

**Improved description of ion stopping power in  
*compounds* in MARS code**

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# Outline

- Motivation
- Formalism
- Comparisons with measurements

# Motivation: Does it really matter?

## Does it matter for $\mu 2e$ ?

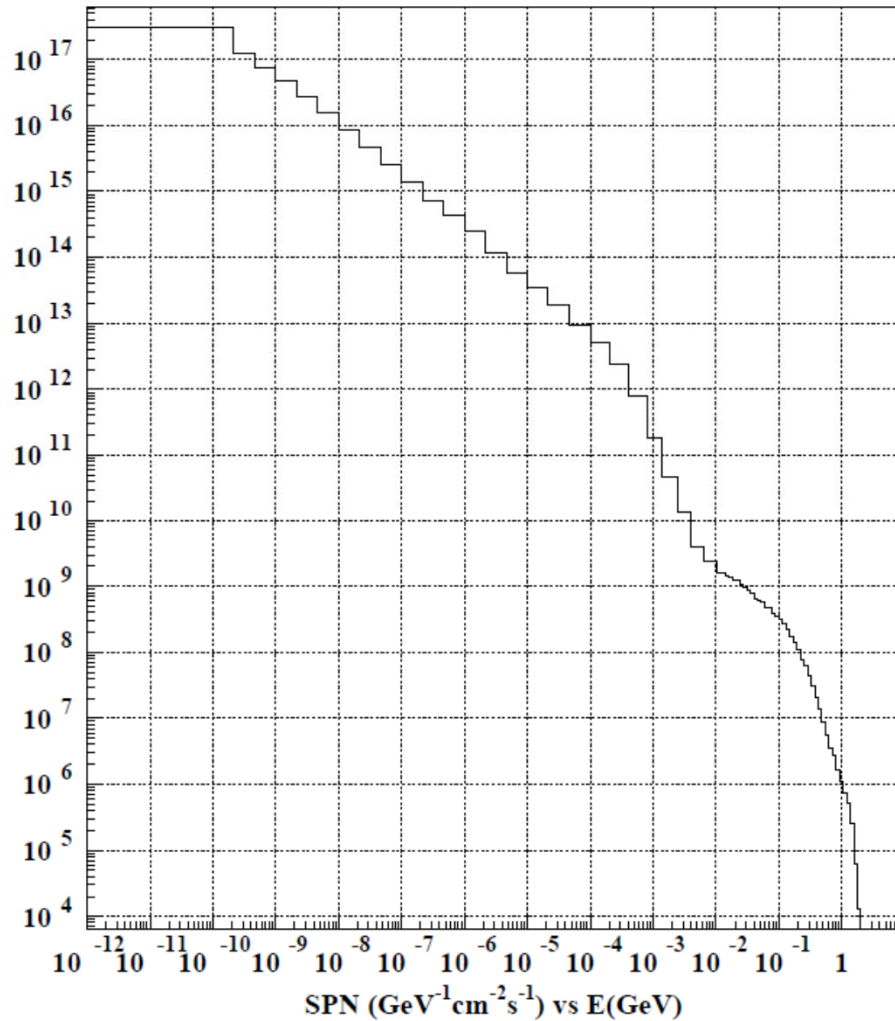
- Accelerator magnets usually contain insulation materials such as **G10, kapton, epoxy** (most radiation prone) → all are compounds.
- Stopping power  $dE/dx$  for compounds is usually described using **Bragg's rule**:

