

# A GPU-based framework for multi-variate analysis in particle physics

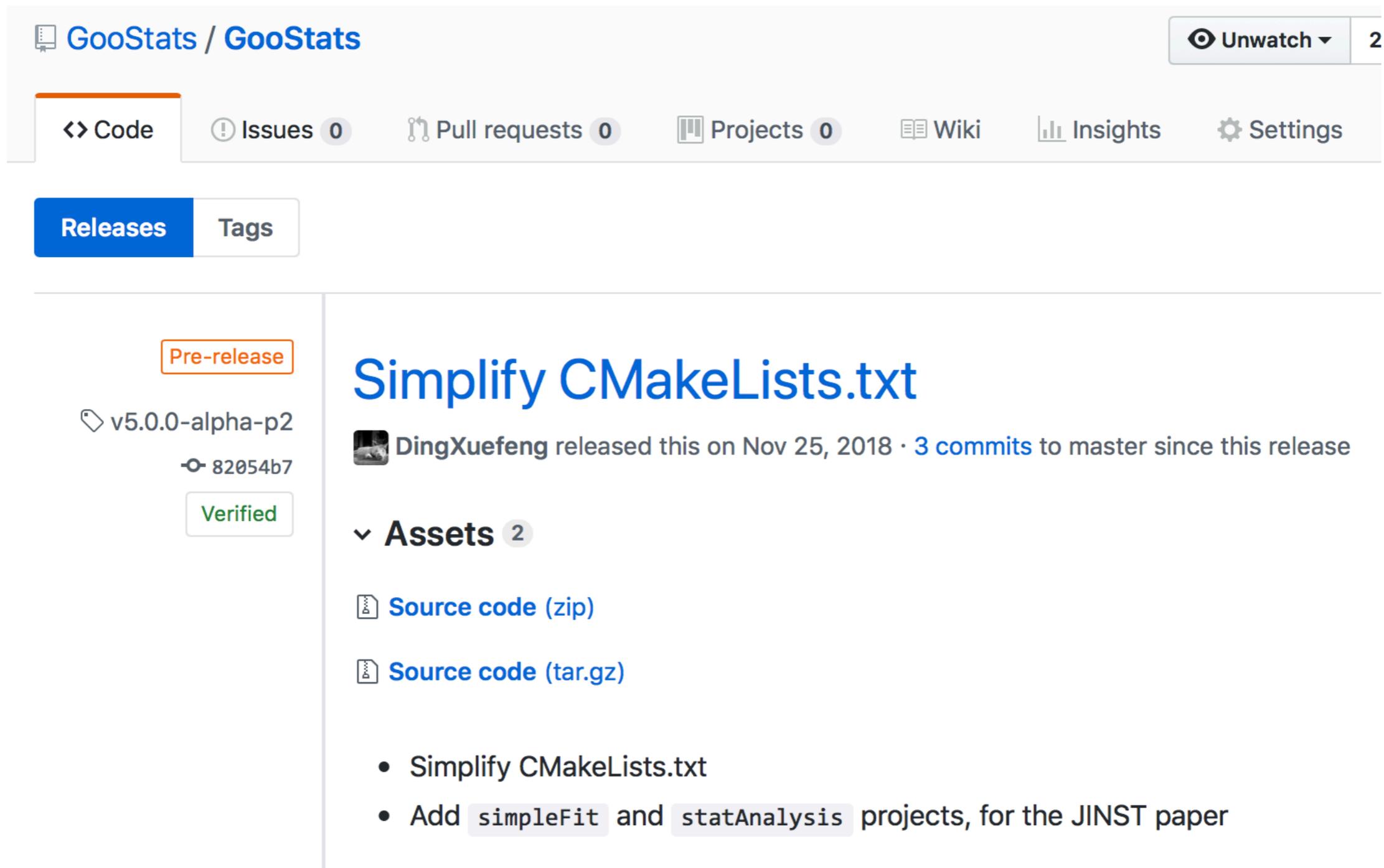
25 Jan 2019, 12:10 (CET) @503-1-001, 12'+3'

**Xuefeng Ding<sup>1,2</sup>**

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- The GooStats framework
- Motivation: Borexino spectrum fit
- Package design
- Systematic uncertainty estimation
- Conclusion

<https://github.com/GooStats/GooStats.git>



GooStats / GooStats Unwatch 2

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**Pre-release**

v5.0.0-alpha-p2  
82054b7

**Verified**

## Simplify CMakeLists.txt

DingXuefeng released this on Nov 25, 2018 · [3 commits](#) to master since this release

Assets 2

- Source code (zip)
- Source code (tar.gz)

- Simplify CMakeLists.txt
- Add `simpleFit` and `statAnalysis` projects, for the JINST paper

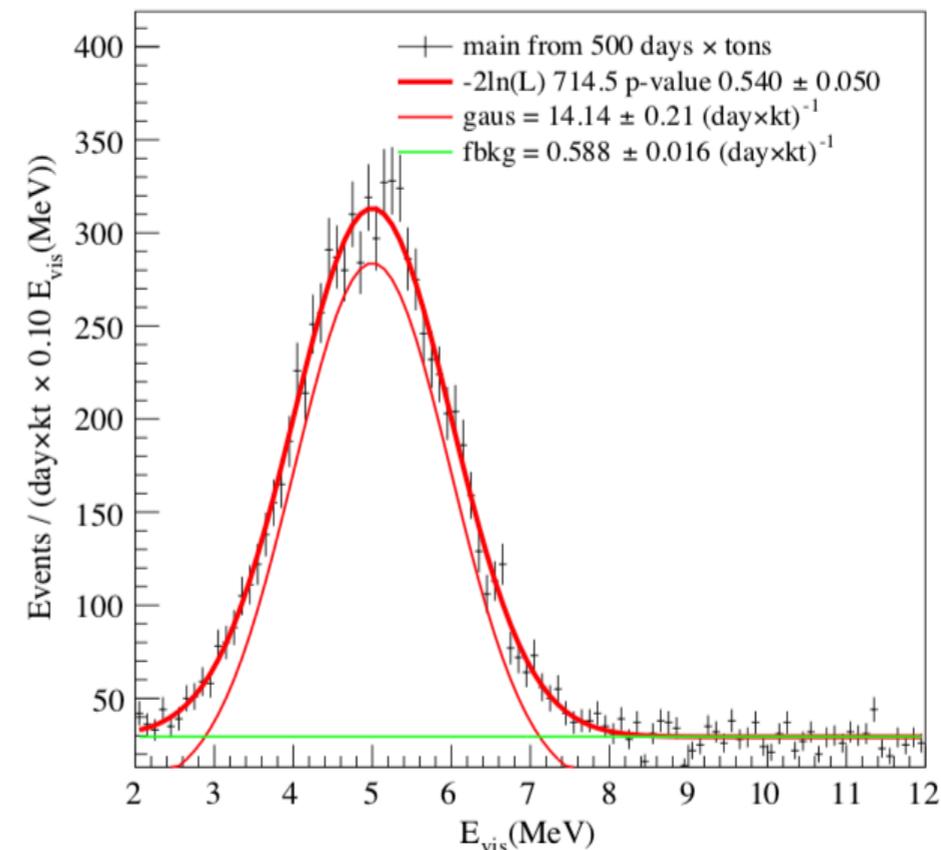
# What and Why?

