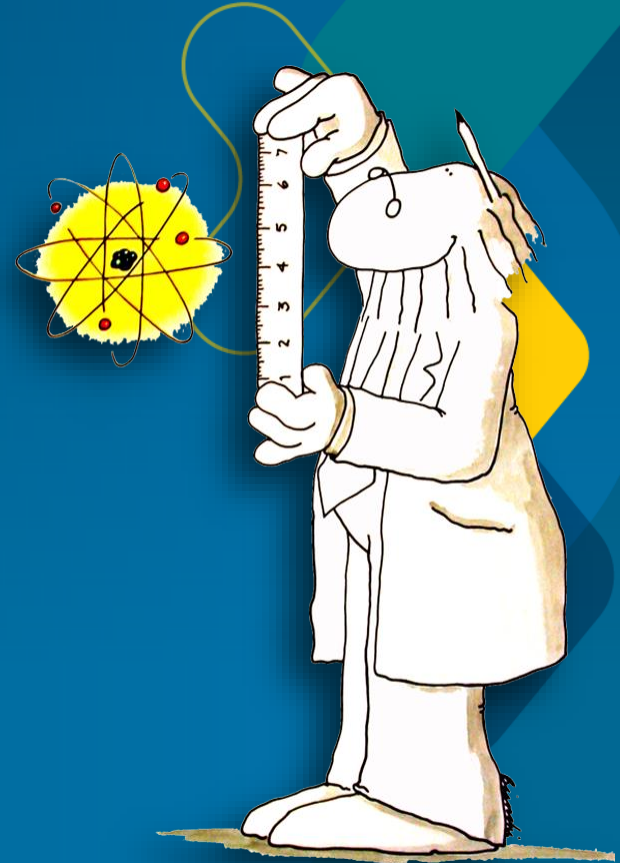


EGSnrc update and Monte Carlo simulation verification

Frédéric Tessier

Measurement Science and Standards
National Research Council Canada





EGSnrc update and Monte Carlo simulation verification

Ernesto Mainegra-Hing

Atomic relaxation and PE cross sections

Frederic Tessier

Reid Townson

Measurement Science and Standards

Radionuclide decay modelling

National Research Council Canada

Dave Rogers

Improved kerma calculations



***“This talk is almost, but not quite,
entirely unlike a scientific presentation.”***

— paraphrasing Douglas Adams

Team

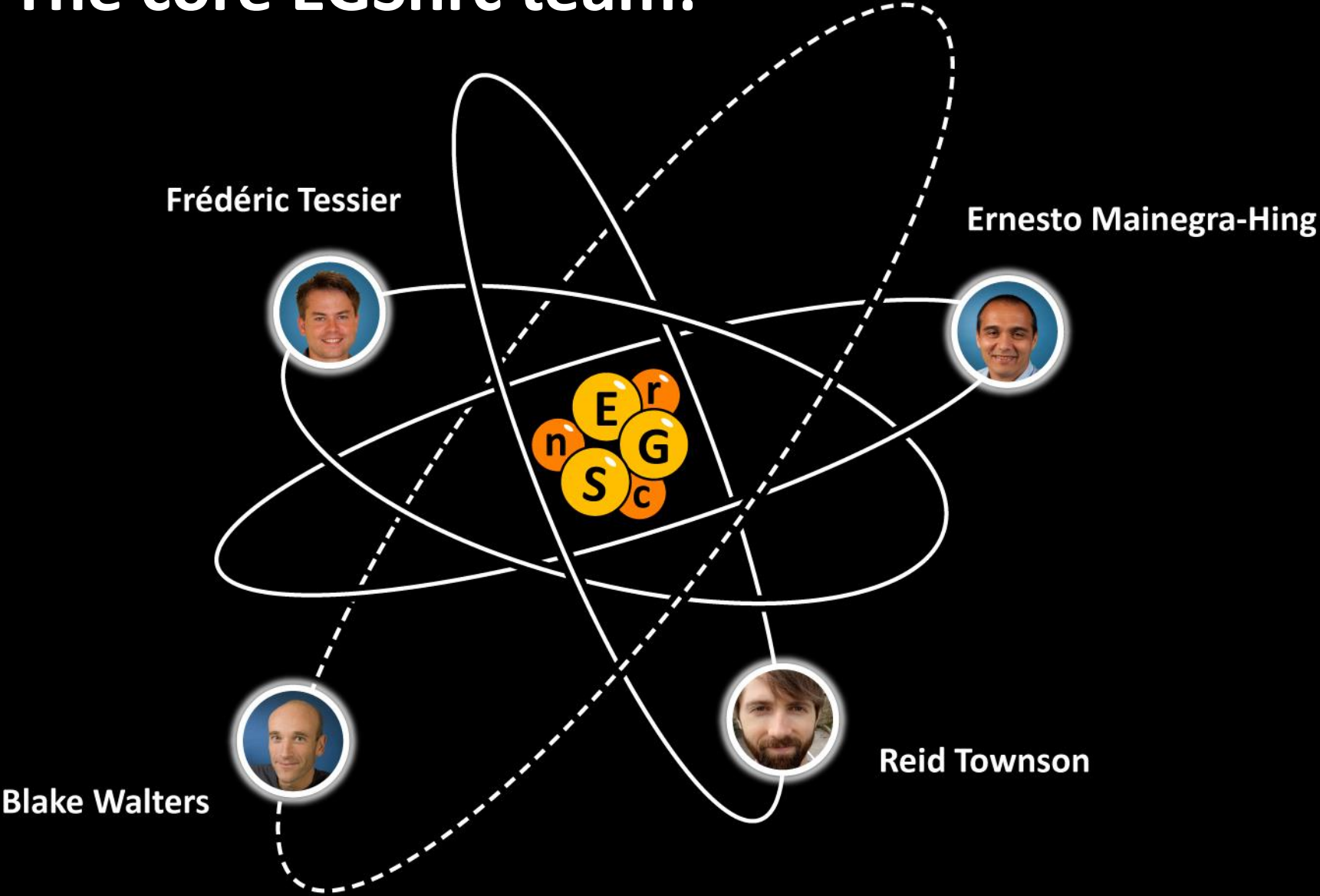
Code

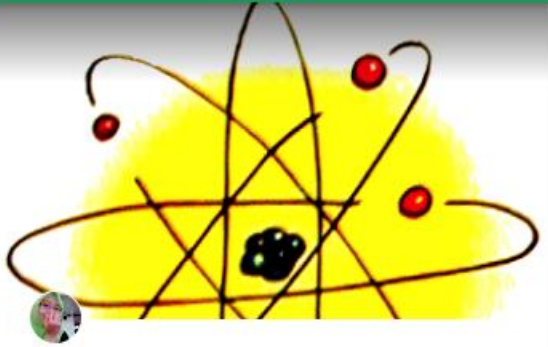
News

Bugs

Test

The core EGSnrc team:





807 members - Public

EGSnrc

Monte Carlo simulation of radiation transport

JOIN

Search Community

About Community

EGSnrc is a software package for the Monte Carlo simulation of coupled electron and photon radiation transport. Its current range of applicability in energy is 1 keV to 10 GeV.

EGSnrc

Filter

Frederic Tessier Moderator
Course

Wrapping up a successful two-week EGSnrc course at ICTP in Trieste, Italy. Congratulations to all the students! Thank you to the organizers and sponsors, on behalf of all lecturers. +1 if you enjoyed the course!

Grignano



+1 17

Marc Chamberland: Great group of students! It was a blast and I can't wait to hear what y...

55karlpdf Discussion 2d

Hi all, Where can I find download EGS_WINDOWS? I'm not looking for the

dian kurniawati Discussion 18h

Hi all,
i have problem, how to combine *.3ddose data on phyton or matlab (CERR) ?
please help me.
thanks

+1 1

Oanh Oanh: Hi,
You have downloaded and run the CERR on...

Pete Watson Bug report 4d

I had been noticing some infinite loop behaviour when using a beam shared library as source in egs_chamber (see: <https://plus.google.com/100288585807196289271/posts/hPgndfQjMw5>).

With the help of Marc-Andre Renaud, we determined that the cause was in egs_beam_source.cpp, where particles in the source were being rejected beca...

Hi all, I have an electron beam model for which I have built, compiled and...

Hi all, I have an electron beam model for which I have built, compiled and made a shared library in \$EGS_HOME. Whenever I use this beam model as a beam ... - Julien...
plus.google.com

Team.

Code

EGSnrc is now in the public domain

Since 2016, the EGSnrc software is distributed under the **GNU Affero GPL v3.0** open source licence.

BEAMnrc is now integrated in the EGSnrc installation.

Permissions

- ✓ Commercial use
- ✓ Modification
- ✓ Distribution
- ✓ Patent use
- ✓ Private use

Limitations

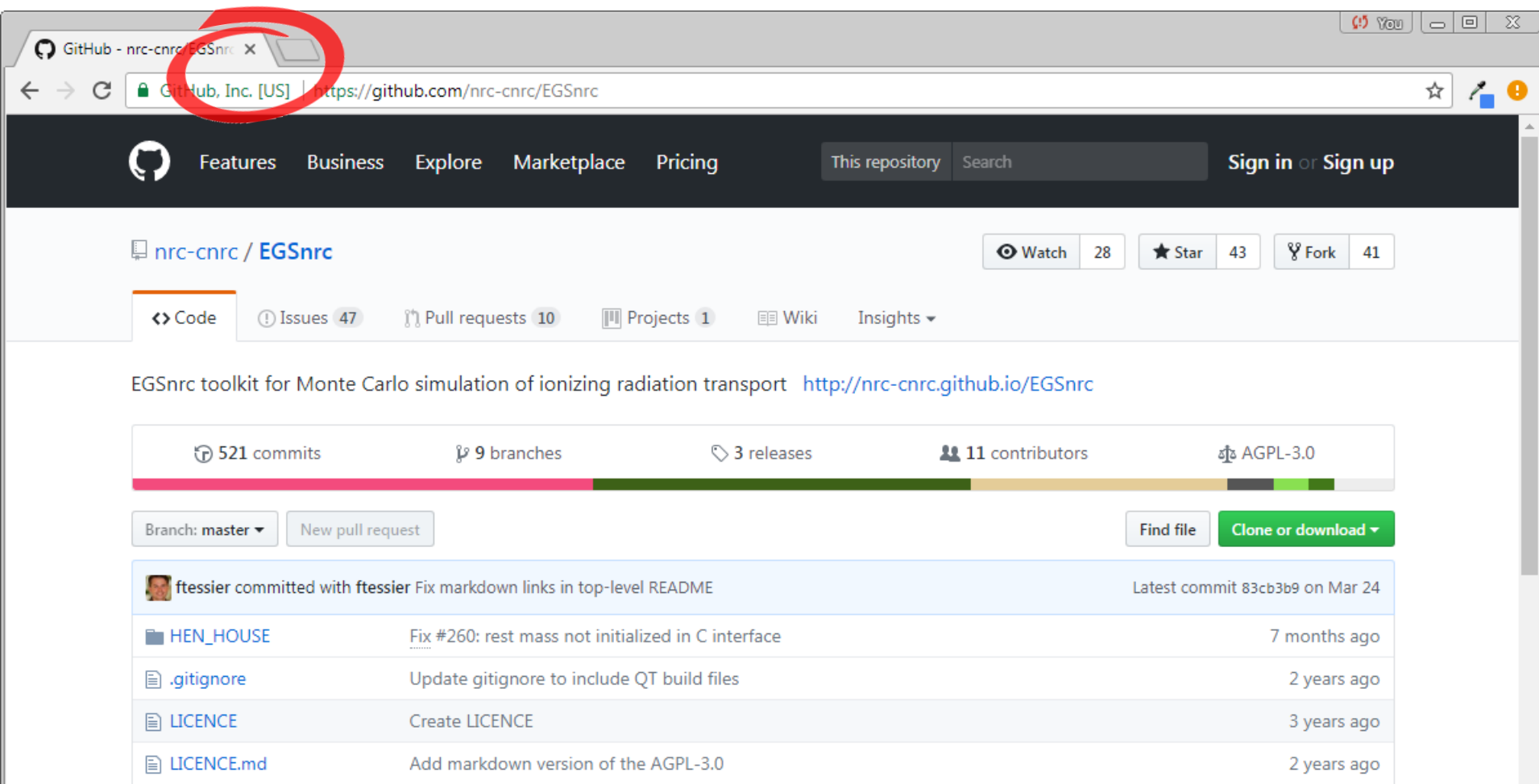
- ✗ Liability
- ✗ Warranty

Conditions

- ⓘ License and copyright notice
- ⓘ State changes
- ⓘ Disclose source
- ⓘ Network use is distribution
- ⓘ Same license

EGSnrc is now hosted on github.com

<https://github.com/nrc-cnrc/EGSnrc>



GitHub - nrc-cnrc/EGSnrc

GitHub, Inc. [US] | <https://github.com/nrc-cnrc/EGSnrc>

Features Business Explore Marketplace Pricing This repository Search Sign in or Sign up

nrc-cnrc / EGSnrc

Watch 28 Star 43 Fork 41

Code Issues 47 Pull requests 10 Projects 1 Wiki Insights

EGSnrc toolkit for Monte Carlo simulation of ionizing radiation transport <http://nrc-cnrc.github.io/EGSnrc>

521 commits 9 branches 3 releases 11 contributors AGPL-3.0

Branch: master New pull request Find file Clone or download

ftessier	ftessier committed with ftessier Fix markdown links in top-level README	Latest commit 83cb3b9 on Mar 24
HEN_HOUSE	Fix #260: rest mass not initialized in C interface	7 months ago
.gitignore	Update gitignore to include QT build files	2 years ago
LICENCE	Create LICENCE	3 years ago
LICENCE.md	Add markdown version of the AGPL-3.0	2 years ago



nrc-cnrc / EGSnrc

Watch 28

Star 43

Fork 41

Code

Issues 47

Pull requests 10

Projects 1

Wiki

Insights

Branch: develop

Commits on Sep 12, 2017



Update source directory path in EXphantom.egsinp

blakewalters committed with fteissier 26 days ago



d3d95a3



Add iaea_phsp build target for Linux and Windows

mainegra committed with fteissier 28 days ago



7793cff



Update default spectra location in egs_inprz

mainegra committed with fteissier 29 days ago



79aeb69



Commits on Sep 5, 2017



Set egs_gui compilation flags properly (#340)

mainegra committed with fteissier 29 days ago



15cb6ca



Fix rest mass and energy in DOSXYZnrc source 20

blakewalters committed with fteissier 29 days ago



ca35d90



Fix radionuclide relaxation input (#336)

rtownson committed with fteissier 29 days ago



7b517f8



Commits on Sep 3, 2017



Adjust examin application to recent changes (#337)

mainegra committed with fteissier on Sep 3



0f3caf9



nrc-cnrc / EGSnrc

report problems

Watch 28 Star 43 Fork 41

Code Issues 47 Pull requests 10 Projects 1 Wiki Insights

is:issue is:open Labels Milestones [New issue](#)

