

**Leftovers**

**Odds and Ends**

**Bits and Bobs**

**Cleaning out the Fridge**

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# Time Conversion

- Timekeeping is complicated. **Time formats** are independent of **time scales**!
- E.g. TT = terrestrial time does not have leap seconds, UTC does.
- MET overview: [https://fermi.gsfc.nasa.gov/ssc/data/analysis/documentation/Cicerone/Cicerone\\_Data/Time\\_in\\_ScienceTools.html](https://fermi.gsfc.nasa.gov/ssc/data/analysis/documentation/Cicerone/Cicerone_Data/Time_in_ScienceTools.html)
- Online conversion tool: <https://heasarc.gsfc.nasa.gov/cgi-bin/Tools/xTime/xTime.pl>
- DIY: **MJDREFI=51910; MJDREFF=7.428703703703703e-4** (86400 seconds/day)
- Use [astropy.time](#) to convert between scales, e.g. TT to calendar time (UTC).

- **MET**: Mission elapsed time (s)
- **MJD**: Modified Julian date (d)
- **TT**: Terrestrial time
- **UTC**: Coordinated Universal Time

```
from astropy.time import Time

myTT = 675702050 / 86400 + (51910+7.428703703703703e-4)
t = Time(myTT, format='mjd', scale='tt')

print("TT: ", t, " = ", t.isot)
print("UTC:", t.utc, " = ", t.utc.isot)

TT: 59730.62632157407 = 2022-05-31T15:01:54.184
UTC: 59730.62552083333 = 2022-05-31T15:00:45.000
```

## PLSuperExpCutoff4:

Example: [XML Model Definition](#)

For modeling pulsars.

# PLEC4

$$\frac{dN}{dE} = \begin{cases} N_0 \left( \frac{E}{E_0} \right)^{\gamma_0 - \frac{d}{2} \ln \frac{E}{E_0} - \frac{db}{6} \ln^2 \frac{E}{E_0} - \frac{db^2}{24} \ln^3 \frac{E}{E_0}}, & \text{if } \left| b \ln \frac{E}{E_0} \right| < 1e^{-2} \\ N_0 \left( \frac{E}{E_0} \right)^{\gamma_0 + d/b} \exp \left( \frac{d}{b^2} \left( 1 - \left( \frac{E}{E_0} \right)^b \right) \right) & \text{otherwise} \end{cases}$$

where

- Prefactor =  $N_0$
- IndexS =  $\gamma_0$
- Scale =  $E_0$
- ExpfactorS =  $d$
- Index2 =  $b$

Note:

- $N_0$  is the normalization (flux density) at  $E_0$ .
- $\gamma_0$  is the local spectral index at  $E_0$ .
- $d$  is the local curvature at  $E_0$

