



multimessenger/supernova*

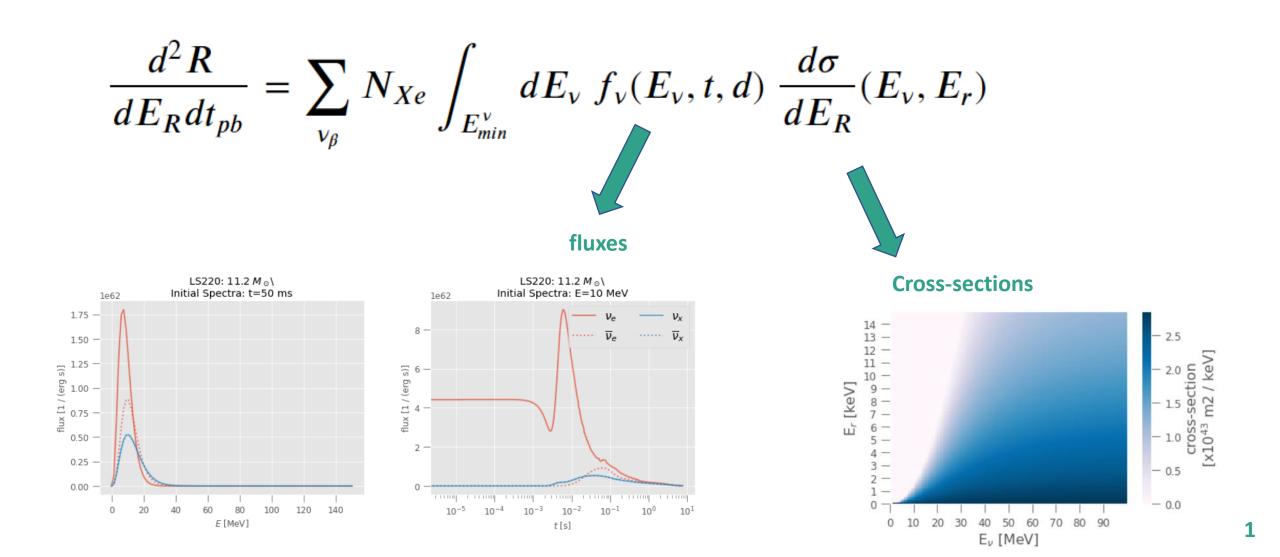
Tool for calculating CEvNS interactions

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SNEWS Collaboration Meeting 05.08.2022 Thursday



* cause we are running out of abbreviations



A = sn.Models(model_name='Fornax_2021')

- [0] lum spec 12M r10000 dat.h5 lum_spec_13M_r10000_dat.h5 [1]lum spec 14M r10000 dat.h5 [2] lum spec 15M r10000 dat.h5 [3] [4] lum spec 16M r10000 dat.h5 lum spec 17M r10000 dat.h5 [5] [6] lum spec 18M r10000 dat.h5 [7] lum_spec_19M_r10000_dat.h5 lum spec 20M r10000 dat.h5 [8] lum_spec_21M_r10000_dat.h5 [9] [10] lum spec 22M r10000 dat.h5 lum spec 23M r10000 dat.h5 [11] lum_spec_25M_r10000_dat.h5 [12] lum spec 26.99M r10000 dat.h5 [13] [14] lum spec 26M r10000 dat.h5
- Fluxes can be fetched from any snewpy model

- Object oriented
- Flexible (change the input composite)
- Scalable by the distance
- Computes rates per time,

 $\frac{d^2 R}{dE_R dt_{pb}} = \sum_{v_a} N_{Xe} \int_{E_{min}^v} dE_v f_v(E_v, t, d) \frac{d\sigma}{dE_R}(E_v, E_r)$

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- Fluxes can be fetched from any snewpy model

- Object oriented
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- Computes rates per time, rates per recoil energy

 $\frac{d^2 R}{dE_R dt_{pb}} = \sum_{v_{\theta}} N_{Xe} \int_{E_{min}^v} dE_v f_v(E_v, t, d) \frac{d\sigma}{dE_R}(E_v, E_r)$

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- Fluxes can be fetched from any snewpy model

- Object oriented
- Flexible (change the input composite)
- Scalable by the distance
- Computes rates per time, rates per recoil energy per neutrino flavor

 $\frac{d^2 R}{dE_R dt_{pb}} = \sum_{\nu_{\beta}} N_{Xe} \int_{E_{min}^{\nu}} dE_{\nu} f_{\nu}(E_{\nu}, t, d) \frac{d\sigma}{dE_R}(E_{\nu}, E_r)$