47 Open ✓ 123 Closed	Author	Labels	Projects	Milestones	Assignee	Sort
Infinite looping when using beam source in egs_chamber #351 opened 23 hours ago by pgfwatso						1
DOSXYZnrc Source 20 simulations in pegsless mode do not work #349 opened 13 days ago by ojalaj						1
bremsstrahlung correction=NRC missing for RMISW521 in material.dat #348 opened 13 days ago by ojalaj						1
Incorrect density for RW-1 water substitute in density corrections file #347 opened 13 days ago by ojalaj						3
EGSnrc gui "Go" button fails when make_prog variable includes arguments #346 opened 16 days ago by ftessier						
Fuzzy boundaries #345 opened 19 days ago by Kawrakow						1
Warn when a geometry name is not unique #344 opened 20 days ago by mchamberland						
Inconsistent terminology in iaea_phsp_source #334 opened on Aug 25 by mchamberland						
Feature request: start phase-space from random particle enhancement #333 opened on Aug 23 by skmartin						2
Issue compiling egs view on macOS Sierra compilation						15

nrc-cnrc / EGSnrc

submit code

Watch 28 Star 43 Fork 41

Code Issues 47 Pull requests 10 Projects 1 Wiki Insights

is:pr is:open Labels Milestones New pull request

10 Open ✓ 172 Closed	Author	Labels	Projects	Milestones	Reviews	Assignee	Sort
Fix #351: Infinite loop using beam source in egs++ ✓ bug #352 opened an hour ago by rtownson • Review required							
OpenMP support in EGSnrc ✓ enhancement #341 opened 28 days ago by edoerner • Review required						5	
Add option to output .3ddose files to EGS_DoseScoring ✓ enhancement #316 opened on Jul 25 by blakewalters • Approved						1	
Add EGS_DynamicSource class ✓ enhancement #308 opened on Jun 23 by blakewalters • Approved						11	
Implement RANLUX random number generator in egs++ ✓ enhancement #292 opened on May 19 by seirios • Review required						3	
QT5 migration for QT based GUI apps ✓ enhancement #263 opened on Mar 7 by crcrewso • Review required Release 2018						4	
Enhanced EM transport enhancement #211 opened on Jan 13 by victorMalkov • Review required Release 2018						3	
Add egs++ volume calculation enhancement #184 opened on Sep 21, 2016 by rtownson • Review required Release 2018						2	
Additional egs++ geometries and shapes ✓ enhancement #103 opened on Jun 8, 2016 by randlet • Approved Release 2018						45	

Installing EGSnrc in a nutshell

(but preferably in a Linux shell)

```
$ git clone https://github.com/nrc-cnrc/EGSnrc.git  
$ cd EGSnrc  
$ HEN_HOUSE/scripts/configure
```

Installing EGSnrc in a nutshell

(but preferably in a Linux shell)

```
$ git clone https://github.com/nrc-cnrc/EGSnrc.git
$ cd EGSnrc
$ git checkout develop # use the develop branch
$ HEN_HOUSE/scripts/configure
```

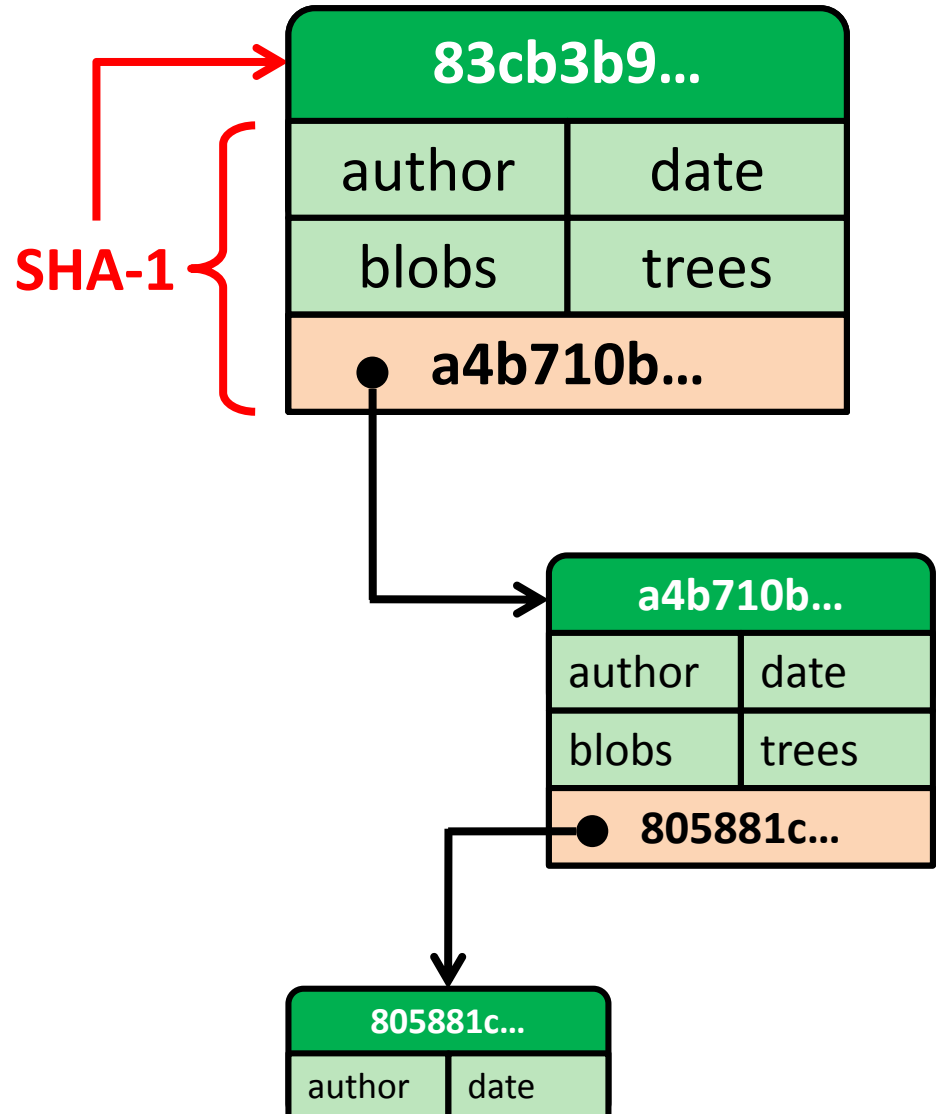
There are two main branches:

- 1. master:** updated yearly, versioned by **year** (EGSnrc 2017).
- 2. develop:** ongoing changes, versioned by **commit** (d3d95a3).

Cloning provides the entire commit history (try `git log`)

git is a robust version control system

- distributed, decentralized
- offline repository
- no repository setup
- atomic commits
- commit staging
- fast, efficient
- flexible and safe
- lightweight branches
- github, bitbucket, etc.



Code.

Ernesto Mainegra-Hing

Atomic relaxation and PE cross sections

Reid Townson

Radionuclide decay modelling

Dave Rogers

Improved kerma calculations

News

EGSnrc can model magnetic fields, *again!*


air

10 MeV
electrons

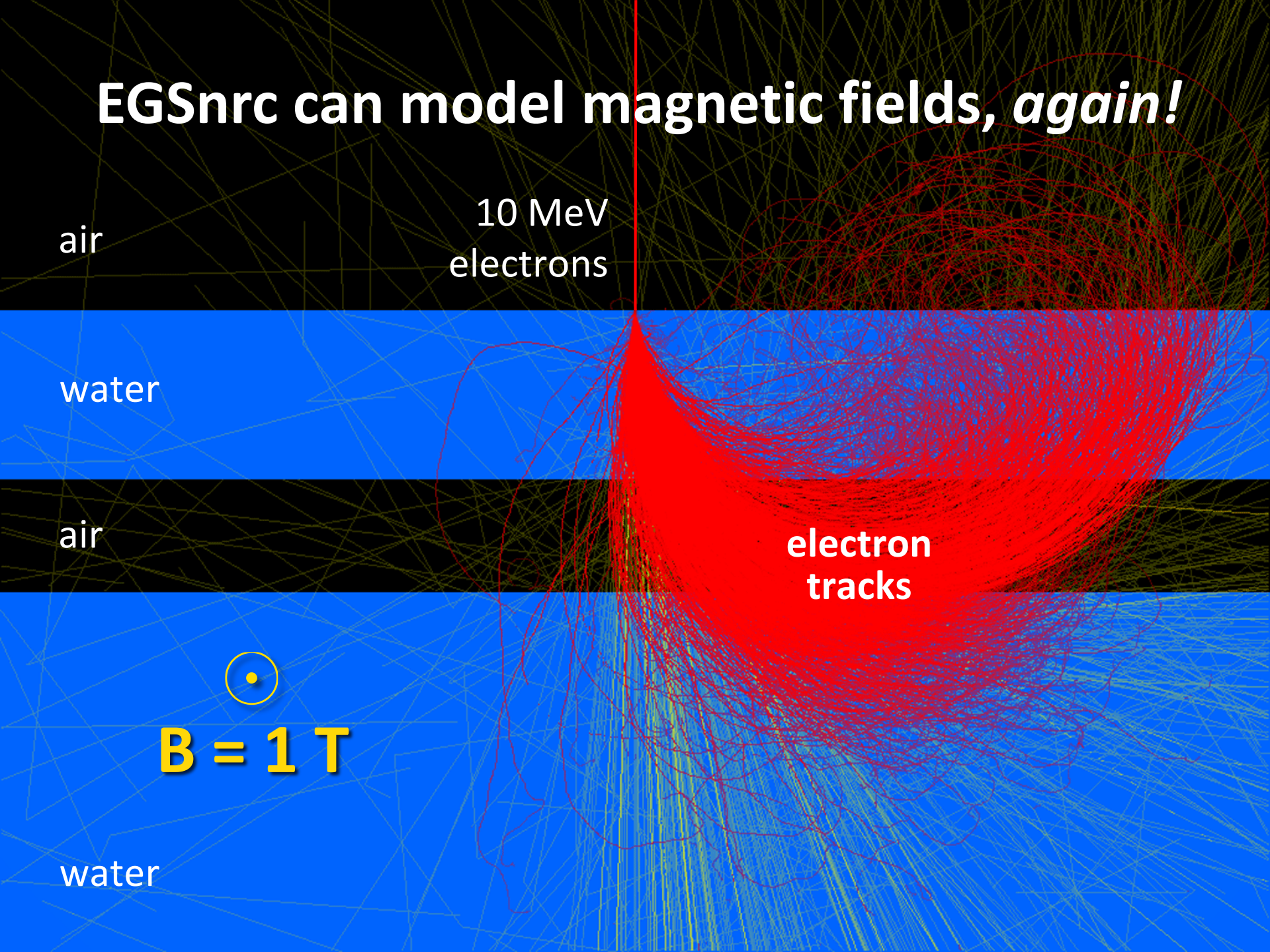
water

air

electron
tracks


B = 1 T

water



EM fields requires `emf_macros.mortran`

Electromagnetic fields are not included by default; you have to include the EMF macros in the compilation chain, e.g.,

```
EGSPP_USER_MACROS = cavity.macros \  
    $(EGS_SOURCEDIR)emf_macros.mortran
```

Fields are defined in the input file:

```
:start MC transport parameter:
```

```
Magnetic Field = 0 0 1 # Bx By Bz (in T)  
Electric Field = 0 0 0 # Ex Ey Ez (in V/cm)  
EM ESTEPE      = 0.02
```



ϵ_B

```
:stop MC transport parameter:
```

Malkov proposed a higher-order method

Charged particle transport in magnetic fields in EGSnrc

V. N. Malkov^{a)} and D. W. O. Rogers^{a)}

*Carleton Laboratory for Radiotherapy Physics, Physics Department, Carleton University,
Ottawa, Ontario K1S 5B6, Canada*

(Received 7 January 2016; revised 30 May 2016; accepted for publication 8 June 2016;
published 29 June 2016)

Purpose: To accurately and efficiently implement charged particle transport in a magnetic field in

Sensitive volume effects on Monte Carlo calculated ion chamber response in magnetic fields

Victor N. Malkov^{a)} and D. W. O. Rogers

Department of Physics, Carleton Laboratory for Radiotherapy Physics, Carleton University, Ottawa, ON, Canada

(Received 17 May 2017; revised 8 June 2017; accepted for publication 13 June 2017;
published 19 July 2017)

Purpose: The development of magnetic resonance-guided radiation therapy (MRgRT) necessitates accurate Monte Carlo (MC) models of ion chambers for computing ion chamber corrections to compensate for the presence of the magnetic field. This study evaluates the sensitivity of the ion chamber dose response in a magnetic field on the collection volume used in the MC simulation.

Methods: The EGSnrc system's egs_chamber application is used with a recently developed and validated magnetic field transport code. The calculated dose to the sensitive volume of the chamber per unit incident photon fluence, normalized to that at 0 T, is evaluated as a function of magnetic field for the PTW 30013, PTW 31006, PTW 31010, Exradin A12S, and Exradin A1SL chambers. The sensitive region is varied by excluding the volume corresponding to either 0, 0.5, or 1 mm of distance from the stem. The photon field, magnetic field, and ion chamber are all oriented perpendicular

News.

Bugs

Zero electron rest mass for 30 days!

 Closed

ftessier opened this issue on Mar 1 · 0 comments

Thank you to **Shahid Naqvi**



ftessier commented on Mar 1 • edited

Owner



Description

A critical bug has been uncovered in EGSnrc which causes the electron energy to be offset by the value of the electron rest mass in some applications. This bug was introduced in commit [1eaf898](#) and is caused by the `prm` variable (precise rest mass) being used before it is initialized via `call hatch`.


Scope

Affects versions of EGSnrc downloaded between **3 February 2017 and 2 March 2017, inclusively**. For those working off the `develop` branch, the corrupted date range is 24 January to 2 March.

Remediation

The fix involves patching three source files, as detailed in commit [805881c](#). The EGSnrc repository has been patched accordingly as of **1 March 2017**, on both branch `master` (commit [805881c](#)) and branch `develop` (commit [1c1cdb2](#)).

Wrong MS coefficients for 17 years!