- A convenient GPU multivariate analysis framework
- Easy spectrum fitting

```
./fit toyMC.cfg out exposure=500
```

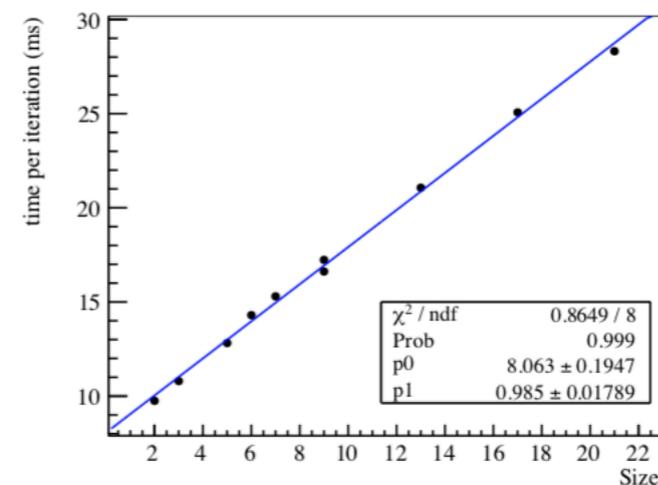
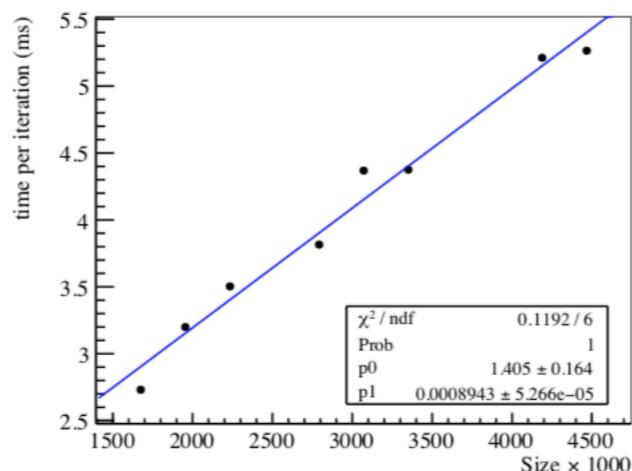
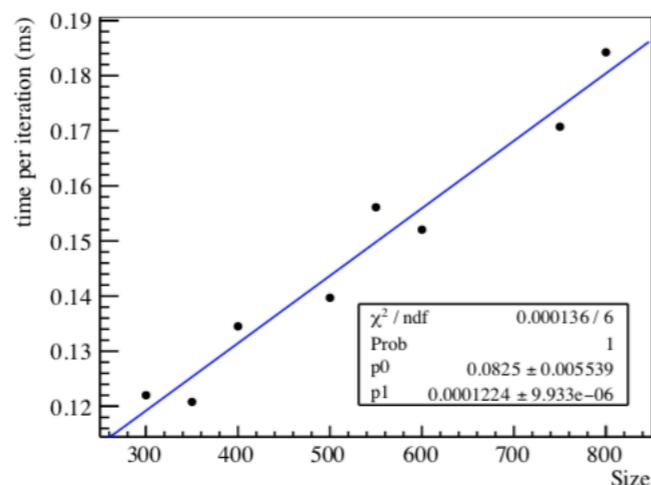
```
FIT PARAMETERS
  gaus = 14.14 #pm 0.21 (day#timeskt)^{-1}
  fbkg = 0.588 #pm 0.016 (day#timeskt)^{-1}
  sdn = 0.991 #pm 0.014
  ^{1}v = 0 [Fixed]
  sigmaT = 0 [Fixed]
  gaus Epeak = 5.000 #pm 0.014 [p] 4.90 #pm 0.50

  chi^2 = 112.7
  chi^2/N-DOF = 1.1739
  p-value = 0.117
  Minimized -2Ln(Likelihood) = 714.53
  Likelihood p-value = 0.540 ± 0.050
```



**Figure 7.** Left: screen shot of fit result summary. Right: produced figure in the file format of pdf.

- Speed up: around 200 times

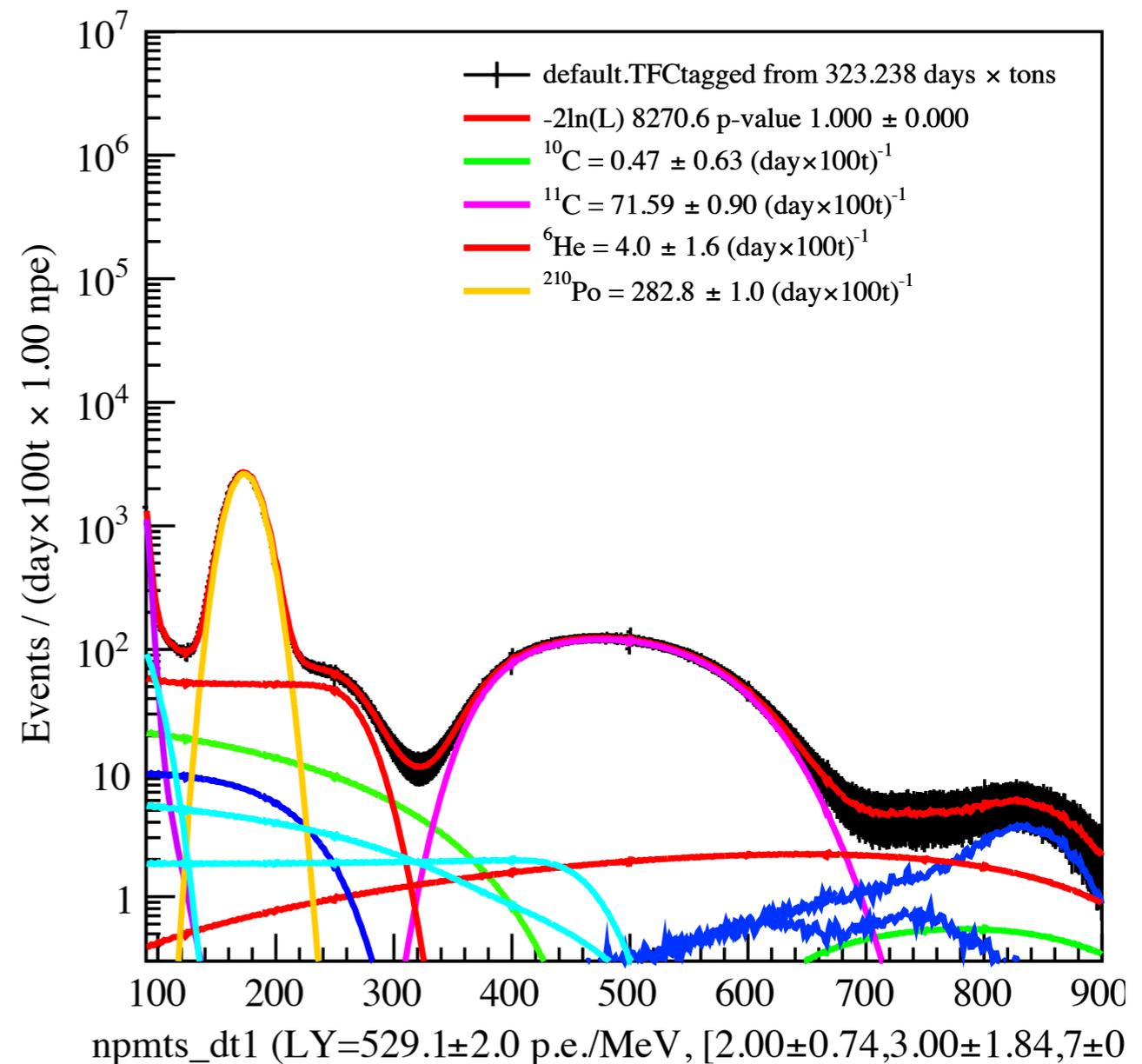
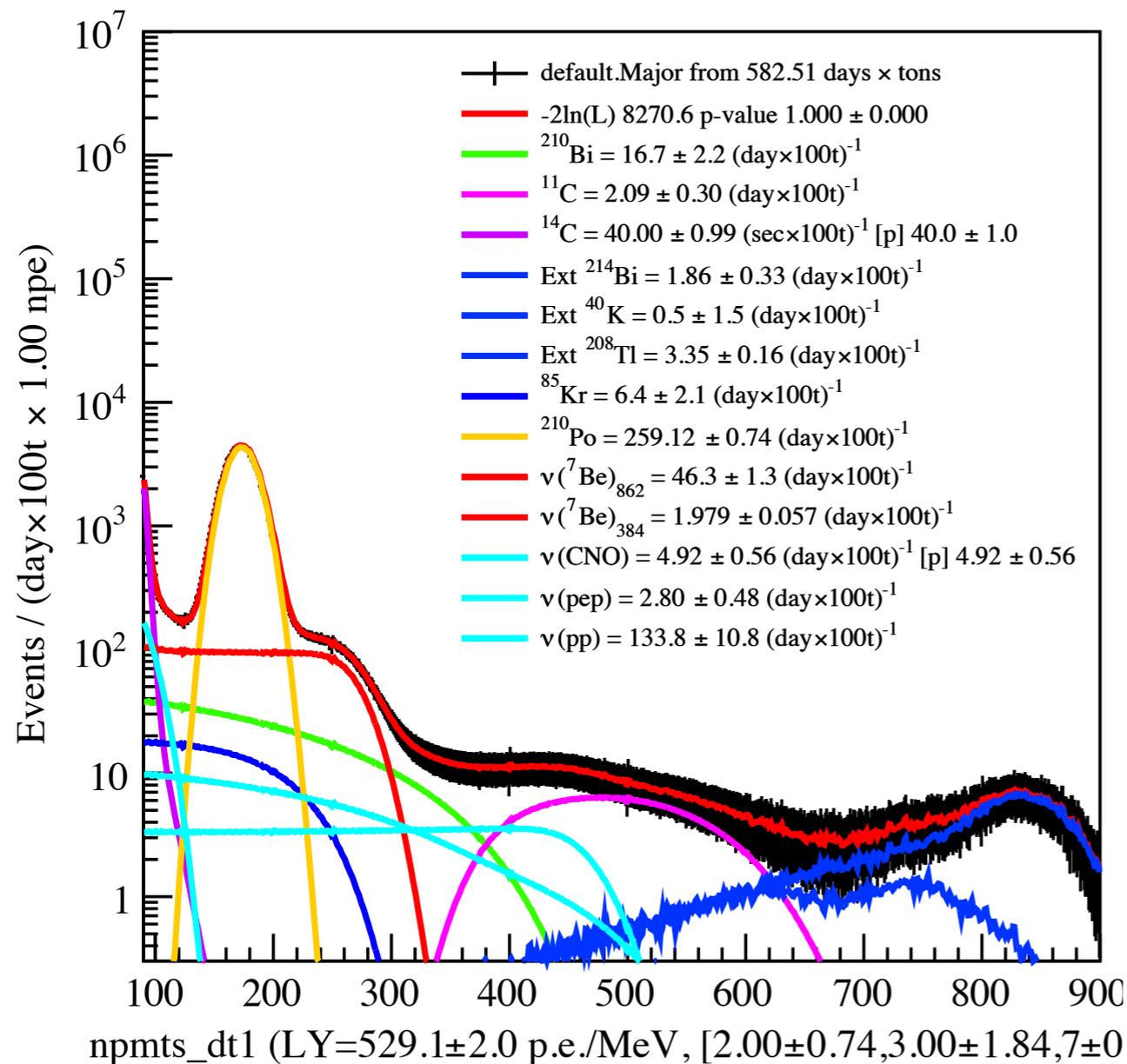