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- Fluxes can be fetched from any snewpy model

- Object oriented
- Flexible (change the input composite)
- Scalable by the distance
- Computes rates per time, rates per recoil energy per neutrino flavor, per isotope

 $\frac{d^2 R}{dE_R dt_{pb}} = \sum_{\nu_o} N_{Xe} \int_{E_{min}^{\nu}} dE_{\nu} f_{\nu}(E_{\nu}, t, d) \frac{d\sigma}{dE_R}(E_{\nu}, E_r)$

CEvNS int

A = sn.Models(model_name='Forna

> Available files for this mode

[0] lum spec 12M r10000 dat [1]lum_spec_13M_r10000_dat lum spec 14M r10000 dat [2] lum_spec_15M_r10000 dat [3] [4] lum spec 16M r10000 dat lum spec 17M r10000 dat [5] [6] lum spec 18M r10000 dat [7] lum spec 19M r10000 dat lum spec 20M r10000 dat [8] lum_spec_21M_r10000_dat [9] [10] lum spec 22M r10000 dat lum spec 23M r10000 dat [11] lum_spec_25M_r10000_datated dat [12] lum spec 26.99M r10000 [13] [14] lum spec 26M r10000 dat

Fluxes can be fetched fron

Atom.py [🛃 Supernova_Models.py [📥 Recoil_calculations.py 🔾 Plotter.pv Xe134 = { 'Type' : 'Xe134', : 134, 'AtomicNum' : 54, : 133.9053945, 'Spin' 'Fraction' : 0.104357 Xe136 = { 'Type' : 'Xe136', : 136, 'AtomicNum' : 54, : 135.907219, 'Spin' 'Fraction' : 0.088573 $ATOM_TABLE = {$ 'Xe124' : Xe124, 'Xe126' : Xe126, 'Xe128' : Xe128, 'Xe129' : Xe129, 'Xe130' : Xe130, 'Xe131' : Xe131, 'Xe132' : Xe132, 'Xe134' : Xe134, 'Xe136' : Xe136

composite)

rates per recoil energy pe

 $dE_{\nu} f_{\nu}(E_{\nu}, t, d) \frac{d\sigma}{dE_{R}}(E_{\nu}, E_{r})$ E_{min}^{ν}

CEvNS int

A = sn.Models(model name='Forna

> Available files for this mode

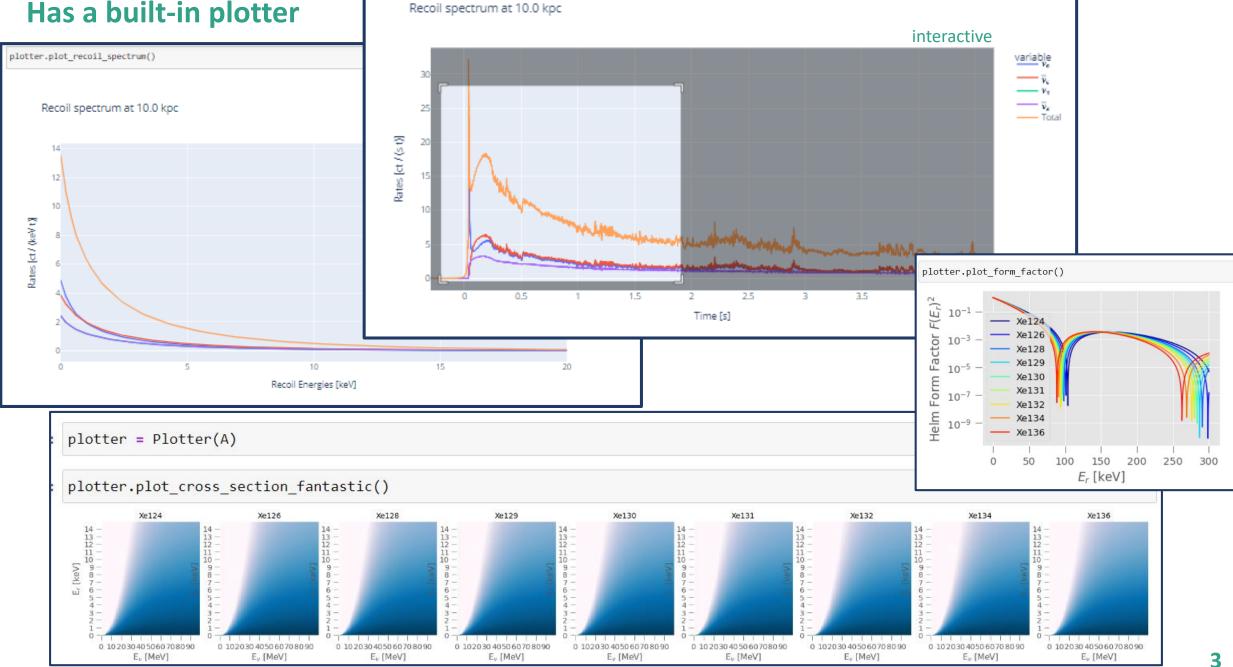
[0] lum_spec_12M_r10000_dat lum_spec_13M_r10000_dat [1] [2] lum_spec_14M_r10000_dat lum_spec_15M_r10000_dat [3] lum spec 16M r10000 dat [4] lum spec 17M r10000 dat [5] [6] lum_spec_18M_r10000_dat lum_spec_19M_r10000_dat [7] lum_spec_20M_r10000_dat [8] lum_spec_21M_r10000_dat [9] [10] lum spec 22M r10000 dat lum_spec_23M_r10000_dat [11] lum_spec_25M_r10000_datated dat [12] lum spec 26.99M r10000 [13] lum spec 26M r10000 dat [14]

