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# **pydirtygrid Documentation**

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This is an affiliated package for the AstroPy package. The documentation for this package is here:



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## PyDirtyGrid

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Code to access and work on the DirtyGrid set of stellar population/dust RT models.

Uses the astropy affiliated package template





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## Repository

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Github: <<https://github.com/karllark/pydirtygrid>>



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## DirtyGrid Data

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The DirtyGrid is provide as a large hd5f file. This file provides photometry in a number of bands.

Location: <<https://stsci.box.com/v/dirtygrid>>



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## Reference API

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### pydirtygrid.PhotDG Module

#### Classes

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<code>PhotDG([datafile])</code>	Read in the DirtyGrid cube from a HDF5 file
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#### PhotDG

**class** pydirtygrid.PhotDG.**PhotDG**(*datafile*='data/dirtygrid\_29mar17.hdf5')  
 Read in the DirtyGrid cube from a HDF5 file

##### Returns

(self.)seds: a list that grows with the SED you choose to extract

(self.)dgrid: the HDF5 file: with attributes corresponding to the parameters values; with each dataset corresponding to each band

#### Methods Summary

<code>findIndexFromParams</code> (grain, geom, sf_type, ...)	Give a set of parameters and find where they will be stored in the
<code>photAddNew</code> (wave0, band_name, newcube[, datafile])	Create the new dataset and add it to the file
<code>photGet</code> (grain, geom, sf_type, metal, age, ...)	Save photometry points given a set of parameters.
<code>photPlot</code> ([ind])	Plot photometry points, either giving a specific set of bands, or all
<code>print_parameters</code> ()	

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#### Methods Documentation

**findIndexFromParams**(*grain, geom, sf\_type, metal, age, sfr, tau*)

Give a set of parameters and find where they will be stored in the DirtyGrid cube

##### Parameters

**grain:** string

type of grain

**geom:** string

geometry

**sf\_type: string**

star formation type

**metal: float**

metallicity

**age: float**

age of the stellar population

**sfr: float**

star formation rate

**tau: float**

optical depth

#### Returns

indgt, indgm, indst, indmt, indsa, indsr, indta: integers

the corresponding indices

**photAddNew**(*wave0*, *band\_name*, *newcube*, *datafile*= 'data/dirtygrid\_29mar17.hdf5')

Create the new dataset and add it to the file

#### Parameters

**wave0: float**

central wavelength of the new filter

**band\_name: string**

name of the new filter

**newcube: array(3, 6, 2, 5, 50, 29, 25) (float)**

new photometry to be added to the file as a new dataset

**photGet**(*grain*, *geom*, *sf\_type*, *metal*, *age*, *sfr*, *tau*, *bands=-1*)

Save photometry points given a set of parameters. The function allows for:

- values not in the DGrid parameter sampling, and finds the closest point
- specific bands only

#### Parameters

**grain: string**

type of grain

**geom: string**

geometry

**sf\_type: string**

star formation type

**metal: float**

metallicity

**age: float**

age of the stellar population

**sfr: float**

star formation rate

**tau: float**

optical depth

**Returns**

seds: save the corresponding SED

**photPlot**(*ind=-1*)

**Plot photometry points, either giving a specific set of bands, or all**

of them either giving a specific index of which if more that one saved, or all of them

**Parameters**

**ind: integer(s) (Optional)**

indices of the SED you wish to plot

**print\_parameters()**

## Class Inheritance Diagram

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graph TD; PhotDG[PhotDG];
```

## pydirtygrid.SpecDG Module

### Classes

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<code>SpecDG()</code>	Read in the mapping file for all spectra
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### SpecDG

**class** pydirtygrid.SpecDG.**SpecDG**

Read in the mapping file for all spectra

**Returns**

(self.)plan: Table

table containing the identification of each spectrum, as well as its parameters

**(self.)seds: list**  
empty list to save the spectra if you want to

## Methods Summary

<code>findFile(file_id)</code>	Returns the full filename for a given GID from the spectrum mapping.
<code>findGidFromParam(grain, geom, sf_type, ...)</code>	Returns the GID given a set of parameters
<code>findParamsFromGid(thisgid)</code>	Returns the parameter values for a given GID
<code>spec2Phot(trans_curve, trans_waves, wave0[, ...])</code>	Compute the new photometry
<code>specGet(filename)</code>	Save a spectrum from the filename
<code>specPlot([ind])</code>	Plot SEDs, either giving a specific index of which if more that

## Methods Documentation

### `findFile(file_id)`

Returns the full filename for a given GID from the spectrum mapping.

#### Returns

filepath+filename: string  
the absolute path and name of the .fits file

### `findGidFromParam(grain, geom, sf_type, metal, age, sfr, tau)`

Returns the GID given a set of parameters (If you want to plot or something)

#### Parameters

**grain: string**  
type of grain  
**geom: string**  
geometry  
**sf\_type: string**  
star formation type  
**metal: float**  
metallicity  
**age: float**  
age of the stellar population  
**sfr: float**  
star formation rate  
**tau: float**  
optical depth

#### Returns

file\_gid: string  
identification number of the spectrum with the given set  
of parameters



**findParamsFromGid**(*thisgid*)

Returns the parameter values for a given GID (Used later to update the cube)

**Parameters**

**thisgid:** string

identification number of the spectrum

**Returns**

thisgt, thisgm, thissf, thismt, thissa, thissr, thista: integers

indices of the parameter values for the given spectrum,

in the multidimensional photometry cube

**spec2Phot**(*trans\_curve*, *trans\_waves*, *wave0*, *energy=1*)

Compute the new photometry

**Parameters**

**trans\_curve:** array

transmission curve of the new filter /!Alread read, as an array

**trans\_waves:** array

wavelengths of the new filter

**wave0:** float

center/effective wavelength of the new filter

**energy:** integer (Optional)

specification for integration with energy(=1, default) or photons

**Returns**

newcube: array(3, 6, 2, 5, 50, 29, 25) (float)

new photometry cube

**specGet**(*filename*)

Save a spectrum from the filename

**Parameters**

**filename:** string

absolute path to the spectrum you want to read

**Returns**

(self.)seds: list

update the list to add the SED

(self.)waves: array

the wavelengths to the SED

**specPlot**(*ind=-1*)

Plot SEDs, either giving a specific index of which if more that one saved, or all of them

**Parameters**

**ind:** integer(s) (Optional)

indices of the SED you wish to plot

## Class Inheritance Diagram



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## Indices and tables

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