Updates on the sensitivity study to BSM physics.

New and improved multi-dimensional frequentist limits & improved Bayesian models

Maros Petro

Outline

- 1. Background model
- 2. Methodology
 - 1. Feldman Cousins sensitivity
 - 2. Bayesian sensitivity

3. Results

- 1. 0
 uetaeta
- 2. $0
 uetaeta\chi^0$
- 3. $0
 uetaeta\chi^0\chi^0$
- 4. $\nu_R \nu_L \beta \beta$



The first step in calculating the sensitvity is to have a proper background model, which defines how sensitive our experiment can be!

Obtaining the background model

To obtain the background model used in this analysis I performed a series of **simulations** of various background **sources** (identified mostly from NEMO-3).

Simulation setup:

- 1. Falaise version 5.1.5
- 2. Reconstruction pipeline: MockCalibration -> Cimrman -> ChargedParticleTracker ->
 GammaClusterizer -> SNCuts -> MiModule
- 3. Magnetic field off
- 4.8% FWHM energy resolution
- 5. Foil geometry: RealisticFlat

Data-cuts (using SNCuts)

- 2 tracks
- 2 foil vertices
- 2 **distinct** associated calo-hits
- ullet $E_{sum} \in (300, 3500)$ keV
- ullet max vertex distance on foil r < 50 mm
- ullet ToF: $P_{int} \geq 4\%$ & $P_{ext} \leq 1\%$

Background sources

Simulated 100M events each:

1.
$$2
uetaeta$$
 $ightharpoonup T_{1/2}^{2
u}=9.4 imes 10^{19}$ yr

2.
214
Bi \rightarrow A = $10~\mu Bq/kg$

3.
$$^{208}\text{TI} \rightarrow A = 2~\mu Bq/kg$$

4.
40
K \rightarrow A = $58~mBq/kg$

5.
$234m$
Pa -> A = $17 \; mBq/kg$

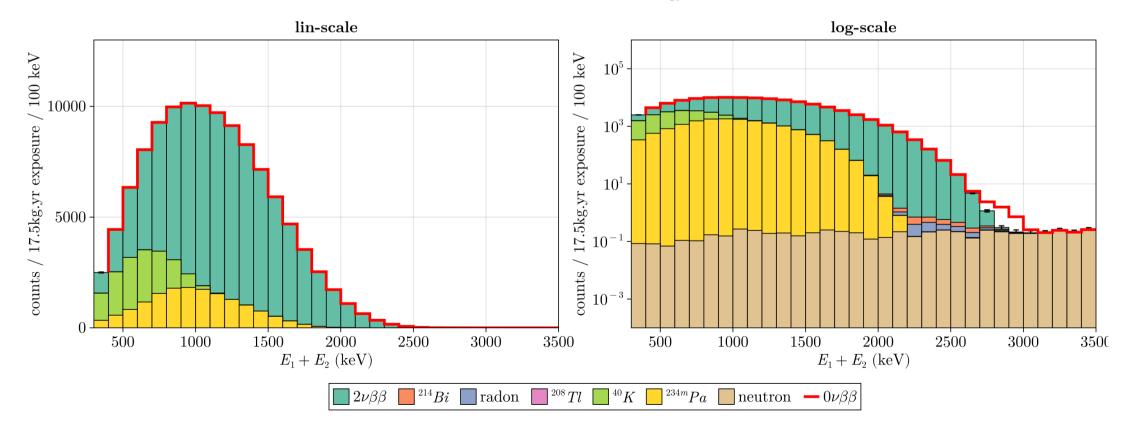
6. Radon
$$\rightarrow$$
 A = $150 \mu Bq/m^3$

Neutron data is obtained from Sam but not used everywhere cause I wasn't able to do it properly in time.

Background Model



Total background model summed 2-electron energy



Background by source

process	bkg counts
2nubb	87663.0±29.0
Bi214	8.903±0.022
Radon	10.62±0.15
Tl208	1.0227±0.0034
K40	12155.0±63.0
Pa234m	15578.0±39.0
neutron	5.91
total	115422.91±80.0

- Largest contribution due to $2\nu\beta\beta$ almost 90k events (+10k compared to old tracking)
 - Good for 2nu physics
 - Bad for onu
- K40 and Pa234m relevant **only** at lower energies
- Radon will highly depend on whether or not we have anti-Rn
- For now, "simple" neutron data



2.1. Feldman Cousins sensitvity

$$T^{1/2} = ln(2)rac{N_A\cdot m\cdot t}{W}rac{arepsilon}{\mathcal{S}(ar{b})}$$

The one you know: 1D approach!

