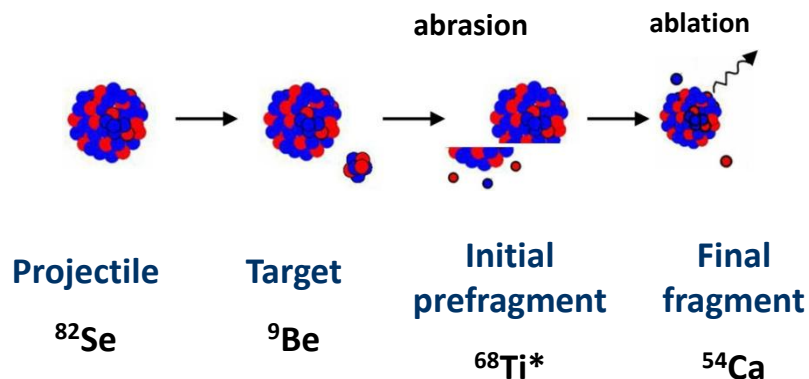
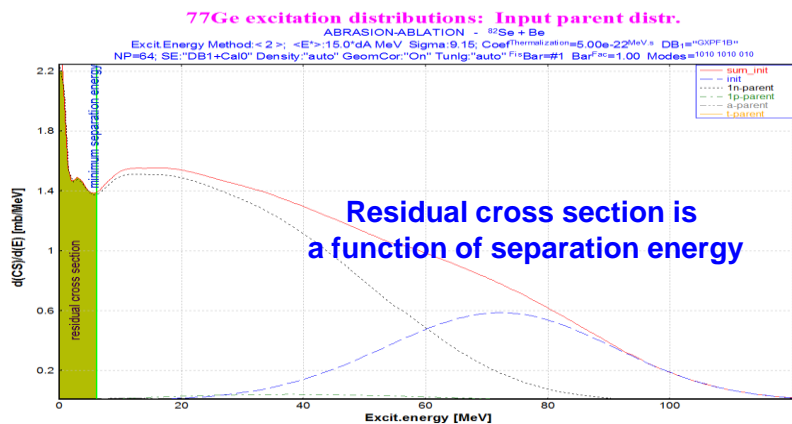


Projectile fragmentation

Abrasion-ablation model



Ablation step (Evaporation cascade) uses a mass table to obtain separation energies!!



Nuclear Physics A531 (1991) 709-745
 North-Holland

A REEXAMINATION OF THE ABRASION-ABLATION MODEL FOR THE DESCRIPTION OF THE NUCLEAR FRAGMENTATION REACTION*

J.-J. GAIMARD¹ and K.-H. SCHMIDT

We propose the following approach for calculating the excitation energy of the prefragment: The nucleons are bound in the potential well of the nucleus. During the abrasion, the orbits of the nucleons not removed are preserved. This is suggested by the short time span of the abrasion, in the order of $(2-5) \times 10^{-23}$ s. By the abrasion, a certain number of single-particle levels is vacated, and the excitation energy is given by the sum of the energies of these holes with respect to the Fermi surface. For a quantitative estimate we take a Woods-Saxon potential with an average depth of -47.4 MeV for neutrons and protons²⁹. We neglect that the density reduction caused by the abrasion decreases the potential depth because it is reestablished after the contraction to normal nuclear density. The energy generated by one hole varies between 0 and 40 MeV depending on its position if we assume a Fermi energy around -7.4 MeV. In order to calculate the mean energy induced by one hole in the potential well below the Fermi surface, we use the single-particle level density $g(\epsilon)$ of the Woods-Saxon potential which can be described approximately by $g \sim \epsilon$ [ref.³⁰] if ϵ is the single-particle energy counted from the bottom of the potential well, and we assume that the probability for generating the hole is the same for each level. This statistical hole-energy model gives an average excitation energy of 13.3 MeV per hole.

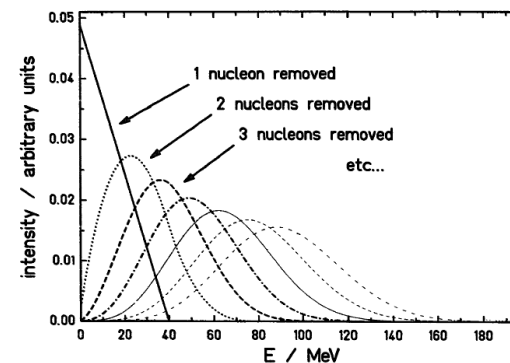


Fig. 2. Excitation-energy distributions as calculated by the diabatic model (this work) for different prefragments, after the abrasion of 1-8 nucleons from the projectile nucleus.

v. 9.4.44 06/09/12
thermalization
of excitation energy

In the following we will investigate whether the basic assumption of our proposed model is justified, namely that the orbits of the nucleons of the prefragment are untouched by the abrasion process, that means whether the abrasion process at relativistic energies is diabatic. For this purpose we will estimate the part of the excitation energy which is thermalized during the abrasion, that means the part which is equilibrated and modifies the orbits of the nucleons of the prefragment. According to calculations of Bertsch³¹), Nörenberg³²) estimated an intrinsic thermalization time, also called intrinsic equilibration time, by the following relation:

$$\tau_{intr}(t) = 2 \times 10^{-22} \text{ MeV} \cdot s / e^*(t), \quad (3.1)$$

User cross-section analysis
using Abrasion-Ablation model

http://lise.nsci.msu.edu/7_5/lise++_7_5.pdf#page=85

$$Final = w_1 \cdot Lo\chi^2_{local} + w_2 \cdot LoD_{local} + w_3 \cdot Lo\chi^2_{total} + w_4 \cdot LoD_{total}$$

$$\text{where } Lo\chi^2 = \ln(\chi^2) \text{ and } LoD = \sum_{i=1}^N \left| \log_{10}(y_{exp}) - \log_{10}(y_{calc}) \right| / N$$

Global fit by AA for all CSs obtained in the experiment allows to deduce the excitation energy parameters

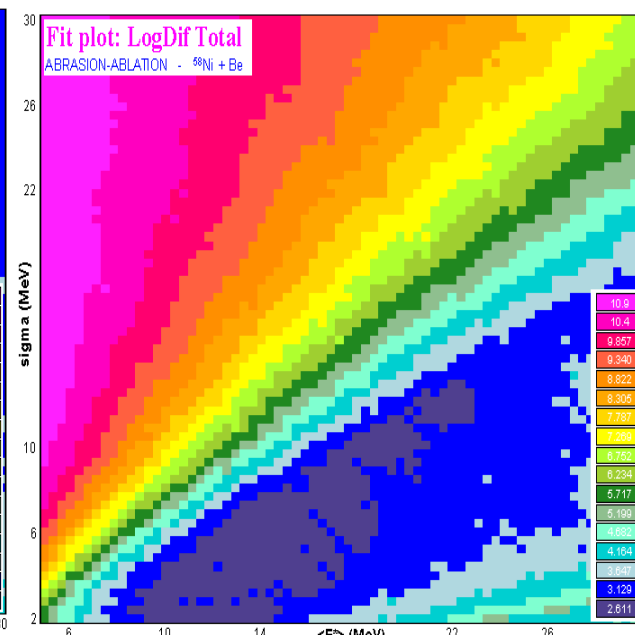
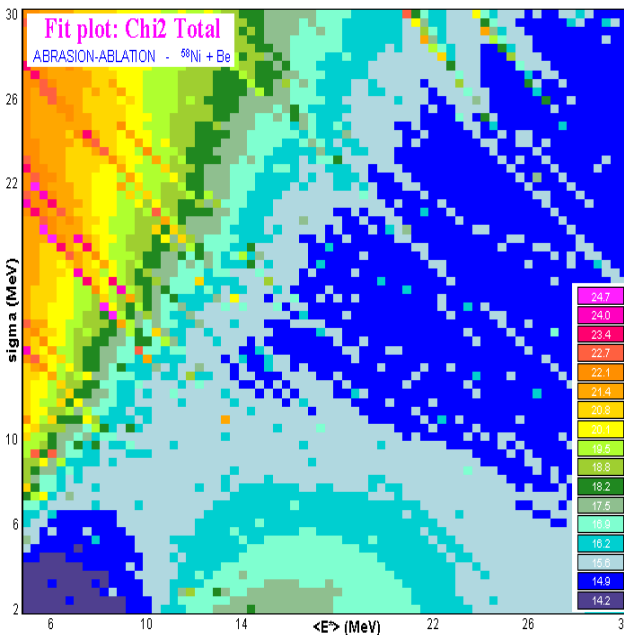
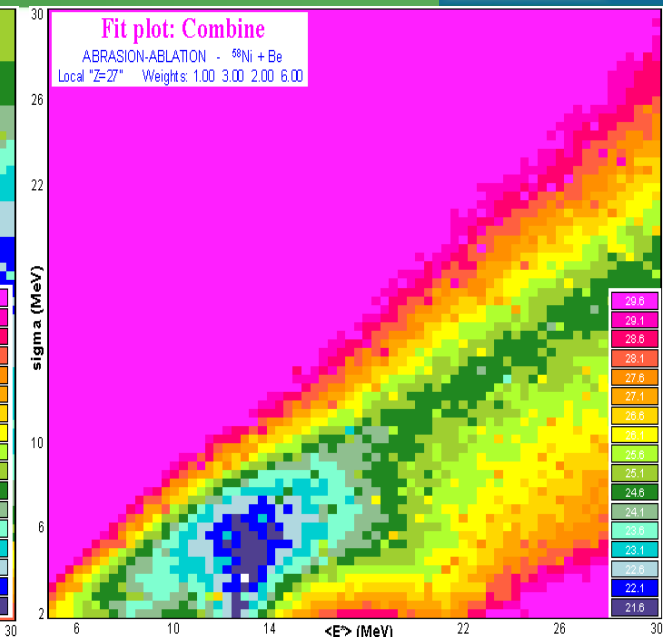
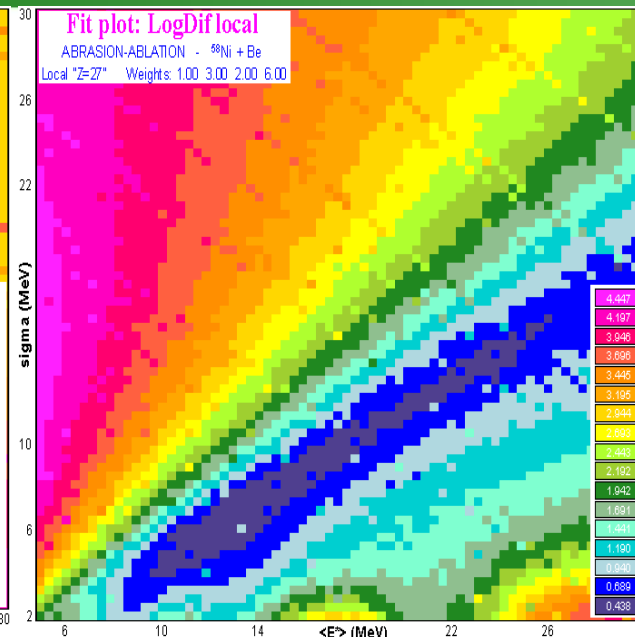
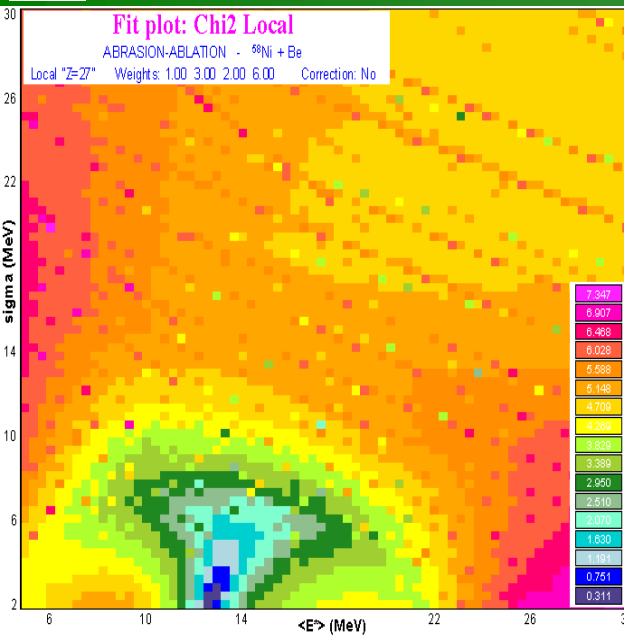