 Closed jantolak opened this issue on Jun 24, 2016 · 1 comment

Thank you to **John Antolak**



jantolak commented on Jun 24, 2016 · edited by fressier

Contributor

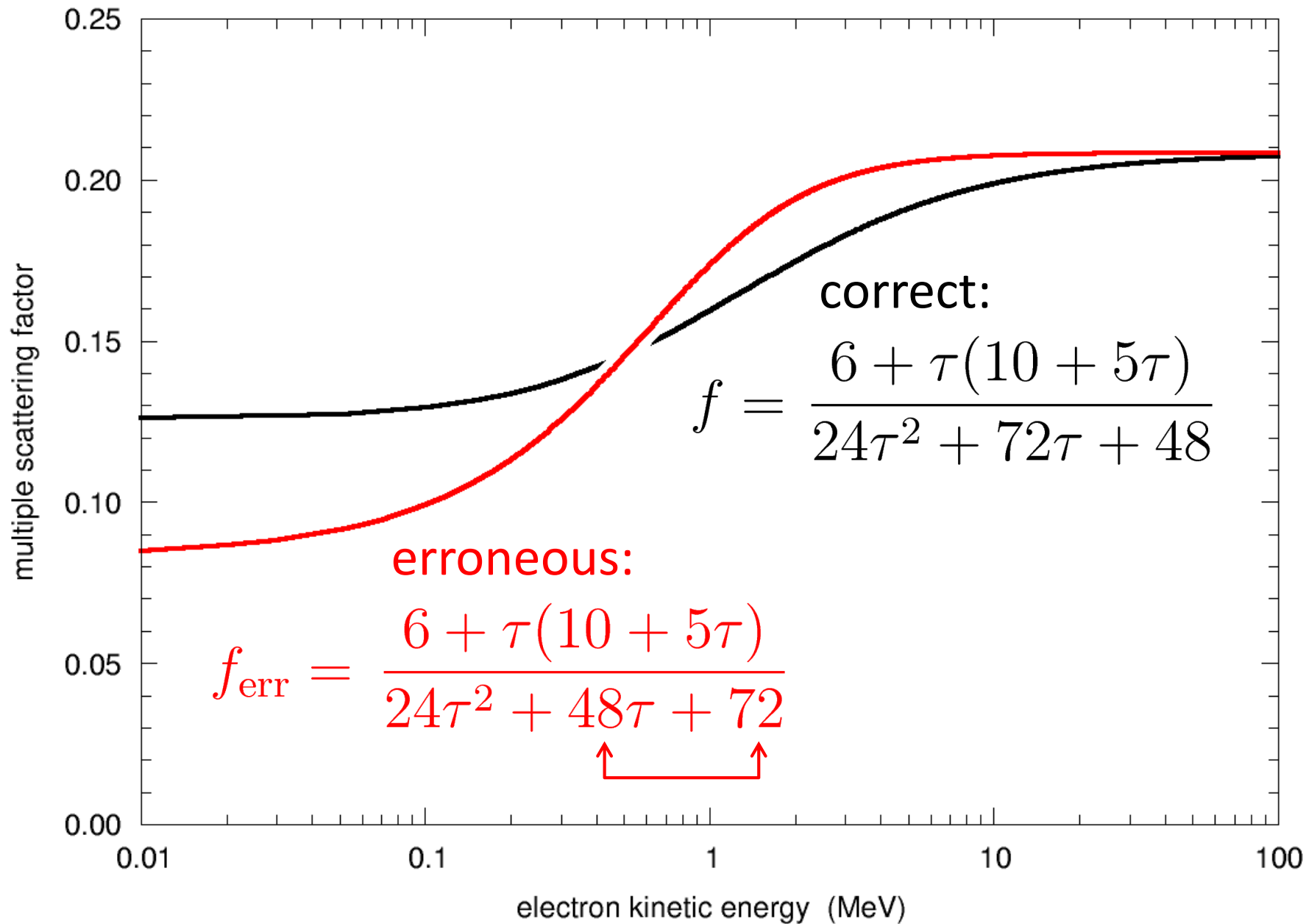
In `egsnrc.mortran`, subroutine `msdist_pII` at line 4092, it appears that the polynomial in the denominator is incorrect. Here is the current code at those lines.

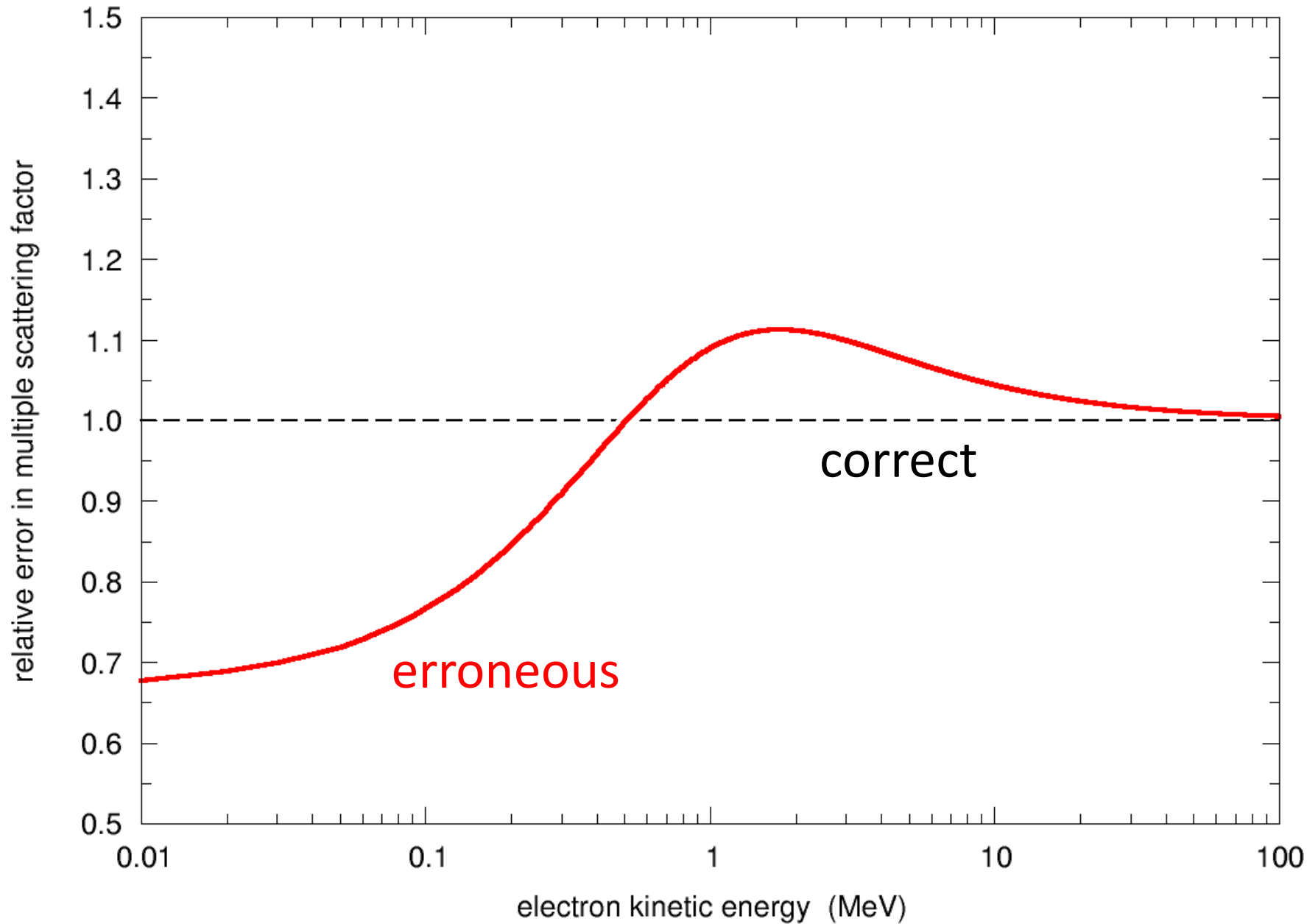
```
4091  "e      = e * (1 - epsilon*epsilon*((6+tau*(10+5*tau))/(tau+1)/((tau+2))/24);
4092  e      = e * (1 - epsilon*epsilon*(6+10*tau+5*tau2)/(24*tau2+48*tau+72));
```

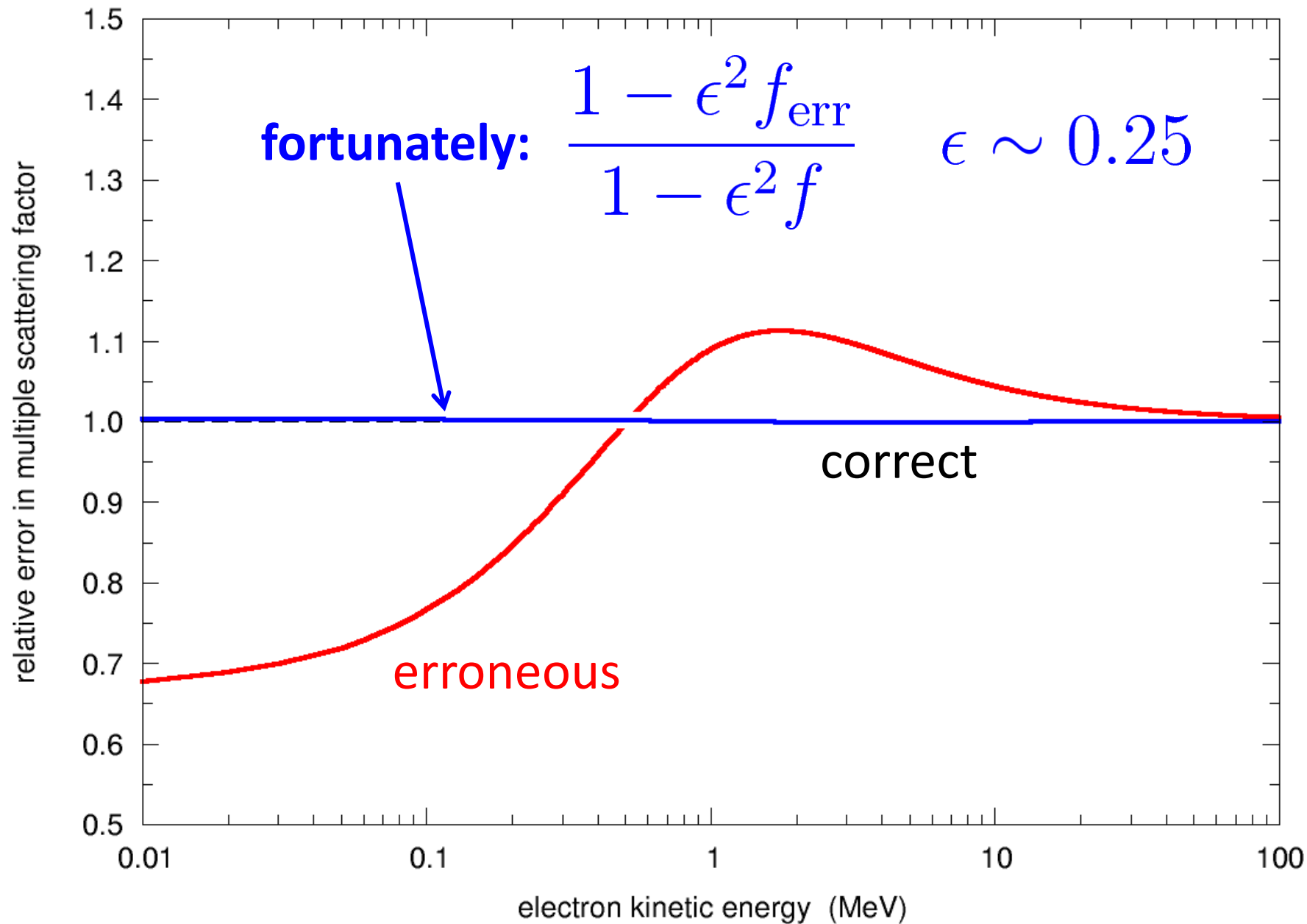
Assuming that the expression on line 4091 is correct, the denominator in the last term should be `24*(tau+1)*(tau+2) = 24*(tau2+3*tau+2) = (24*tau2+72*tau+48)`: the 72 and 48 are not in the right place in the code. Therefore, line 4092 should read

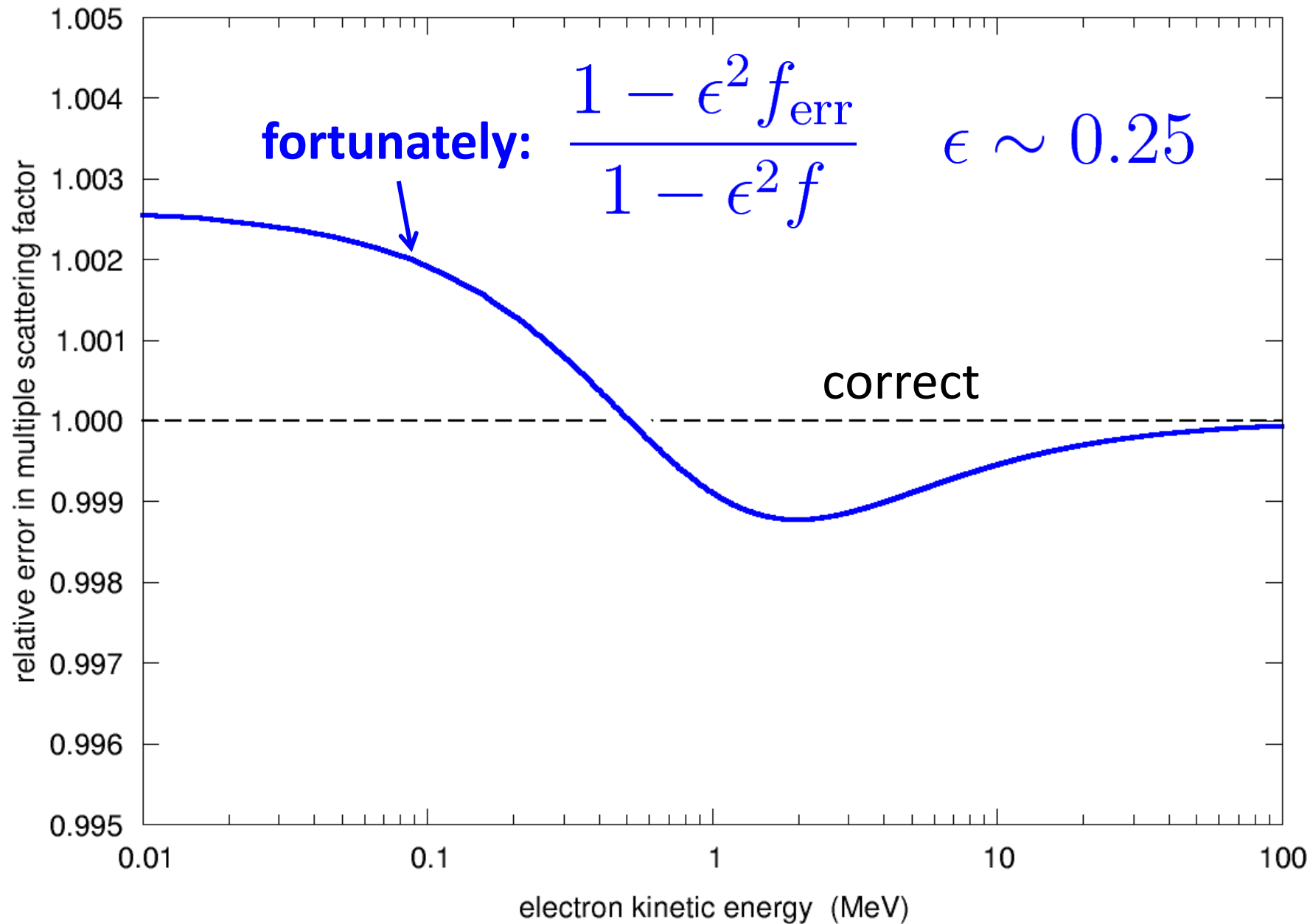
```
4092  e      = e * (1 - epsilon*epsilon*(6+10*tau+5*tau2)/(24*tau2+72*tau+48));
```