$$\frac{dE}{dx} = \sum_i \left( \frac{dE}{dx} \right)_i$$

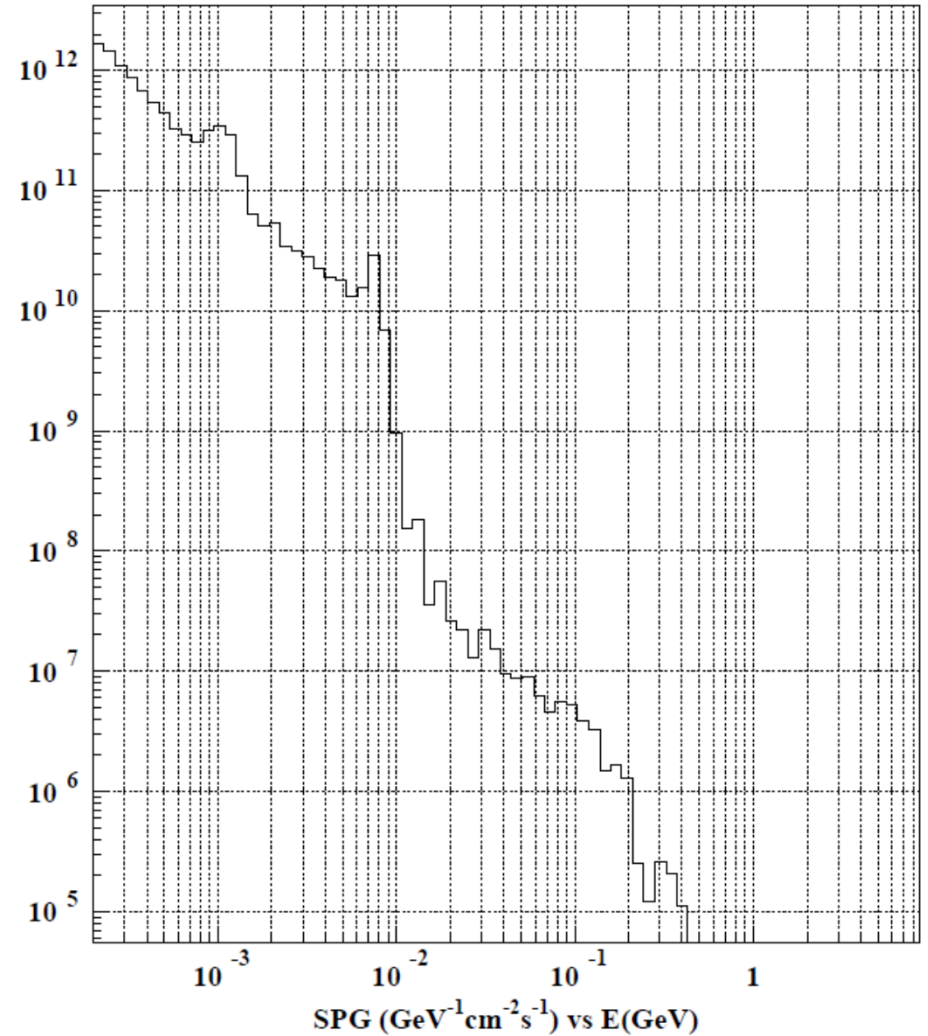
- At low energies,  $\sim 100$  keV/amu, for **low-Z materials** ( $Z < 12$ , especially **hydrocarbons**) the recipe introduces errors, and the difference between theory and experiment can be as large as **20%**.
- Let us see what energy spectra we get in Production Solenoid coil.

Vitaly Pronskikh provided particle energy spectra in the **hottest spot** of the Production Solenoid coil calculated at  $\sim 1/3$  nominal power ( $6.7 \times 10^{12}$  p/s)