Figure. Comparisons of experimental data [Bla94] and Abrasion-Ablation calculations for the $^{58}\text{Ni}(650\text{AMeV})+\text{Be}$ reaction. Left top and middle top plots represent comparisons for a Local line, which was set to $Z=27$. Left bottom and middle bottom plots are done for the global data set. The right top plot shows the sum of all four models, using coefficients 1,3,2,6 (see insertions in plots). Using this analysis it is possible to conclude that the best parameters to use in the AA model to describe these experimental data are $\langle E^ \rangle = 13$ and $\text{sigma} = 4$ MeV.*

During analysis of GSI's ^{238}U , RIKEN's ^{238}U , MSU's ^{82}Se experiments there was significant modification of LISE⁺⁺ AA:

- Improving/Fixing problems (interpolation, new methods),
- new properties (excitation energy thermalization and etc),
- new utility: Initial prefragments plot, Decay Analysis utility update,
- new mass tables (AME2011, GXPF1B), unknown masses extrapolation procedure update and so on

Evaporation calculator

Initial nucleus: 82Se
Excitation energy window: Lower = [] MeV (gaussian), Upper = [] MeV (rectangle)
Initial nucleus production cross-section = [] mb
make calculations down to Z = [1]

Modes:
 Fragmentation of beam (Abrasion-Ablation)
 Excited nucleus evaporation
 Load initial conditions from file

2D-plots:
 Final Evap. Residue CS
 Decay channel analysis
 Fission channel CS
 Temperature
 Break-up channel CS
 Fission Excitation Energy

CALCULATE Evaporation settings

Final nucleus: A=54, Element=Ca, Z=20, stable
 Table of Nuclides
 Excitation energy plot

Final fragment production cross-section: 6.14e-8 mb
 Initial production CS of Final fragment: 1.04e+0 mb
 Cross section from EPAX 2.15: 2.79e-6 mb

Minimum separation energy (SE): 3.59 MeV
 Minimum sum of (SE + deduced effective Coulomb barrier): 3.59 MeV
 Fission barrier at L=0: 46.89 MeV

Average values: < E_x > = 220.69, < T > = 5.39

PARENT	3.49e+0	4.1e-1	1.78e-1	2.01e-1			4.28e+0	55Ca
Decay modes	1n	1p	alpha	t			Break-up	sum
DAUGHTER	4.95e+0	1.64e-1	5.03e-2	1.61e-1			5.32e+0	53Ca

N° of all calculated nuclei: 1331

Sum	1.09e+4	2.34e+3	7.84e+2	1.97e+2			Initial	Residues	Fission	Break-up
							1.66e+3	1.66e+3		1.16e-2

Output cross-section file: [] 08234_00904.lcs [Browse] [Show]
 Output file of parent-daughter: [] 08234_00904.lpd [Browse] [Show]
 Fission CS output file: [] 08234_00904.lcs [Browse] [Show]

Initial Prefragments Plot for final nucleus [] [Help] [Quit]

54Ca excitation distributions: Input parent distr.

ABRASION-ABLATION - $^{82}\text{Se} + \text{Be}$

Excit.Energy Method: < 2 >; < E* >: 15.0*dA MeV Sigma: 9.15; Coef^{Thermalization}=5.00e-22 MeV.s DB₁="GXPF1B"
 NP=64; SE:"DB1+Cal0" Density:"auto" GeomCor:"On" Tunlg:"auto" FisBar=#1 BarFac=1.00 Modes=1010 1010 010

Decay channel analysis

Channels: Current mode: Initial CS -> [S residue] / [Sr total]

Mode:

Absolute value

Ratio

Value (selected channel):

S parent

S residue

S evap

(So total) evap

Take ratio to:

S parent

Si total

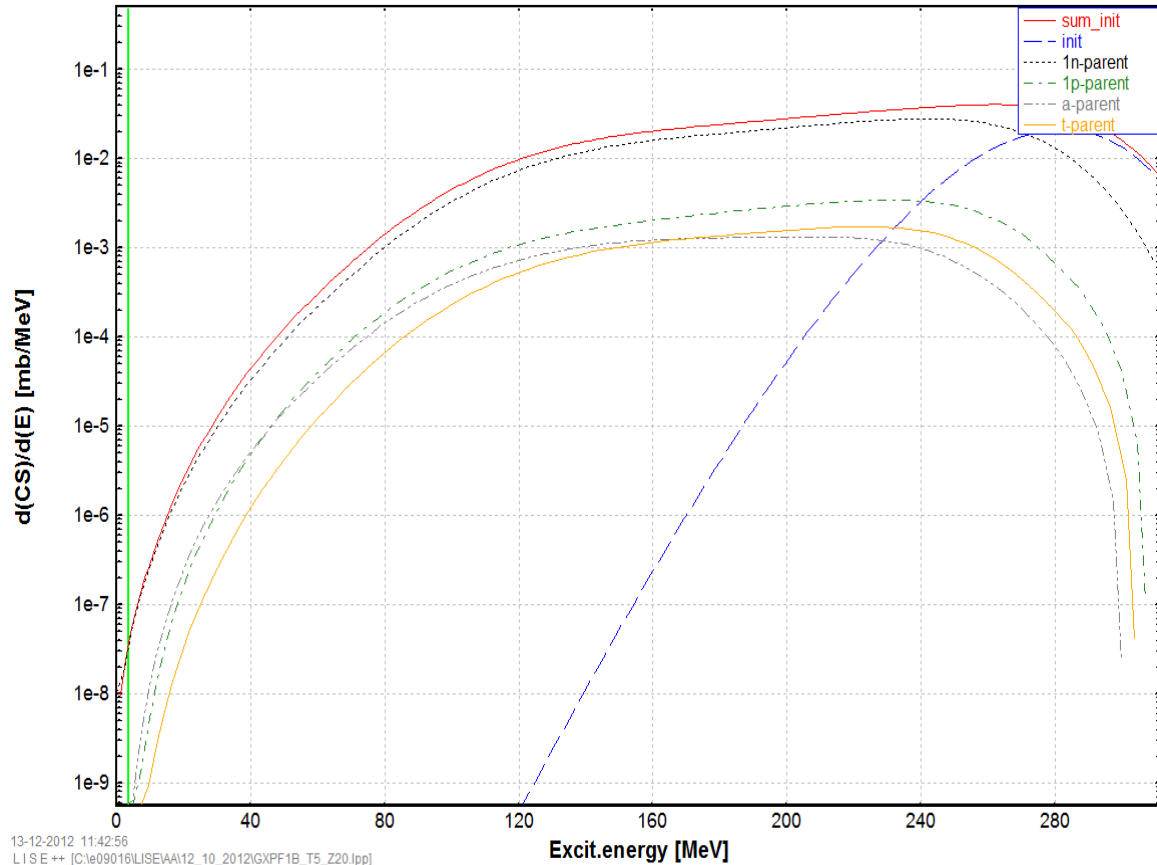
Sr total

So total

Description:

SE min = minimum separation energy
 S residue = residue cross-section from this channel
 S evap = decay channel in the daughter nucleus from this parent channel
 S parent = excitation function incoming from the parent nucleus
 Si total = sum of all incoming (from parents) channels including Init CS
 Sr total = sum of all residue cross-sections
 So total = sum of all outgoing (into daughters) channels
 (So total) evap = decay channel into the daughter from ALL parent incoming channels

Fission and Break-up are only output channels; Init (AA) is only input channel

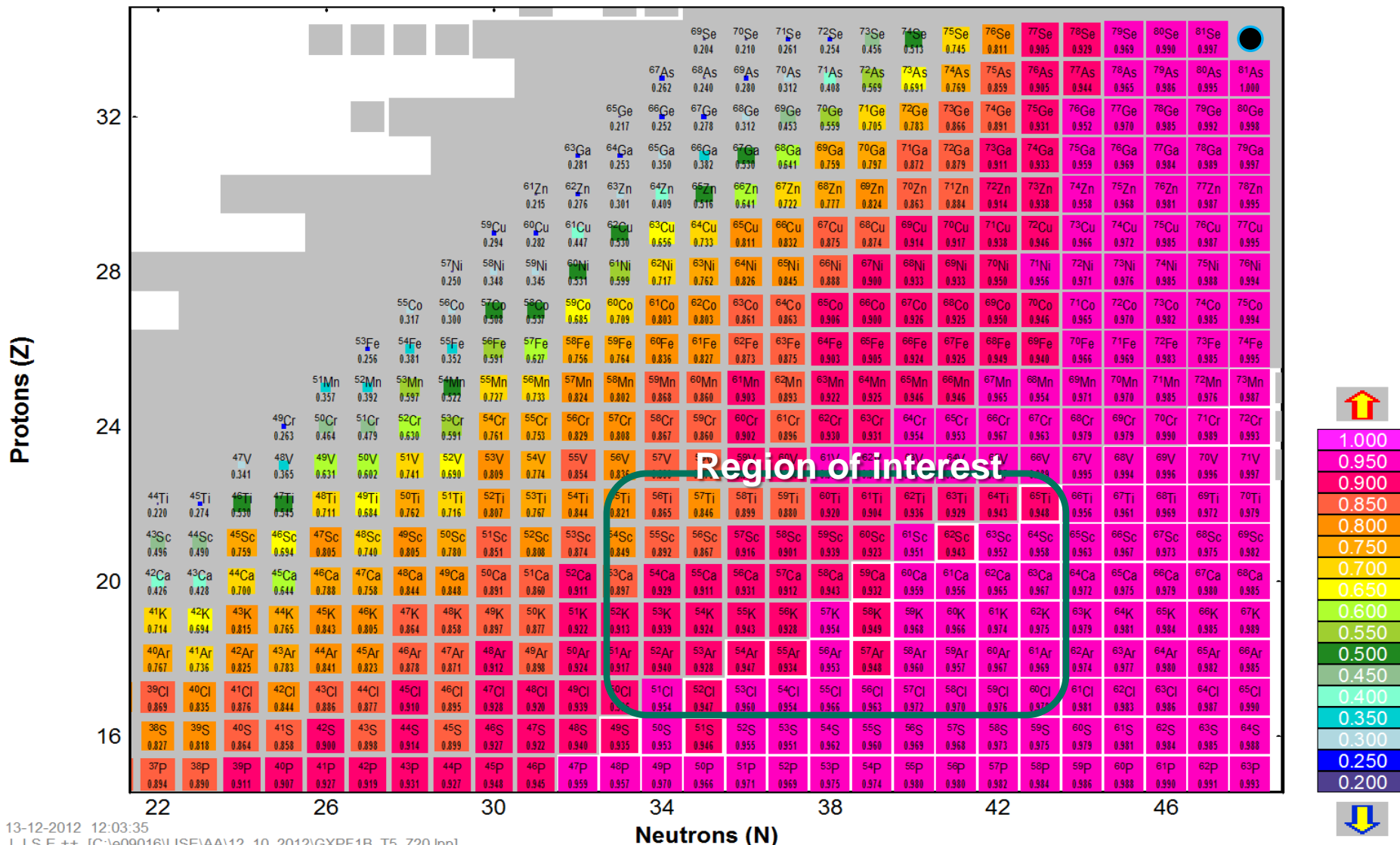


GXPF1B5 + LDM0, $E^* = 15.0$ & 9.15
 (deduced from ^{82}Se experimental data)

Current mode: 1n -> [(So total) evap] / [So total]

ABRASION-ABLATION - ⁸²Se + Be

Excit.Energy Method:< 2 >; <E*>:15.0*dA MeV Sigma:9.15; CoefThermalization=5.00e-22MeV.s DB₁="GXPF1B"
 NP=64; SE:"DB1+Ca10" Density:"auto" GeomCor:"On" Tunlg:"auto" FisBar=#1 BarFac=1.00 Modes=1010 1010 010