I already gave several talks on this: DocDB#5943 and DocDB#5833

The general idea

- 1. Simulate relevant signal process
- 2. Chose 1 variable which is tracked (i.e. E_{sum} , E_i , ϕ ..)
- 3. For each possible combination of ROI calculate signal arepsilon and $\mathcal{S}(ar{b})^*$ as signal-to-background ratio $r(ROI)=rac{arepsilon}{\mathcal{S}(ar{b})}$
- 4. Find which ROI maximizes r(ROI) —> from the max(r) calculate sensitvity (simulation)
- 5. Once you've measured data -> fit in ROI to get $ar{b}$ -> calculate sensitivity
 - \circ use ε from simulation, \bar{b} from data
- * $\mathcal{S}(ar{b})$ is a Fledman Cousins limit calculated at lpha CL for total expected number of background counts in ROI $ar{b}$

- ullet ROI in this case is E^l_{sum}, E^u_{sum} , where l,u are the lower and upper boundaries of ROI
- Then sensitivity is a function of $oldsymbol{E}_{sum}$ ROI:

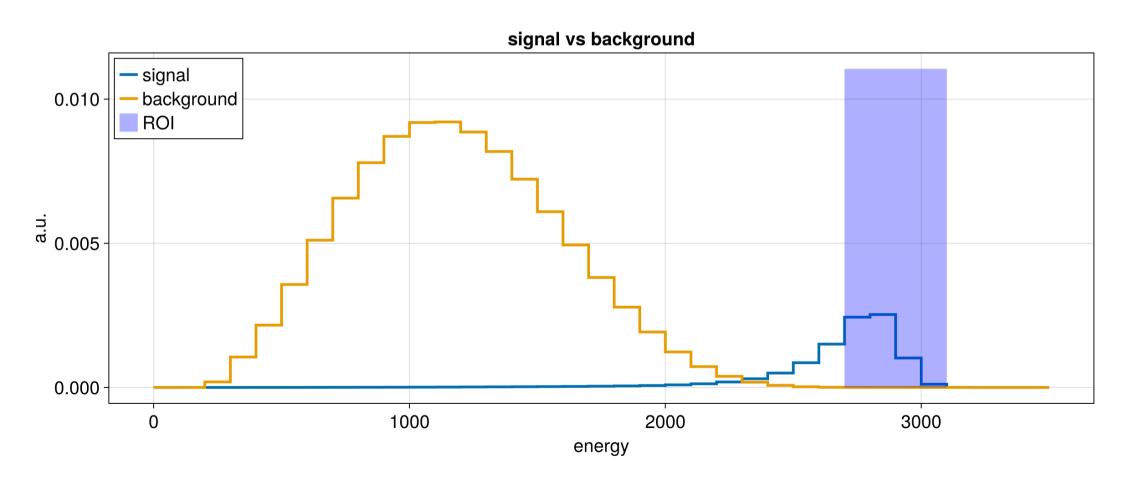
$$T^{1/2}(ROI) = T^{1/2}(E^l_{sum}, E^u_{sum}) = const. \ rac{arepsilon(E^l_{sum}, E^u_{sum})}{\mathcal{S}(ar{b}(E^l_{sum}, E^u_{sum}))}$$

ullet We calculate $T^{1/2}$ for each combination and create a 2D map:

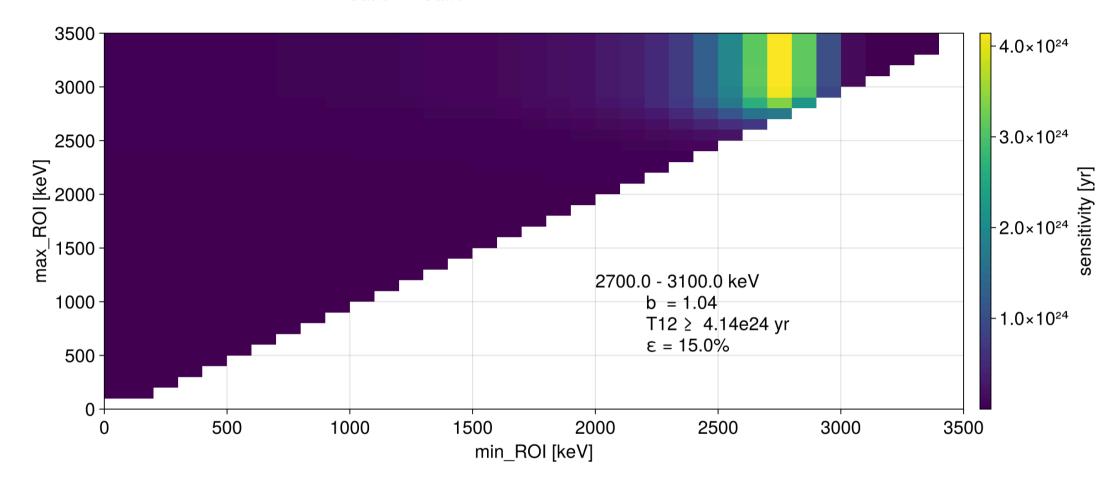
Boundaries of ROI affect signal + background differently!