**Table 1.** Comparison of fitting time between GooStats and original software used by the Borexino collaboration.  $T_{\text{tot}}$ : total execution time.  $N$ : the number of iterations taken to converge in MINUIT.  $T_{\text{it}}$ : average execution time per iteration. Speed up:  $T_{\text{it}}(\text{CPU})/T_{\text{it}}(\text{GPU})$ .

Type	CPU			GPU			speed up
	AMD Opteron(TM) Processor 6238			nVidia Tesla K20m			
Size	$T_{\text{tot}}$ (s)	$N$	$T_{\text{it}}$ (ms)	$T_{\text{tot}}$ (s)	$N$	$T_{\text{it}}$ (ms)	
400	27.6	1128	24.4	0.181	1346	0.135	181
350	29.4	1331	22.1	0.156	1294	0.121	183
300	22.5	1239	18.2	0.243	1995	0.122	149

# Motivation: Borexino spectrum fit

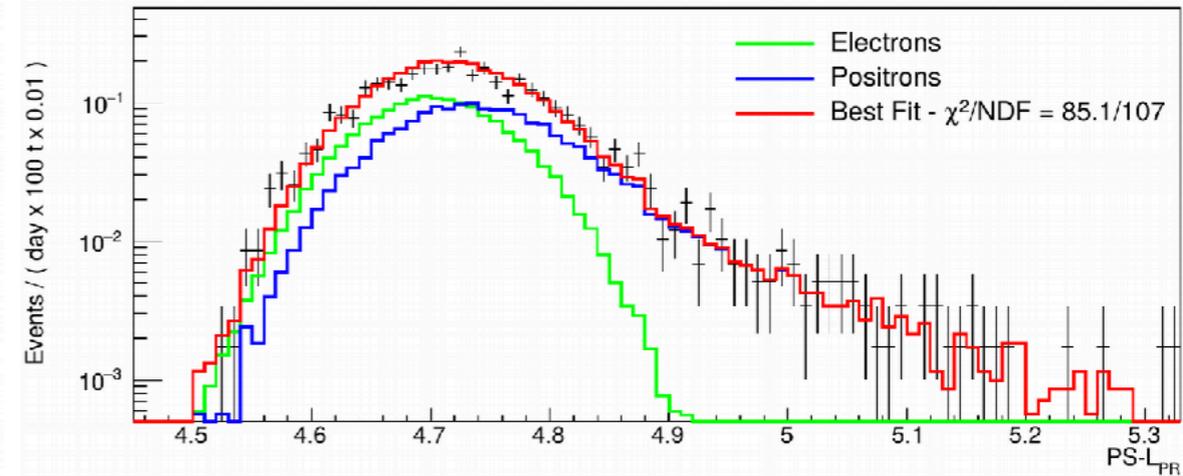
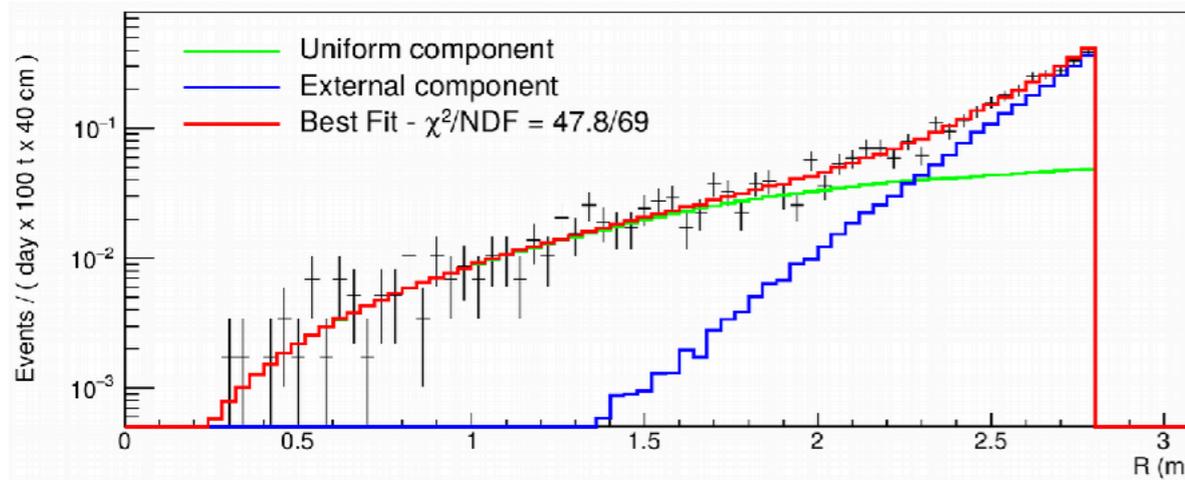
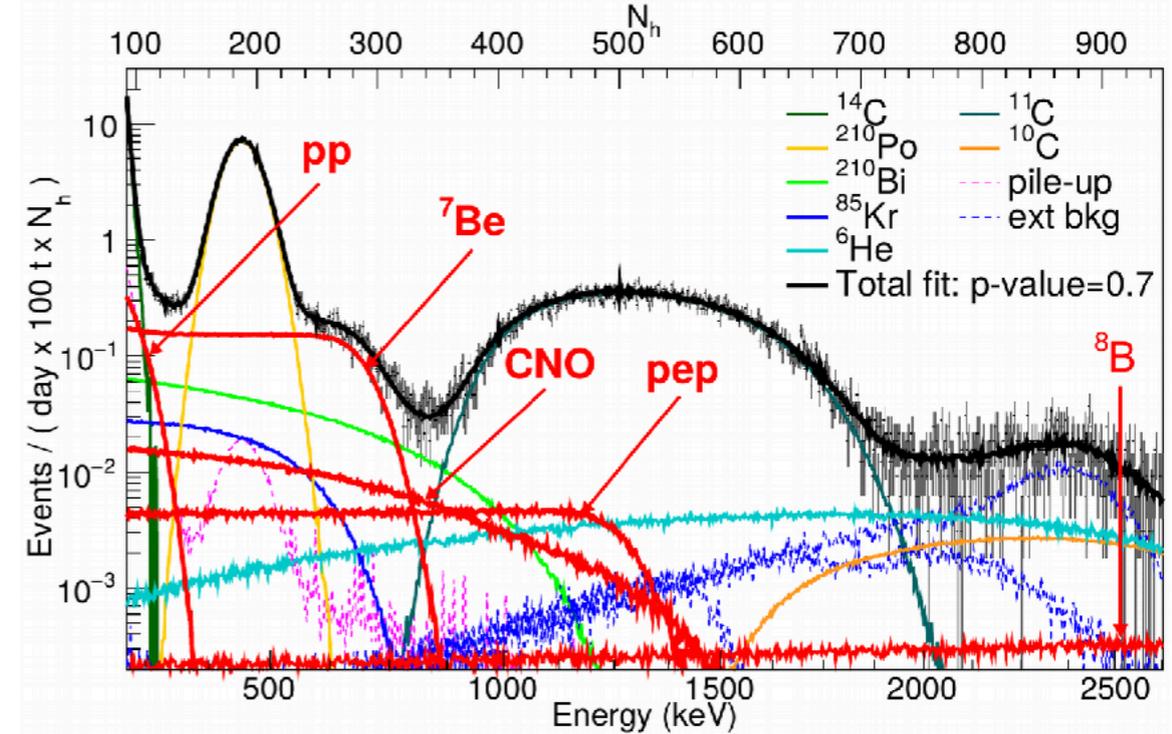
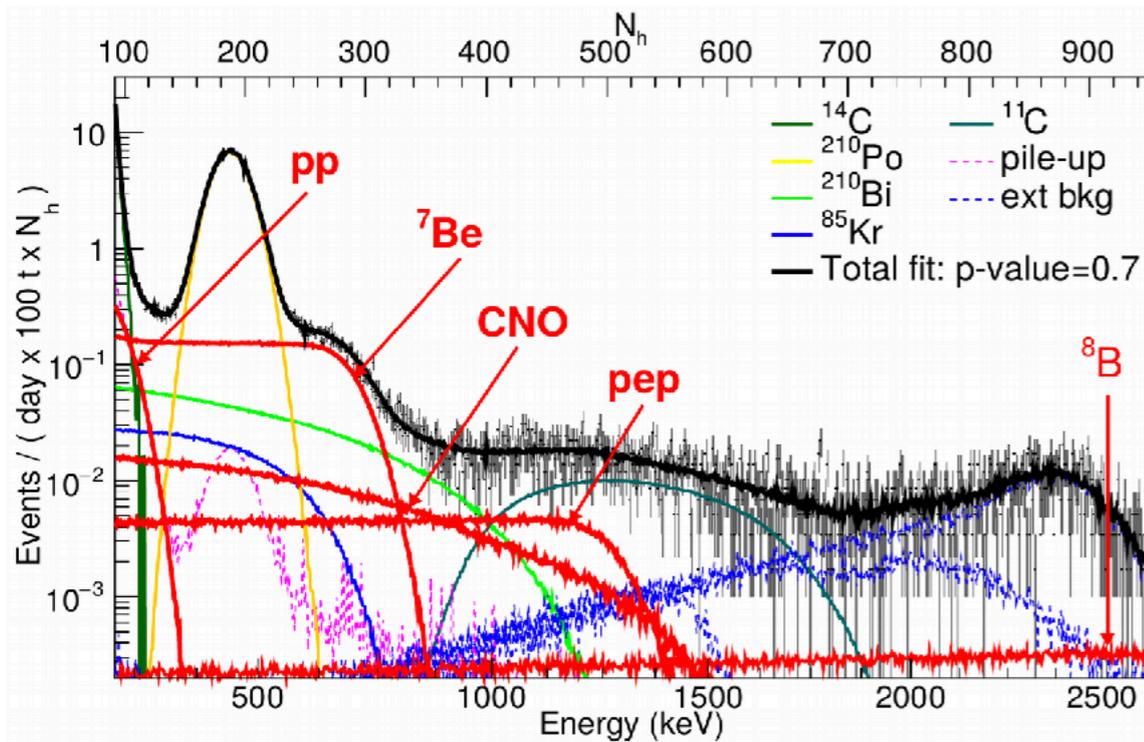
- Motivation: spectral-fitting tool for Borexino analysis