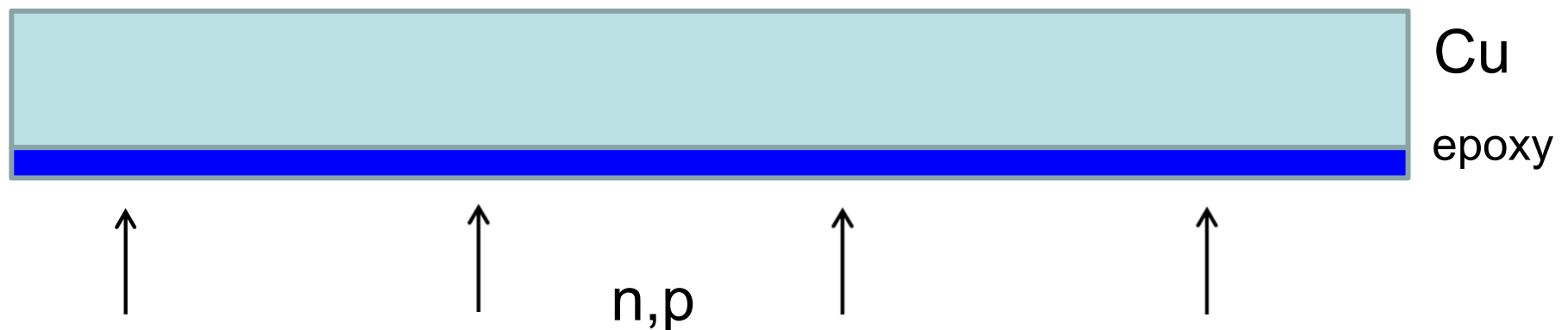
### Neutrons



### Gammas

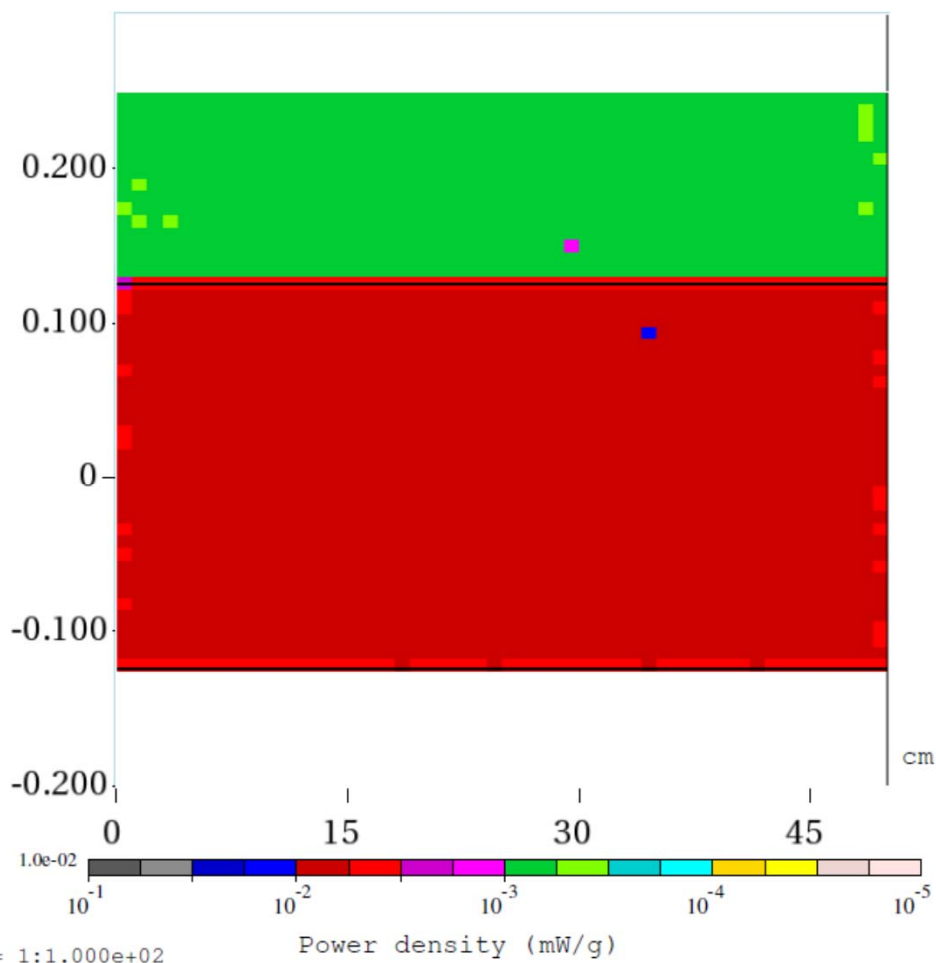


- Comparison between partial contributions to **energy deposition** (absorbed dose) in an epoxy layer (ethylene oxide  $C_2H_4O$ ) 2.5 mm thick.
- $\phi_n = 5.5 \times 10^9 \text{ cm}^{-2}\text{s}^{-1}$ ,  $\phi_p = 4.7 \times 10^5 \text{ cm}^{-2}\text{s}^{-1}$
- Separate calculation for a simple model in order to get a more reliable result in the thin epoxy layer.