That we are interested in is:

- signal efficiency arepsilon(ROI)
- ullet expected background counts $ar{b}(ROI)$



For each combination of ROI (E^l_{sum}, E^u_{sum}) we get a different value of $T^{1/2}$ —> pick maximum



get_bkg_counts(): passed isotope bb0nu_foil_bulk is a signal process!!

- best ROI is 2700-3100 keV
- ullet with $ar{b}=1.04$ and arepsilon=15 we get sensitivity of $T^{1/2}\geq 4.14\cdot 10^{24}$ yr

What if we look at another variable? Another signal process?

- For different signal processes (signal shapes!) we can be more sensitive in different channels: i.e. ϕ should be a better channel for RH spectra (will get to this later)
- Do we really need to chose only one? What if we can maximize all of them at once!

N-Dim ROI Search: the next step!

The general idea

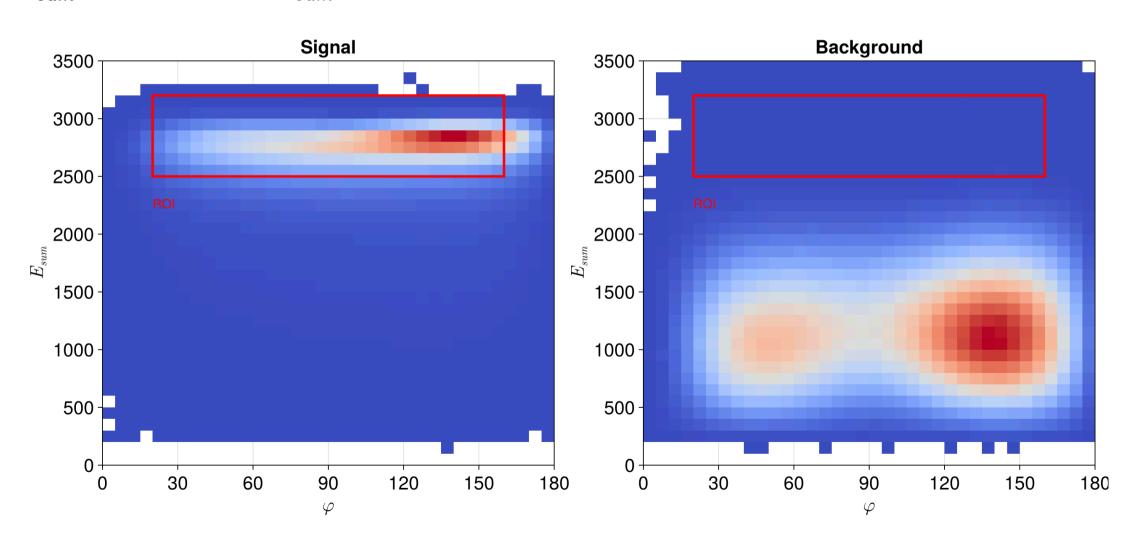
- 1. Simulate relevant signal process (again...)
- 2. Chose n variables, i.e. E_{sum} , ϕ , r
 - That means we have 6 parameters (2 for each ROI bound)
 - \circ n can be essentially any number, but computational requirements grow fast (Can be used to study optimal vertex-distance, E_{max} , E_{min} , 2D angle etc.)
- 3. For each possible combination of ROI calculate signal arepsilon and $\mathcal{S}(ar{b})^*$ as signal-to-background ratio $r(ROI)=rac{arepsilon}{\mathcal{S}(ar{b})}$
 - o •• This takes way too long and grows way too fast, when we add new dimensions ••
- 4. Find which ROI maximizes r(ROI) —> from the max(r) calculate sensitivity (simulation)
- 5. Once you've measured data \rightarrow fit in ROI to get $ar{b}$ \rightarrow calculate sensitivity
 - \circ use arepsilon from simulation, $ar{m{b}}$ from data

Extending to multiple dimensions:

Let's take the example methodology above, but use 2 variables for best ROI: E_{sum} and arphi

Again, we look at the spectrum for signal and background (now a 2D histogram):

$$E^l_{sum}$$
 = \sim 2500, E^u_{sum} = \sim 3200, φ^l = \sim 20, φ^u = \sim 160



We can see that taking a 2D ROI (basically a data-cut), we can keep a lot of signal while reducing the amount of background. In this example however, it is not exactly clear which is the **best** choice!

This approach can be extended to many dimensions!

Example for $0\nu\beta\beta$ N-Dim:

- Choose n = 3, with variables φ, E_{sum}, r
 - \circ ROI consists of 6 parameters: (E^l_{sum}, E^u_{sum}), (φ^l, φ^u) and (r^l, r^u), where l, u are the lower and upper boundaries of ROI
 - \circ let's collectively call them Θ
- Choose optimization algorithm
 - there's many libraries written for this purpose, the goal is to choose one that works well for the particular problem
 - I tried multiple, the best results seem to be found with Evolutionary Centers Algorithm ECA¹
- Then sensitivity is a function of Θ :

$$T^{1/2}(\Theta) = const. \, rac{arepsilon(\Theta)}{\mathcal{S}(ar{b}(\Theta))}$$

1https://doi.org/10.1007/978-981-13-0860-4_6

Example for $0\nu\beta\beta$ N-Dim:

Basics of ECA approach

- 1. The algorithm starts with N randomly placed initial guesses Θ_i
- 2. Calculates the values of $T^{1/2}(\Theta_i)$ for each
- 3. Identifies "best performing" Θ_i
- 4. Creates new set of N around the best performers
- 5. Repeat until convergence (or failure...)