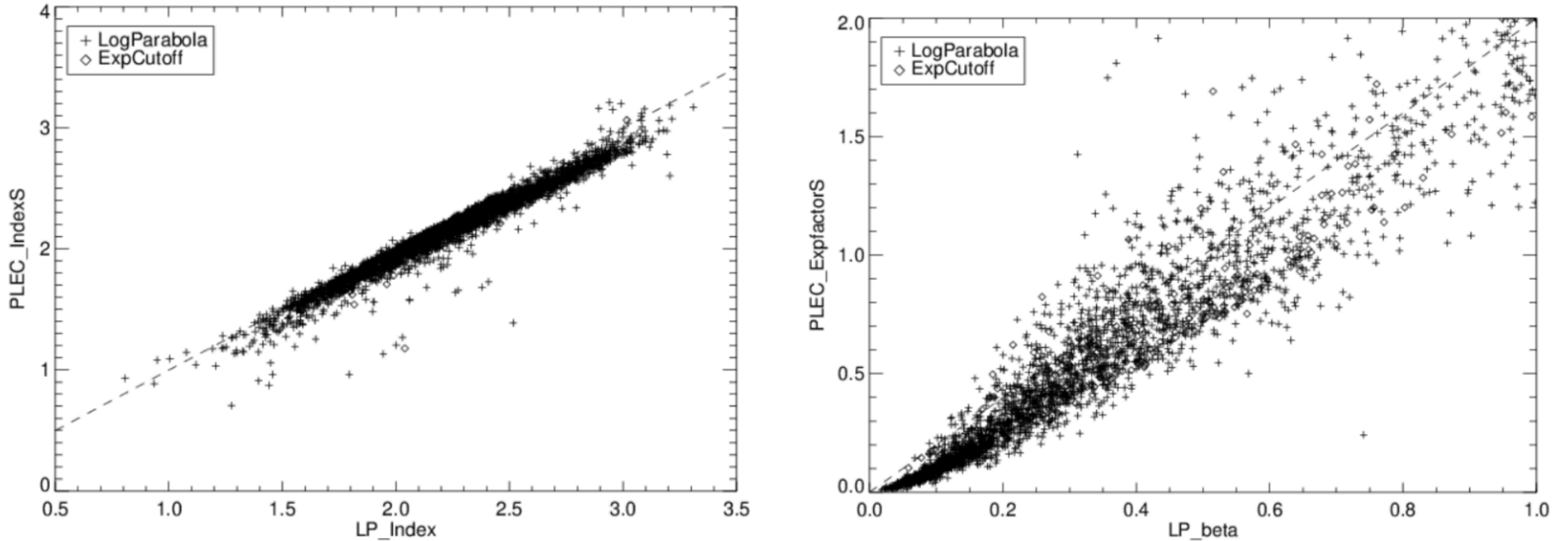
To provide more information on curved spectra, we now report systematically in the catalog the peak energy in  $\nu F_\nu$  and its uncertainty for all sources (including those not significantly curved) and both models as (Unc\_)LP\_EPeak and (Unc\_)PLEC\_EPeak:

$$E_{\text{peak}}(\text{LP}) = E_0 \exp \left( \frac{2 - \alpha}{2\beta} \right) \quad (4)$$

$$E_{\text{peak}}(\text{PLEC}) = E_0 \left( 1 + \frac{b}{d} (2 - \Gamma_S) \right)^{1/b} \quad (5)$$

The  $b = 0$  case corresponds to a LogParabola with  $\alpha = \gamma_0$  and  $\beta = d/2$ .

# PLEC4



**Figure 4.** Left: comparison of  $\alpha$  of the LP model and  $\Gamma_S$  of the PLEC model (LP\_Index and PLEC\_IndexS in the FITS file), showing that these parameters are largely similar. Right: comparison of LP  $\beta$  and PLEC  $d$  (LP\_beta and PLEC\_ExpfactorS in the FITS file), showing that  $d$  is well correlated to  $2\beta$ . The dashed lines show a one-to-one correlation. The outliers have large errors on both plots (none is farther off than  $2\sigma$ ).



# Spectral models in 4FGL-DR3

The new PLEC parameterization allows a 4-parameter fit (with **free b**) in fainter sources. We have applied it to **all pulsars with TS > 10,000** (28, up from 6 in DR1 and DR2). At  $TS < 10,000$   $\Delta b$  becomes larger than 0.15, which is the natural scatter on b in the brightest pulsars, so freeing b is no longer beneficial. For all other significantly curved pulsars, **b is fixed to 2/3 as in 4FGL**. For comparison, the median b over the 28 pulsars with free b is 0.51, its weighted (with  $1/\sigma^2$ ) average is 0.55, and its intrinsic dispersion is 0.16.

The **Small Magellanic Cloud has b = 1**, as before (Abdo et al. 2010). Besides 3C 454.3, the five other **AGN with TS > 80,000** were modeled with PLEC and free b as well: CTA 102, Mkn 421, S5 0716+71, 3C 279 and PKS 1424–41.