$$M : f(E) \mapsto g(\text{charge}) = \int_0^{E_{\text{end}}} dE \cdot f(E) \cdot \text{RPF} [\text{charge}; \mu(E), \text{var}(\mu)]$$

- Analytical shape of spectrum of mono-energetic events
  - **Momentum based approximation**
  - Match the average ( energy scale + non-linearity model )
  - Match the variance ( energy resolution model )
  - ... (—> simplified)
  - More: “Mask”, “pile-up” etc...
- We can simplify because
  - Borexino response is simple: small FV in center, low energies => no irregular tail
  - We are not sensitive.. => small systematics
  - **Fit full MC to get the bias introduced in simplification**

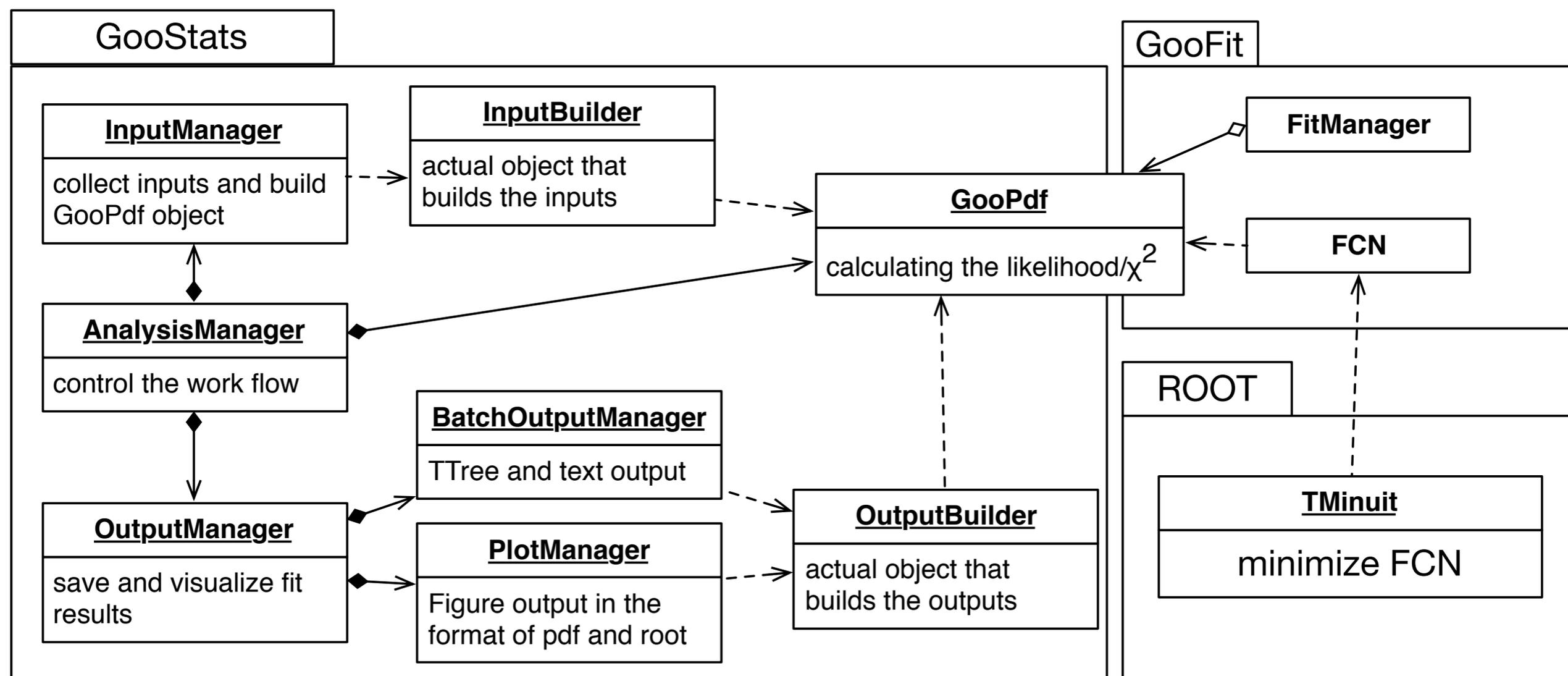
# Multivariate Likelihood (Davini method)



$$\mathcal{L}_{MV}(\vec{\theta}) = \mathcal{L}_{\text{TFC-sub}}(\vec{\theta}) \cdot \mathcal{L}_{\text{TFC-tagged}}(\vec{\theta}) \cdot \mathcal{L}_{\text{RD}}(\vec{\theta}) \cdot \mathcal{L}_{\text{PS}}(\vec{\theta})$$