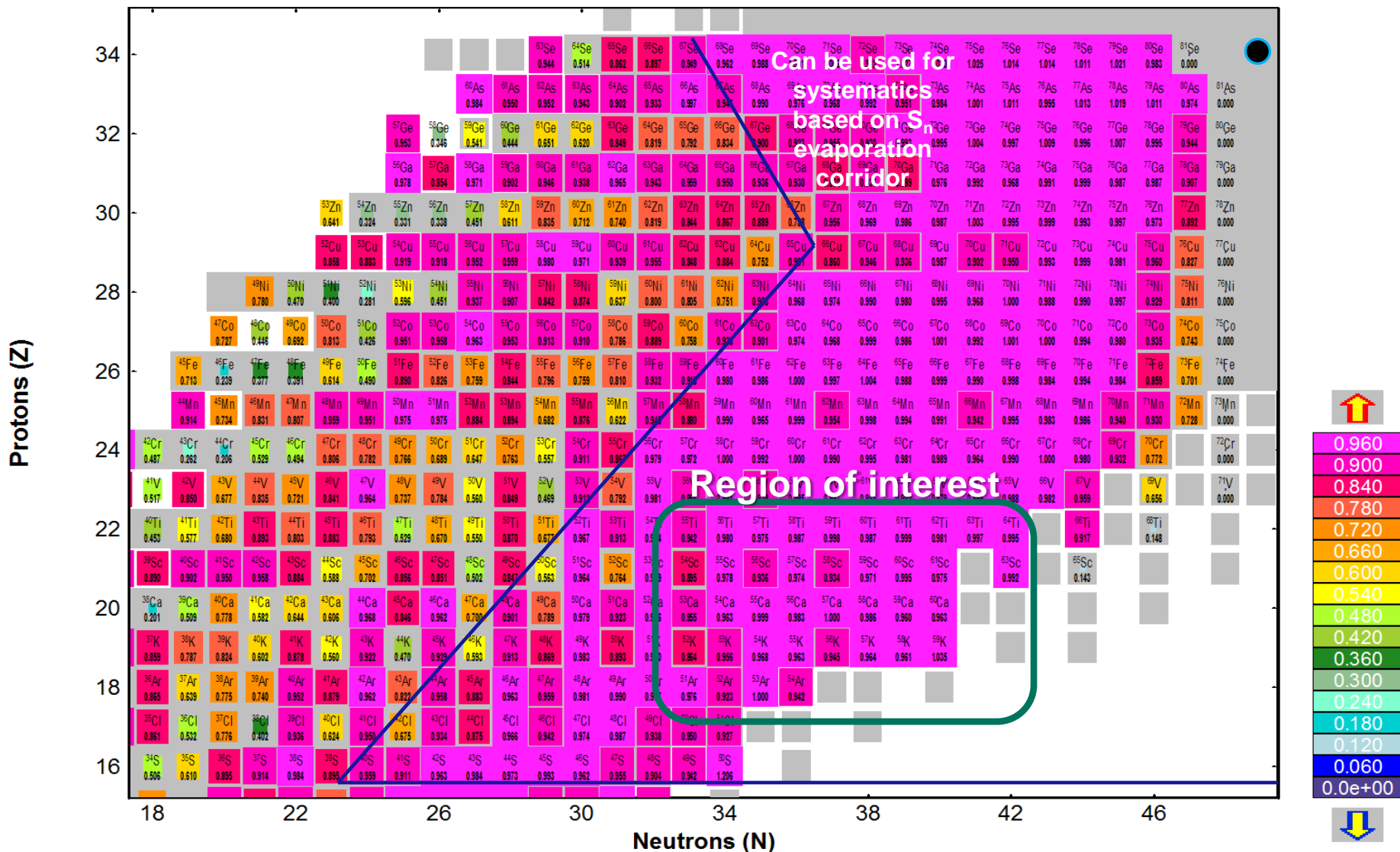


Current mode: 1n -> [S residue] / [Sr total]

ABRASION-ABLATION - ⁸²Se + Be

Excit.Energy Method:< 2 >; < E* >:15.0*dA MeV Sigma:9.15; Coef^Thermalization=5.00e-22MeV.s DB₁="GXPF1B"

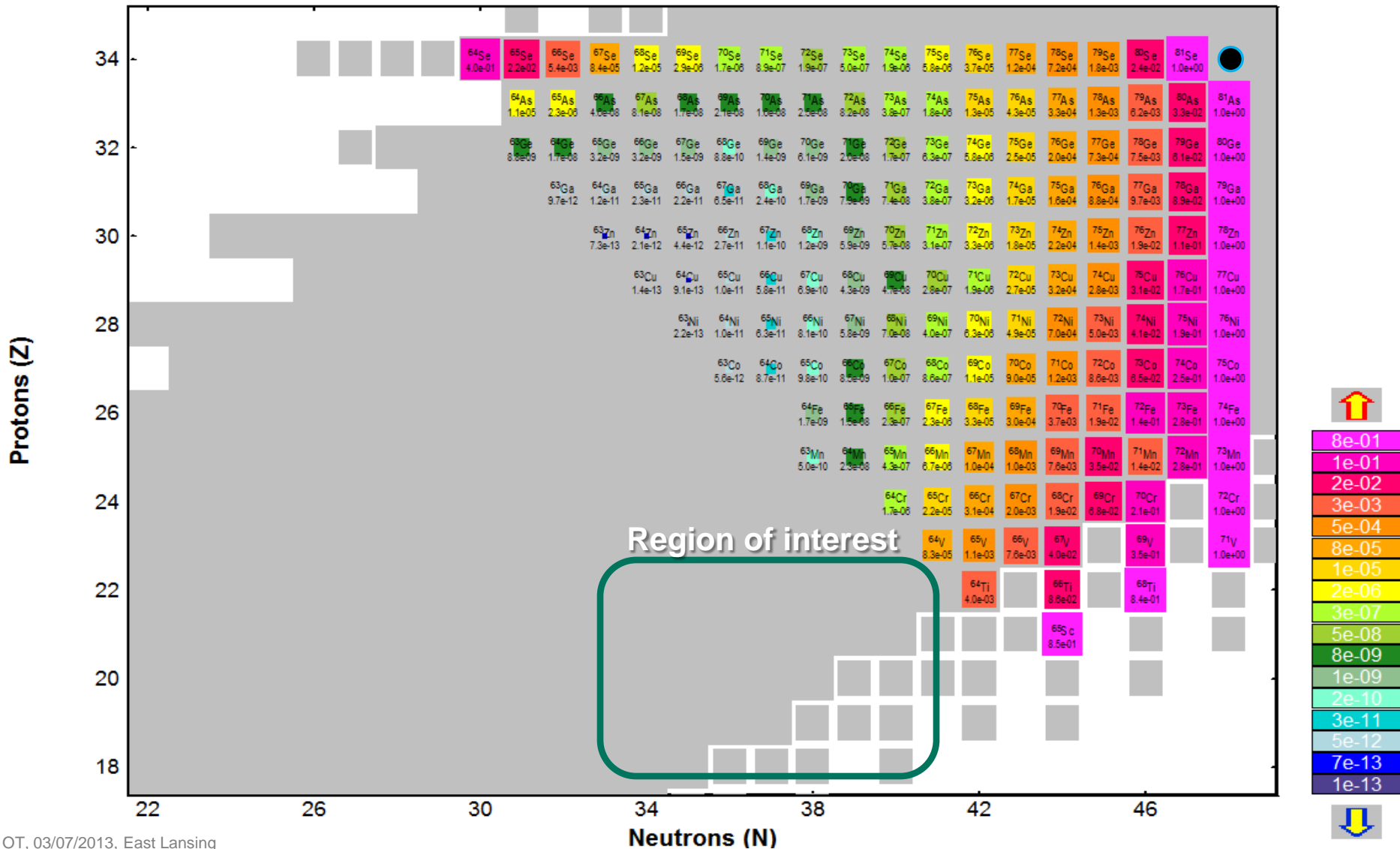
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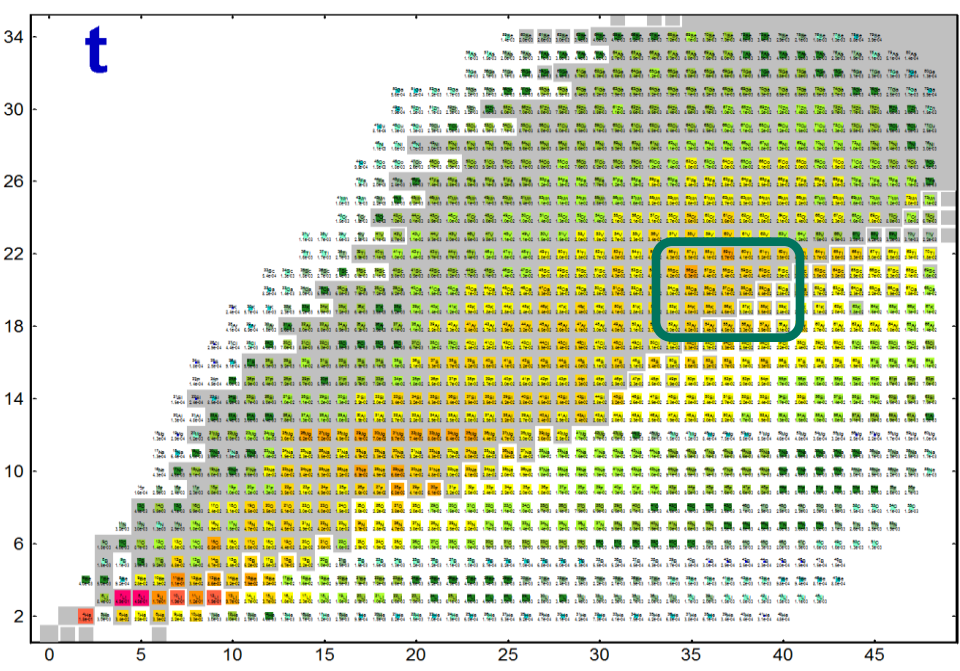
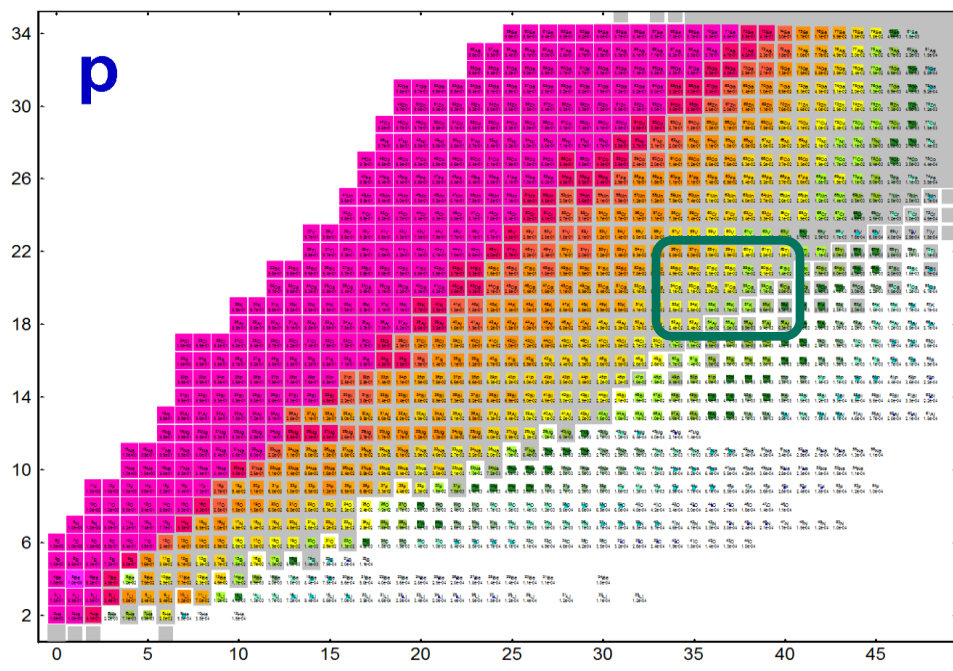
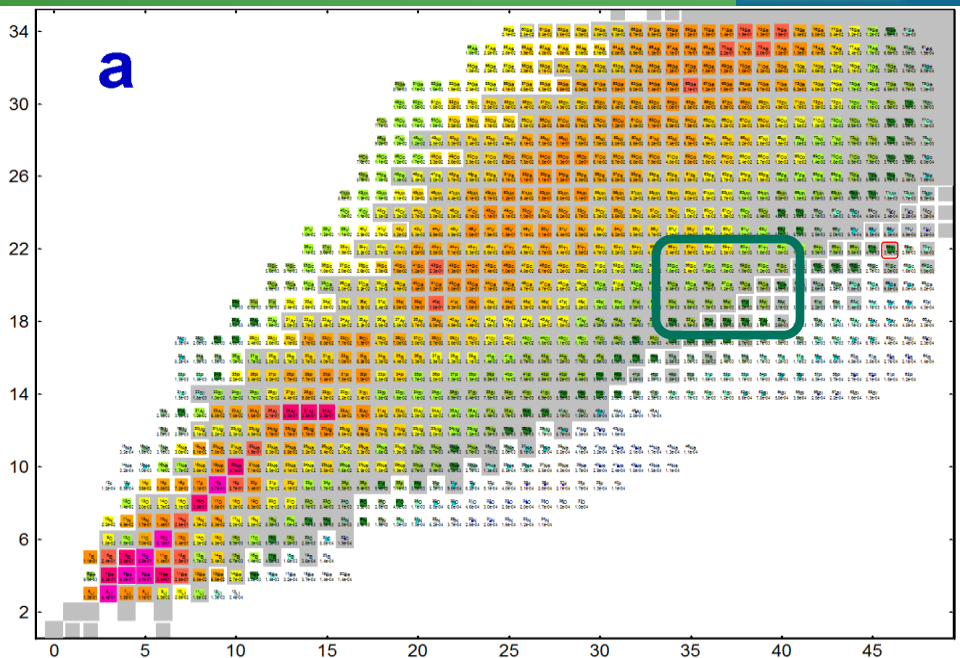
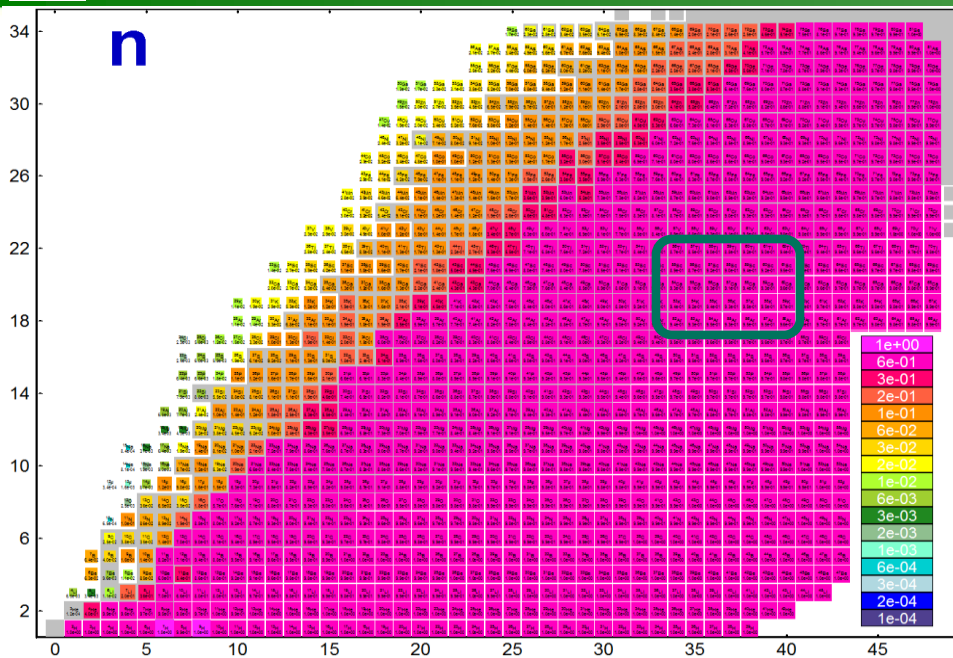