The new parameterization **contains the LP model** (for  $b=0$ ) so there is no risk of a worse fit, but non-convergence can occur when the model is not constrained well enough. The TS threshold for free b is higher in AGN than pulsars, because the curvature is much less in AGN, so  $\Delta b$  is larger. The main objective of fitting PLEC with free b remains to **improve the modeling of the surroundings of very bright sources at low energy**.

In the DR1 and DR2 catalogs, the sources were represented with a curved spectral model (LP or PLEC) rather than a power law (PL) when  $TS_{\text{curv}} = 2 (\log L(\text{curved spectrum}) - \log L(\text{PL}))$  was larger than 9 ( $3 \sigma$ ). **For DR3 we have lowered that threshold to 4 ( $2 \sigma$ ) [...].**

# Parameter bounds in fermipy

```
def set_parameter(self, name, par, value, true_value=True, scale=None, bounds=None,
error=None, update_source=True):
    """
    Update the value of a parameter. Parameter bounds will automatically be adjusted to
    encompass the new parameter value.

    Parameters
    -----
    name : str - Source name.
    par : str - Parameter name.
    value : float - Parameter value. By default this argument should be the unscaled
    (True) parameter value.
    scale : float - Parameter scale (optional). Value argument is interpreted with
    respect to the scale parameter if it is provided.
    error : float - Parameter error (optional). By default this argument should be the
    unscaled (True) parameter value.
    update_source : bool - Update the source dictionary for the object.
    """
```

# Examples

```
gta.roi["4FGL J0425.6+5522e"].spectral_pars["norm"]  
  
{'name': 'norm',  
 'value': 1.236583491,  
 'error': 0.06524381672,  
 'min': 1e-05,  
 'max': 1000.0,  
 'free': True,  
 'scale': 1e-13}
```

```
gta.set_parameter("4FGL J0425.6+5522e", "norm",  
1.236583491, bounds=[0.01, 100], scale=1e-13,  
true_value=False)
```

```
gta.roi["4FGL  
J0425.6+5522e"].spectral_pars["norm"]
```

```
{'name': 'norm',  
 'value': 1.236583491,  
 'error': 0.06524381672,  
 'min': 0.01,  
 'max': 100.0,  
 'free': True,  
 'scale': 1e-13}
```

```
gta.set_parameter("4FGL J0425.6+5522e",  
"norm", 1.236583491e-13, bounds=[0.01, 100],  
scale=1e-13, true_value=True)
```

```
gta.roi["4FGL  
J0425.6+5522e"].spectral_pars["norm"]
```

```
{'name': 'norm',  
 'value': 1.236583491,  
 'error': 0.06524381672,  
 'min': 0.01,  
 'max': 100.0,  
 'free': True,  
 'scale': 1e-13}
```

# Fermipy configuration

- Configuration options for `config.yaml` are described in the [documentation](#)
- Can be overwritten at run time when you call the relevant functions.

## roiopt

The options in `roiopt` control the default behavior of the `optimize` method. For more information about using this method see the [ROI Optimization and Fitting](#) page.

### roiopt Options

Option	Default	Description
<code>max_free_sources</code>	5	Maximum number of sources that will be fit simultaneously in the first optimization step.
<code>npred_frac</code>	0.95	
<code>npred_threshold</code>	1.0	
<code>shape_ts_threshold</code>	25.0	Threshold on source TS used for determining the sources that will be fit in the third optimization step.
<code>skip</code>	None	List of str source names to skip while optimizing.



