

v.16.4.3
07/12/22

Many thanks to Marc and Dave for comments and suggestions

- Account for material lengths in optics
- Update of gas mixture use
- Exponential Abrasion Model (v.16.3)
- LISE $^{++}$ site : secure connection
- Important updates (Reaction mechanism, ETACHA, bugs)
- List of all updates (v.16.2.12-16.4.3)

https://lise.nscl.msu.edu/16/16_4_MaterialAsDrift.pdf

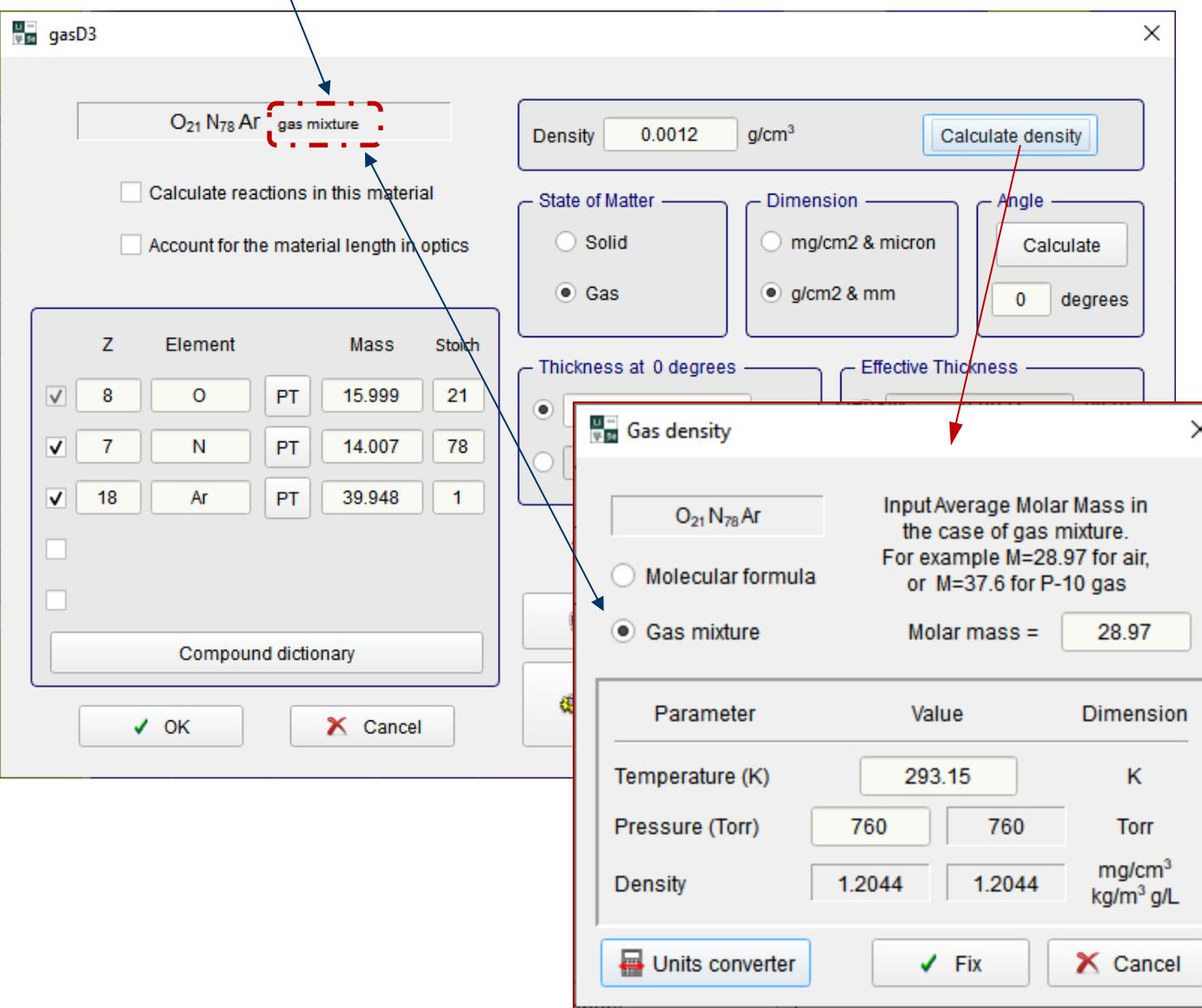
New version allows to account for the material length in optics and as well in time of flight calculations.

The Block compound in this case is considered as an optical drift block in optical matrix calculations.

Pay attention to use this option for compounds inside of a separator!!!

