

Geometric Transformation

Story

I want a way to quickly and efficiently go from un-dispersed to dispersed images reference frame. It is therefore an $(x,y) \rightarrow (dx,dy)$ transformation but wavelength factors in this, as well as where on the detector we are. It is therefore, in a more general way a transformation from $(x,y,\lambda) \rightarrow (dx,dy,\lambda)$, where (x,y) is the location on the detector to use to account for any field dependency.

I can think of several things I would use this for:

- Simulating: In this case I want to be able to compute the offsets (dx,dy) coordinates in the dispersed image where light originating from (x,y) in the undispersed image and at a particular wavelength λ ends up.
- Extracting: This is a reverse operation. I know the coordinates in a dispersed spectrum (x',y') and therefore (dx,dy) and I want to know what the wavelength of that light is since I know it originates from the pixel (x,y) in the direct image
- In between cases: Some times, I want to be able to go back and forth from (x,y,λ) to (dx,dy,λ) with missing values of x , y or λ , or dx,dy,λ so I need functions that give me the flexibility to solve any of the in between problems, equally as fast.

I envision that I will need to compute this several millions of times so I want to keep things as fast as possible. It could be a 2D polynomial solution, which can be inverted either analytically or numerically in the case of higher polynomial orders (as in UVIS).

I want this system to be flexible and extendable to highly distorted spectra so I do not want the calibration to be dependent on the x or y coordinates, or the pathlength along the trace (which is difficult to compute and invert) or λ directly. The choice of x or y is a bad one for example in cases where the spectra curve dramatically and either of these coordinates become a poor way to measure where the trace is. Instead I want a system where each of this transformation is given as a function of an intermediate variable t which can then be defined as best fit a particular disperser. Hence the functions will be $dx = f_x(x,y,t)$, $dy = f_y(x,y,t)$, $\lambda = f_\lambda(x,y,t)$ and the inverse functions $x = f_x'(dx,dy,t)$, $y = f_y'(dx,dy,t)$, $\lambda = f_\lambda'(dx,dy,t)$. The variable t can be defined as $t = x$ in the case of a simple x -direction disperser. It can also be defined as $t = \lambda$, or $t = (\lambda - \lambda_0)/(\lambda_1 - \lambda_0)$ where λ_0 and λ_1 are the minimum and maximum wavelength in the dispersed spectrum. The latter results in $0 < t < 1$ which could be beneficial to some.

Inputs

- A grism configuration file that described the transformations
- A set of coordinates and wavelengths, etc...

Outputs

- As set of coordinates and wavelengths, etc...

Computations

- Computations are handled by a small module which simply implements the equation described above, using parameters values determined during the course of the instrument calibration.

Drawbacks

- The use of the variable t provides more flexibility but some bad choices can still be made. How to parametrize t should be defined by balancing adequacy, speed, as well as the ability of the instrument team to calibrate the grism in that manner