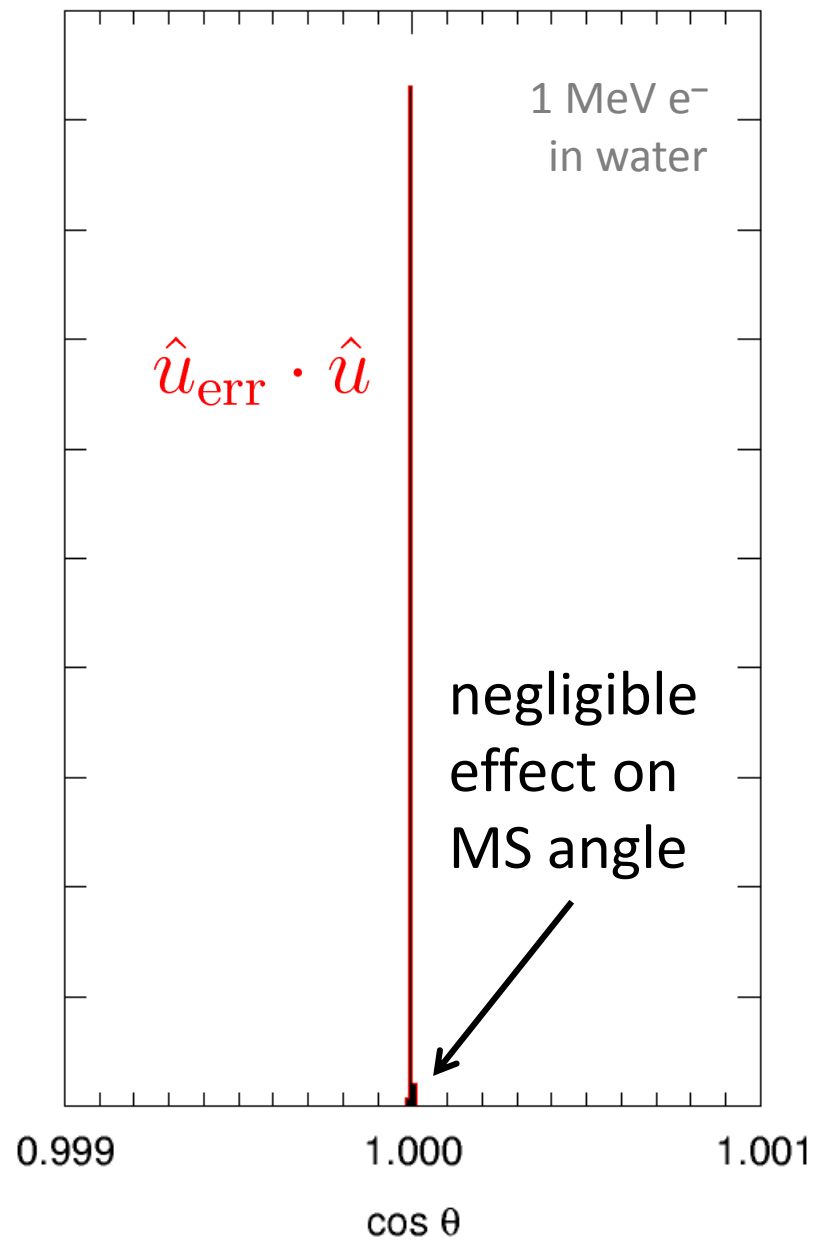
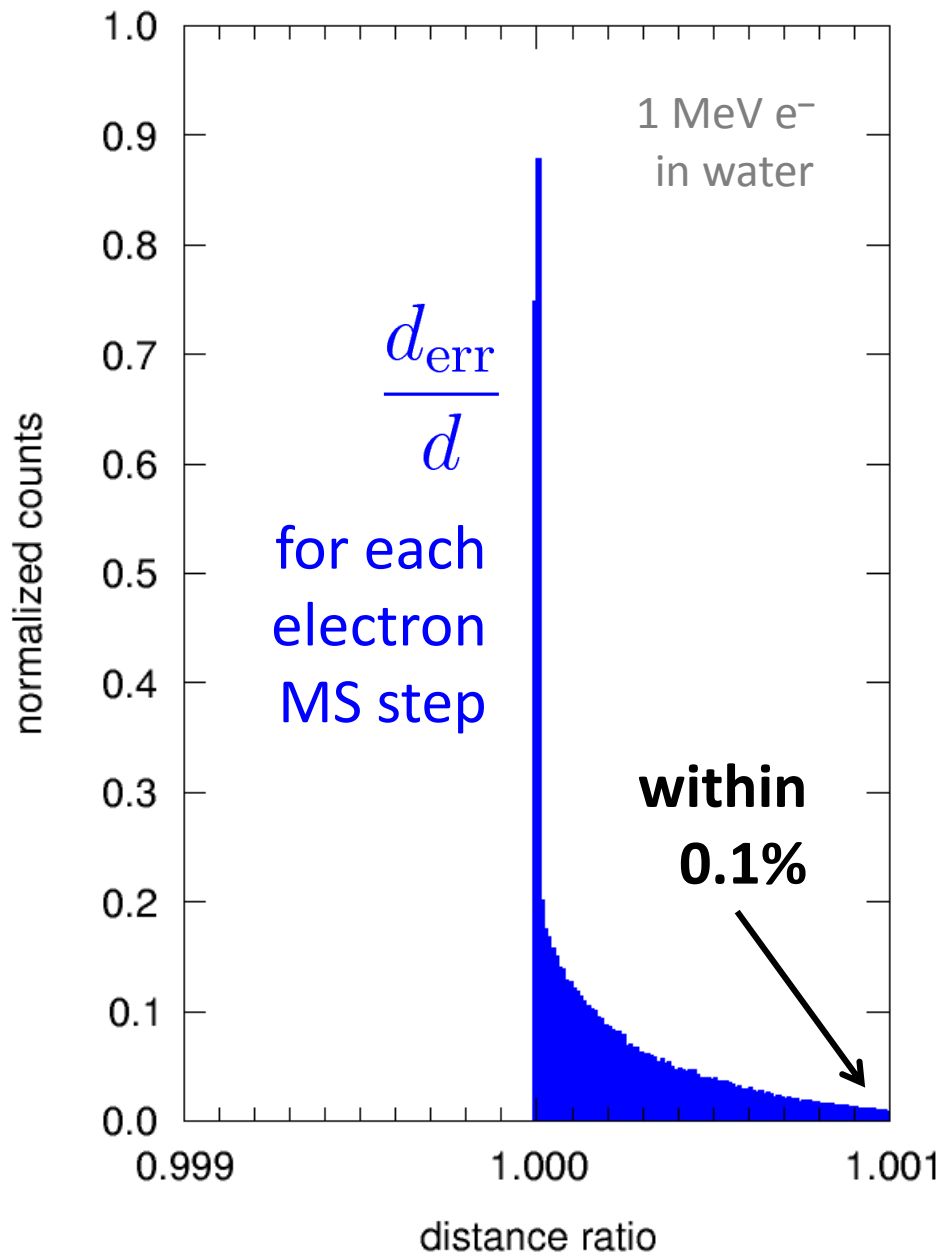
It appears that this error makes very little difference, which is likely why it has not been noticed. The error was actually pointed out to me by a trainee who was working on his own MC code for electron scattering and wanted to compare to an existing code as a baseline.











Someone else's bug: ESTAR I-value

NIST National Institute of Standards and Technology Physical Meas. Laboratory

stopping-power and range tables for electrons

The ESTAR program calculates stopping power, density effect parameters, range, and radiation yield tables for electrons in various materials. Select a material and enter the desired energies. Energies are specified in MeV, and must be in the range from 0.001 MeV to 10000 MeV.

[Help](#) [Text version](#) [Material composition data](#)

Select a common material: 1: Hydrogen
or enter a [unique material](#)

Graph stopping power:
 Total Stopping Power
 Collision Stopping Power
 Radiative Stopping Power

Graph density effect parameter

Graph CSDA range

Graph radiation yield

No graph

Additional Energies (optional):
Use energies from a file*
Choose File No file chosen

or
Use energies entered below (one per line)

 Include default energies

Note: Only stopping powers and the density effect parameter will be calculated if additional energies are used.

Submit Reset

* Your browser must be file-upload compatible.

[contents](#)

Someone else's bug: ESTAR I-value

The screenshot shows the ESTAR web interface. At the top, there is a banner with the ESTAR logo and the text "stopping-power and range tables for electrons". Below the banner, a message states: "The information below will be used to determine ESTAR's output for *graphite*." The main form area is purple and contains several sections. A text input field for "mean excitation energy" is highlighted with a red circle and contains the value "81.0". A red arrow points from this field to a callout box on the right. Below the input field, there are several radio button options for graphing: "Graph stopping power:" (with sub-options for Total, Collision, and Radiative Stopping Power), "Graph density effect parameter", "Graph CSDA range", "Graph radiation yield", and "No graph". To the right of these options is the "Additional Energies (optional)" section, which includes a "Choose File" button and a text area for entering energies. A red arrow points from the "Choose File" button to the same callout box. At the bottom of the form, there are "Submit" and "Reset" buttons. A note at the bottom of the form reads: "Note: Only stopping powers and the density effect parameter will be calculated if additional energies are used." Below the form, there is a footer with "back" and "contents" links. A yellow box at the bottom of the page contains the text "Validate against ESTAR.f program".

ESTAR calculated the mean excitation energy to be: 81.0 eV.
You may change this value now.

Help

Graph stopping power:
 Total Stopping Power
 Collision Stopping Power
 Radiative Stopping Power

Graph density effect parameter

Graph CSDA range

Graph radiation yield

No graph

Additional Energies (optional):
Use energies from a file*
Choose File No file chosen

or
Use energies entered below (one per line)

Include default energies

Note: Only stopping powers and the density effect parameter will be calculated if additional energies are used.

Submit Reset

* Your browser must be file-upload compatible.

back contents

Validate against ESTAR.f program

custom I-value is not taken into account when custom energies are supplied

Bugs.

Test

Are Monte Carlo simulations *traceable*?



Mass measurements are in principle traceable to the BIPM kilogram in Paris (until 2018).

Are we doing everything we can to ensure the validity of Monte Carlo simulation?

Monte Carlo simulation results are widely trusted, for example in dosimetry protocols.

Clients have started to ask for official Monte Carlo simulation calibration certificates!

What is software traceability anyway?

1. robust versioning, robust source code

- migrate to git version control system ✓
- port the EGSnrc core code to C++

2. automated, continuous integration testing

- compilation test on every commit (Travis CI) ✓
- run standard simulation set for numerical comparison

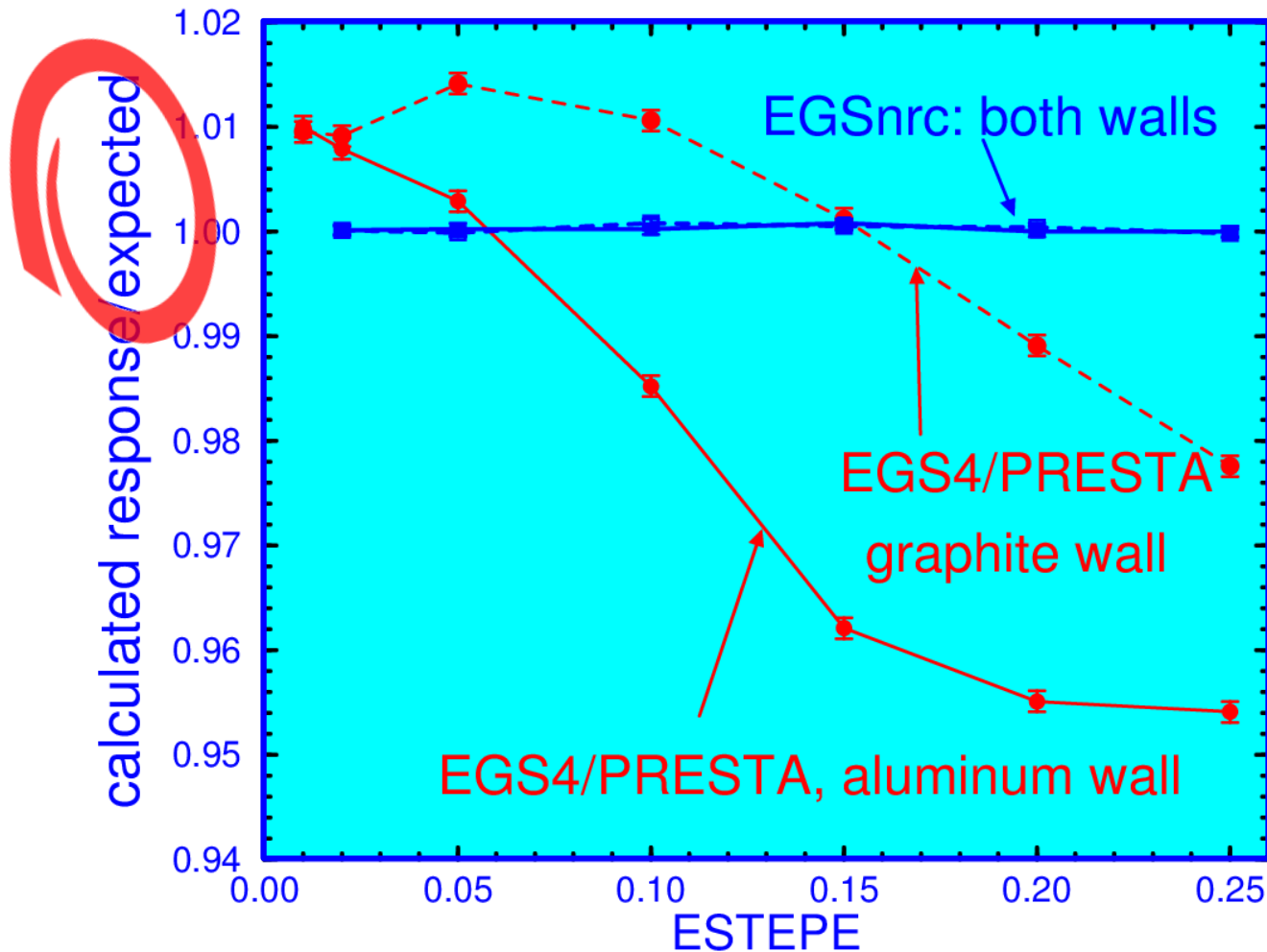
3. automated, ongoing key comparisons between codes

- agree on key data and key scenarios
- develop a common simulation description language?

4. Monte Carlo simulation verification

Kawrakow's famous Fano test graph

ion chamber response in ^{60}Co beams



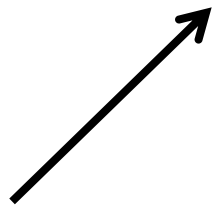
“EGSnrc is accurate to within 0.1%, with respect to its own cross sections.”

This remains a distinguishing feature of EGSnrc today!

Fano theorem provides a rigorous test

A Monte Carlo simulation algorithm is essentially solving the Boltzmann transport equation, numerically:

$$\mathbf{u} \cdot \nabla_r \phi = \rho (s + I[\phi])$$



change in fluence

$$0 = \rho (s + I[\phi])$$

source

atomic interactions

If the atomic properties are identical everywhere, **per unit mass**, a uniform fluence implies a uniform source (per unit mass).

Since the solution to the Boltzmann equation is presumed unique: **turn this around to verify the Monte Carlo algorithm.**

Reference dosimetry in the presence of magnetic fields: conditions to validate Monte Carlo simulations

Hugo Bouchard¹, Jacco de Pooter², Alex Bielajew³ and Simon Duane¹

¹ Acoustics and Ionising Radiation Team, National Physical Laboratory, Hampton Road, Teddington TW11 0LW, UK

² VSL B.V., Thijsseweg 11, NL-2629 JA Delft, The Netherlands

³ Department of Nuclear Engineering and Radiological Sciences, The University of Michigan, Ann Arbor, MI 48109, USA

E-mail: hugo.bouchard@npl.co.uk

Received 3 February 2015, revised 22 June 2015

Accepted for publication 26 June 2015

Published 13 August 2015



CrossMark

Abstract

With the advent of MRI-guided radiotherapy, reference dosimetry must be thoroughly addressed to account for the effects of the magnetic field on absorbed dose to water and on detector dose response. While Monte Carlo plays an essential role in reference dosimetry, it is also crucial for determining

Fano theorem within a magnetic field

The magnetic field adds a Lorentz force term in the Boltzmann transport equation:

$$\mathbf{u} \cdot \nabla_r \phi = \rho \left(-s + \frac{q\mathbf{u}}{c} \cdot \left[\frac{B(\hat{\mathbf{u}})}{B} \times \hat{\mathbf{b}} \right] \right) \cdot \nabla_p \phi$$

There are **two choices** to recover a testable *Fano condition*:

1. scale the magnetic field with density
2. make this gradient parallel to velocity

$$B \rightarrow B/\rho$$

The condition $\nabla_p \phi \sim \hat{\mathbf{u}}$ implies that the magnetic term vanishes: a **uniform isotropic source** yields a uniform fluence!