Current mode: Initial CS -> [S residue] / [Sr total]

ABRASION-ABLATION - $^{82}\text{Se} + \text{Be}$

Excit.Energy Method:< 2 >; < E* >:15.0*dA MeV Sigma:9.15; Coef^{Thermalization}=5.00e-22 MeV.s DB₁="GXPF1B"
 NP=64; SE:"DB1+Cal0" Density:"auto" GeomCor:"On" Tunlg:"auto" FisBar=#1 Bar^{Fac}=1.00 Modes=1010 1010 010



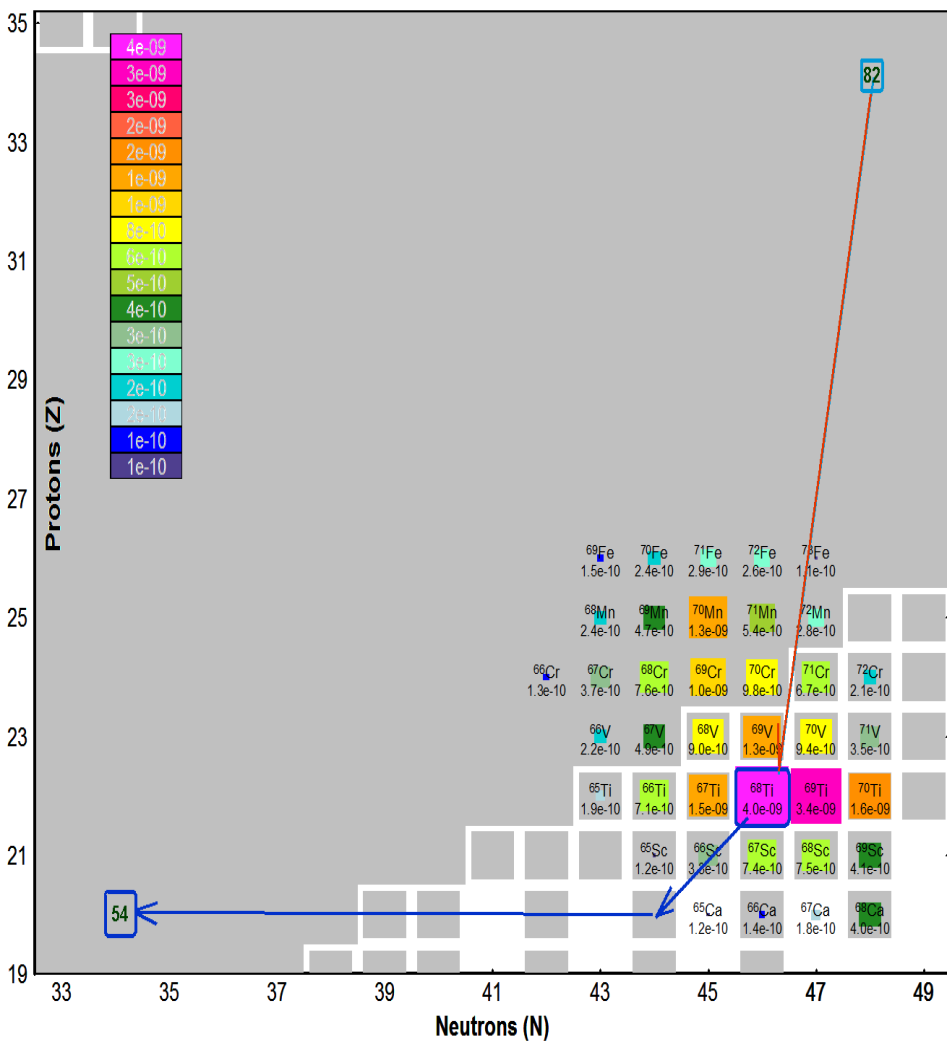


Initial Prefragments Plot for ⁵⁴Ca (2.78e-08 mb)

ABRASION-ABLATION - ⁸²Se + Be: more probable ⁶⁸Ti(4.02e-09 mb); <-dZ>=2.88 <-dN>=11.78

Excit.Energy Method:< 2 >; <E*>:15.0*dA MeV sigma:9.20; Thermal.Intr.Coef. = 5.00e-22 MeV*s

NP=64; SE:"DB1+CaI0" Density:"auto" Geom.Corr:"On" Tunlg:"auto" FisBar=#1 BarFac=1.00 Modes=1010 1010 010

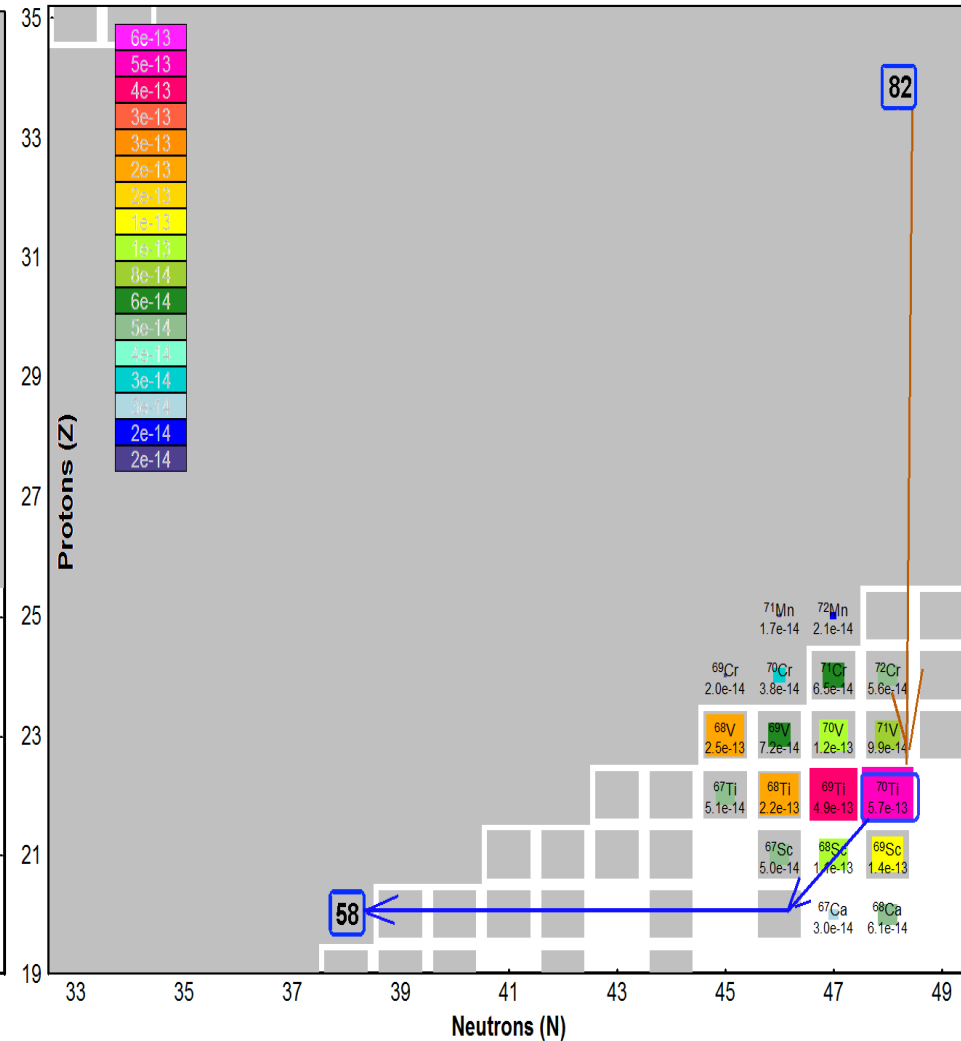


Initial Prefragments Plot for ⁵⁸Ca (2.57e-12 mb)

EVAPORATION - Compound nucleus ⁶⁸Ti: more probable ⁷⁰Ti(5.73e-13 mb); <-dZ>=2.27 <-dN>=8.89