Fluxes can be fetched fron

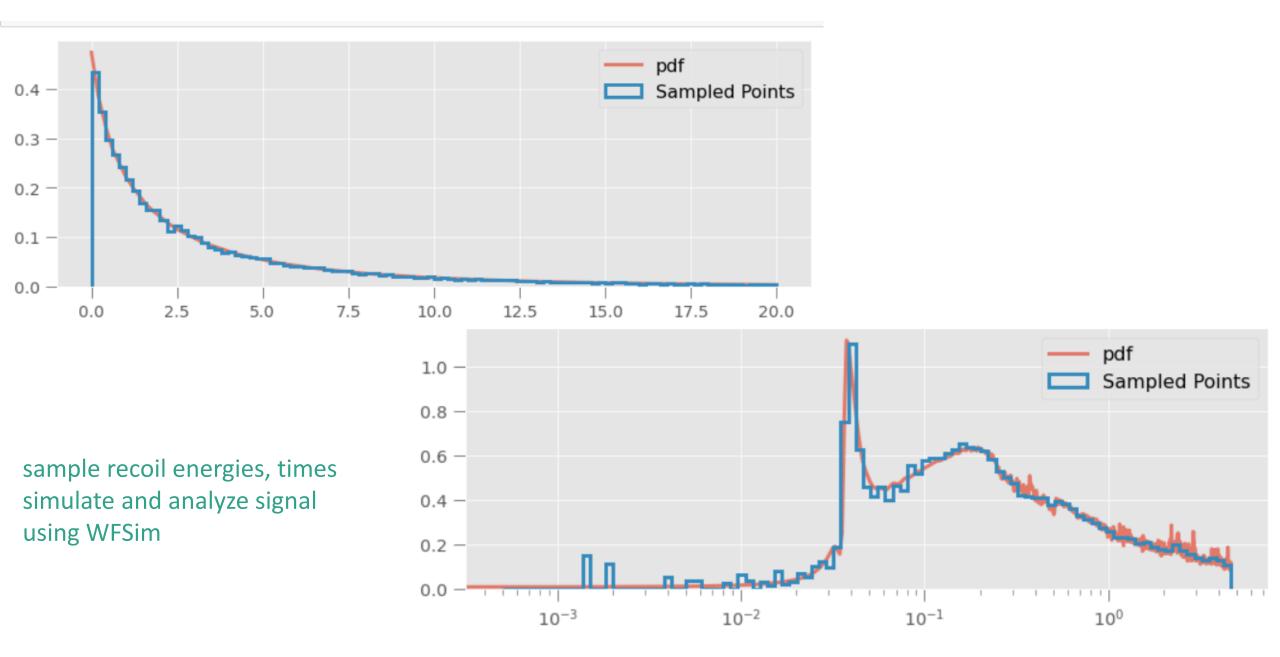
enon_Atom.py 🗡 🐔 Supernova_Models.py 🗡 🐔	Recoil_calculations.py × 🛛 👸 Plotter.py
-Xe134 = {	
'Type' : 'Xe134', 'MassNum' : 134,	
'AtomicNum' : 54,	
'Mass' : 133.9053945,	
'Spin' : 0,	
'Fraction' : 0.104357	
₽ }	
	computed, can always be fetched in < 1s
A = sn.Models(model_name='Forna	ax_2021', index=5)
A.compute_rates();	
Computing for all isotopes: 100%	9/9 [00:00<00:00, 457.87it/s]
100%	9/9 [00:00<00:00, 46.67it/s]
100%	9/9 [00:00<00:00, 46.67it/s] 9/9 [00:00<00:00, 44.24it/s]
100%	
100%	
100% ///////////////////////////////////	
100% 'Xe120 : Xe120, 'Xe128' : Xe128, 'Xe129' : Xe129,	
100% ///////////////////////////////////	
100% 'Xe120 : Xe120, 'Xe128' : Xe128, 'Xe129' : Xe129, 'Xe130' : Xe130,	9/9 [00:00<00:00, 44.24it/s]
100% 'Xe128' : Xe128, 'Xe129' : Xe128, 'Xe130' : Xe130, 'Xe131' : Xe131,	9/9 [00:00<00:00, 44.24it/s]
100% 'Xe120 . Xe120, 'Xe128' : Xe128, 'Xe129' : Xe129, 'Xe130' : Xe130, 'Xe131' : Xe131, 'Xe132' : Xe132,	