Fano testing requires 3 ingredients

1. uniform atomic interaction cross sections:

set all regions to the same material, vary the density.

2. a uniform, isotropic, density-scaled source of particles:

before: parallel photon beam, regenerate photons.

now: use the `egs_fano_source` class.

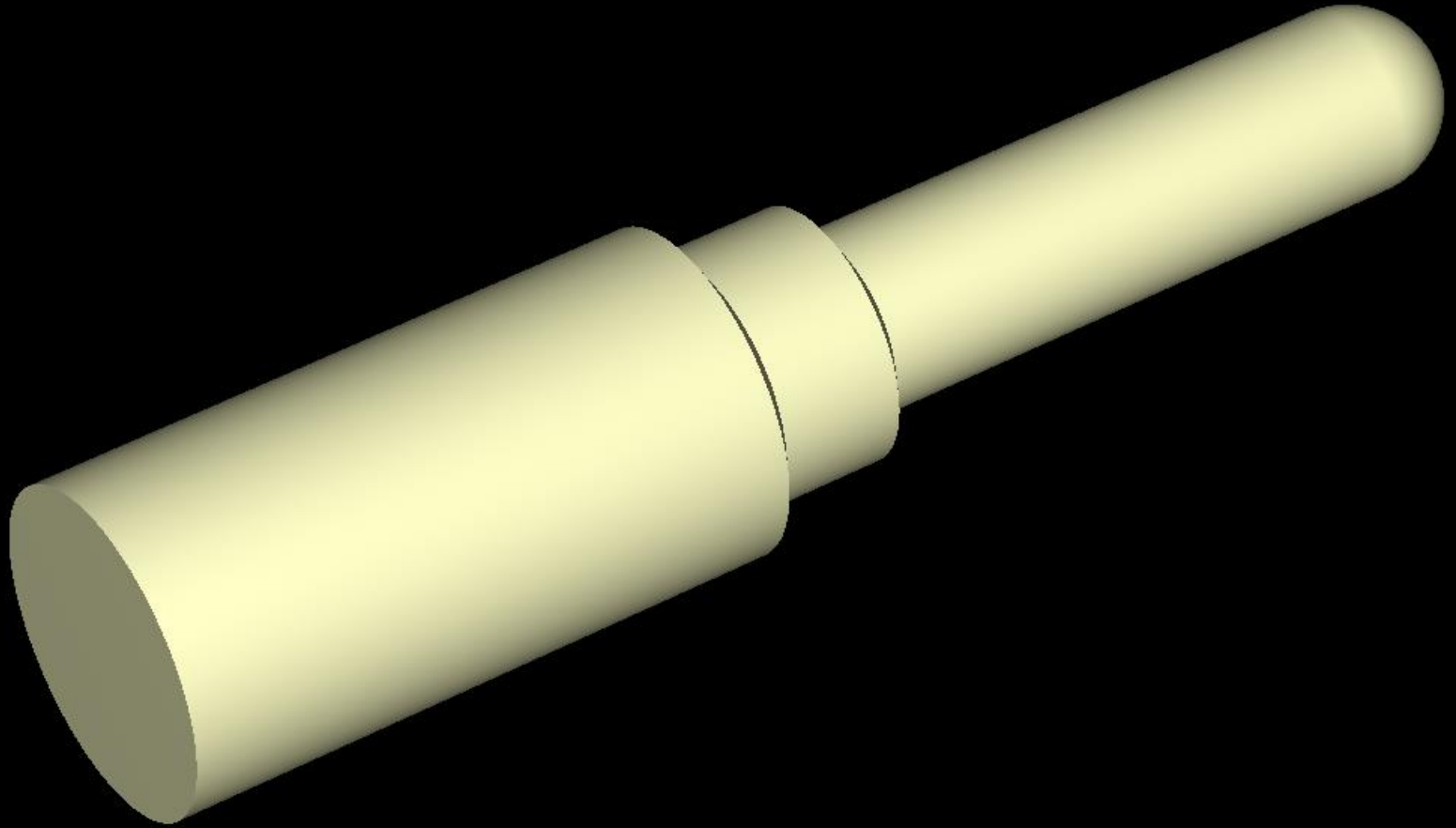
3. an infinite simulation space:

before: discard photon, worry about electron range...

now: use an *infinite* simulation space!