Excit.Energy: 149.0-207.0 MeV; Fus.CS: 0.0 mb; Fus.Barrier: 10.82 fm; h_omega = 2.0 MeV

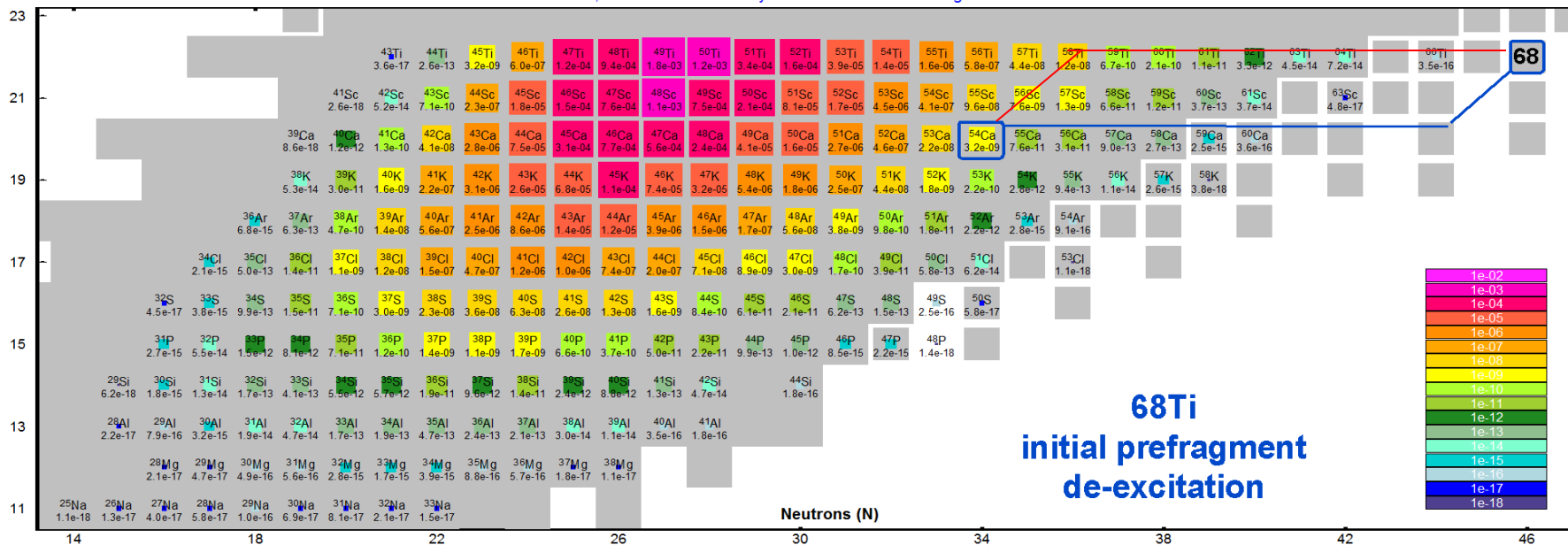
NP=64; SE:"DB1+CaI0" Density:"auto" GeomCor:"On" Tunlg:"auto" FisBar=#1 BarFac=1.00 Modes=1010 1010 010



More probable prefragments are Ti-isotopes (dZ=2)

Final Evaporation Residue cross-sections (LisFus)

EVAPORATION - Compound nucleus ^{68}Ti
 Excit.Energy: 149.0-207.0 MeV; Fus.CS: 0.0 mb; Fus.Barrier: 10.82 fm; $h_\omega = 2.0$ MeV
 NP=64; SE:"DB1+Cal0" Density:"auto" GeomCor:"On" Tunlg:"auto" FisBar=#1 BarFac=1.00 Modes=1010 1010 010



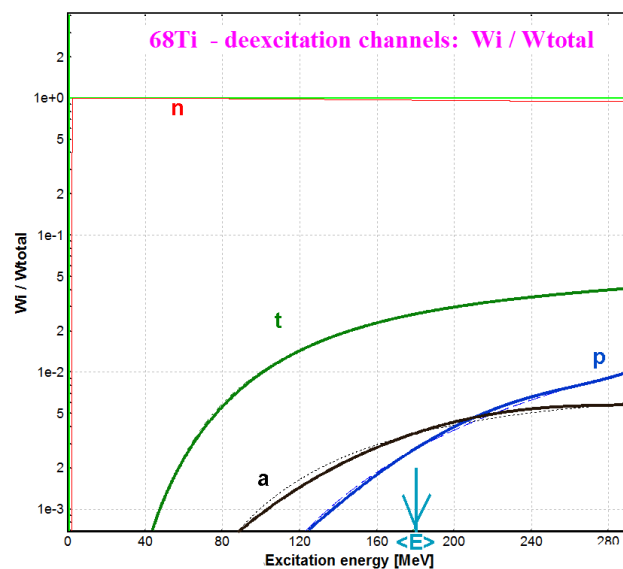
**^{68}Ti
initial prefragment
de-excitation**

Probability for $^{68}\text{Ti}^*(\text{Ex}=180\text{MeV})$

$t = 2.6\text{e-}2$
 $a = 3.6\text{e-}3$
 $p = 9.3\text{e-}3$

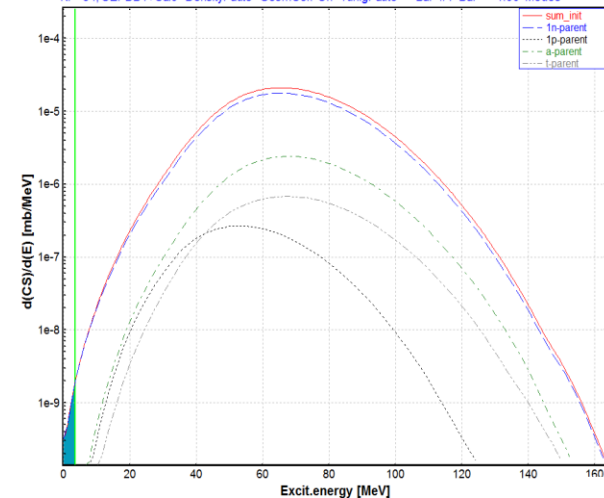
Probability ($dZ=2$)
 $t^{\wedge}2 = 6.8\text{e-}4$
 $a = 3.6\text{e-}3$
 $p^{\wedge}2 = 8.7\text{e-}5$

It is necessary to create the MC version to gate for ^{54}Ca residual in order to answer where ($^{68}\text{Ti} \rightarrow ^{54}\text{Ca}$) alpha de-excitation is more probable



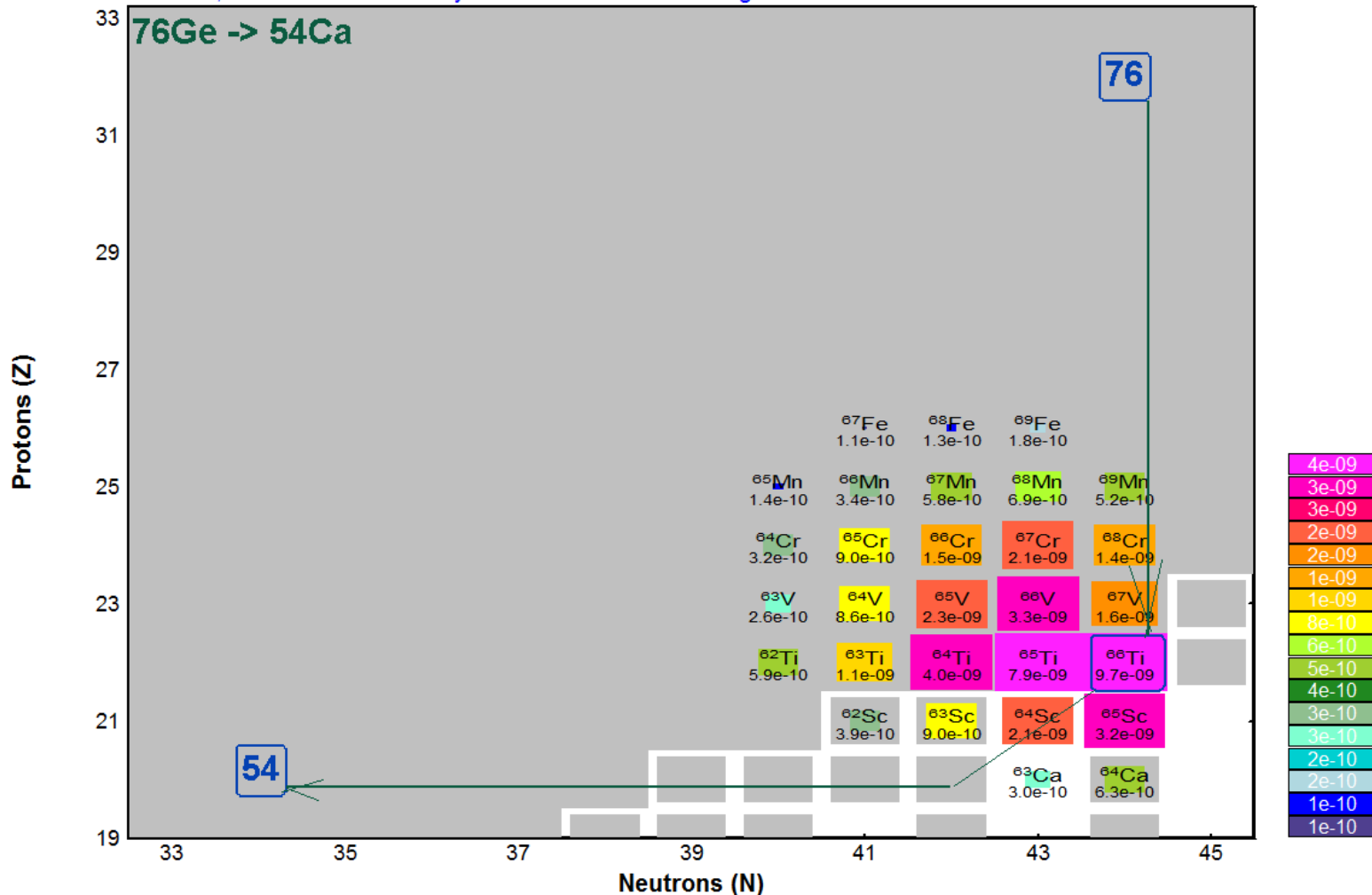
^{54}Ca excitation distributions: Input parent distr.