Oleg Tarasov @ MSU 04/06/2022

New feature: “gas mixture” message



Some compounds in
..../lisecfg/compound.dat
file contain the molar
mass value that tells
LISE++ that this gas is a
mixture

Xylene, H₁₀ C₈, 0.8611
* Gases
Acetylene, H₂ C₂, 0.0010825
Air (gas mixture), O₂₁ N₇₈ Ar₁, 0.001205, 28.97
Allene Propadiene, H₄ C₃, 0.0016656
Ammonia, H₃ N₁, 0.00070804
Butane, H₁₀ C₄, 0.0024164

http://lise.nscl.msu.edu/16/16_3_ExpAbrasion.pdf

LISE ⁺⁺_{cute} : Exponential Abrasion Model

v.16.3.1
04/06/22

- Exponential Abrasion Model implementation to LISE ⁺⁺_{cute}
- Use of the Excitation energy fissile nuclei parameters from the BeAGLE calculations in LISE ⁺⁺_{cute}
- LISE ⁺⁺_{cute} Abrasion-Fission 3 EER with the EIC settings
- Comparison of calculation results of different models
- Fission Fragment Kinematics

This PC > Documents > LISEcute > files > examples > afission

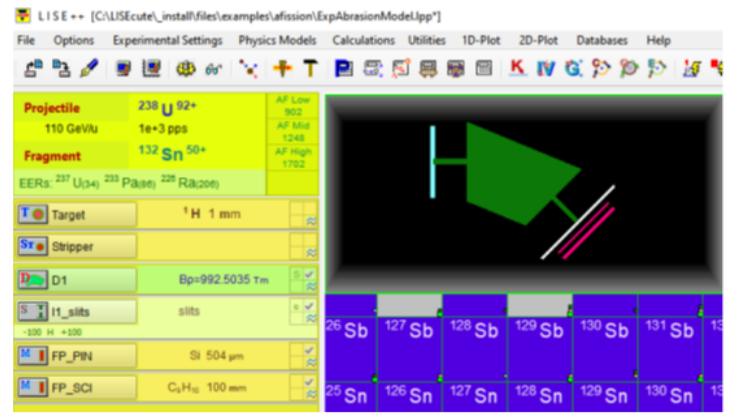
| Name | Status | Date modified |
|-------------------------------|--------|-------------------|
| AF_208Pb_Be.ipp | | 10/1/2021 2:27 PM |
| AF_238U_Be_CS_fromBigRIPS.ipp | | 10/1/2021 2:27 PM |
| AF_238U_Be_highZ.ipp | | 10/1/2021 2:22 PM |
| AF_238U_C.ipp | | 10/1/2021 2:24 PM |
| AF_238U_Pb.ipp | | 10/1/2021 2:27 PM |
| ExpAbrasionModel.ipp | | 4/6/2022 3:07 AM |

LISE ++ [C:\LISEcute\install\files\examples\afission\ExpAbrasionModel.ipp]
File Options Experimental Settings Physics Models Calculations Utilities 1D-Plot 2D-Plot Databases Help

Projectile 238 U 92+ AF Low 902
110 GeV/u 1e+3 pps AF Mid 1248
Fragment 132 Sn 50+ AF High 1702
EERs: 237 U(34) 233 Pa(86) 228 Ra(206)

Target ¹H 1 mm
Stripper
D1 Bp=992.5035 Tm
IT_slits slits
FP_PIN Si 504 μm
FP_SCI C₆H₆ 100 mm

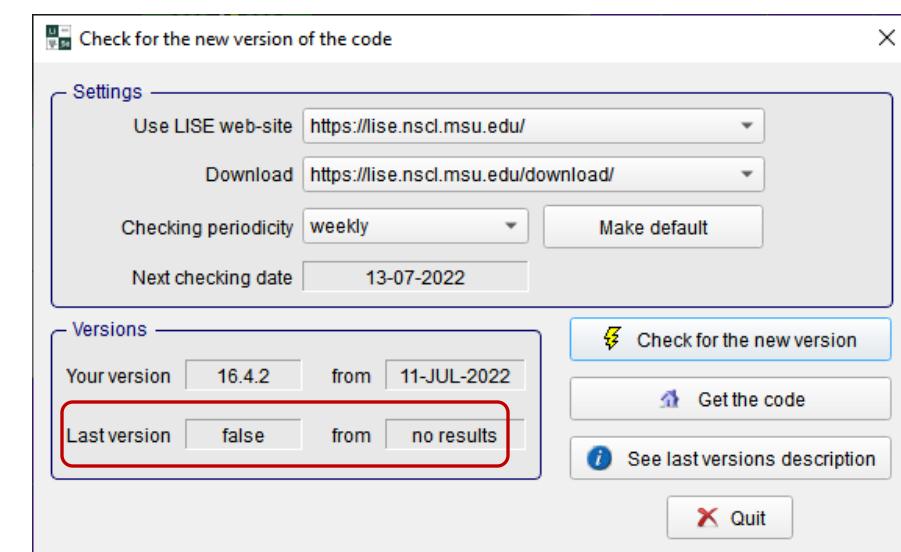
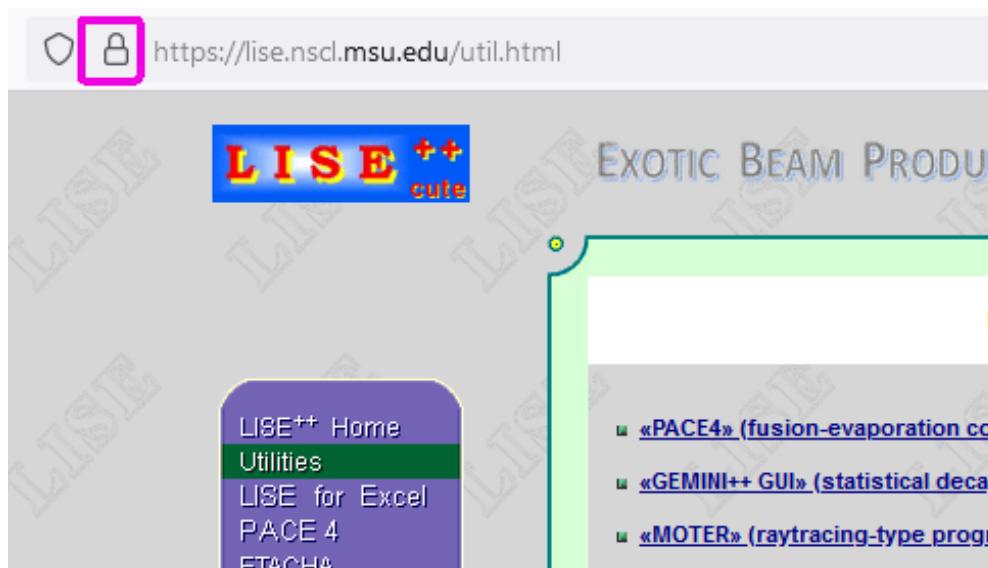
26 Sb 127 Sb 128 Sb 129 Sb 130 Sb 131 Sb
25 Sn 126 Sn 127 Sn 128 Sn 129 Sn 130 Sn 131 Sn