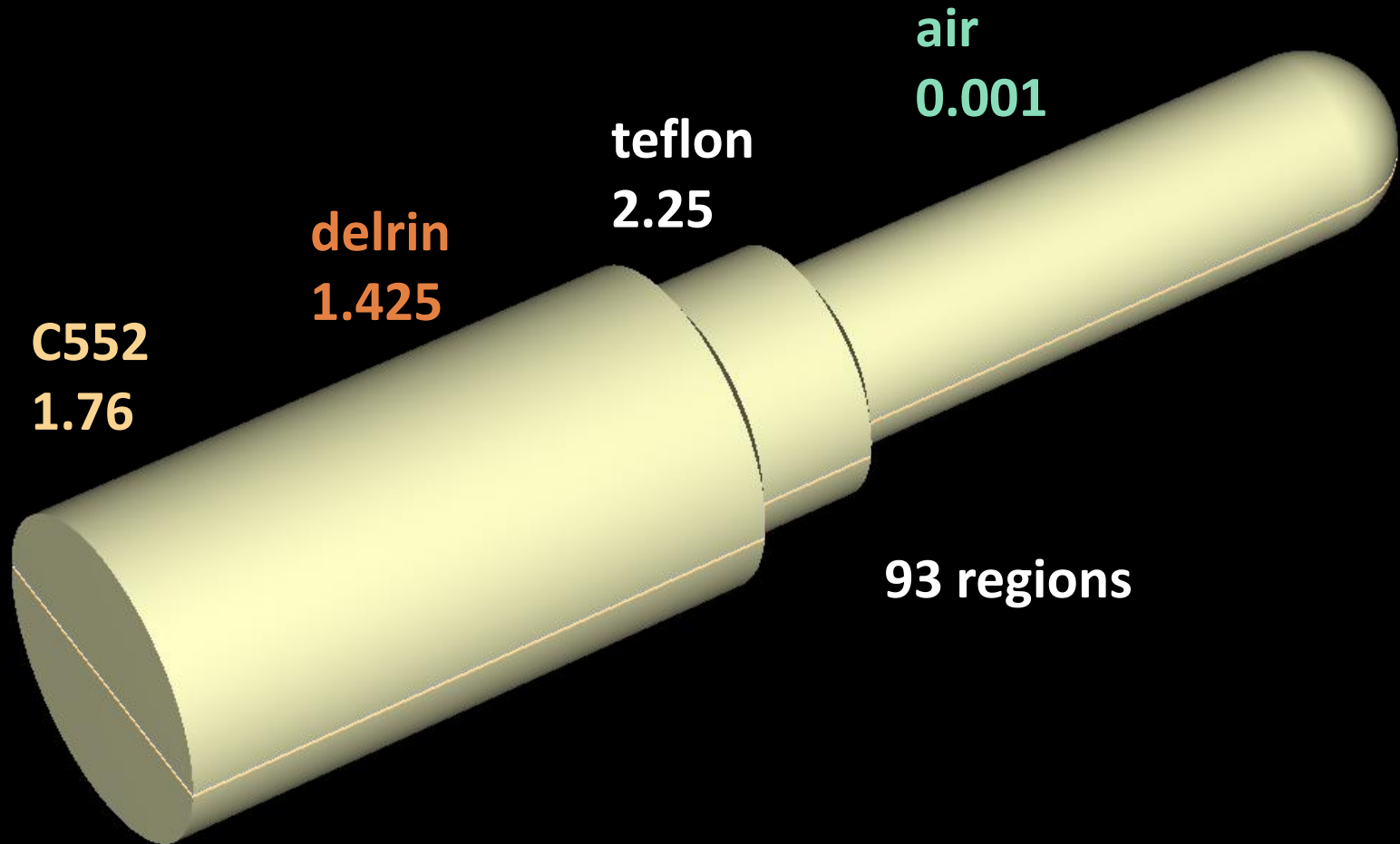
Fano testing an ion chamber

Exradin A12, 0.6 cm³ chamber



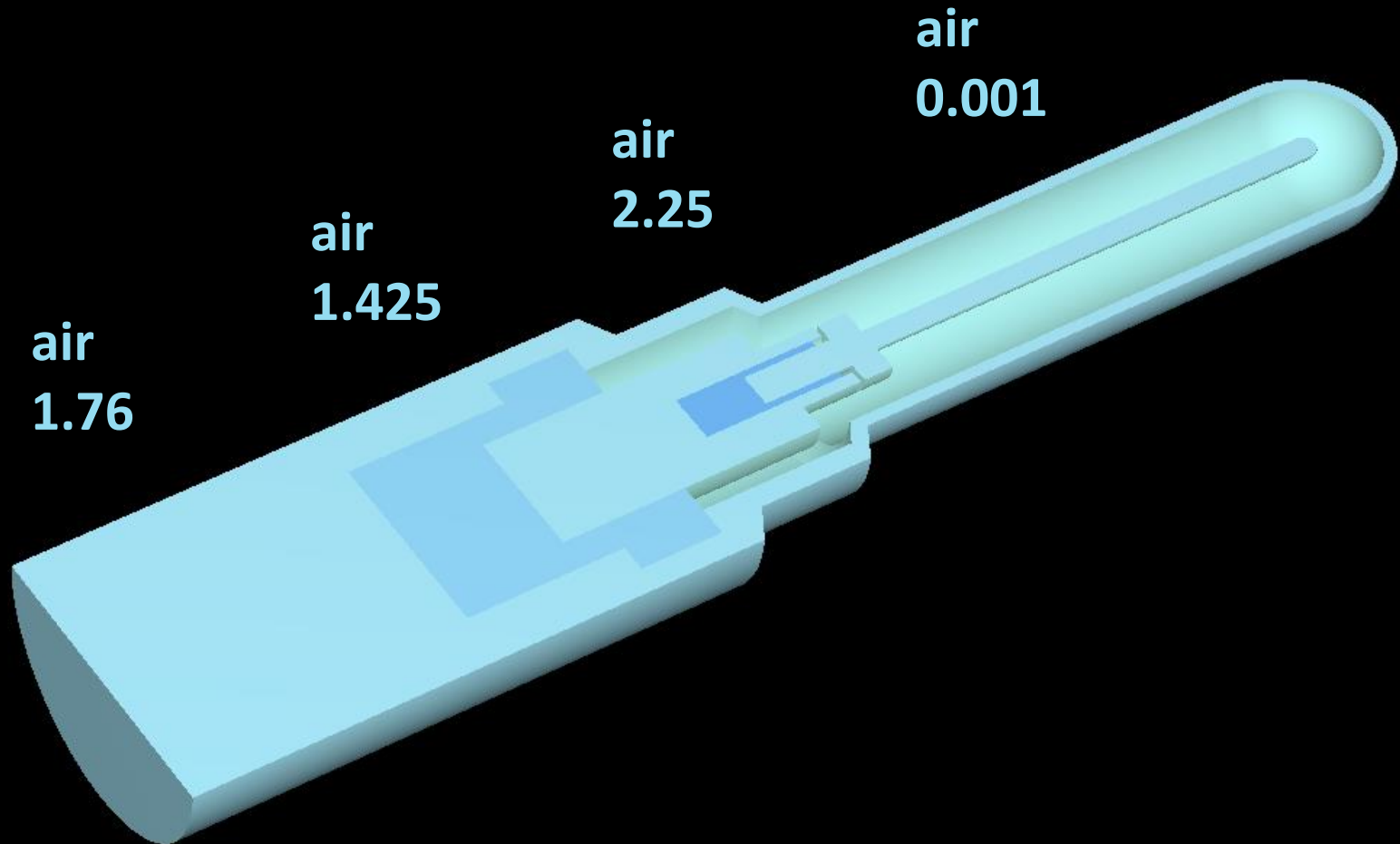
Fano testing an ion chamber

Exradin A12, 0.6 cm³ chamber



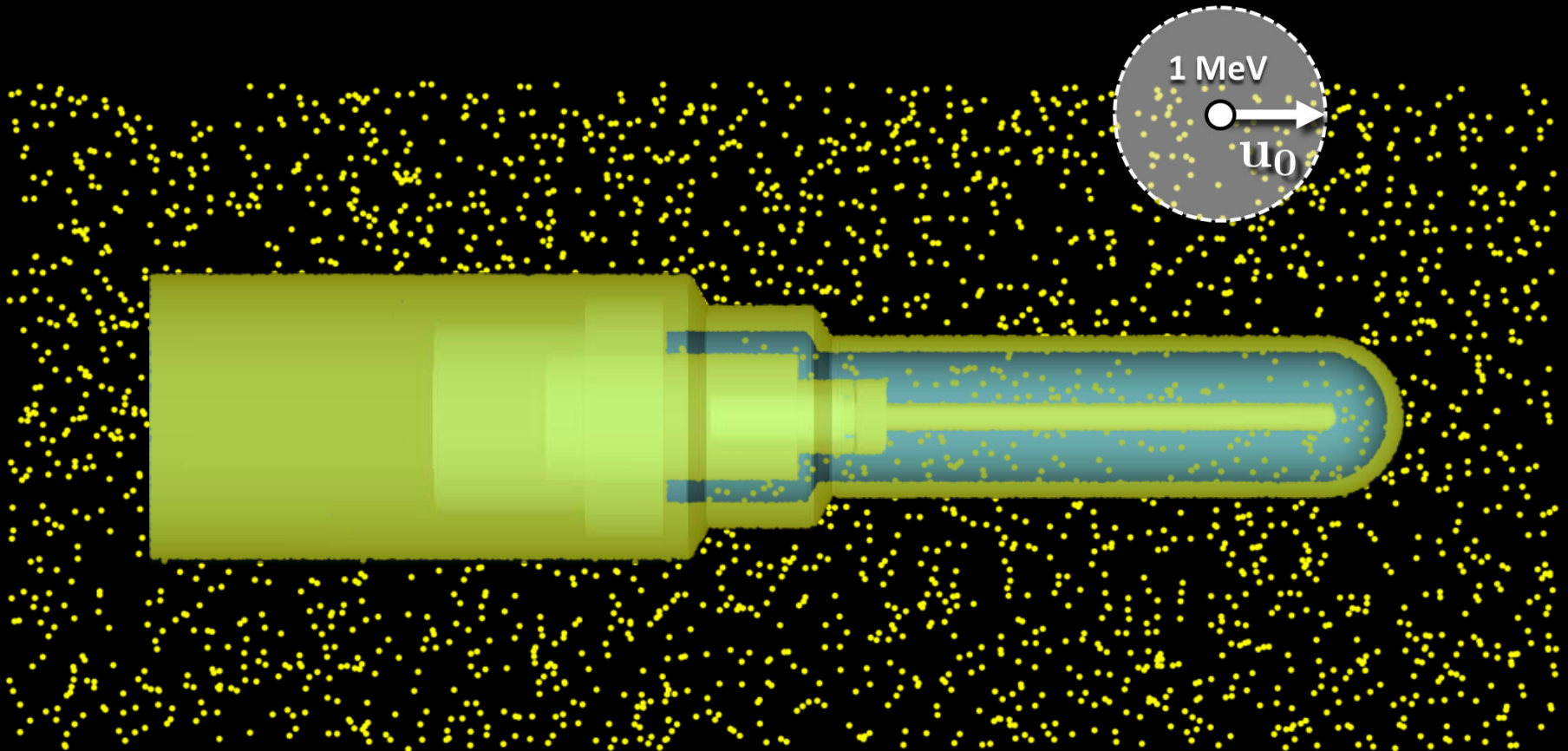
Fano testing an ion chamber

1. uniform atomic interaction cross sections



Fano testing an ion chamber

2. a uniform, isotropic, density-scaled source of particles

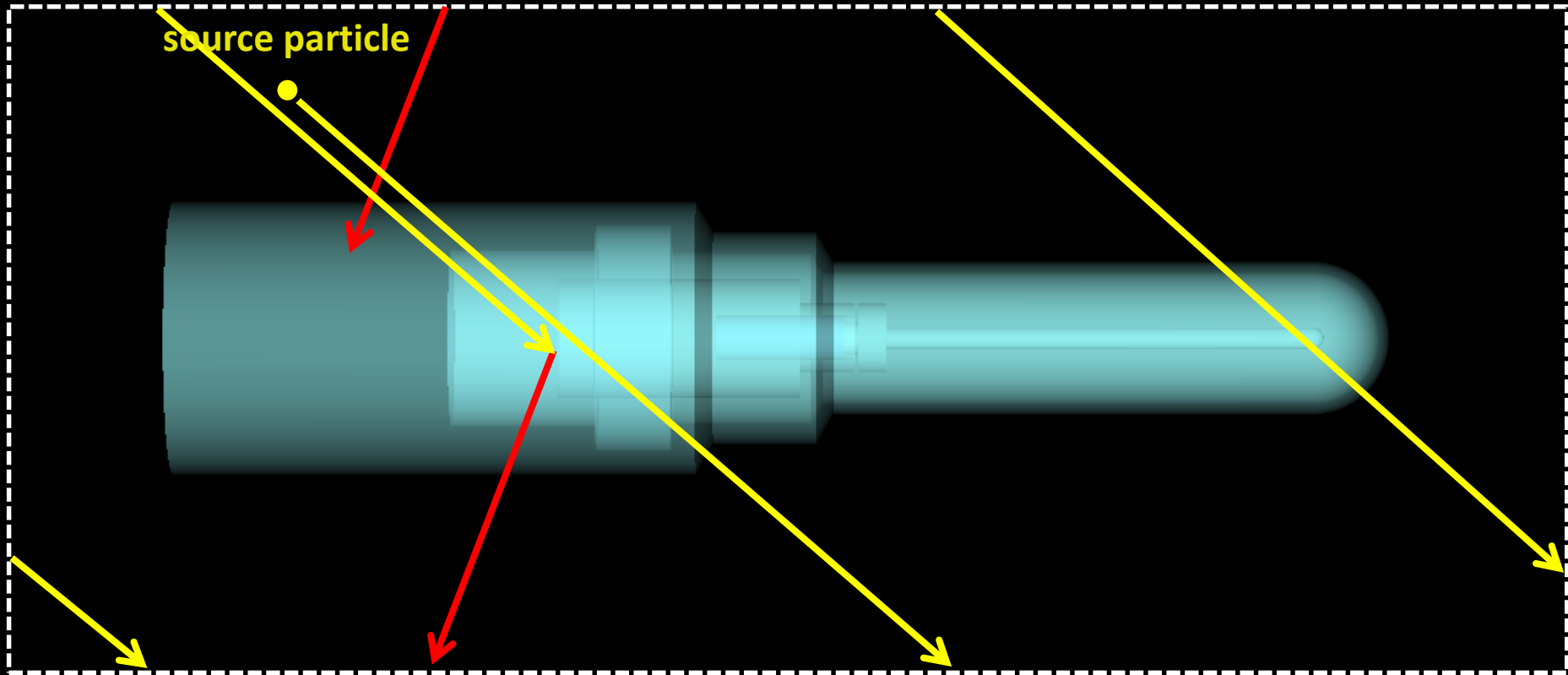


egs_fano_source

Fano testing an ion chamber

3. an infinite simulation space

periodic boundary conditions

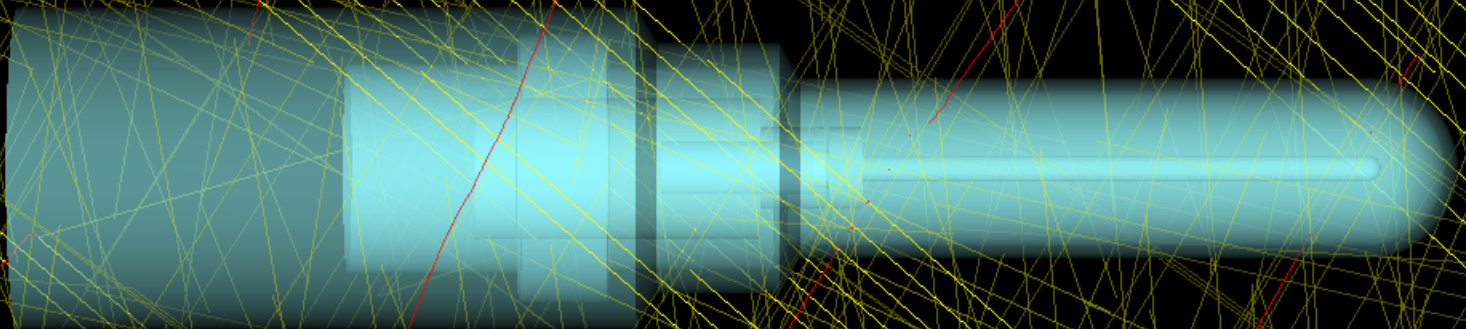


Fano testing an ion chamber

3. an infinite simulation space

periodic boundary conditions

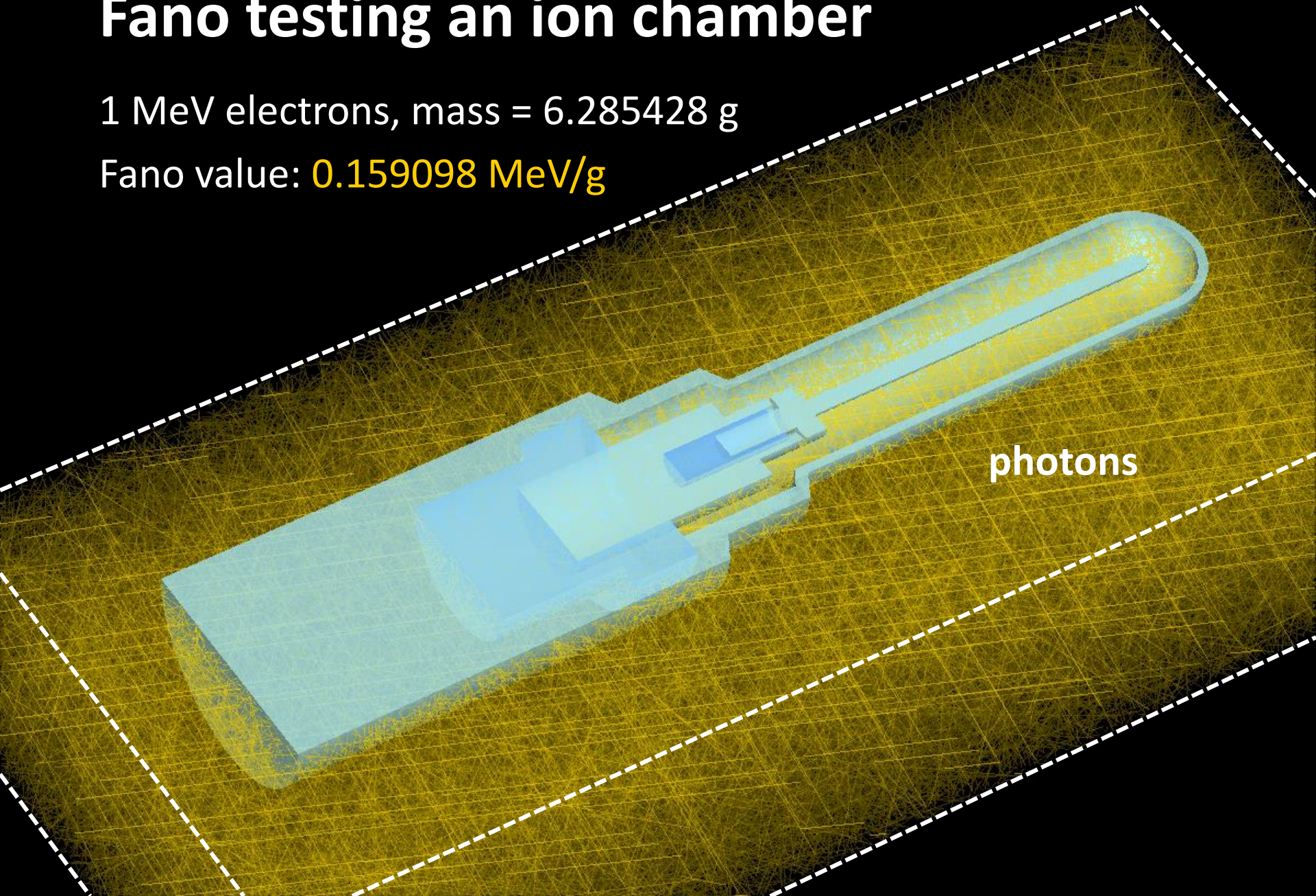
source particle



Fano testing an ion chamber

1 MeV electrons, mass = 6.285428 g

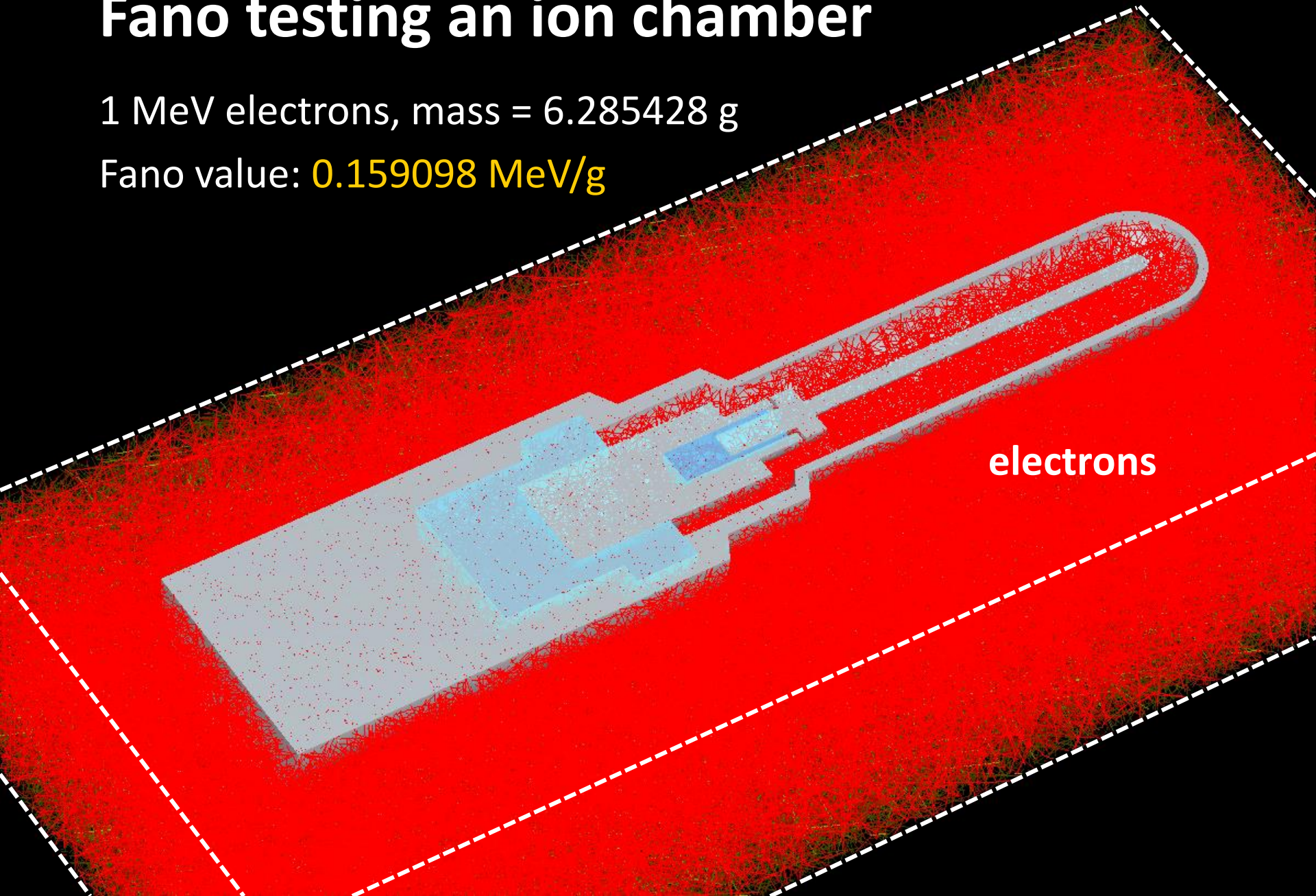
Fano value: **0.159098 MeV/g**