Example for $0\nu\beta\beta$ N-Dim:

Best
$$\Theta =$$

variable	ROI		
$\boldsymbol{E_{sum}}$	(2710, 3350) keV		
arphi	(10, 180)°		
r	(0, 50) mm		

Resulting sensitivity

variable	value
arepsilon	0.145
$ar{b}$	0.73
$T^{1/2}$	$\geq 4.22 yr$

2.2 Bayesian sensitivity

$$T^{1/2} = ln(2) rac{N_A \cdot m \cdot t}{W} rac{arepsilon}{\mu_S}$$

where μ_S is the posterior 90% CI on the signal!

2.2 Bayesian sensitivity

The general idea

- 1. Simulate relevant signal and background processes
 - form truth pdfs on signal and background processes
- 2. Chose 1 variable which is tracked (i.e. E_{sum} , E_i , ϕ ..)
- 3. Create appropriate model which describes the data
 - \circ One of the model's parameters is n_S -> the expected number of signal events
- 4. Generate a pseudo-data from the simulated pdfs
- 5. Extract **pdf** of the signal counts **given data**: $pdf(n_S|data)$
- 6. Extract the 90% CI from $pdf(n_S|data)$ —> Calculate $T^{1/2}$
- 7. Repeat steps 4.-6. N times to mitigate statistical fluctuations and obtain median sensitivity
- 8. Profit 🍾 🍾

More detailed description in: DocDB#5943

Example, generic Bayesian inference

Let's say we have some Gaussian signal and Exponential background and have measured some data.

We have:

- ullet μ_{sig} mean of signal Gaussian
- Θ_{sig} proportion of signal in data
- $oldsymbol{\lambda}_{bkq}$ decay rate of background
- Θ_{bkg} proportion of background in data
- ullet For simplicity, let's say we know μ_{sig} and λ_{bkg} .
- We just want to find: Θ_{sig} , Θ_{bkg} .

 $^{^*}$ let's ignore σ for now..

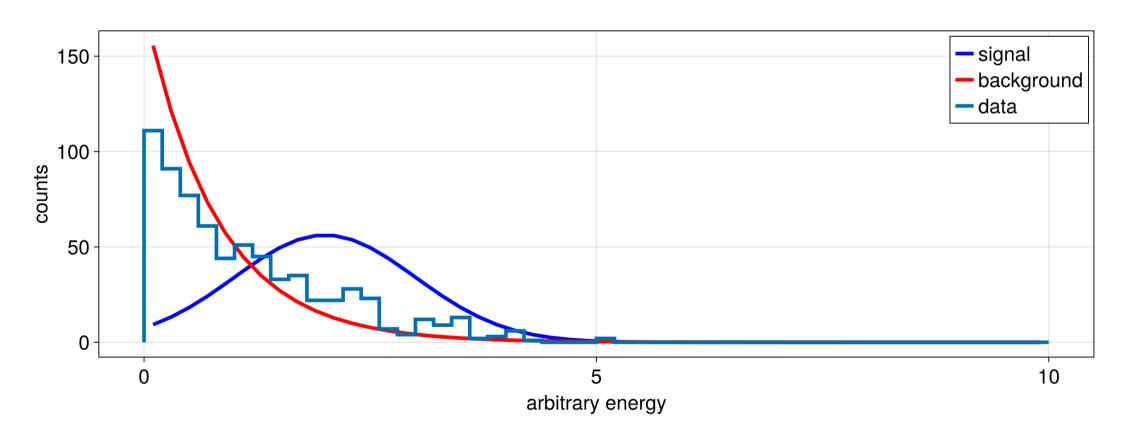
Example, generic Bayesian inference

Our likelihood:

$$\mathcal{L}(data|\Theta_{sig},\Theta_{bkg}) = \prod_{i}^{N^{obs}} \left(\Theta_{bkg}\lambda_{bkg}e^{-\lambda_{bkg}E_i} + \Theta_{sig}rac{1}{\sigma\sqrt{2\pi}}e^{-(rac{E_i-\mu_{sig}}{\sigma})^2}
ight)$$

Example experiment with n data-points:





Parameters of the experiment:

$$\mu_{sig}$$
 = \bigcirc 2.0

$$\Theta_{sig} =$$
 0.301

$$\lambda_{bkg} =$$
 0.8

$$\Theta_{bkg}$$
 = 1 - Θ_{sig} = 0.7

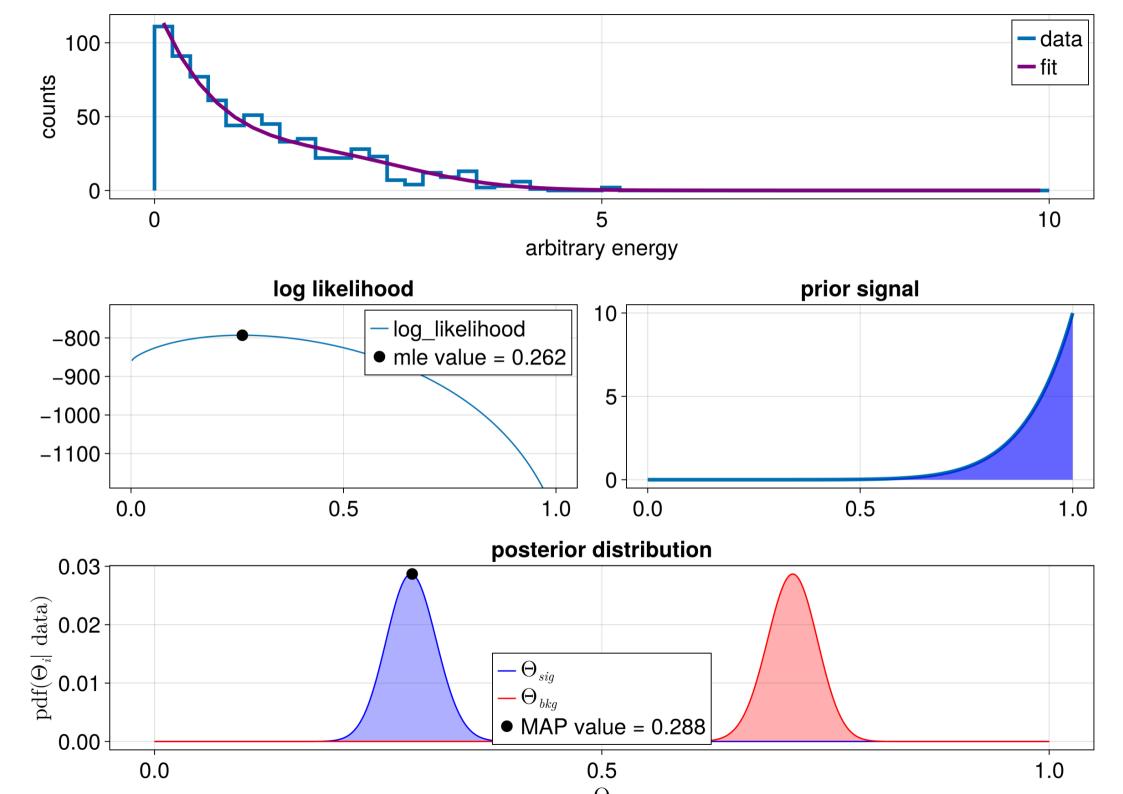
For a Bayesian inference we must chose the so-called **prior** (describes our knowledge of the paramaters)

$$pdf(\Theta) \sim ext{some distribution}$$

What we really want as the result is the so-called **posterior**: **The distribution of the parameter given data!**

Since our parameter of interest is Θ_{sig} :

$$pdf(\Theta_{sig}|data) \propto \mathcal{L}(data|\Theta_{sig},\Theta_{bkg}) \cdot pdf(\Theta_{sig})$$



What was...bad!

In DocDB#5943 I showed general model Exponential background + gaussian signal

$$\mathcal{L}(data|n_S,n_B,\lambda) = rac{1}{n_S+n_B} \prod_i^{N^{obs}} \left(n_B \lambda e^{-\lambda E_i} + n_S rac{1}{\sigma \sqrt{2\pi}} e^{-(rac{E_i-\mu}{\sigma})^2}
ight)$$

Where $\mathcal{L}(data|n_S, n_B, \lambda)$ is the likelihhod of data **given** parameters, n_S, n_B, λ are expected number of signal, background counts and decay constant. N_{obs} are the observed events.

This description works for most experiments that are not SuperNEMO!!

- Our ROI is very wide:
 - o cannot be described well with and exponential, or flat (as is the case for CUPID and Gerda)
- Our Signal is not gaussian:
 - due to energy losses in gas, the signal is more Landau shaped!

What is...better!

Detailed likelihood model with each process having its own pdf!

$$\mathcal{L}(data|ec{\Theta}) = rac{1}{\sum{\Theta_i}}\prod_i^{N^{obs}} \left(\Theta_1 pdf(sig, E_i) + \sum_{j=2}^{n_{pars}} \Theta_j pdf(bkg_j, E_i)
ight)$$

Where $pdf(sig, E_i)$ and $pdf(bkg_j, E_i)$ are the signal and background normalized spectra evaluated at E_i .

• This way we have better description of the shapes of the spectra (if we consider simulation to be correct...)