Thanks to Greg Comins!

New approach for visitor statistics.
It allows to get the code downloading statistics.

http → https



Unfortunately, the “New version” dialog is currently not operating outside the FRIB net. It will be solved soon.

v.16.3.7 :

OT's manual correction for light charge particles production cross section can be applied for all PF models
(before it was used only for EPAX2.15-user)

Projectile fragmentation

$^{48}\text{Ca} (140.0 \text{ MeV/u}) + \text{Be} \rightarrow ^{42}\text{S}$

Fragment velocity ✓ Momentum distribution ✓ Cross section, Excitation energy and etc.

Prefragment and Evaporation options Excitation energy for Abrasion-Ablation model

Cross Sections

5 - FRACS 1.1: B.Mei,Phys.Rev.C95(2017)034608 [E=300MeV/u]

Use O.T.'s manual corrections for light charge particles (H, He, Li) production cross sections

"FAST" mode for Abrasion-Ablation calculations* * use this mode only for heavy projectiles as Uranium. Evaporation distribution dimension is equal to 8.

Coefficients for modified EPAX 2.15

| | |
|----------------|----------------------------|
| normalization | p-rich slope |
| U_norm [1.0] 1 | U1 [1.7808] 1.79 |
| n-rich slope | U2 [4.721e-3] 0.00472 |
| Un [1.65] 1.65 | U3 [-1.303e-5] -1.3000e-05 |

Make default OK Cancel Help

v. 16.2.27 04/01/22

The Convolution model : each option has its own σ_0 value

Convolution of Gaussian (Fragmentation) and Exponent (Friction) distributions

$^{48}\text{Ca} (140.0 \text{ MeV/u}) + \text{Be} \rightarrow ^{42}\text{S}$

$$f(p) \approx \exp\left(\frac{p}{\tau}\right) \cdot \left[1 - ferr\left(\frac{p - p_0 + \frac{\sigma_{II}^2}{\sqrt{2}} - shift \cdot \tau}{\sqrt{2} \sigma_{II}}\right)\right]$$

$$\sigma_{II}^2 = (\sigma_0^{conv} \sqrt{\beta_p})^2 \frac{A_F^*(A_p - A_F^*)}{A_p - 1} \quad \tau = \frac{coeff}{\beta} \sqrt{A_F^* E_S}$$