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Has a built-in plotter



In XENONnT ...

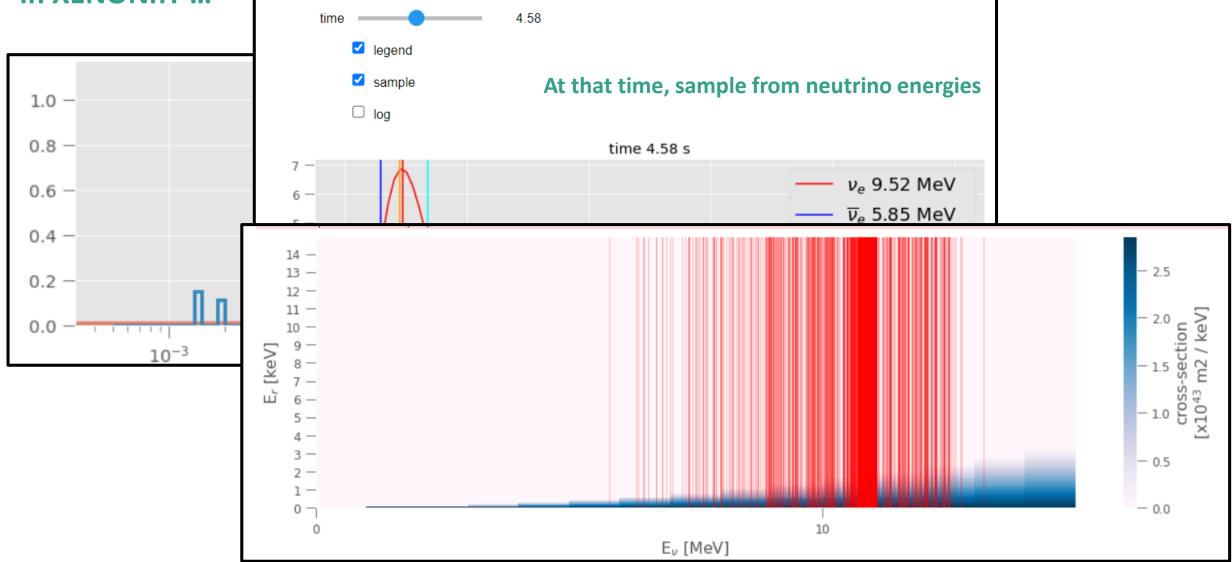


Summary:

- An efficient way of computing interaction rates For any snewpy model at any given time for any given recoil energy from any given <u>neutrino energy</u> for any given atom/isotope for all neutrino flavors



In XENONnT ...



At that neutrino energy sample a recoil energy based on cross-section probability

This method actually gives you a terrible spectra Regardless if you sample from time first (almost only returns first second) Or sample time uniformly and look at the energy distributions.

Although you get those neutrinos more, they don't recoil energetic enough. So you only see the high energy ones (and there are still a lot 10^12)

It makes more sense to study what we can see, and energy neutrino can detect those. Then looking into flux distribution to see what models can generate enough of those events, and at what times

Sampling from time integrated recoil energy distribution is probably the best approximation to this.