Priors and posetrior

Priors

- ullet the parameters are set up as a "proportion of the total spectrum", so each process contributes some percentage of the total spectrum: $\Sigma_i\Theta_i=1$
- ullet signal $p(\Theta_1|data) \sim Uniform(0,10^{-4})$
 - $_{\circ}~10^{-4}$ represents roughly $T^{1/2} \geq 10^{24} yr$ for 17.5 kg.~y

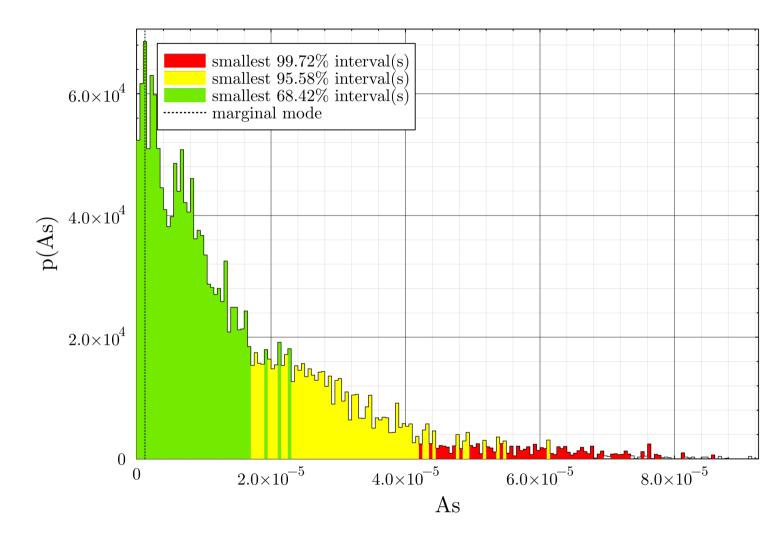
ullet backgrounds $\Theta_i \sim Uniform(0,1)$

The only condition is that $\sum \Theta_i = 1.0$.

Posterior

The posterior distributions represent the **possible distributions of the parameter** given data. That means, that for each parameter (proportion of the spectrum), we get a pdf. For example, if $0\nu\beta\beta$ represented say 30% of our measured data, we'd get a pdf with peak around 30%. If there's no signal present, the peak would be expected at 0. The shape of the pdf is how we calculate confidence interval.

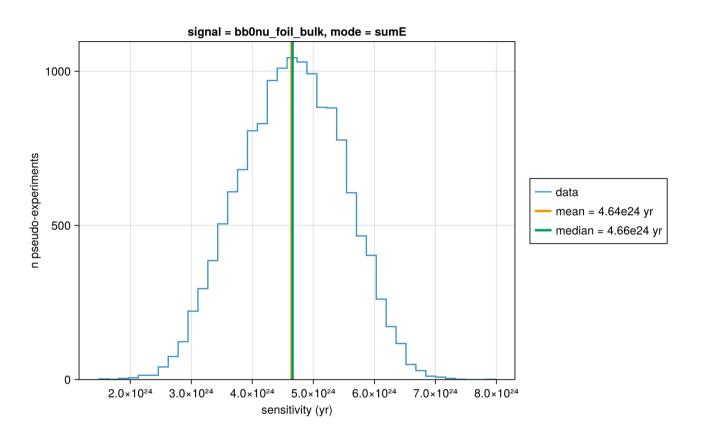
Results: For a single Bayesian inference we get posterior distribution of Θ_{sig} :



This means, the most probable value for Θ_{sig} is very close to zero, with 90% CI around 3×10^{-5} , which corresponds to 3.5 events. This is the 90% CI on the number of signal events that should be plugged into the sensitivity equation.

Danger!

The Bayesian inference method shown can fluctuate a lot! Both due to the statistical nature of the Bayesian MCMC sampling and due to how the data fluctuates (this is true for Frequentist as well). To limit the effects, we create **n pseudo experiments** to gather a large enough statistics for drawing conclusions.



I ran this algorithm in 20 parallel jobs on CC-LYON cluster for 2 days each. The following steps were within each job:

- 1. For each iteration generate pseudo-data based on the background model
- 2. Perform bayesian inference and extract **Sensitivity**
- 3. Repeat for 2 days
- 4. Save all sensitivities into a file

The resulting median sensitivity for 0
uetaeta is:

$$T_{1/2}^{0
u} \geq 4.66 imes 10^2 4 yr$$

Results

Numbers in parenthasis are with **manually** added "total neutrons in given ROI" (without performing a dedicated ND/Bayes analysis)

signal	1D	ND	Bayes
0 uetaeta	$4.14\times 10^{24}y$	$4.22(3.20) imes 10^{24}y$	$4.66 imes10^{24}y$
$0 uetaeta\chi^0$	$1.45 imes10^{23}y$	$1.48(1.45) imes 10^{23}y$	$2.38 imes10^{23}y$
$0 uetaeta\chi^0\chi^0$	$2.31 imes10^{22}y$	$2.32(2.32) imes 10^{22}y$	$1.43 imes10^{22}y$
$ u_L u_R eta eta$	$1.30 imes10^{22}y^*$	$1.30 imes10^{22}y$	$1.09 imes 10^{21} y^{**}$