Settings for Gaussian distribution

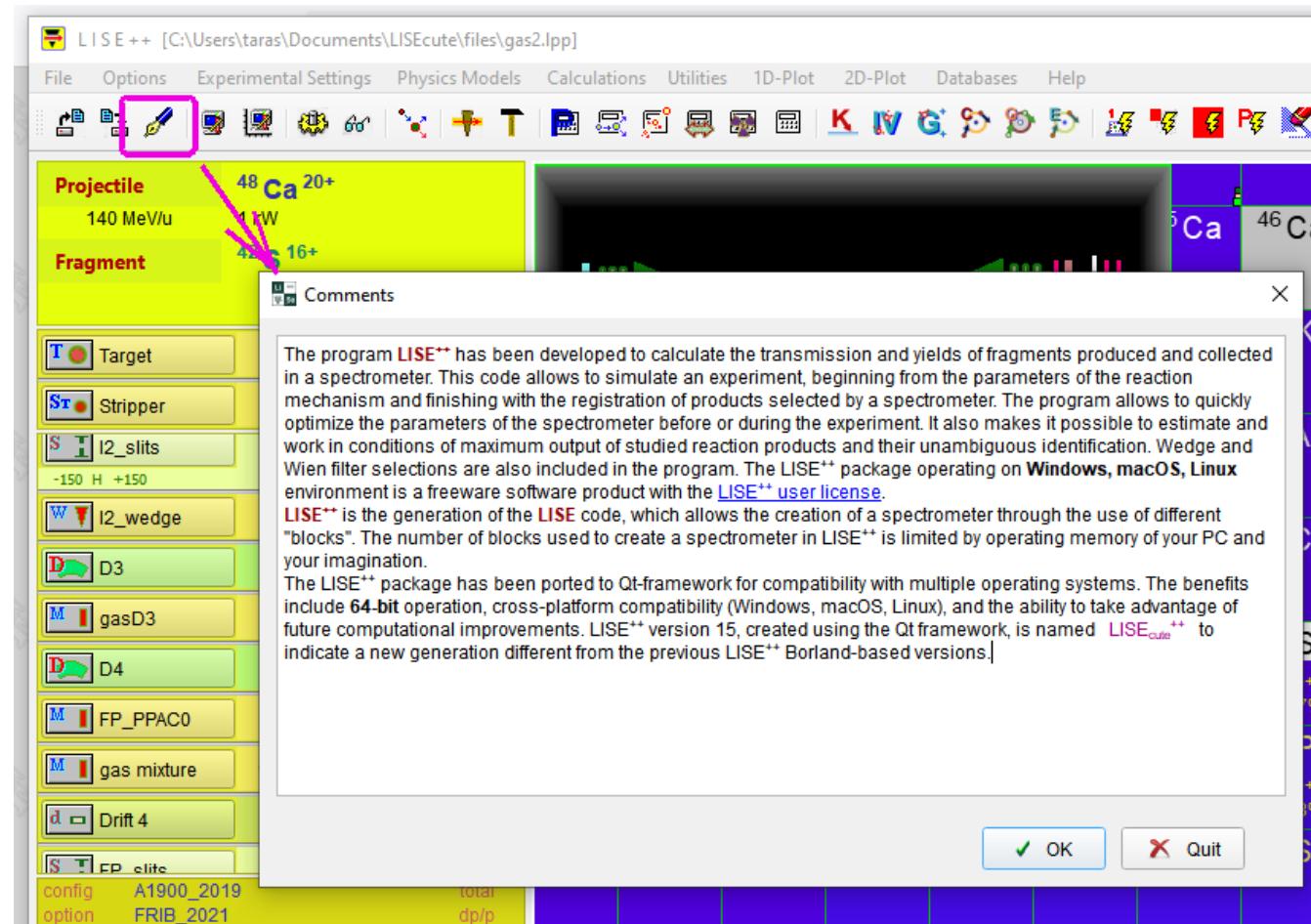
| |
|--------------------------------------|
| $P_0 (\text{MeV/c})$ 22226 |
| V_F / V_B from settings 0.994 |
| Mom.distribution = [1] D.J.Morrissey |
| σ_0 87 MeV/c |
| σ_{II} 244.5 MeV/c (*) |

Settings for convolution

| Separation Energy Model | $E_{separation}$ | σ_0^{conv} | coeff | shift | FWHM / 2.355 (*) | tau | $P(Y_{max})$ | V_F / V_B peak | V_F / V_B mean |
|---------------------------------------|------------------|-------------------|-------|-------|------------------|-------|--------------|------------------|------------------|
| 0. Energy from Qg | 26 | 91.5 | 3.344 | 0.158 | 183.9 | 231.3 | 22079 | 0.997 | 0.993 |
| 1. Excitation from dSurface | 11.7 | 91.5 | 3 | 0.149 | 157.1 | 139.2 | 22137 | 0.998 | 0.996 |
| 2. Excitation from the Abrasion model | 41.2 | 125 | 1 | -1 | 184.2 | 87.1 | 22094 | 0.995 | 0.994 |

$\sigma_0^{conv} = 125 \text{ MeV/c}$ $g = 0.95 \text{ MeV/fm}^2$ (*) - with y-factor

