/N, we use a larger BWIDTH.
      par=b_median=0,b_mean1=0,b_mean2=0,bdist=30,bwidth=30'
      calstis,7386,par,h0,m0,w0,f0,q0,e0
      calstis,7394,par,h1,m1,w1,f1,q1,e1
      calstis,7388,par,h2,m2,w2,f2,q2,e2
      calstis,7390,par,h3,m3,w3,f3,q3,e3
      calstis,7392,par,h4,m4,w4,f4,q4,e4
; place results in 2-D arrays
      w = [[w0],[w1],[w2],[w3],[w4]]
      f = [[f0],[f1],[f2],[f3],[f4]]
      q = [[q0],[q1],[q2],[q3],[q4]]
      e = [[e0],[e1],[e2],[e3],[e4]]
; merge results from all five observations weighted by the exposure times given in vector t
      dbopen, 'stislog'
      dbext, [7386,7394,7388,7390,7392], 'integ', t
      hrs_merge, w,f,q,e, wave,flux,eps,err, weight=t
      plot,wave,flux
```

## **B.6** Calibrate an Echelle Spectrum Including a Scattered Light Correction

```
; Calibrate an E140M observation of CPD-59D2603 in MAMA Hi-Res (2048x2048) mode
; using the associated wavecal for zero point wavelength adjustment.
;

calstis,12729,'outimage=def,hires=1',/autowave
;
; Perform echelle scattered light correction

echelle_scat,12729
;
; read spectral output for both before and after scattered light correction

before = mrdfits('spec_12729.fits', 1)
    after = mrdfits('newspec_12729.fits', 1)
;
; compare one of the orders before and after the scattered light correction

!grid = 1
    plot, before(35).wavelength, before(35).flux<1e-12, nsum=11
    oplot, after(35).wavelength, after(35).flux, nsum=11, thick=2
```