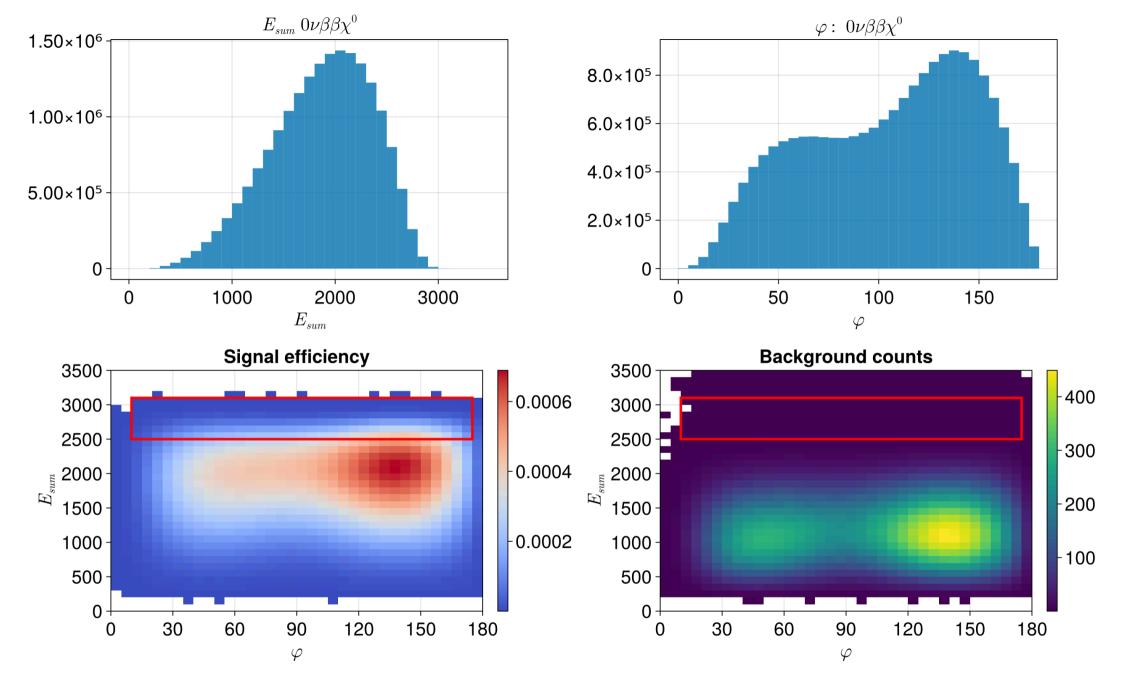
^{*} for angular distribution

^{**} for angular distribution, need to investigate why this is so low.

Results: $0\nu\beta\beta\chi^0$

ROI: $arphi \in (5,175)^\circ, E_{sum} \in (2500,3100) keV, r \in (0,50) mm$

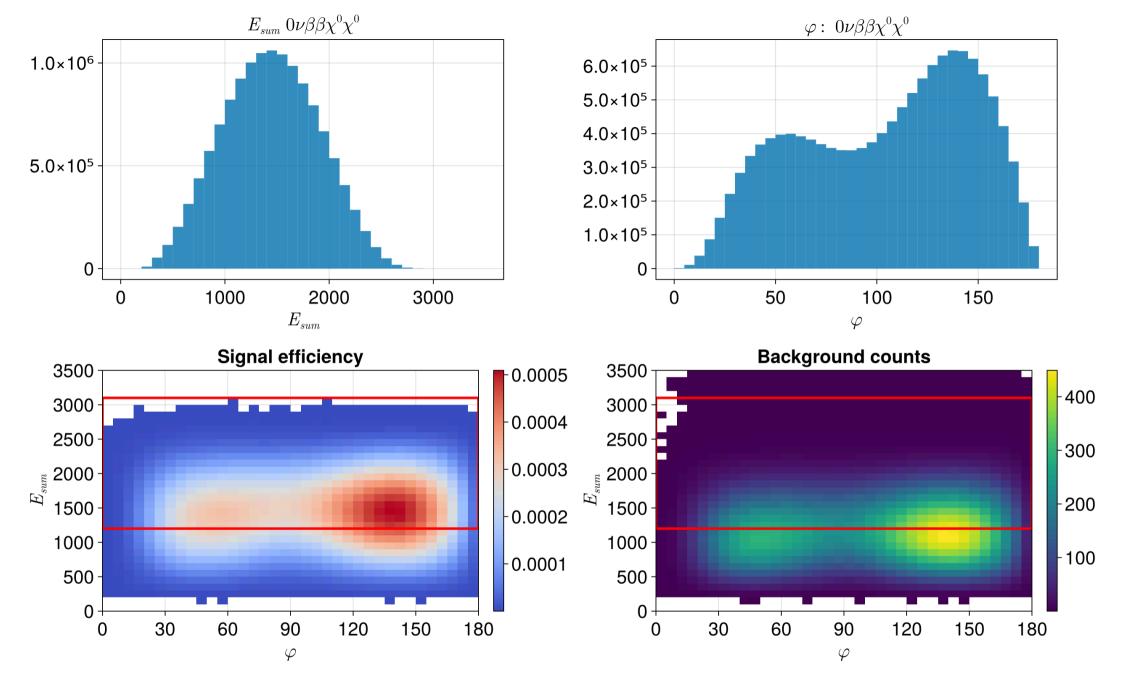
 $T^{1/2} \geq 1.48 imes 10^{23}$ y; arepsilon = 0.017; $ar{b} = 25.944~(+1.2)$