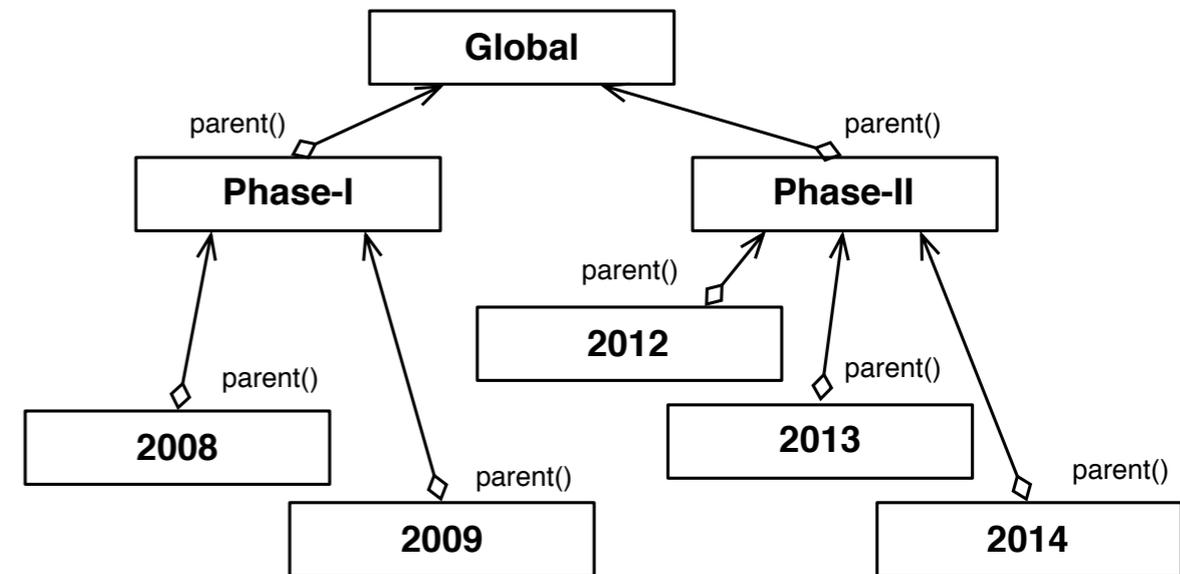
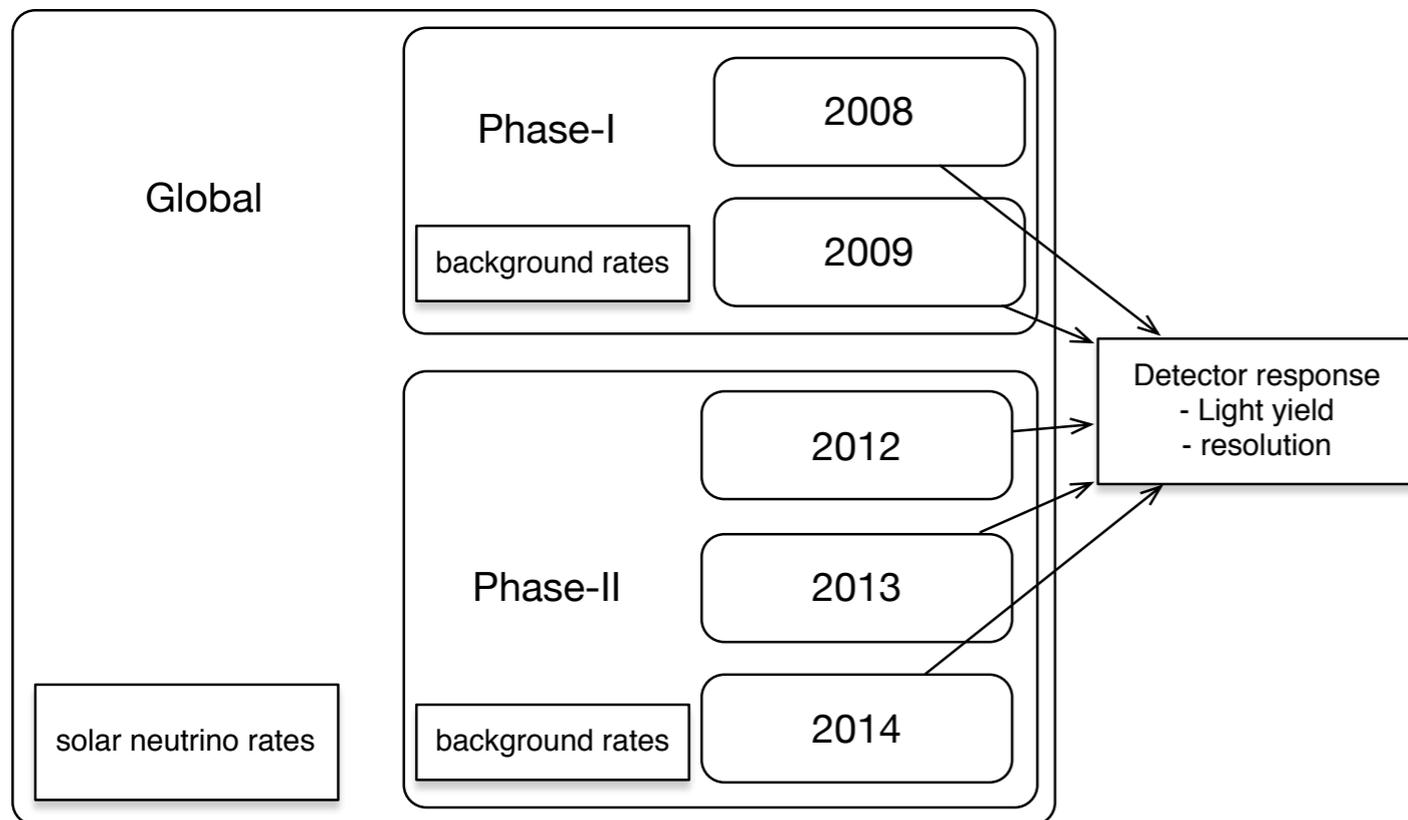
- Scaling factor introduced to remove bias.

- **Middle Layer** between **GooFit** and **User module**



# Parameter Synchronization

- Synchronize fit parameters in joint analysis
- Supported by tree type internal data structure