EVAPORATION - Compound nucleus ^{68}Ti
 Excit.Energy: 149.0-207.0 MeV; Fus.CS: 0.0 mb; Fus.Barrier: 10.82 fm; $h_\omega = 2.0$ MeV
 NP=64; SE:"DB1+Cal0" Density:"auto" GeomCor:"On" Tunlg:"auto" FisBar=#1 BarFac=1.00 Modes=1010 1010 010

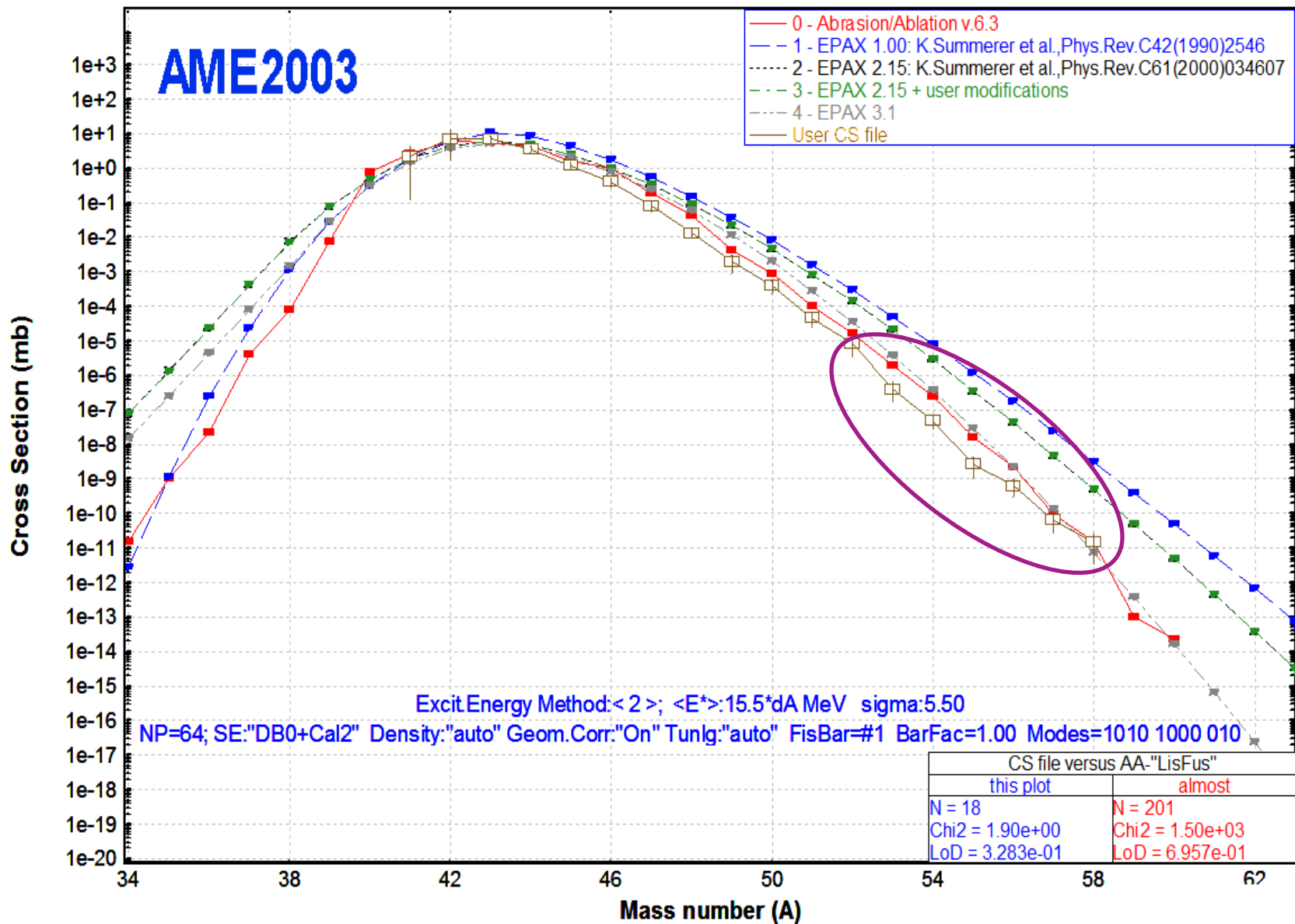


Initial Prefragments Plot for ⁵⁴Ca (4.85e-08 mb)

ABRASION-ABLATION - ⁷⁶Ge + Be: more probable ⁶⁶Ti(9.68e-09 mb); <-dZ>=2.46 <-dN>=8.90
 Excit.Energy Method:< 2 >; <E*>:15.0*dA MeV Sigma:8.60; Coef^Thermalization=5.00e-22MeV.s DB1="GXPF1B"
 NP=64; SE:"DB1+Cal0" Density:"auto" GeomCor:"On" Tunlg:"auto" FisBar=#1 Bar^{Fac}=1.00 Modes=1010 1010 010



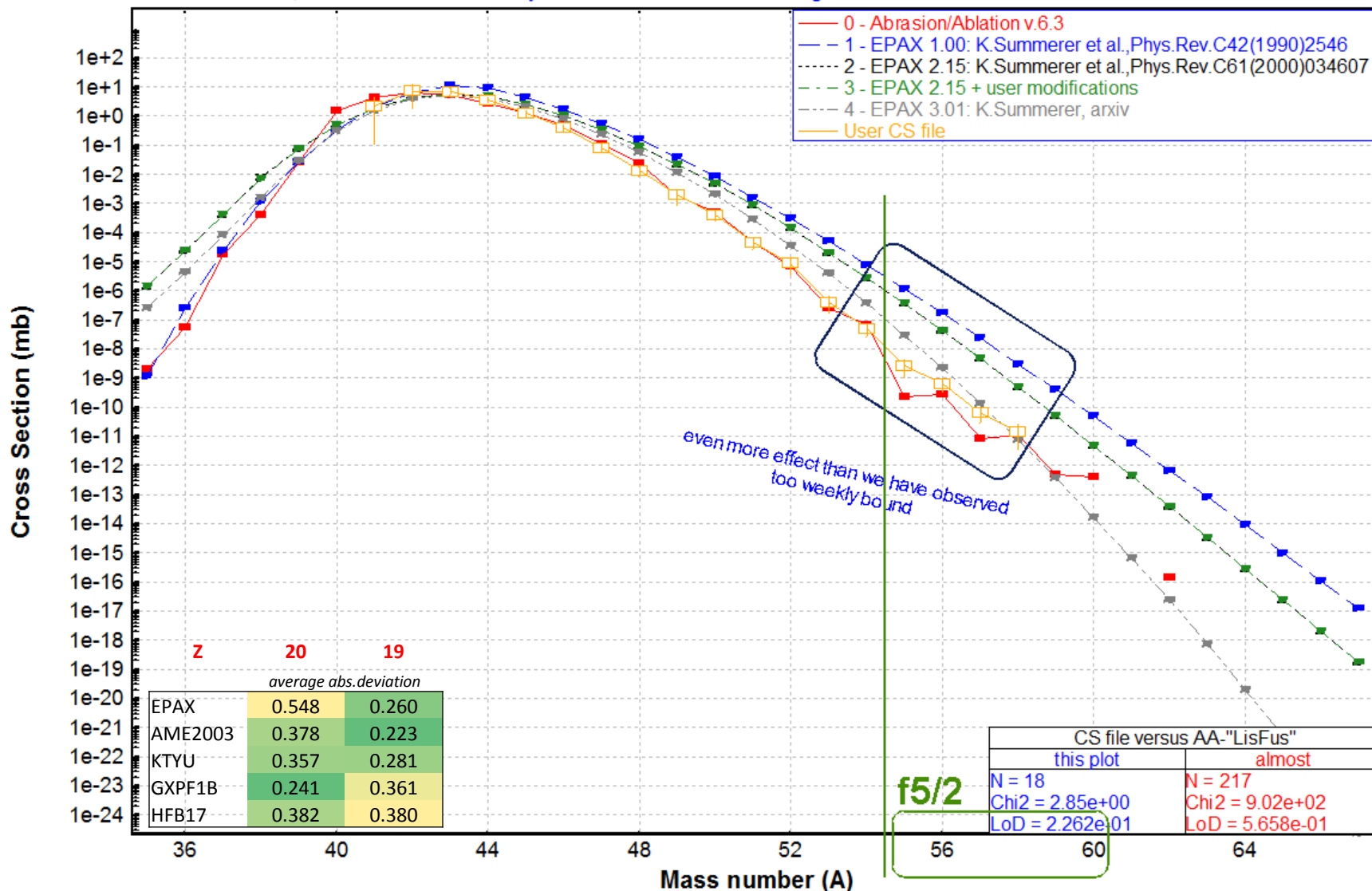
More probable prefragments are Ti-isotopes (dZ=2)



Cross sections (Projectile Fragmentation)

$^{82}\text{Se} + \text{Be} \rightarrow Z=20$

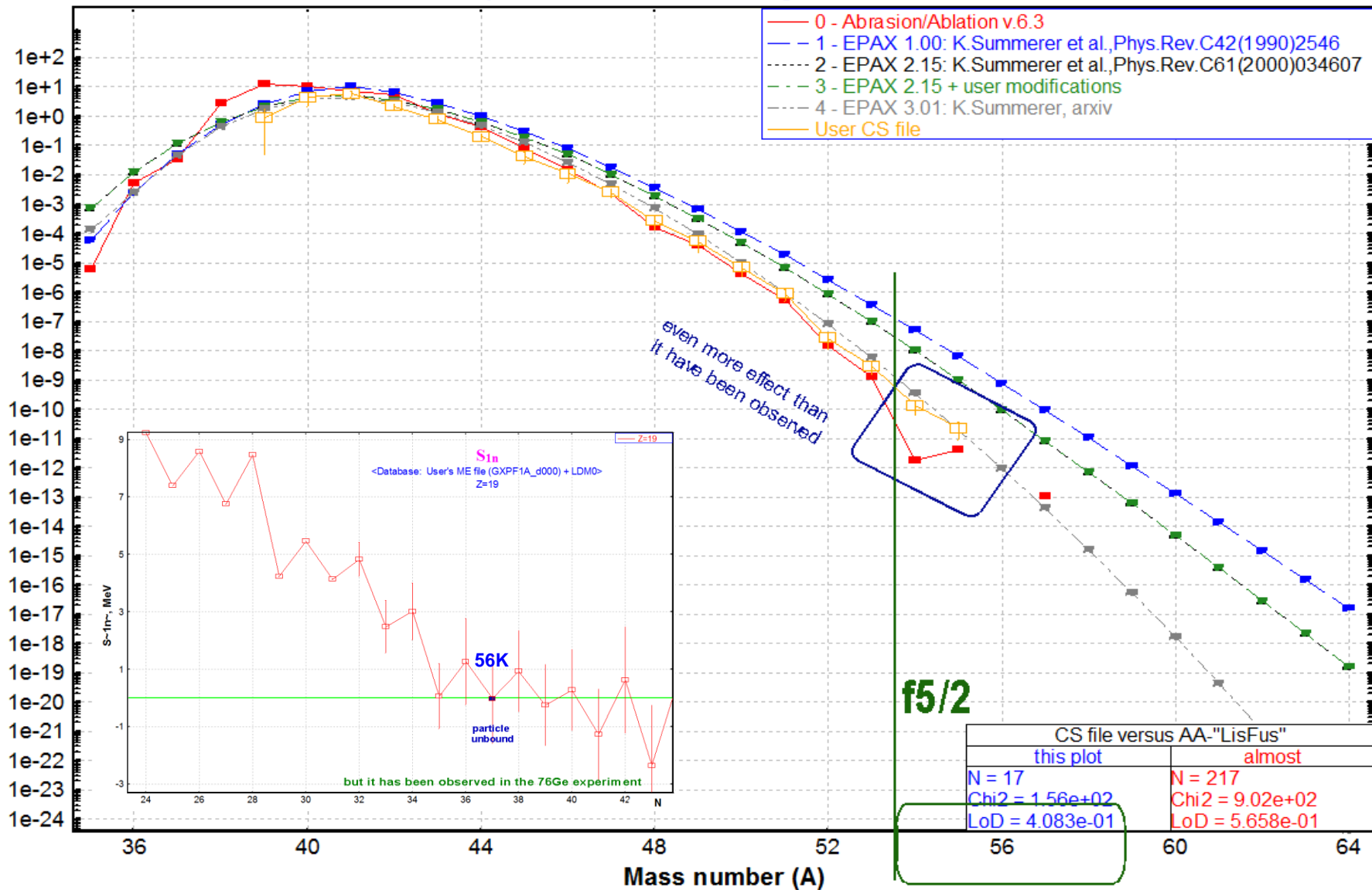
Excit. Energy Method: < 2 >; < E* >: 15.0 * dA MeV Sigma: 9.20; Coef^{Thermalization}=5.00e-22 MeV.s DB₁="GXPF1A_d000"
 NP=64; SE:"DB1+Cal0" Density:"auto" GeomCor:"On" Tunlg:"auto" FisBar=#1 Bar^{Fac}=1.00 Modes=1010 1010 010



Cross sections (Projectile Fragmentation)

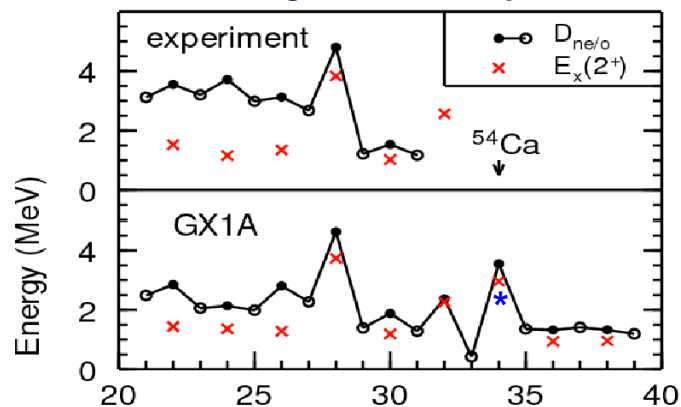
$^{82}\text{Se} + \text{Be} \rightarrow \text{Z}=19$