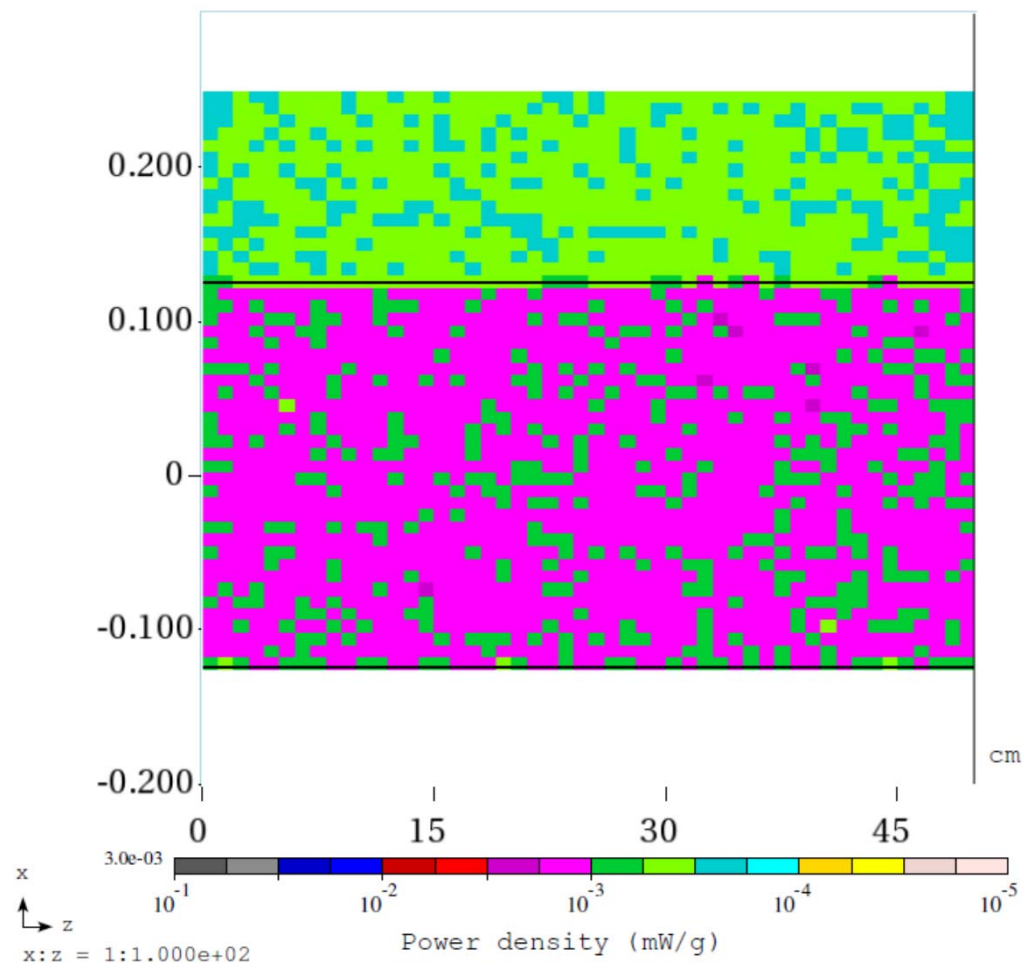


# Power density (arbitrary normalization)

## Total

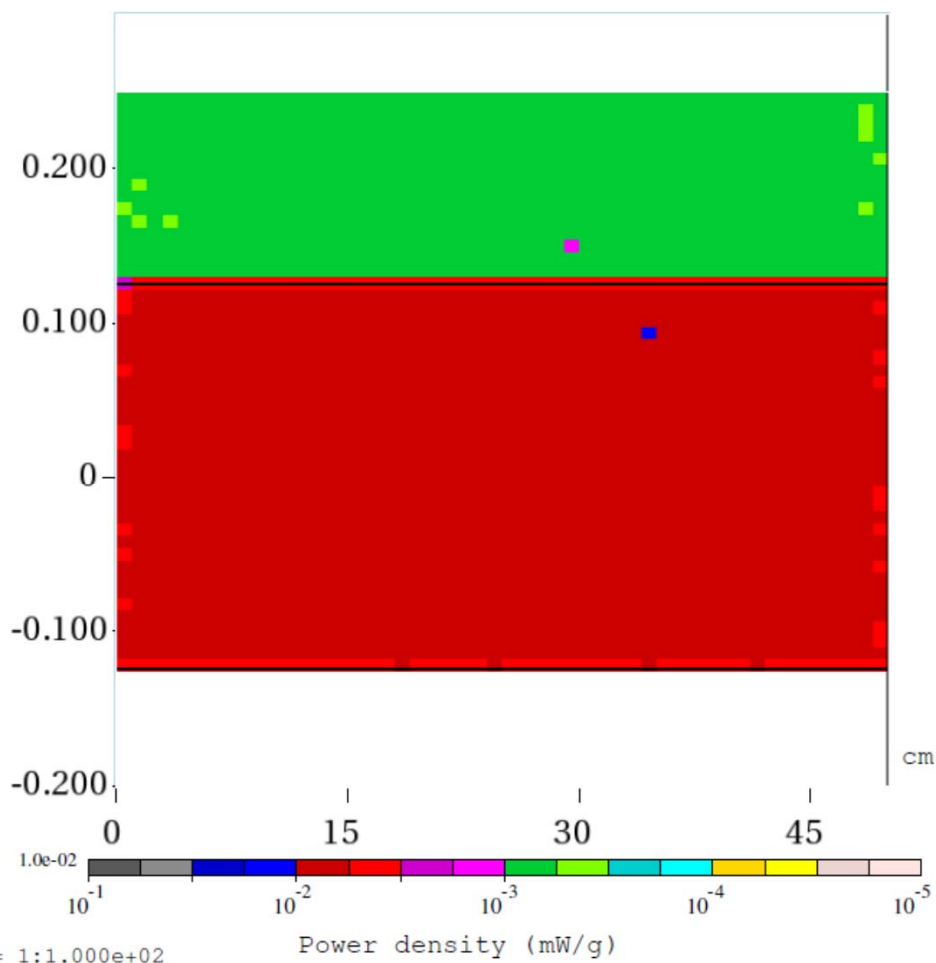


## Neutron contribution

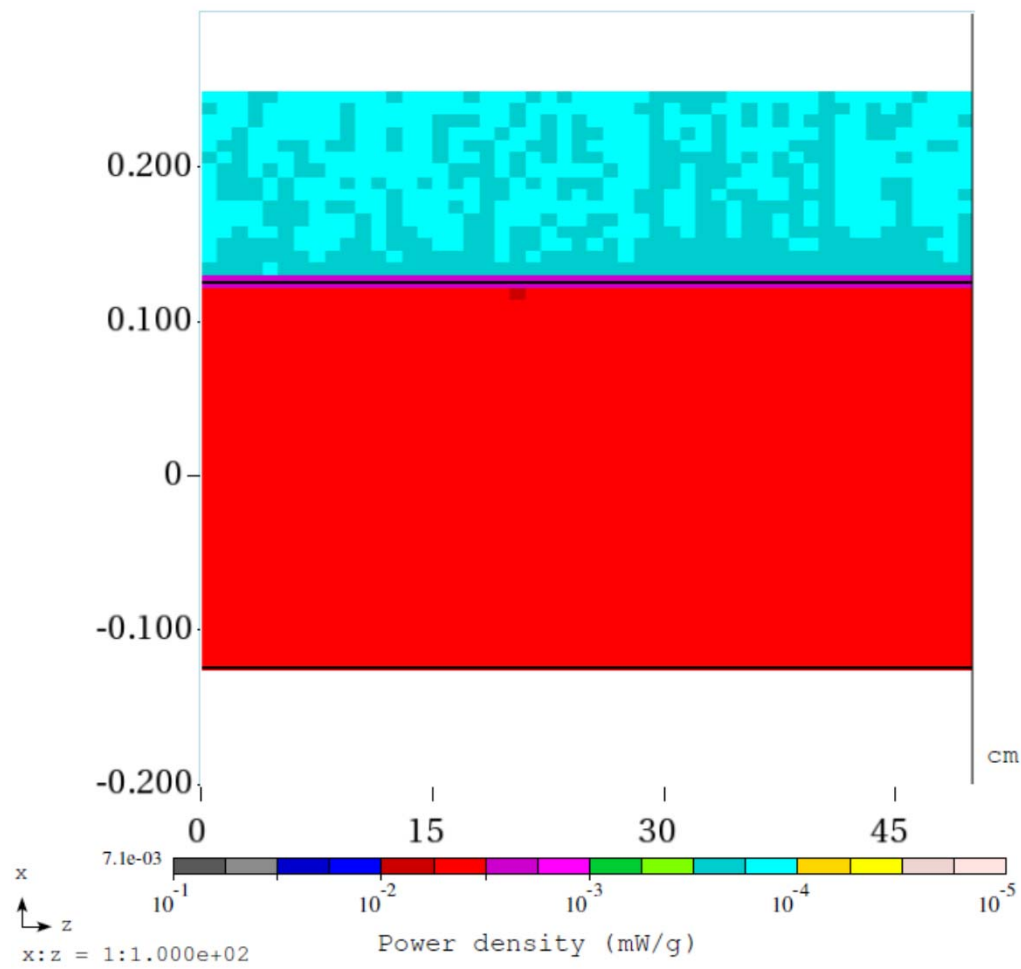


# Power density (arbitrary normalization)

## Total



## Proton contribution



# Numerical data on average power density

- Total  $\sim 8 \times 10^{-3}$  mW/g
- Neutron contribution  $\sim 1 \times 10^{-3}$  mW/g
- Proton contribution  $\sim 7 \times 10^{-3}$  mW/g
  
- Neutrons ( $\sim 90\%$  with energy lower than 0.5 MeV) transfer energy mostly in **elastic collisions** with target nuclei. ***Small fraction*** of neutron kinetic energy is transferred to nuclei like C, O.