Results: $0 u\beta\beta\chi^0\chi^0$

ROI: $arphi \in (0,180)^{\circ}, E_{sum} \in (1200,3100) keV, r \in (0,50) mm$

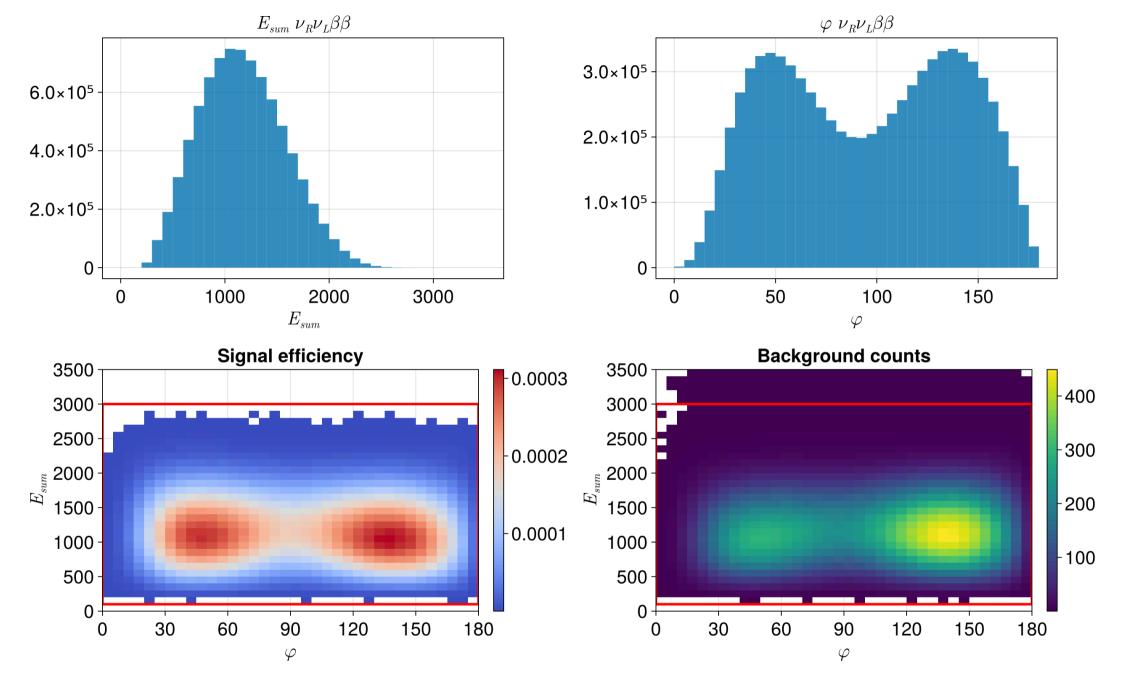
 $T^{1/2} \geq 2.32 imes 10^{22}$ y; arepsilon = 0.09; $ar{b} = 44880.87~(+3.66)$



Results: $\nu_R \nu_L \beta \beta$

ROI: $arphi \in (5,175)^\circ, E_{sum} \in (100,3000) keV, r \in (0,50) mm$. Basically take everything?

$$T^{1/2} \geq 1.30 imes 10^{22}$$
 y; $arepsilon = 0.08$; $ar{b} = 115520.50$



Conclusions

- Using n-dimensional approach (sometimes) leads to slight increase in sensitivities
 - there are issues that need investigation
 - o better optimization should improve the results even further
 - ullet The choice of data-cut on r < 50mm is way too strict!
 - Will try to optimize ToF values as well
 - Choice of variables can be extended, at the cost of computing time
 - Machine learning techniques for MVA could improve it even further
- Bayesian approach is more sensitive than frequentist when signal shape is more different from backgrounds!
 - o not sure what's wrong with RH spectra, but this could be it?
- Need to implement neutron data
- We have an idea to use data-partitioning for angular distribution

Future

Based on research of angular distribution for the electrons, depending on which side of foil they are emitted, the reconstruction is differently efficient. We could use this fact!