Excit.Energy Method:< 2 >; <E*>:15.0*dA MeV Sigma:9.20; Coef^{fThermalization}=5.00e-22^{MeV.s} DB1="GXPF1A_d000"
 NP=64; SE:"DB1+Cal0" Density:"auto" GeomCor:"On" TunJg:"auto" FisBar=#1 Bar^{Fac}=1.00 Modes=1010 1010 010

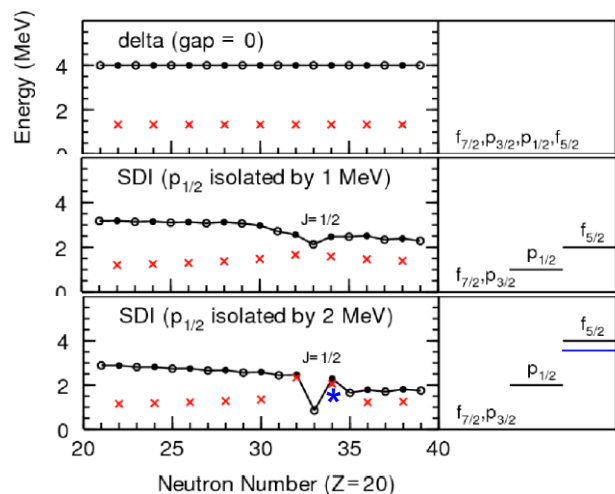


Alex Brown, ENSFN, October 11, 2012

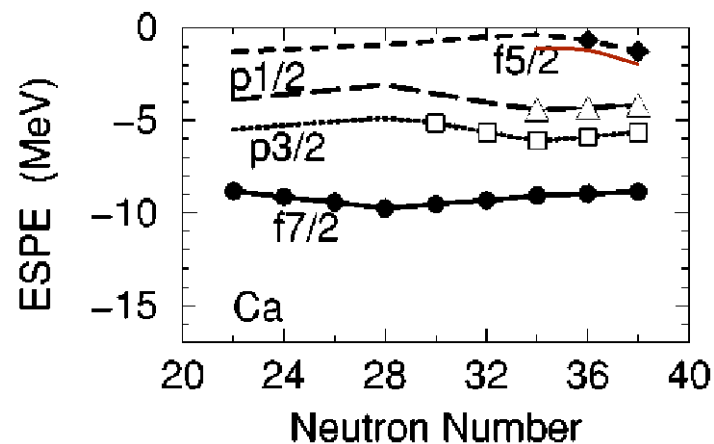
Pairing and Shell Gaps



$$D_n(N) = (-1)^N [S_n(Z, N+1) - S_n(Z, N)]$$



RIKEN Experiment: $E_x(2^+) \rightarrow 0.5$ MeV below predictions

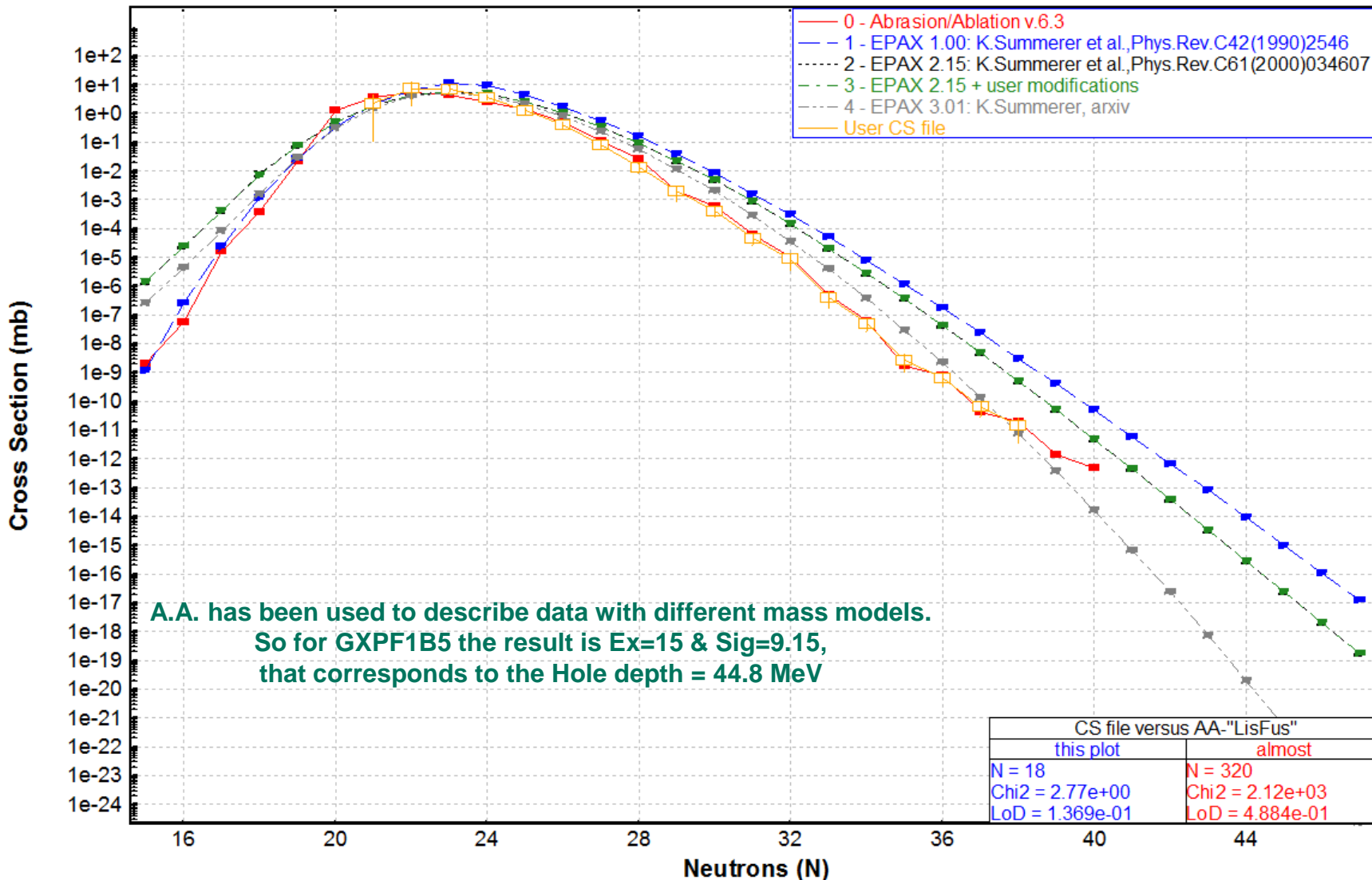


$^{55-60}\text{Ca}$ are more particle bound than GXPF1B predicted

Cross sections (Projectile Fragmentation)

$^{82}\text{Se} + \text{Be} \rightarrow Z=20$

Excit.Energy Method:< 2 >; < E* >:15.0*dA MeV Sigma:9.15; Coef^{Thermalization}=5.00e-22 MeV.s DB₁="GXPF1B"
 NP=64; SE:"DB1+Cal0" Density:"auto" GeomCor:"On" Tunlg:"auto" FisBar=#1 Bar^{Fac}=1.00 Modes=1010 1010 010



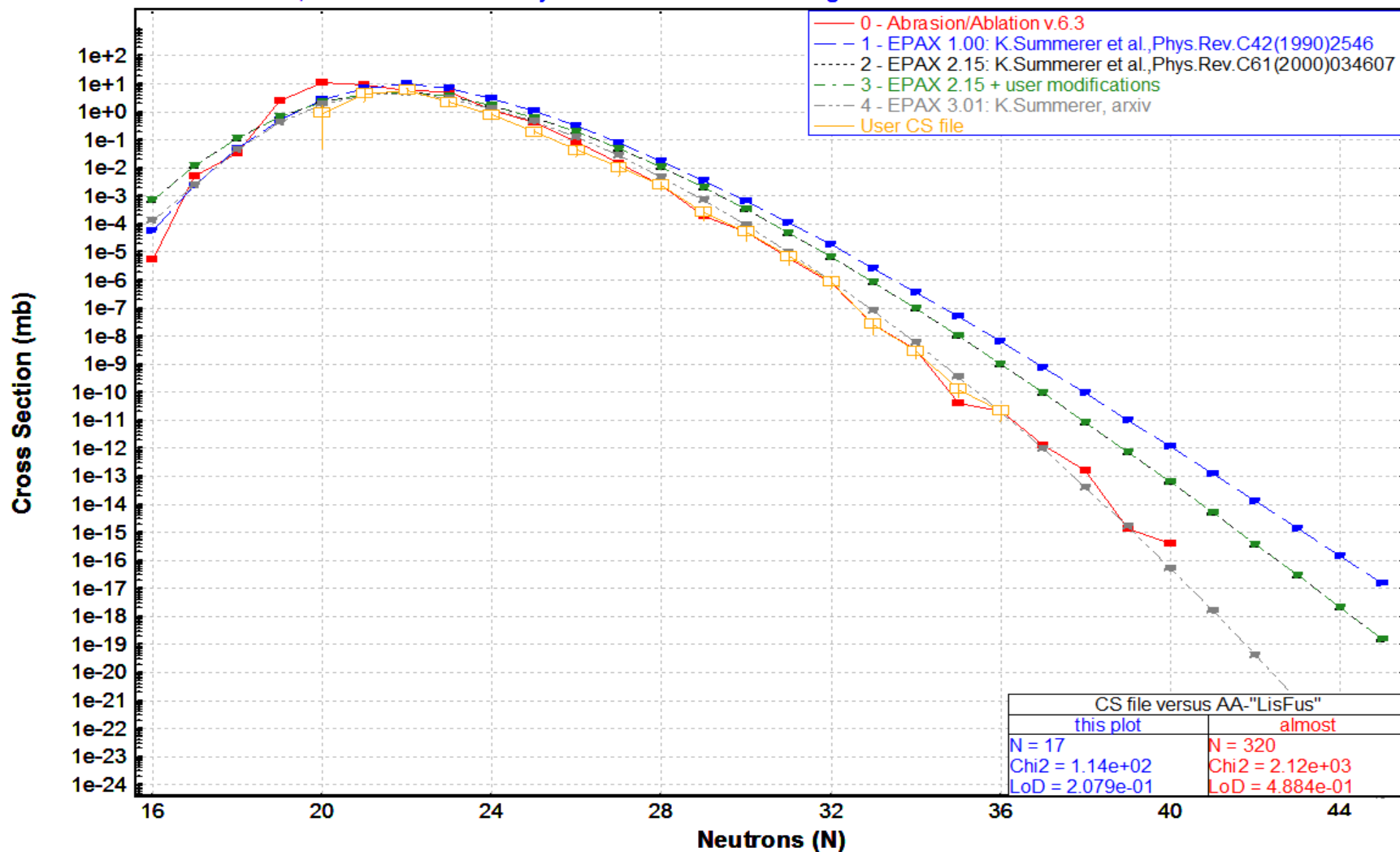
A.A. has been used to describe data with different mass models.
 So for GXPF1B5 the result is Ex=15 & Sig=9.15,
 that corresponds to the Hole depth = 44.8 MeV