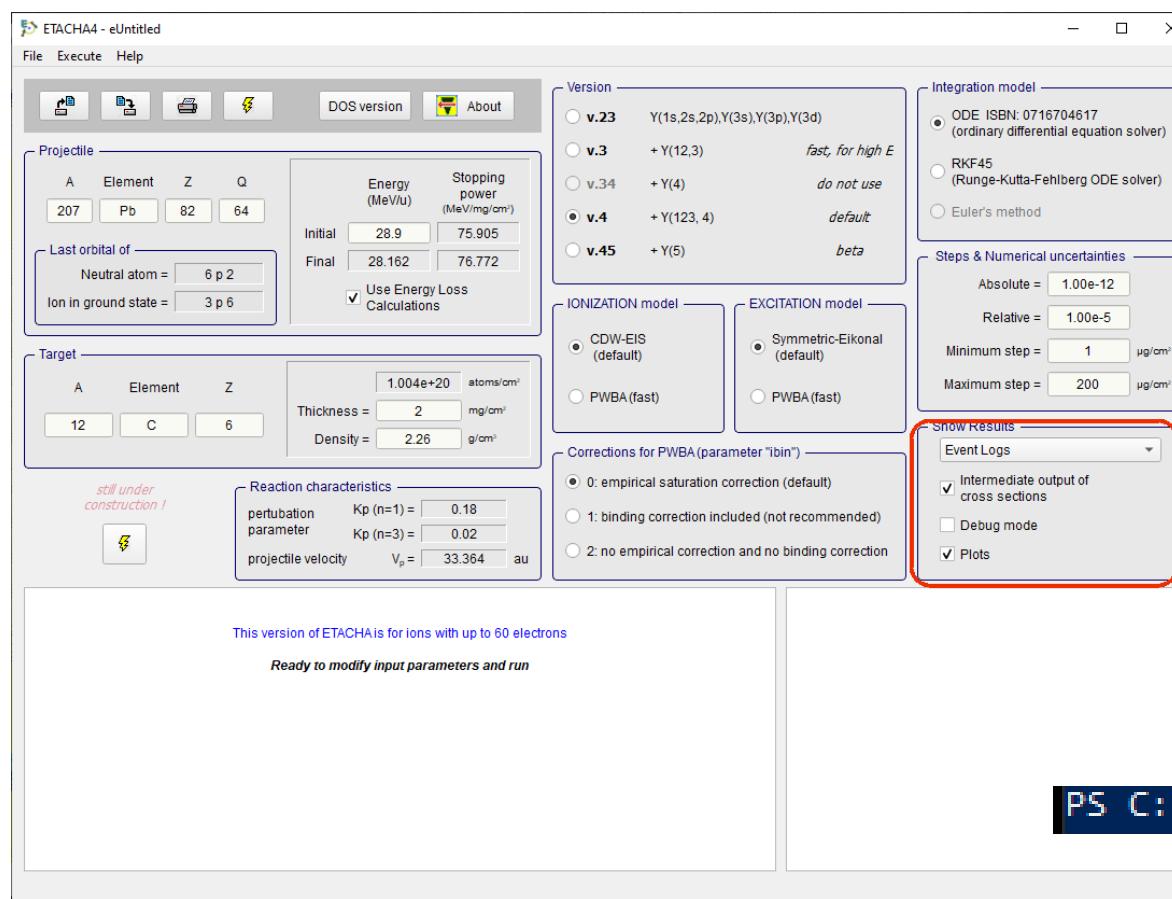
Plot 1D Convolution Analysis Make default OK Cancel Help



-
- 16.2.23 03/29/22
comments action in toolbar; comment icon in file menu
-
- 16.2.22 03/29/22
comments: summarize all comments from files in the case of
append operations
-
- 16.2.21 03/29/22
comments: summarize all comments from files in the **lise files list**
-

- ETACHA 4.4.11 05/30/22 batch mode with command string argument -r
 ETACHA 4.4.10 05/30/22 Correction with a density value input from file
 ETACHA 4.4.09 05/30/22 Plot option in etacha file

The option -r suppresses the graphs and the input window.



