

**NAME**

pyFAI-calib – pyFAI-calib

**DESCRIPTION**

INFO:root:Enter, port=54321. INFO:root:Enter. usage: pyFAI-calib [options] **-w** 1 **-D** detector **-c** calibrant.D imagefile.edf

Calibrate the diffraction setup geometry based on Debye–Sherrer rings images without a priori knowledge of your setup. You will need to provide a calibrant or a "d-spacing" file containing the spacing of Miller plans in Angstrom (in decreasing order). If you are using a standard calibrant, look at <https://github.com/kif/pyFAI/tree/master/calibration> or search in the American Mineralogist database: <http://rruff.geo.arizona.edu/AMS/amcsd.php> The **--calibrant** option is mandatory !

**positional arguments:**

**FILE** List of files to calibrate

**optional arguments:**

**-h, --help**

show this help message and exit

**-V, --version**

**-o FILE, --out FILE**

Filename where processed image is saved

**-v, --verbose**

switch to debug/verbose mode

**-c FILE, --calibrant FILE**

Calibrant name or file containing d-spacing of the reference sample (MANDATORY, case sensitive !)

**-w WAVELENGTH, --wavelength WAVELENGTH**

wavelength of the X-Ray beam in Angstrom. Mandatory

**-e ENERGY, --energy ENERGY**

energy of the X-Ray beam in keV ( $hc=12.398419292\text{keV}\cdot\text{Å}$ ).

**-P POLARIZATION\_FACTOR, --polarization POLARIZATION\_FACTOR**

polarization factor, from **-1** (vertical) to **+1** (horizontal), default is None (no correction), synchrotrons are around 0.95

**-b BACKGROUND, --background BACKGROUND**

Automatic background subtraction if no value are provided

**-d DARK, --dark DARK**

list of comma separated dark images to average and subtract

**-f FLAT, --flat FLAT**

list of comma separated flat images to average and divide

**-s SPLINE, --spline SPLINE**

spline file describing the detector distortion

**-D DETECTOR\_NAME, --detector DETECTOR\_NAME**

Detector name (instead of pixel size+spline)

**-m MASK, --mask MASK**

file containing the mask (for image reconstruction)

**-n NPT, --pt NPT**

file with datapoints saved. Default: basename.npt

**--filter FILTER**

select the filter, either mean(default), max or median

- l DISTANCE, --distance DISTANCE**  
sample-detector distance in millimeter. Default: 0.1m
- poni1 PONI1**  
poni1 coordinate in meter. Default: center of detector
- poni2 PONI2**  
poni2 coordinate in meter. Default: center of detector
- rot1 ROT1**  
rot1 in radians. default: 0
- rot2 ROT2**  
rot2 in radians. default: 0
- rot3 ROT3**  
rot3 in radians. default: 0
- fix-dist**  
fix the distance parameter
- free-dist**  
free the distance parameter. Default: Activated
- fix-poni1**  
fix the poni1 parameter
- free-poni1**  
free the poni1 parameter. Default: Activated
- fix-poni2**  
fix the poni2 parameter
- free-poni2**  
free the poni2 parameter. Default: Activated
- fix-rot1**  
fix the rot1 parameter
- free-rot1**  
free the rot1 parameter. Default: Activated
- fix-rot2**  
fix the rot2 parameter
- free-rot2**  
free the rot2 parameter. Default: Activated
- fix-rot3**  
fix the rot3 parameter
- free-rot3**  
free the rot3 parameter. Default: Activated
- fix-wavelength**  
fix the wavelength parameter. Default: Activated
- free-wavelength**  
free the wavelength parameter. Default: Deactivated
- saturation SATURATION**  
consider all pixel>max\*(1-saturation) as saturated and reconstruct them, default: 0 (deactivated)
- weighted**  
weight fit by intensity, by default not.

- npt** NPT\_1D  
Number of point in 1D integrated pattern, Default: 1024
- npt--azim** NPT\_2D\_AZIM  
Number of azimuthal sectors in 2D integrated images. Default: 360
- npt--rad** NPT\_2D\_RAD  
Number of radial bins in 2D integrated images. Default: 400
- unit** UNIT  
Valid units for radial range: 2th\_deg, 2th\_rad, q\_nm<sup>-1</sup>, q\_A<sup>-1</sup>, r\_mm. Default: 2th\_deg
- no-gui**  
force the program to run without a Graphical interface
- no-interactive**  
force the program to run and exit without prompting for refinements
- r, --reconstruct**  
Reconstruct image where data are masked or <0 (for Pilatus detectors or detectors with modules)
- g** GAUSSIAN, **--gaussian** GAUSSIAN  
Size of the gaussian kernel. Size of the gap (in pixels) between two consecutive rings, by default 100 Increase the value if the arc is not complete; decrease the value if arcs are mixed together.
- square**  
Use square kernel shape for neighbor search instead of diamond shape
- p** PIXEL, **--pixel** PIXEL  
size of the pixel in micron

The output of this program is a "PONI" file containing the detector description and the 6 refined parameters (distance, center, rotation) and wavelength. An 1D and 2D diffraction patterns are also produced. (.dat and .azim files)