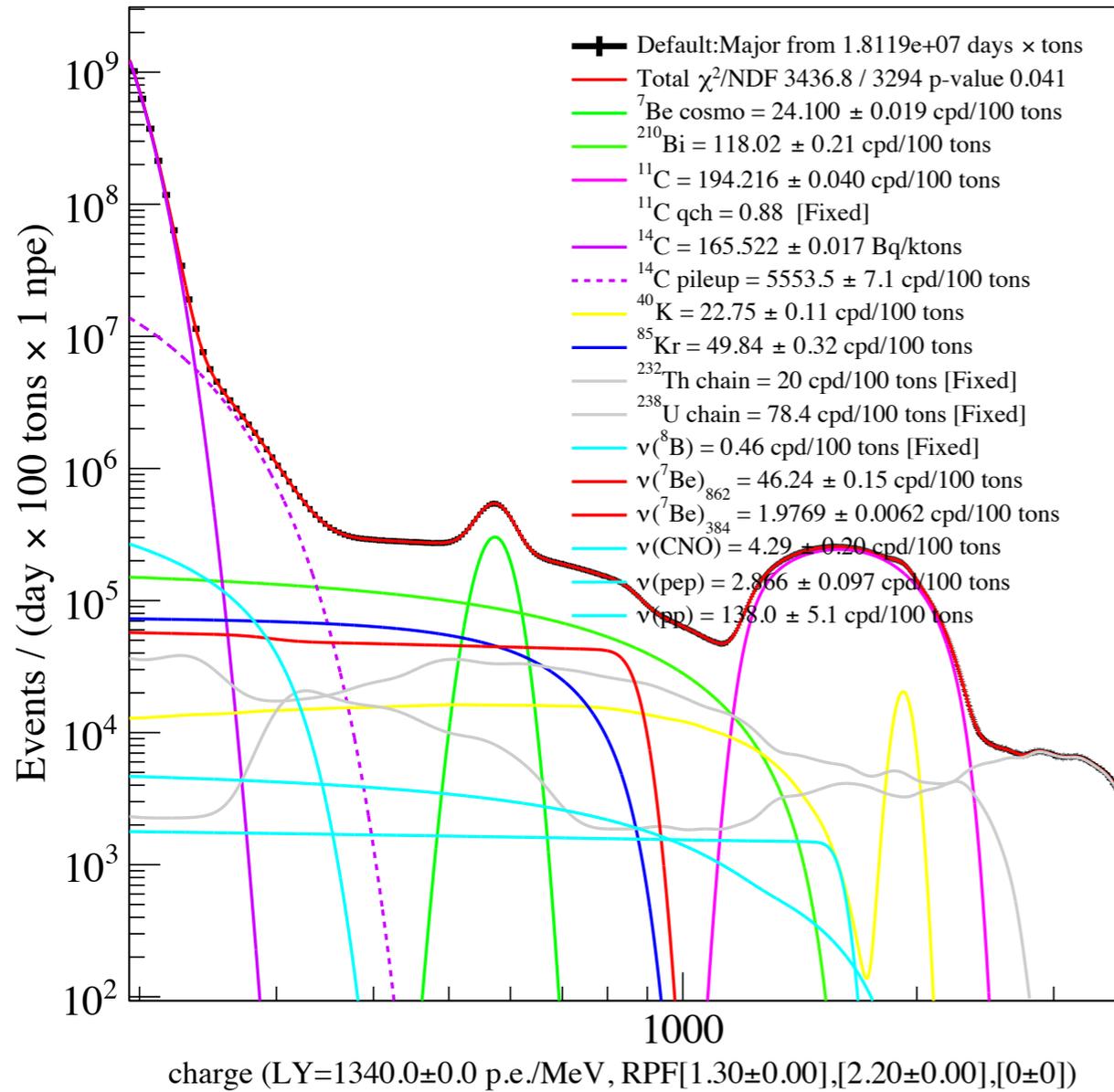
- All fit parameters saved in TTree output automatically
- More quantities can be added as lambda functions in the output builder

```
root [1] .ls
TFile**      test_tree.root
TFile*       test_tree.root
KEY: TTree   fit_result;1   Fit result of GooStats
root [2] fit_result->Show(0)
=====> EVENT:0
default.NReactor = 3.02987
default.NReactor_err = 0.0159146
default.Nbkg      = 1.02405
default.Nbkg_err  = 0.0162459
default.U235      = 0.5
default.U235_err  = 0
default.U238      = 0.2
default.U238_err  = 0
default.Pu239     = 0.1
default.Pu239_err = 0
default.U241      = 0.2
default.U241_err  = 0
default.LY        = 1300
default.LY_err    = 0
default.qc1       = 2.78788
default.qc1_err   = 0
default.qc2       = -0.528003
default.qc2_err   = 0
default.v1        = 0.3
default.v1_err    = 0
default.vT        = 5
default.vT_err    = 0
default.Reactor_dEvis = 1078.28
default.Reactor_dEvis_err = 13.2291
chi2              = 390.448
NDF               = 397
likelihood        = 1883.96
```

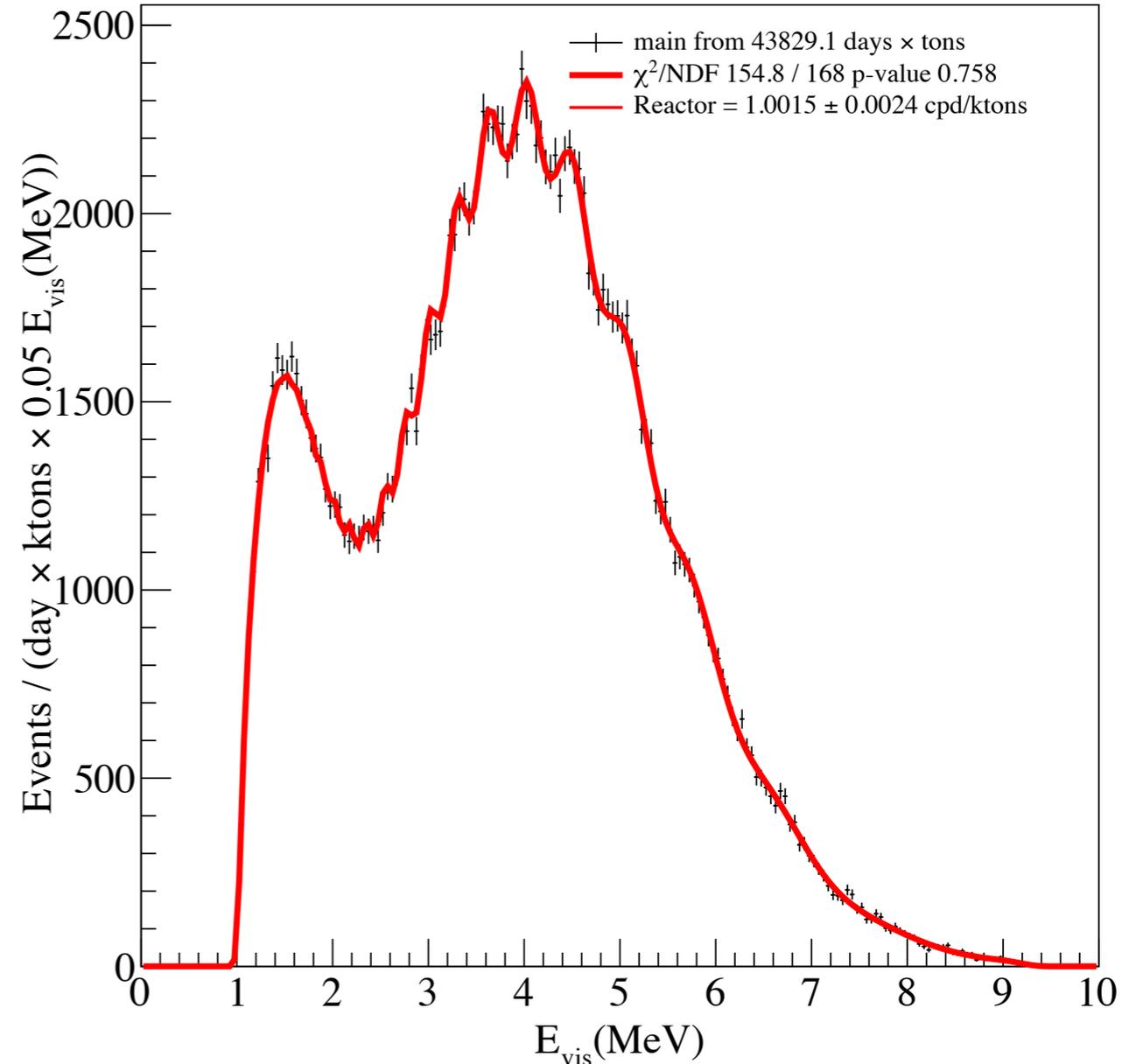
- Adding tasks of statistical analysis by registering new modules inheriting abstract classes
- A few pre-installed modules: ScanPar, DiscoveryTest..

```
PrepareData *data = new PrepareData();  
SimpleFit *fit = new SimpleFit();  
DiscoveryTest *discovery = new DiscoveryTest();  
  
ana->registerModule(inputManager);  
ana->registerModule(data);  
ana->registerModule(fit);  
ana->registerModule(discovery);  
ana->registerModule(outManager);
```