# Stopping power in compounds (a little boring chemistry)

- There is no need to invent the wheel: the **Cores-and-Bonds** approach (CAB) was developed in 80s by G. Both *et al.*, Köln University.

- $S_{\text{ion}}^{Z_{\text{eff}}} \rightarrow S_{\text{p}} \rightarrow S_{\text{p,Bragg}}, S_{\text{p,Bragg}}(125 \text{ keV}) \text{ and } S_{\text{p,CAB}}(125 \text{ keV})$

- $S_{\text{p,CAB}}(125 \text{ keV}) = \sum \text{Cores} + \sum \text{Bonds}$

- Cores: **atoms** from H to Cl.
- Bonds: **single** like H-H, C-H *etc*, **double** like C=C, C=O *etc*, **triple** like C≡C, O≡O.

# Stopping power in compounds (a little boring chemistry)

- All chemical bonds in a compound (may be tricky for some compounds).
- CAB developers claim that such an approach provides accuracy of about 1% (difference between measurement and theory). More than 100 comparisons for a variety of compounds.
- More info at [www.srim.org](http://www.srim.org)  
Computer code by J. Ziegler et al.,  
**Stopping and Range of Ions in Matter**

# Implementation & comparisons

- The CAB method was implemented in MARS code recently for **several** most important **compounds**: water, polyethylene, polystyrene, epoxy.
- Currently MARS users can define composition of compounds using input file. It might be makes sense to provide a possibility to enter the CAB data in the input file if a user wants to do so.