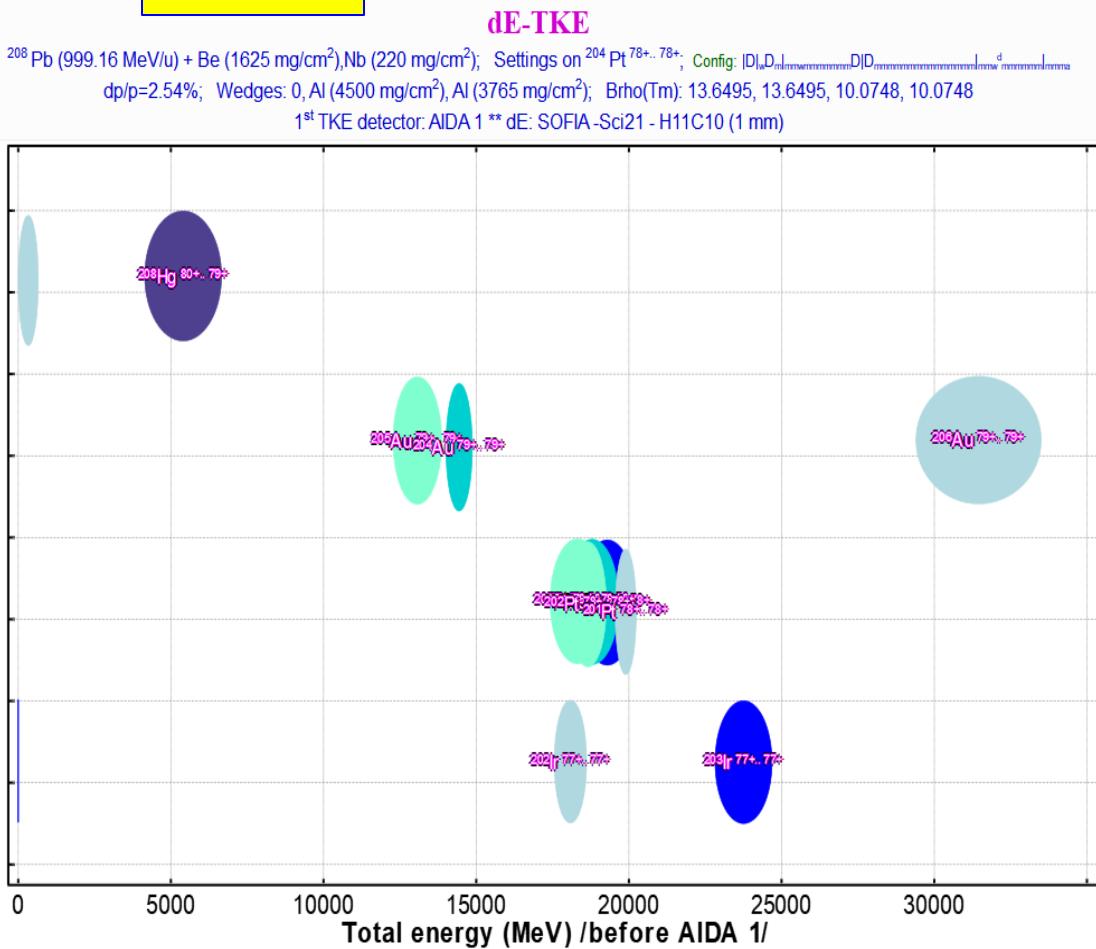
HW's request

```
PS C:\Program Files\LISEcute> .\ETACHA4.exe D:\test.etacha -r
```