# SED configuration

Option	Default	Description
<code>bin_index</code>	2.0	Spectral index that will be use when fitting the energy distribution within an energy bin.
<code>cov_scale</code>	3.0	Scale factor that sets the strength of the prior on nuisance parameters that are free. Setting this to None disables the prior.
<code>free_background</code>	False	Leave background parameters free when performing the fit. If True then any parameters that are currently free in the model will be fit simultaneously with the source of interest.
<code>free_pars</code>	None	Set the parameters of the source of interest that will be freed when performing the global fit. By default all parameters will be freed.
<code>free_radius</code>	None	Free normalizations of background sources within this angular distance in degrees from the source of interest. If None then no sources will be freed.
<code>make_plots</code>	False	Generate diagnostic plots.
<code>ul_confidence</code>	0.95	Confidence level for flux upper limit.
<code>use_local_index</code>	False	Use a power-law approximation to the shape of the global spectrum in each bin. If this is false then a constant index set to <code>bin_index</code> will be used.
<code>write_fits</code>	True	Write the output to a FITS file.
<code>write_npy</code>	True	Write the output dictionary to a numpy file.

# Fermipy Lightcurves: Under the Hood

## Lightcurve fitting routine:

1. Start by freeing target and provided list of sources, fix all else -
  - 1a. If fit fails, **fix all pars except norm** and try again
2. If that fails to converge then try **fixing low TS (<4) sources** and then refit
3. If that fails to converge then try **fixing low-moderate TS (<9) sources** and then refit
4. If that fails then **fix sources out to 1dg away from center of ROI.**  
NB: This comment can be misleading - sources with offset > 1 deg are fixed.
5. If that fails **set values to 0 in output** and print warning message.

<i>lightcurve Options</i>		
Option	Default	Description
<code>binsz</code>	86400.0	Set the lightcurve bin size in seconds.
<code>free_background</code>	False	Leave background parameters free when performing the fit. If True then any parameters that are currently free in the model will be fit simultaneously with the source of interest.
<code>free_params</code>	None	Set the parameters of the source of interest that will be re-fit in each time bin. If this list is empty then all parameters will be freed.
<code>free_radius</code>	None	Free normalizations of background sources within this angular distance in degrees from the source of interest. If None then no sources will be freed.
<code>free_sources</code>	None	List of sources to be freed. These sources will be added to the list of sources satisfying the free_radius selection.
<code>make_plots</code>	False	Generate diagnostic plots.
<code>max_free_sources</code>	5	Maximum number of sources that will be fit simultaneously with the source of interest.
<code>multithread</code>	False	Split the calculation across number of processes set by nthread option.
<code>nbins</code>	None	Set the number of lightcurve bins. The total time range will be evenly split into this number of time bins.
<code>nthread</code>	None	Number of processes to create when multithread is True. If None then one process will be created for each available core.
<code>outdir</code>	None	Store all data in this directory (e.g. "30days"). If None then use current directory.
<code>save_bin_data</code>	True	Save analysis directories for individual time bins. If False then only the analysis results table will be saved.
<code>shape_ts_threshold</code>	16.0	Set the TS threshold at which shape parameters of sources will be freed. If a source is detected with TS less than this value then its shape parameters will be fixed to values derived from the analysis of the full time range.
<code>systematic</code>	0.02	Systematic correction factor for TS:subscript: var . See Sect. 3.6 in 2FGL for details.
<code>time_bins</code>	None	Set the lightcurve bin edge sequence in MET. This option takes precedence over binsz and nbins.
<code>use_local_ltcube</code>	True	Generate a fast LT cube.
<code>use_scaled_srcmap</code>	False	Generate approximate source maps for each time bin by scaling the current source maps by the exposure ratio with respect to that time bin.
<code>write_fits</code>	True	Write the output to a FITS file.
<code>write_npy</code>	True	Write the output dictionary to a numpy file.

# Fermipy as a community contributed tool

- Who's the community?



# Fermipy as a community contributed tool

- Who is the community?
- We all are!
- Many of the original developers have left and/or have little time to implement new features. None are funded to work on fermipy full-time.
- Please let us know if you find a bug, want a new feature, or documentation is lacking.
- <https://github.com/fermiPy/fermipy/issues/>  
(fermipy only - contact FSSC for fermitools or general analysis questions)
- Contribute!
  - Bug reports & feature requests
  - Code (pull request)
  - Documentation updates 🙏
- Please reach out know if you need help to get started.



# Backup