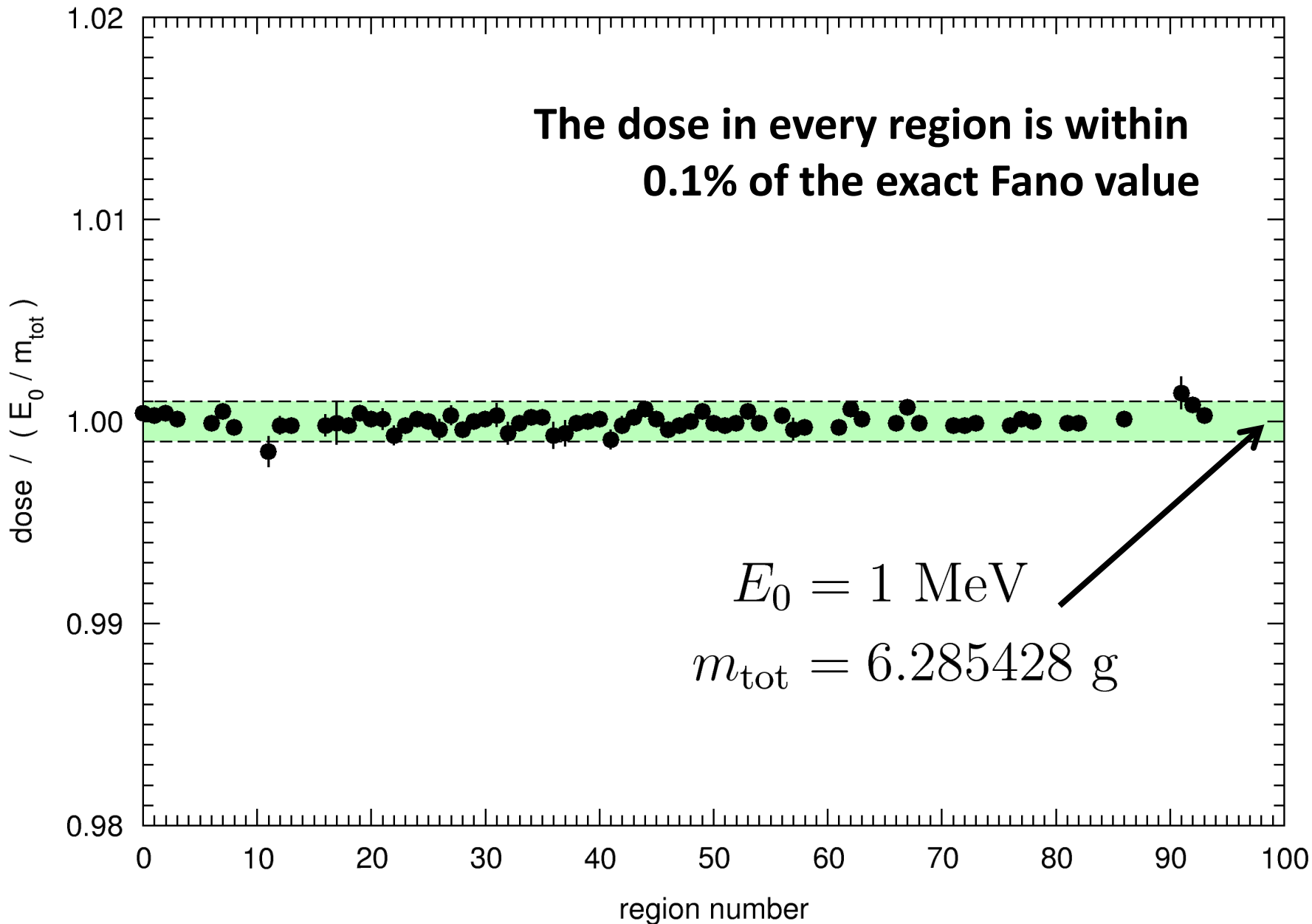
Fano testing an ion chamber

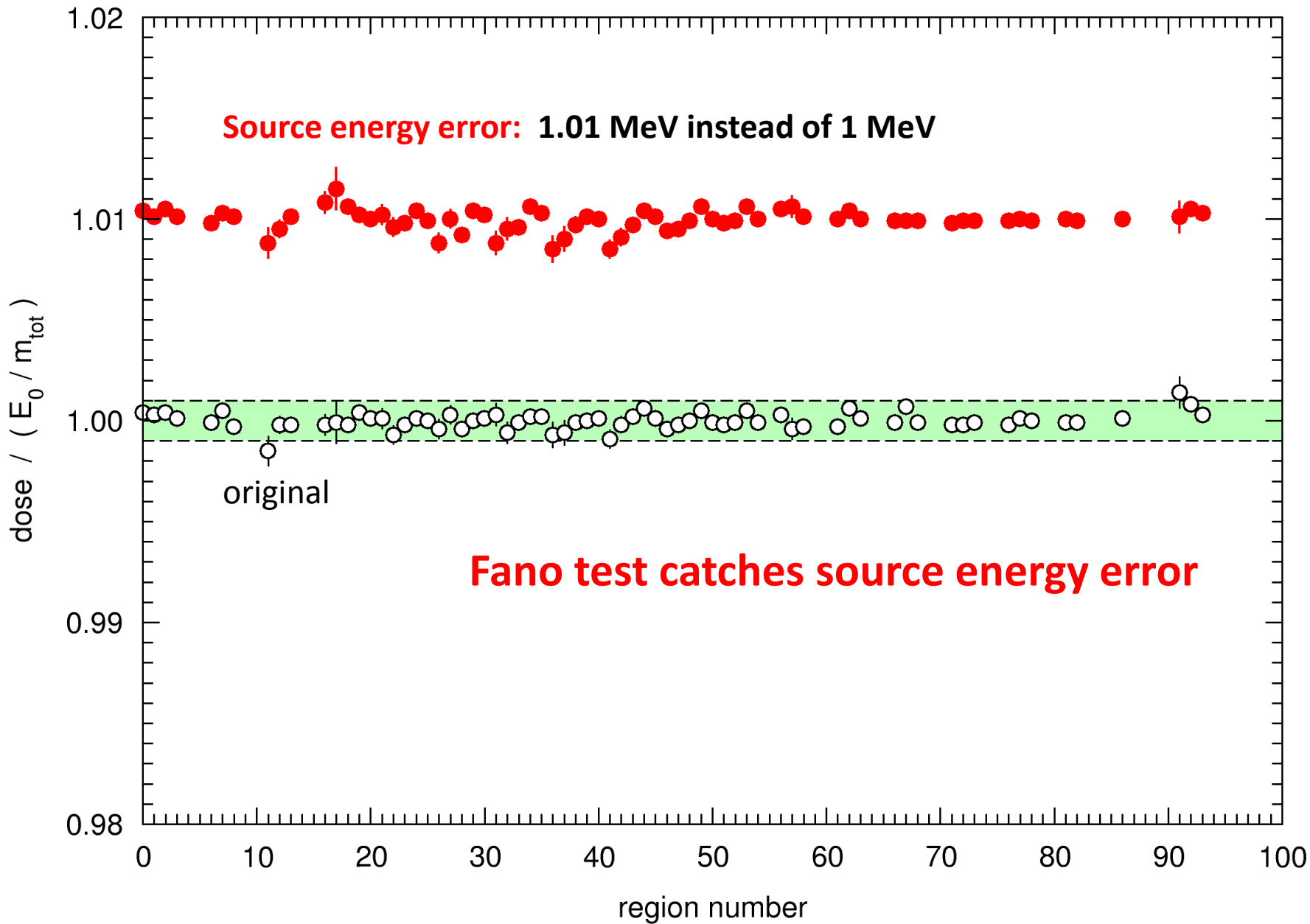
1 MeV electrons, mass = 6.285428 g

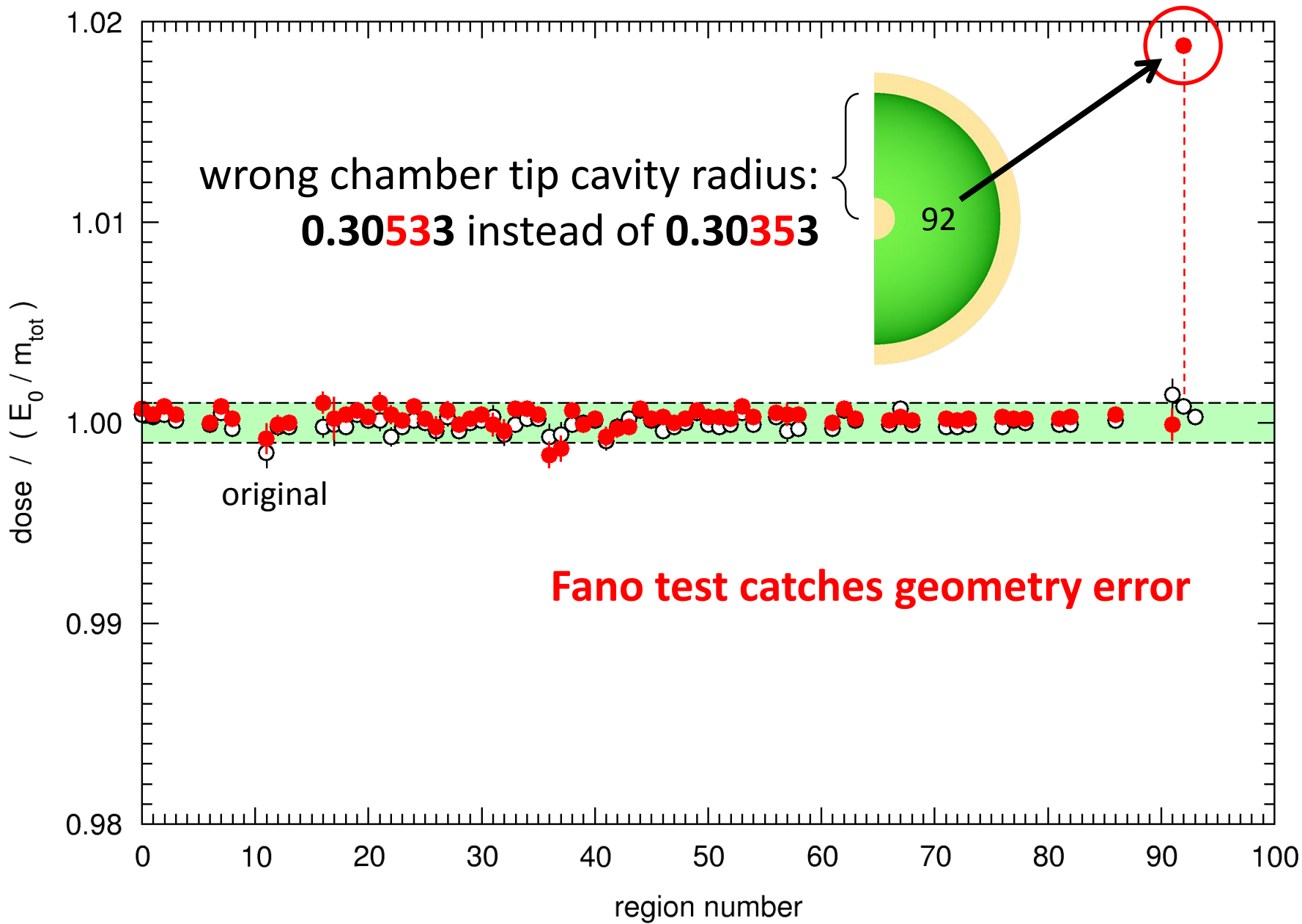
Fano value: **0.159098 MeV/g**



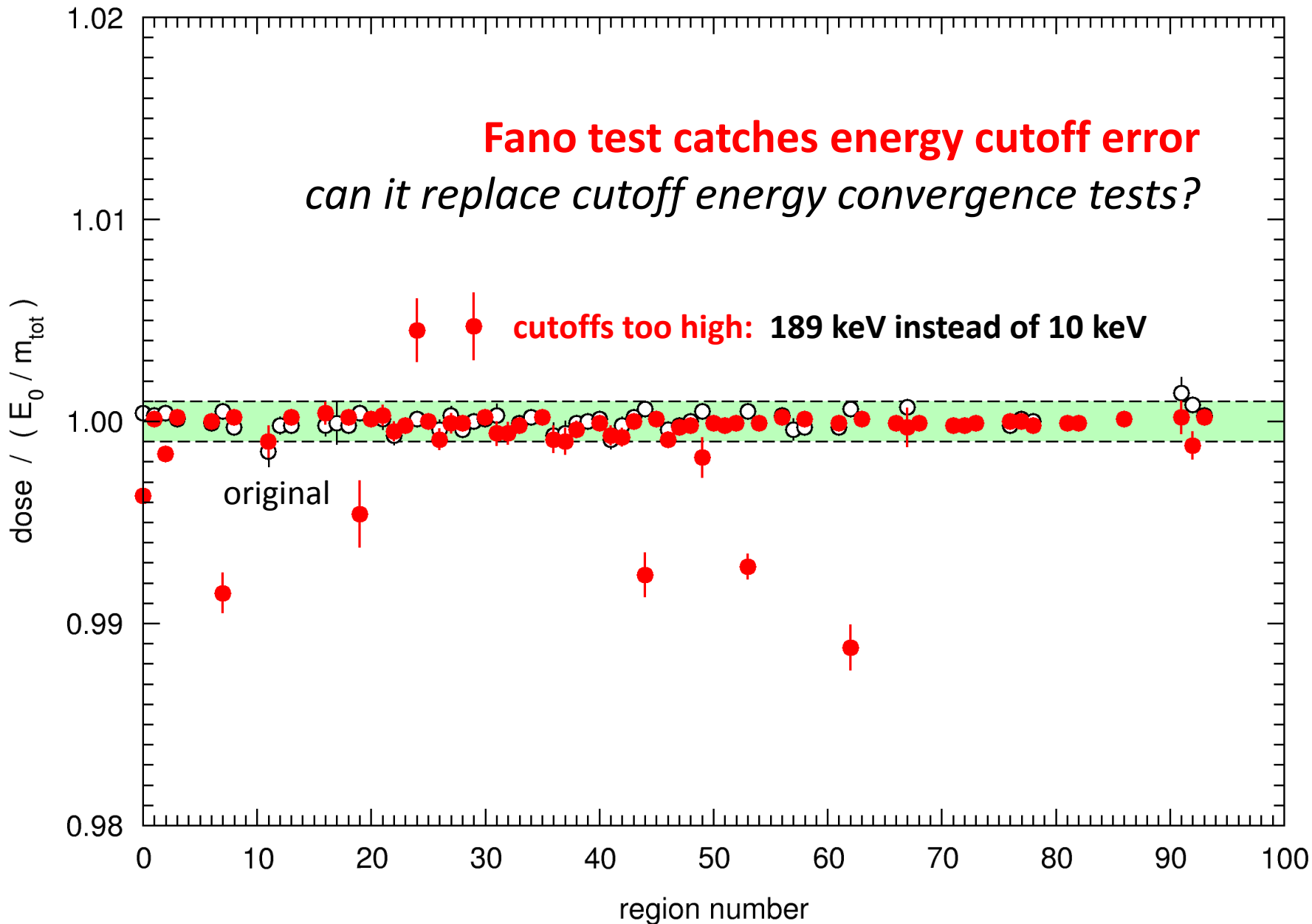
electrons

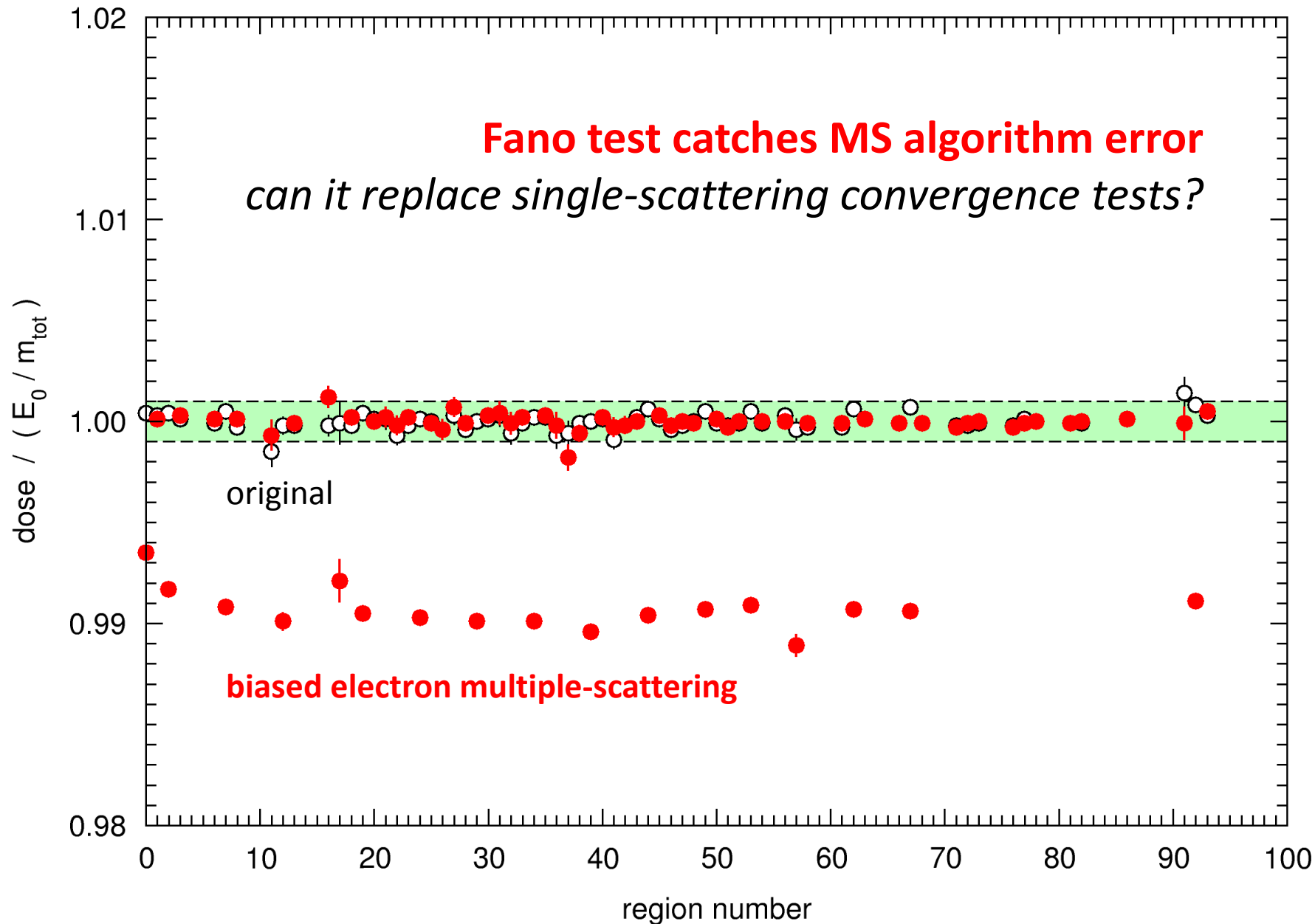


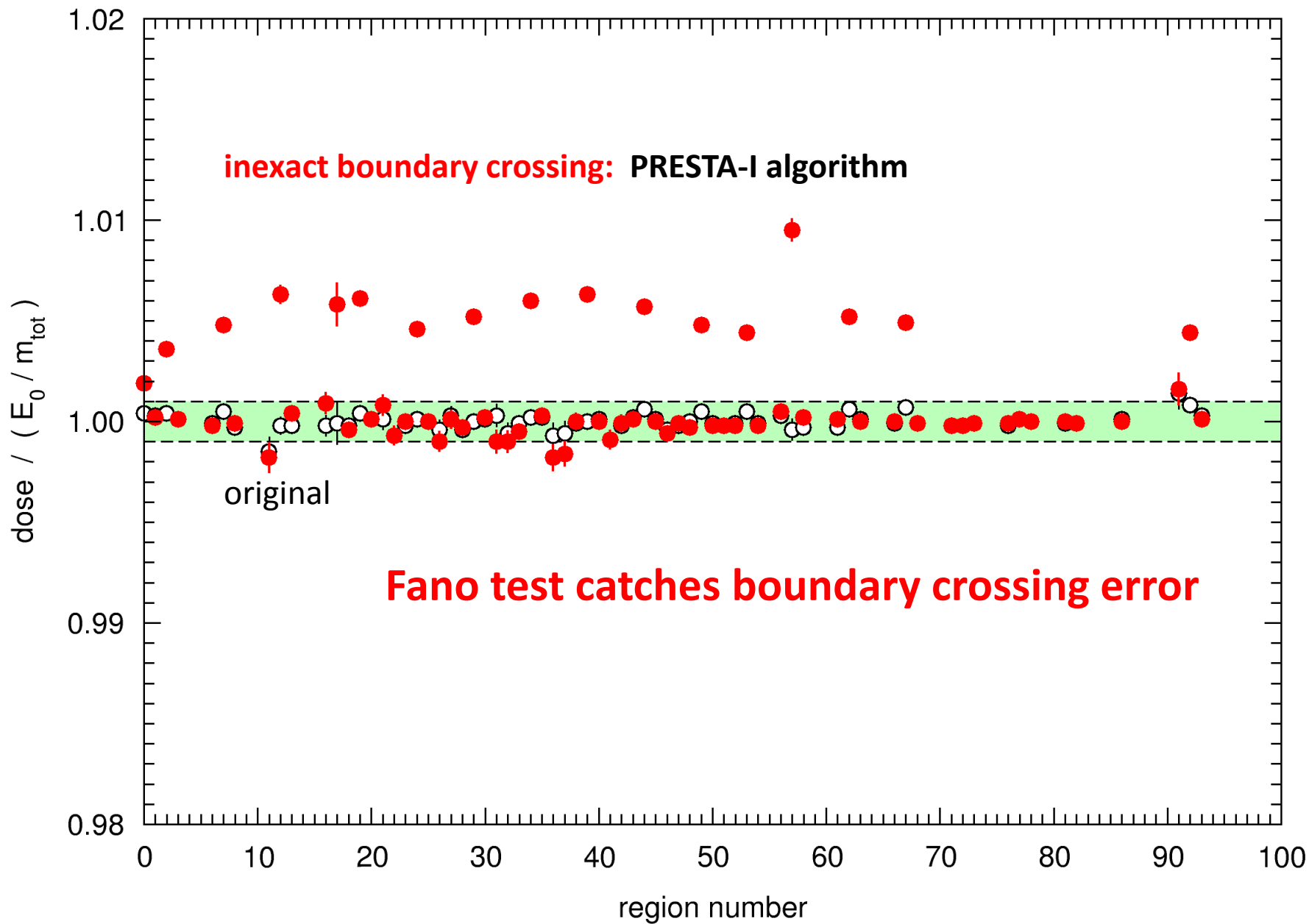


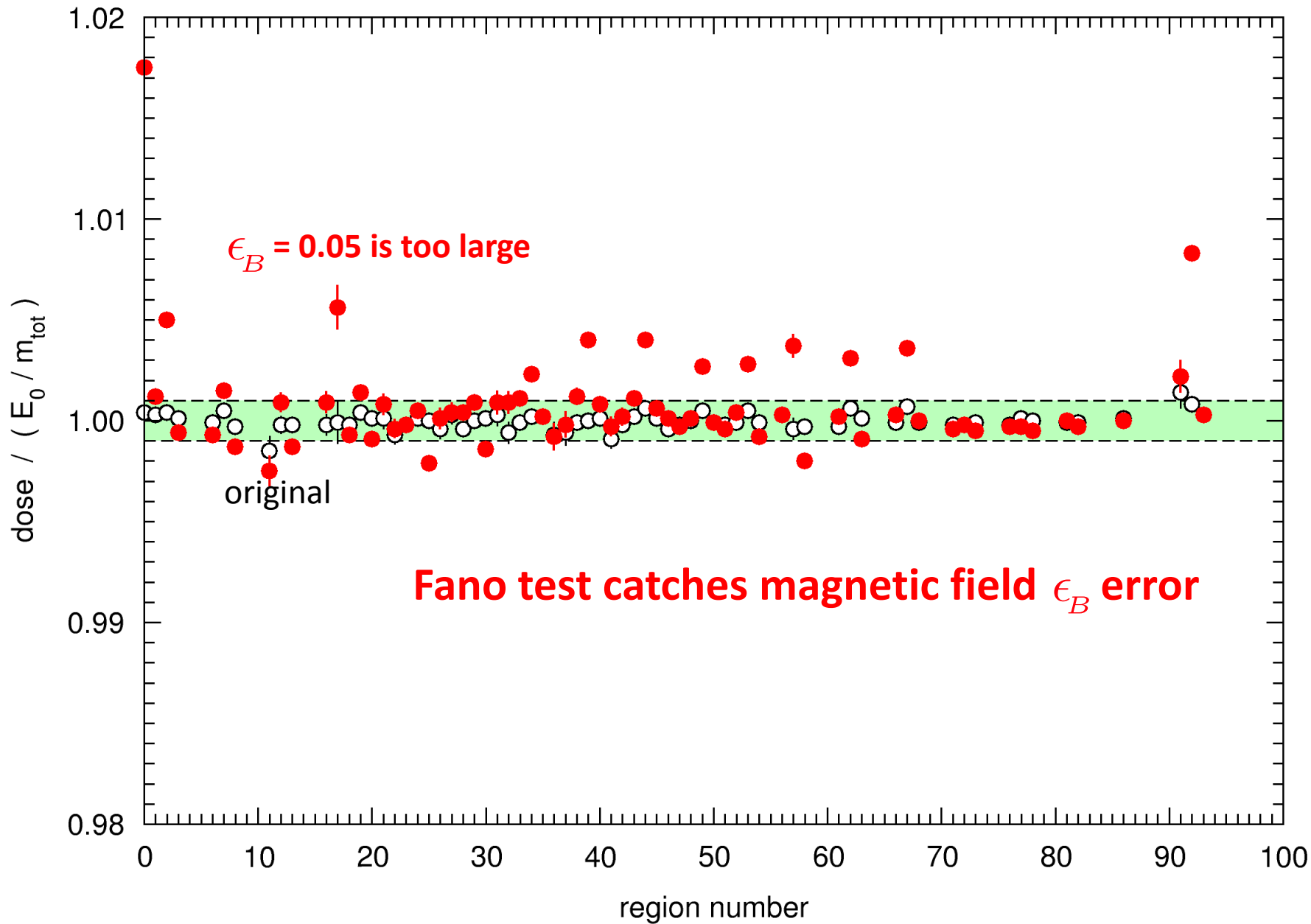


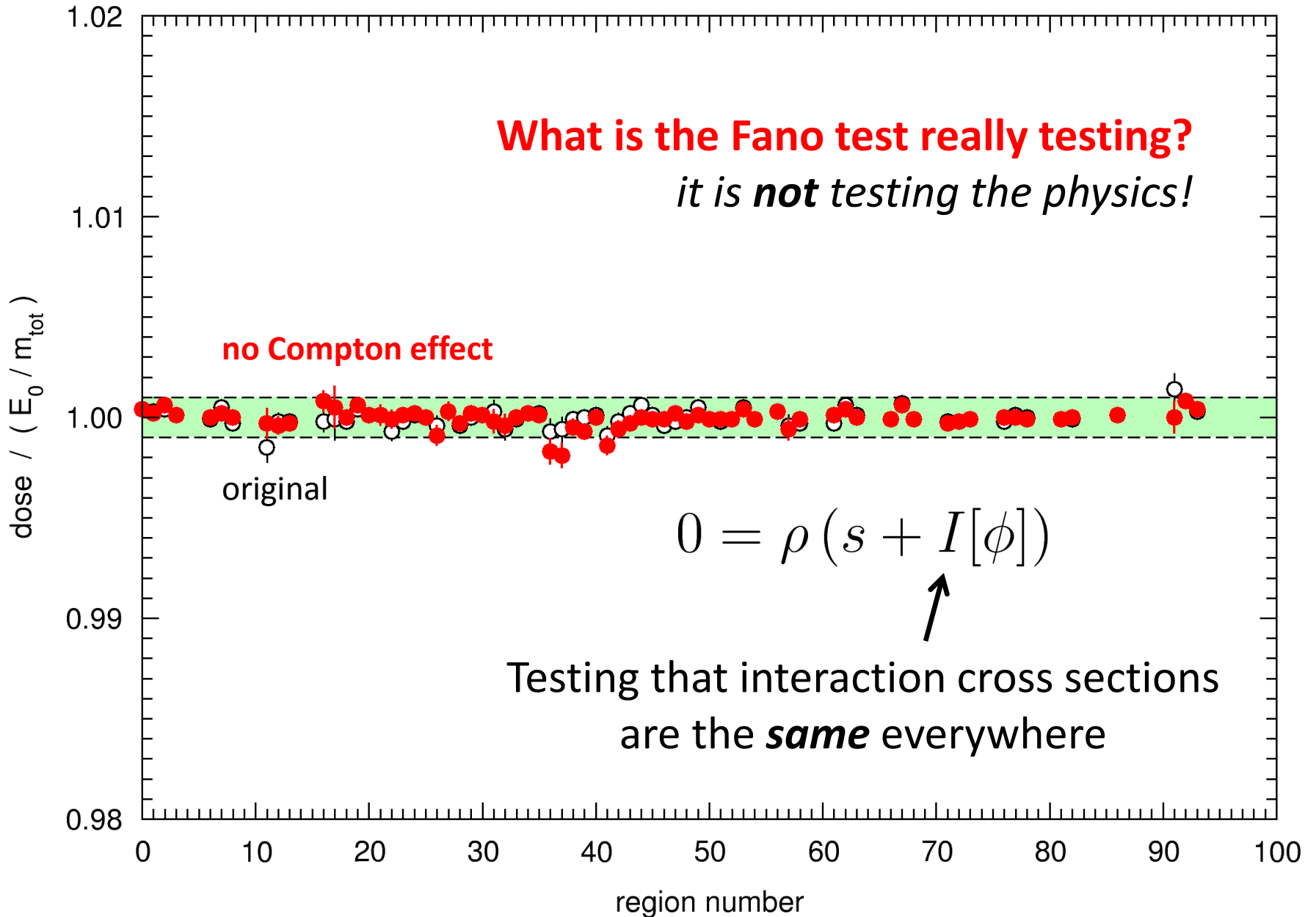
Fano test catches energy cutoff error
can it replace cutoff energy convergence tests?











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Radiation dosimetry in magnetic fields with Farmer-type ionization chambers: determination of magnetic field correction factors for different magnetic field strengths and field orientations

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Appendix. Fano cavity test with the magnetic field macro