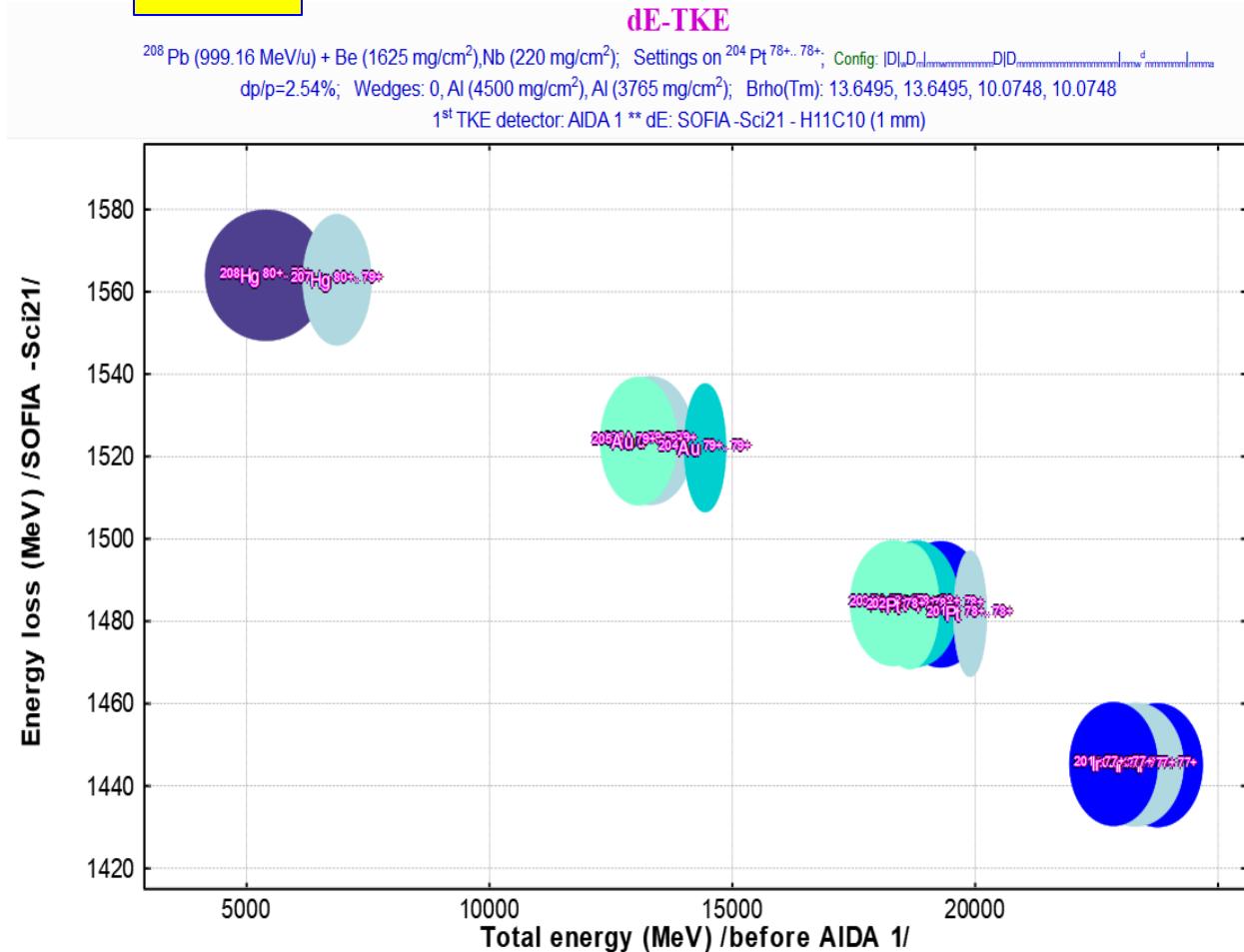
DJM's request

[corrected] Incorrect extrapolation of energy for isotopes completely outside the slits in the case of transmission calculations by the “ellipse” method

before



after



16.3.33 07/09/22

fixed: w_Graph_Envelope crash. (MH)
mostly it has been moved from char* to QStringSize

16.3.11 05/30/22

InitFN_AF_Dlg::CmPrintTitles correction with titles

16.3.9 05/30/22

misspelling corrections in dialogs d_CN0_scanningEnergy and
d_CN0_scanningBrho (MH)

16.3.8 05/11/22

two missing .toString().c.str() in c_Plot_MC.cpp (DB)

16.3.6 04/23/22

correction in angular distribution for the fusion-residual (p,g) case (119I from
118Te beam)

16.3.5 04/23/22

correction in the fusion-residual case for very thin target

16.3.4 04/20/22

* Fixed: crash (only in debug mode) in O-AA_manage.cpp

16.3.2 04/12/22

QSstrcpy and QSstrncpy modifications in win_utilString.cpp (TZ)

16.2.38 04/06/22

initialization beam.Shapes[i] to 1

16.2.30 04/01/22

Fixed: memory leak in read/write blocks

16.2.29 04/01/22

Fixed: bug in file append procedure with high order matrices

16.2.14 03/22/22

Fixed: Results->Transmission A,Z,q1 (summarized by reaction) ==>
bug with stripper charge state value

16.2.13 03/22/22

Fixed: Random wrong initialization of root directory (only in Windows)
(MH)

16.2.12 03/22/22

Fixed: bug in d_Setup_Array after privateName value retyping to
QString (MH)

List of all updates (v.16.2.12-16.4.3) [1]