## Solar Neutrinos



## JUNO style



# Two ways of estimating $\sigma_{\text{sys}}$

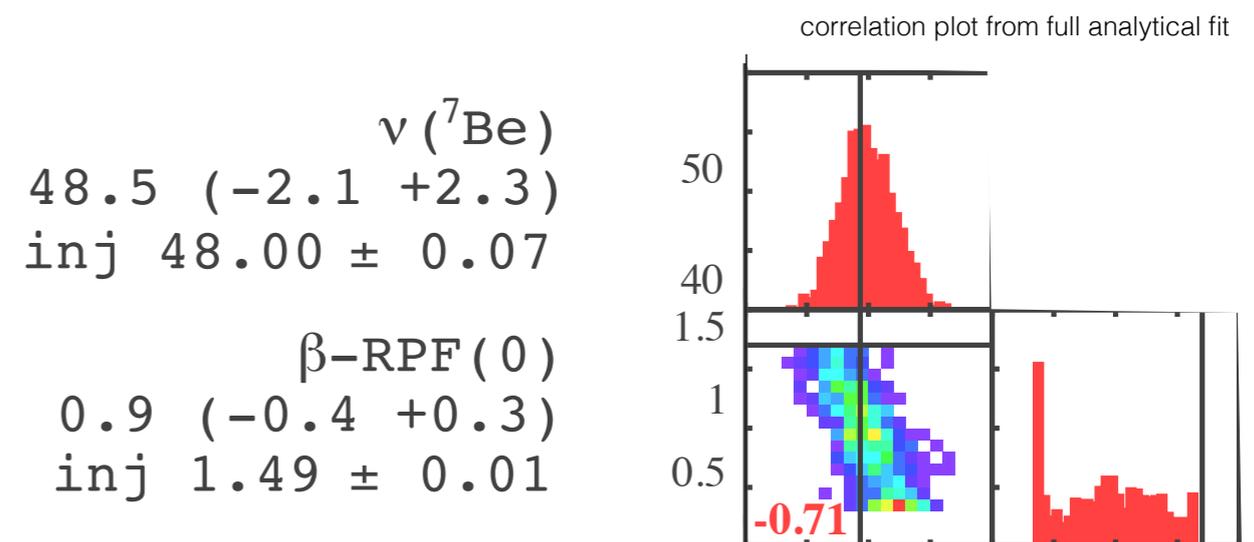
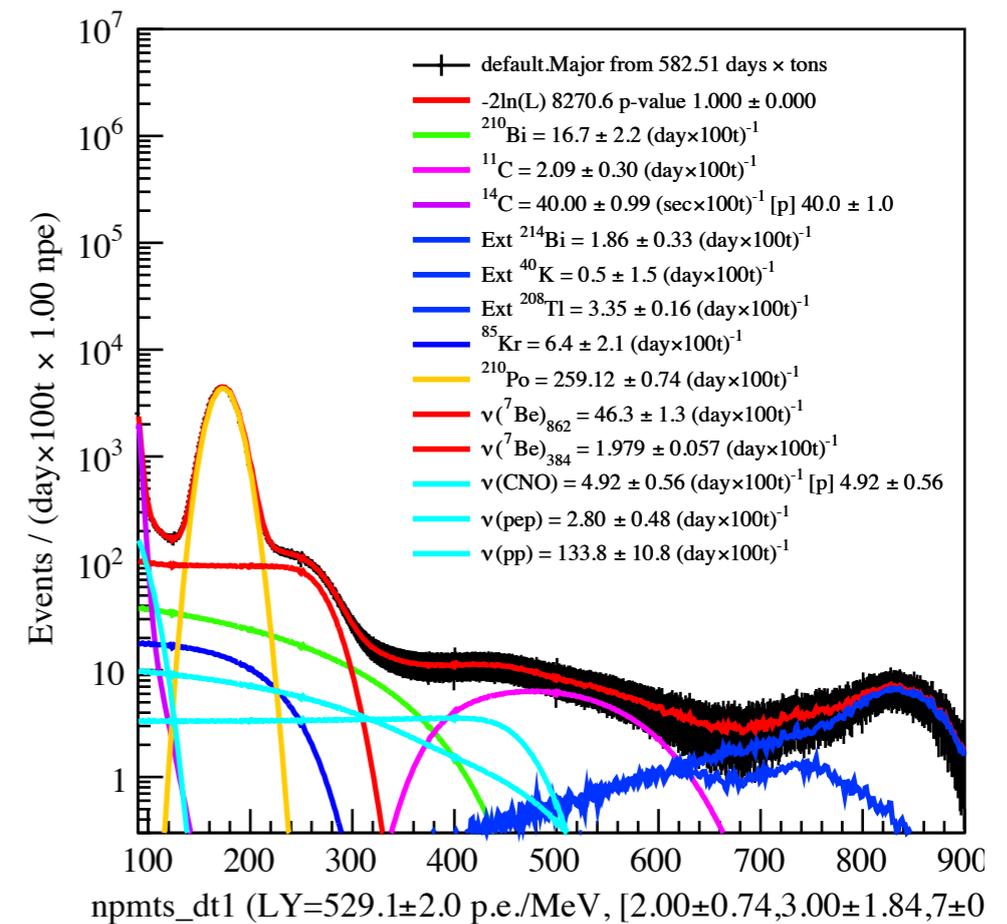


- create an ensemble of models according to experiment precision
- Fit the data with varying models through their coordinates
- Generate pseudo-experiments with varying models and take the width of distribution of best-fit

- **Full MC** during each iteration of the fit, vary **kb / absorption length spectrum etc.** **re-simulate and produce new pdf on the fly** → when one day computer is fast enough
  - ~200, 000 CPU x years per fit
- **Semi-analytical:** analytical non-linearity model + response Matrix
  - ~30 minutes per fit
- **Full analytical**
  - ~2 hours per fit

# Problem with scaling response matrix

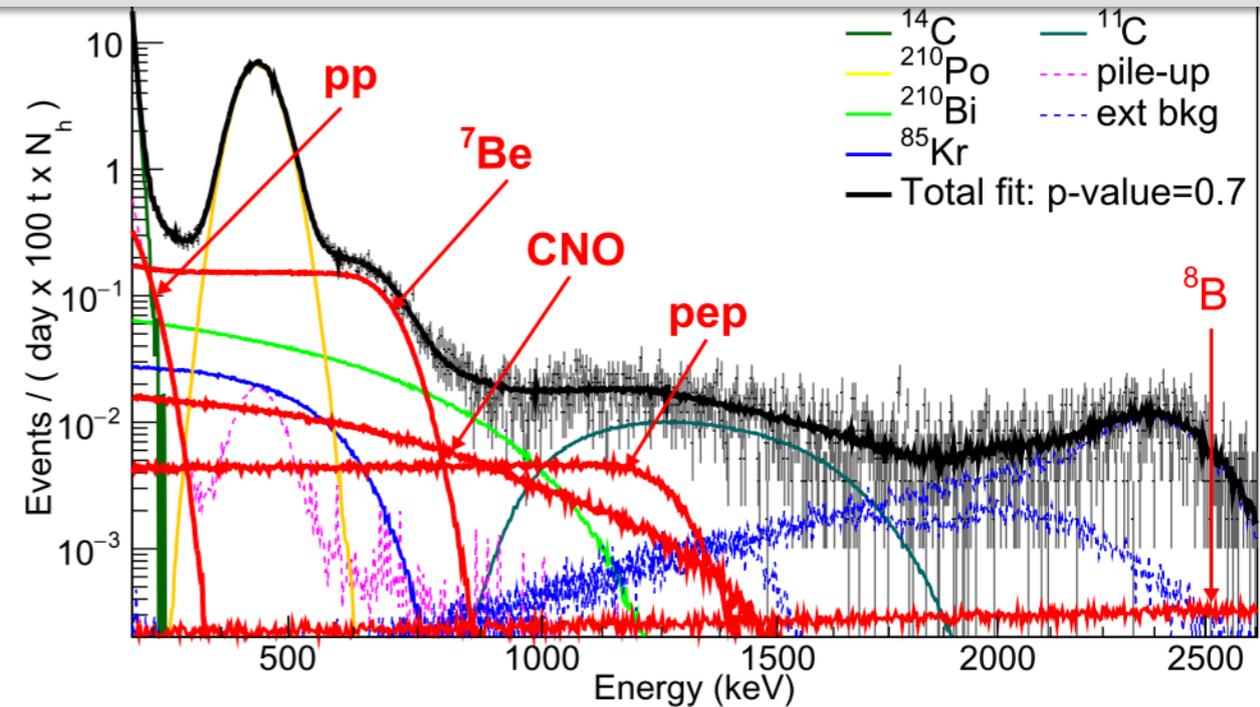
- Systematic uncertainties of LY can be included by scaling the response matrix
- This could be dangerous if the interested parameters is correlated with the resolution parameters
- When LY is scaled, the resolution should also be changed