To test the Monte Carlo simulations with magnetic field macro for consistency, the Fano test in a magnetic field can be applied (Bouchard and Bielajew 2015, Bouchard *et al* 2015, de Pooter *et al* 2015). For this test, a primary source of 1.25 MeV electrons, isotropically distributed and

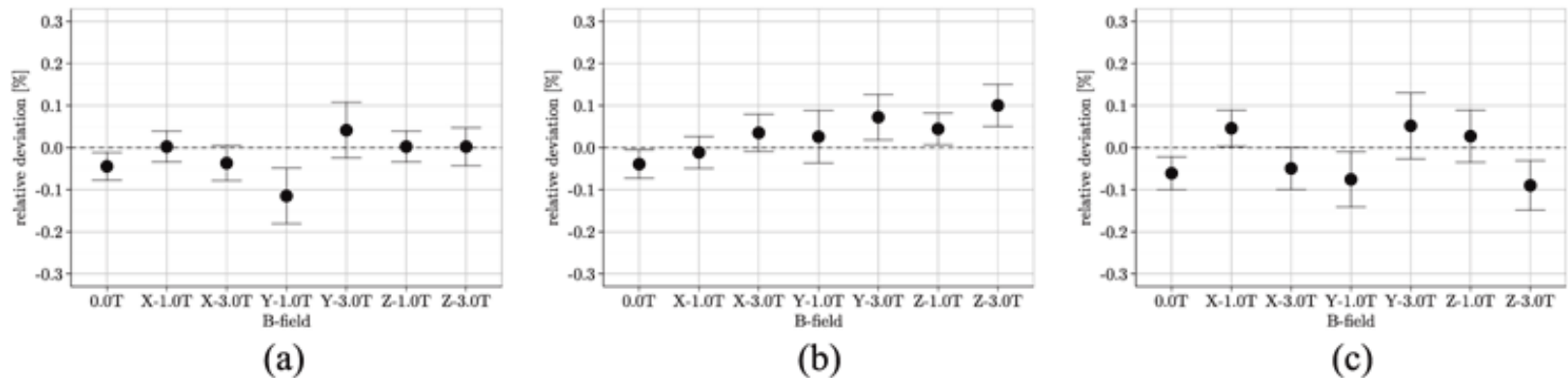


Figure A1. Relative deviation of cavity dose from the expected result for different media and magnetic field geometries. (a) Water. (b) PMMA. (c) Graphite.

All published Monte Carlo simulation results should to be supported by a Fano test calculation.

- Developers should **enable** Fano testing
- Authors should **report** Fano test results
- Reviewers should **request** Fano tests
- Editors should **require** Fano tests

EGSnrc update and Monte Carlo simulation verification

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