| | | |
|--|---|--|
| 16.4.3 07/12/22 PseudoMC update for UseLengthOptics=1 | 16.3.24 07/05/22 Config_read & write updates for block_material length use in optics new function templates in BLOCK_Compound int getUseLengthOptics(){return UseLengthOptics;}; void setUseLengthOptics(int init){ UseLengthOptics = init>0 ? 1 : 0;}; | 16.3.9 05/30/22 misspelling corrections in dialogs d_CNO_scanningEnergy and d_CNO_scanningBrho (MH) |
| 16.4.2 07/11/22 update DrawPlotFile functions for pure ascii output (no Borland, no HTML) (StSc) | 16.3.23 07/05/22 Config_read & write updates for gas mixture molar value | 16.3.8 05/11/22 two missing .toString().c.str() in c_Plot_MC.cpp (DB) |
| 16.4.1 07/11/22 changing middle version | 16.3.22 07/03/22 update d_Thickopt_compound to read gas mixture molar value | 16.3.7 05/06/22 OT's manual correction for light charge particles production cross section can be applied for all PF models |
| 16.3.36 07/09/22 modification LengthNameCompound from 61 to 76 modification LengthNameCompoundPlusMass from 78 to 100 | 16.3.21 07/03/22 Ctarget::input_thickness call in Ctarget | 16.3.6 04/23/22 correction in angular distribution for the fusion-residual (p,g) case (119I from 118Te beam) |
| 16.3.35 07/09/22 update FillTitlePlot for Wedge graph name and Brho greek | 16.3.20 06/30/22 Option Show/Hide Cancel-button in Gauge button. used for initial Table creation | 16.3.5 04/23/22 correction in fusion-residual for very thin target |
| 16.3.34 07/09/22 new BLOCK function bool isLengthBlock() connected to material UseLengthOptics=1, update BLOCK_util for isLengthBlock() | 16.3.19 06/30/22 Material dialog - the Z cell is hidden if this component does not exist | 16.3.4 04/20/22 * attempt to use SSL protocol in the new version dialog * Fixed: crash (debug mode) in O-AA_manage.cpp because of char buff[20]; |
| 16.3.33 07/09/22 fixed: w_Graph_Envelope crash. (MH) mostly it has been moved from char** to QStringSize | 16.3.18 06/30/22 modification of compound.dat file for molecular mass molecular-Mix mass flag in Ctarget class --> later w was erased | 16.3.3 04/15/22 https: instead http: for LISE site links |
| 16.3.30-32 07/08/22 MC envelope update for UseLengthOptics=1 TOF p_Block_util update for UseLengthOptics=1 p_Block_polyном update for UseLengthOptics=1 | 16.3.17 06/30/22 new NNDC isotope link | 16.3.2 04/12/22 QSstrcpy and QSstrncpy modifications in win_utilString.cpp (TZ) |
| 16.3.29 07/08/22 MC_options => always opt->MC_Zaxis_material=1, now opt->MC_Zaxis_material is used as flag in PassMaterialBase for UseLengthOptics=1 | 16.3.16 06/29/22 d_Thick : new CheckBox "Use Length in Optics BLOCK_Compound : int UseLengthOptics | 16.3.1 04/06/22 new middle version http://lise.nscl.msu.edu/16/16_3_ExpAbrasion.pdf |
| 16.3.28 07/08/22 d_Setup_Optics -> CmBeamSigmaPlot -> update for UseLengthOptics | 16.3.15 06/29/22 new icon plotsaveMC, plot_save icons modification (StSc) | 16.2.39 04/06/22 \files\examples\afission\ExpAbrModel.lpp -- settings for Exponential abrasion with EIC parameters |
| 16.3.27 07/08/22 d_Setup_Optics -> CmMatrixEnvelope -> update for UseLengthOptics | 16.3.14 06/29/22 c_Plot_2 -> Actions : Save Plot for mode=20 (Ellipse) | 16.2.38 04/06/22 initialization beam.Shapes[i] to 1 |
| 16.3.26 07/08/22 d_Setup_Optics -> CmMatrixFile -> update for UseLengthOptics correction in p_Block_Compound for om->Unit in the case of drift | 16.3.13 06/29/22 c_Graph2_goodies : PassSlits -> CutSpace -> USeAnalysis=false for isotopes out of slits | 16.2.37 04/06/22 modifications in the Abrasion-Fission dialog and plot |
| 16.3.25 07/05/22 d_Thick dialog update for UseLengthOptics, W_Scheme update for UseLengthOptics, different color and size for materials with lengths o_Optics update for UseLengthOptics in RecalculateGOM, p_Block_Compound for UseLengthOptics in blockLength for optical matrix | 16.3.12 06/26/22 c_Graph2_goodies : PassSlits only for optical blocks | 16.2.36 04/06/22 double Sigma0convolution parameter --> double Sigma0convolution() function |
| | 16.3.11 05/30/22 InitFN_AF_Dlg::CmPrintTitles correction with titles | 16.2.35 04/04/22 ExpSlope abrasion: Nuclide CS map development |
| | 16.3.10 05/30/22 Physical Calculator : increasing output precision (AG) | 16.2.34 04/04/22 ExpSlope abrasion: including AA compare project, ApfExcitation dialog modification |

List of all updates (v.16.2.12-16.4.2) [2]

16.2.33 04/04/22
COSY map default file extensions are "map" and "mat" for loading

16.2.32 04/02/22
Update of FRIB_2022.lopt (default option file)

16.2.31 04/01/22
global revision in read/write blocks with new elements of Zagolovok and Vegas classes

16.2.30 04/01/22
Fixed: memory leak in read/write blocks

16.2.29 04/01/22
Fixed: bug in file append procedure with high order matrices

16.2.28 04/01/22
Read/Write --> struct Zagolovok to class, struct Vegas -> class
zero initialization of char* in these classes
High order class : HOL - new function init(const *str)
no more some functions in Vegas class (alloc, free, read_func)

16.2.27 04/01/22
Convolution model : each option has its own sigma0 value

16.2.26 04/01/22
COSY map default file extension is "map"

16.2.25 04/01/22
new default option file FRIB_2022.lopt
correction in "writing rays to file" order in option file

16.2.24 04/01/22
No more modal property for the File comments dialog
writing high order optics element format increased to %.e5

16.2.23 03/29/22
comments action in toolbar; comment icon in file menu

16.2.22 03/29/22
comments summarize comments in the case of append operations

16.2.21 03/29/22
comments summarize comments of all file in file list

16.2.20 03/27/22
direct multipole dialog edit after clicking on the table in the Optics Setup dialog

16.2.19 03/26/22
global substitution for strReactionRegl function

16.2.18 03/26/22
new Functions QString strReactionRegl(Clement* el, double energy, Compound* compound, Clement* el2, bool);

16.2.17 03/26/22
class Element and Compound -- implementation of QString for name

16.2.16 03/24/22
new Abrasion model : Exponential -- only Dialog and initialization

16.2.15 03/23/22
new function chargeStateTrans (const Ccalc* cc) : product of all charge state transmission values for one ion

16.2.14 03/22/22
Fixed: Results->Transmission A,Z,q1 (summarized by reaction) ==> bug with stripper charge state value

16.2.13 03/22/22
Fixed: Random wrong initialization of root directory (only in Windows) (MH)

16.2.12 03/22/22
Fixed: bug in d_Setup_Array after privateName value retyping to QString (MH)
ARIS.blif file has been updated by MH