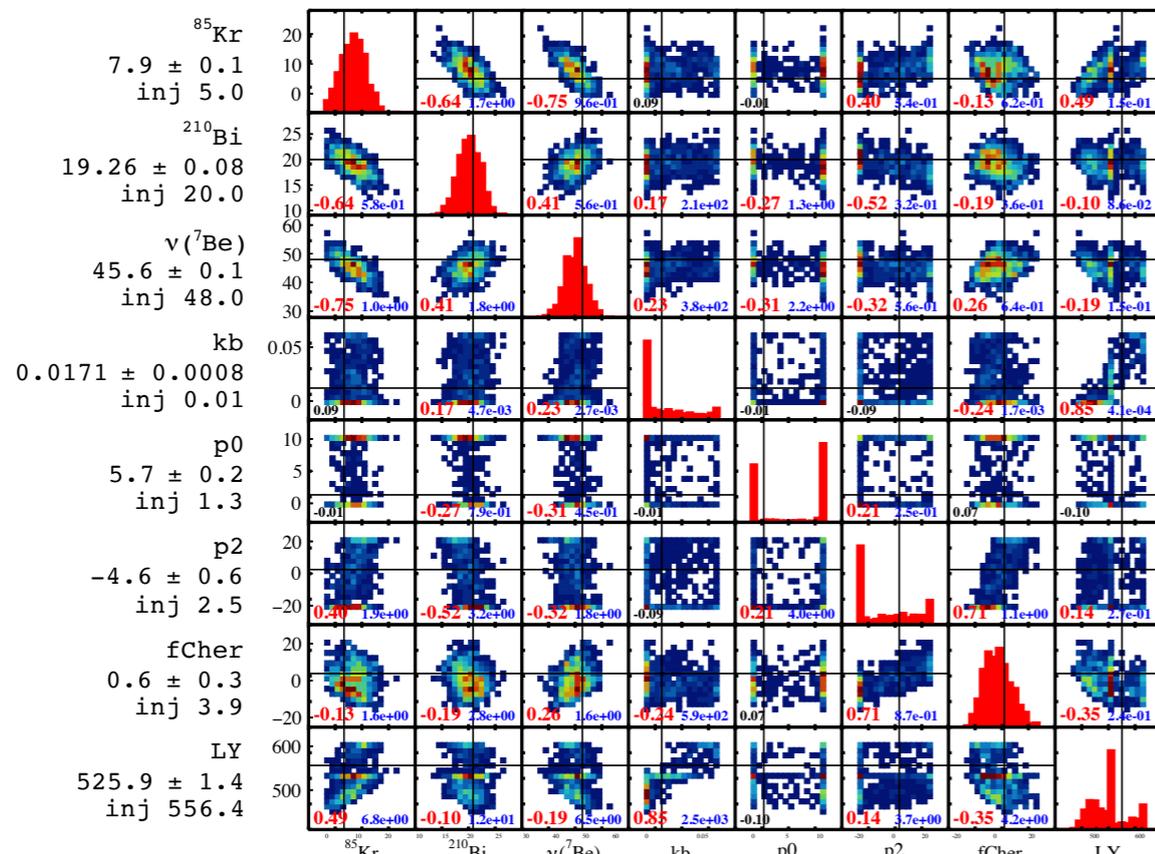
# Correlation with NL/resolution

- Using full analytical response function, we can see the correlation with detector responses

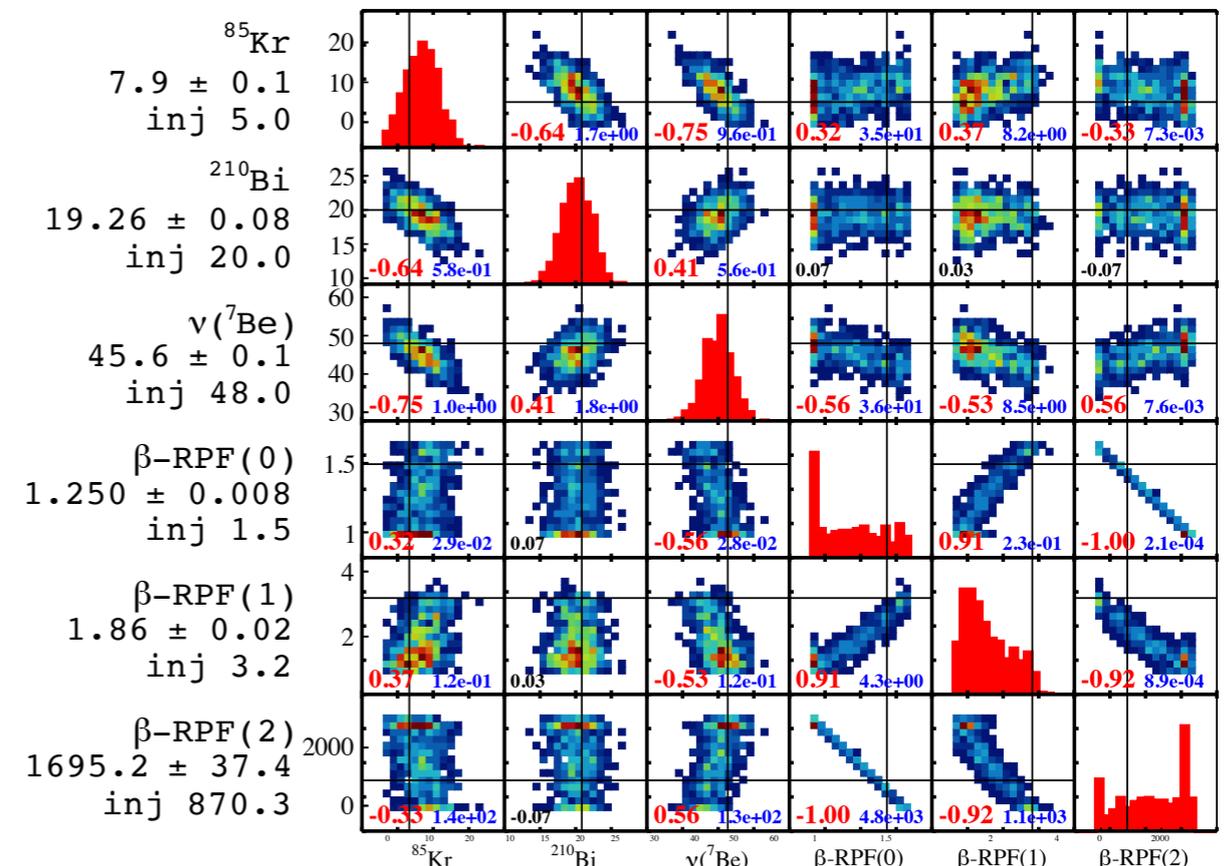
[1] M. Agostini et al., "First Simultaneous Precision Spectroscopy of  $^85\text{Kr}$ ,  $^{210}\text{Po}$ , and  $^{210}\text{Bi}$  Solar Neutrinos with Borexino Phase-II," pp. 1-8, Jul. 2017.



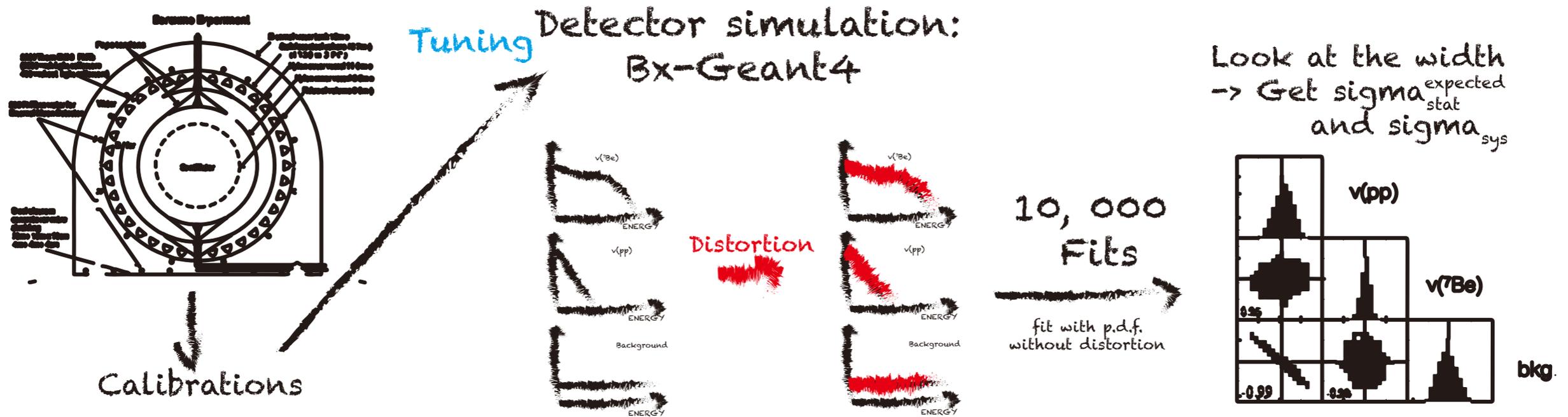
correlation with NL



correlation with res



# Method 2: Monte Carlo method



- pseudo-experiment spectra without distortion —> **statistical sensitivity**
- pseudo-experiment spectra with distortion —> **statistical systematic uncertainty**

- GooStats, a GPU based multivariate analysis framework is introduced.
- Statistical analysis module is easy to be implemented as needed. TTree/figure output provided.
- Full analytical response function has advantage that it treated the NL and resolution in a coherent way when evaluating the systematic uncertainties
- The systematic uncertainties can also be evaluated using Monte Carlo method by looking at change of width of best fit with/without distortion