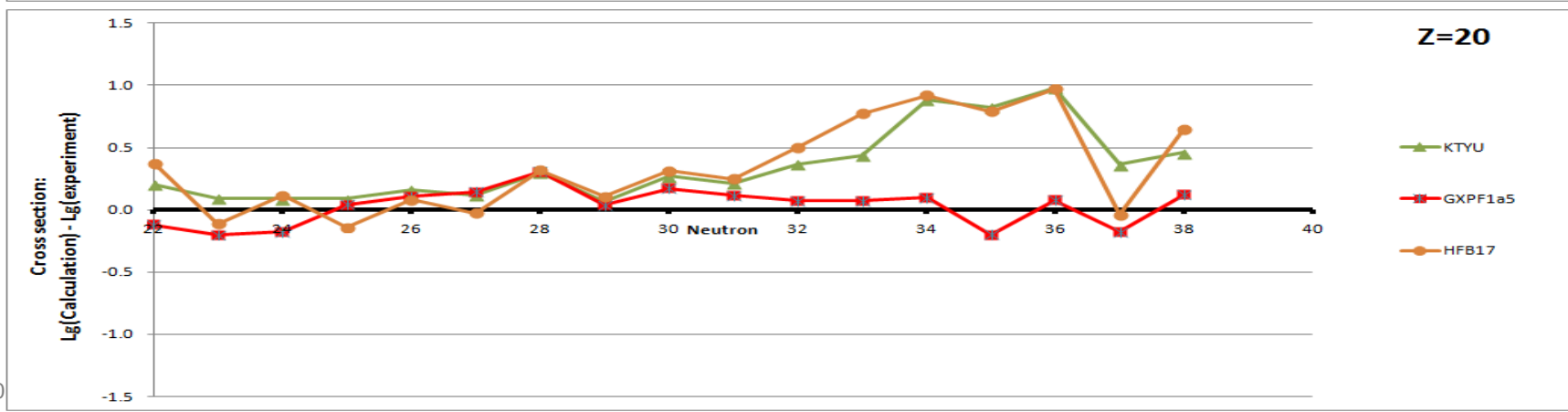
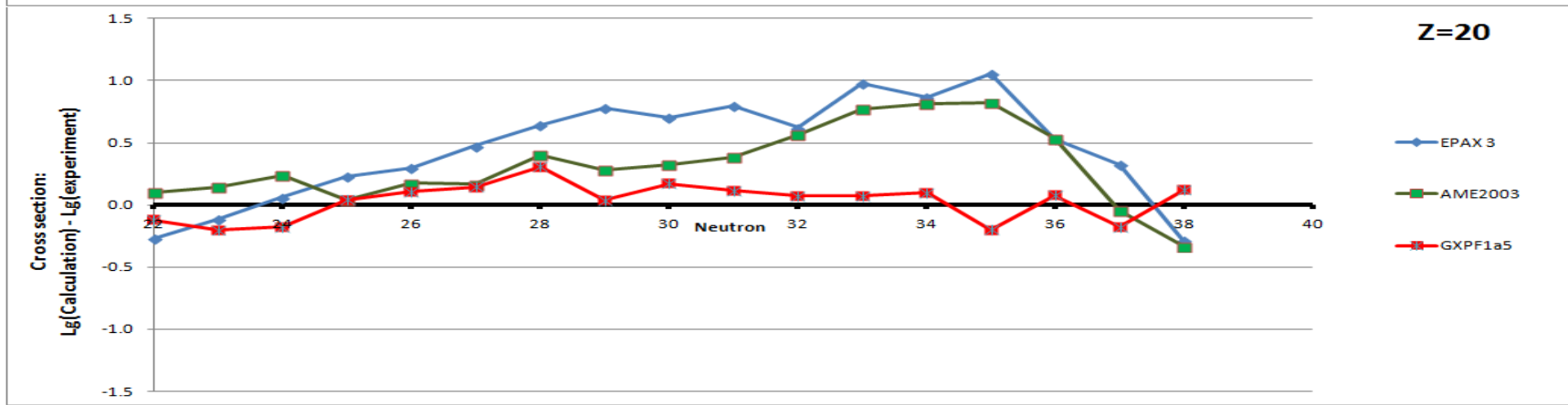
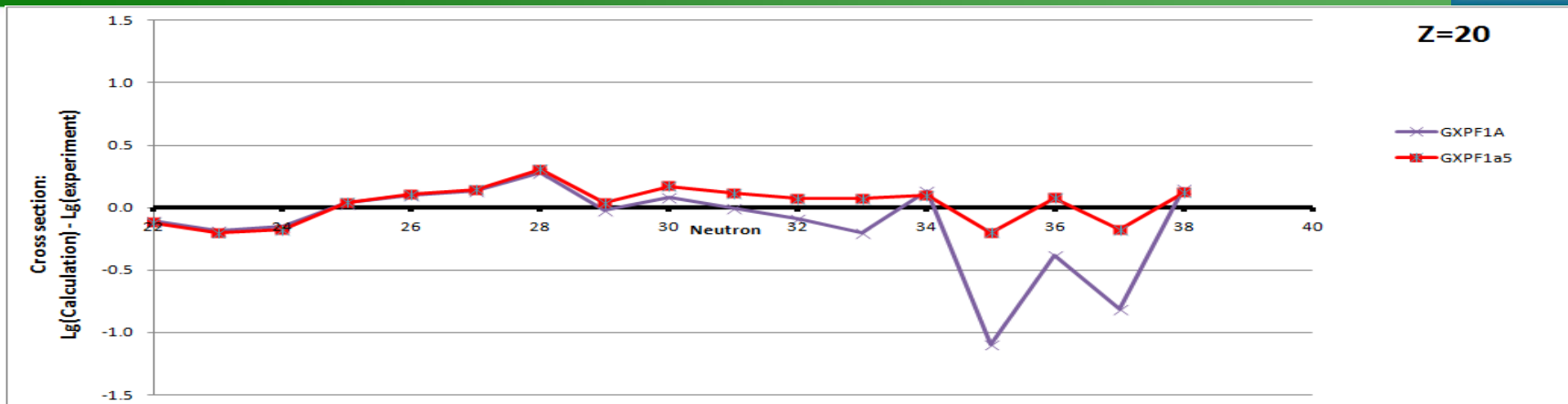
CS file versus AA-"LisFus"	
this plot	almost
N = 18	N = 320
Chi2 = 2.77e+00	Chi2 = 2.12e+03
LoD = 1.369e-01	LoD = 4.884e-01

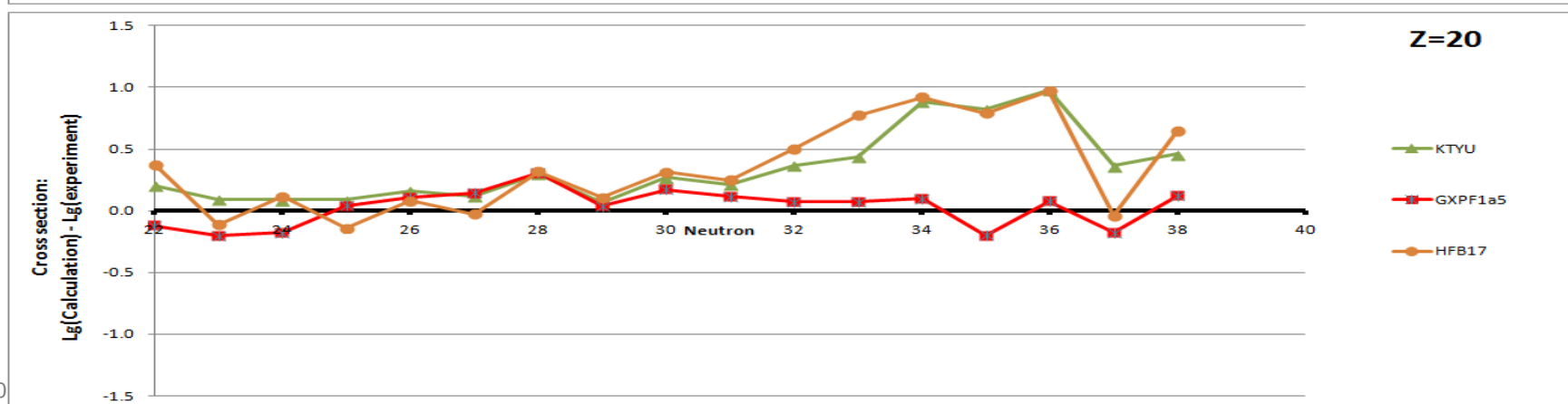
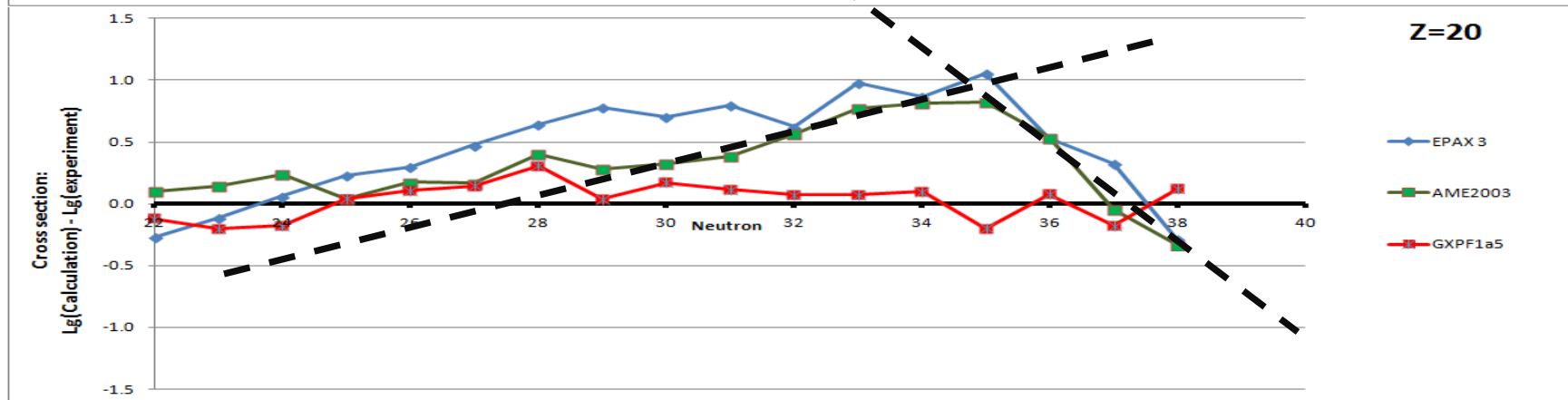
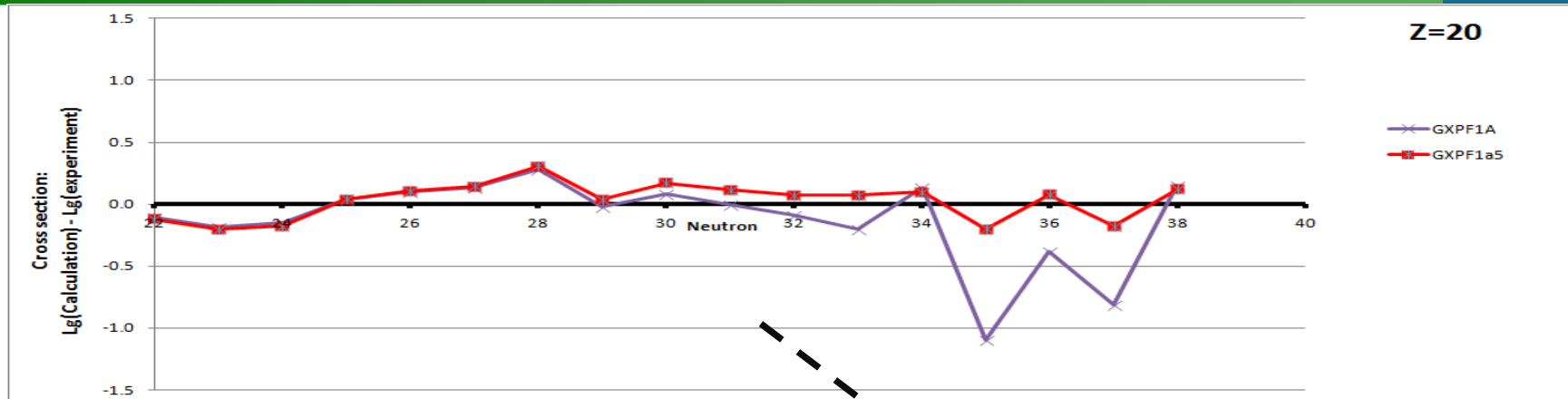
Cross sections (Projectile Fragmentation)

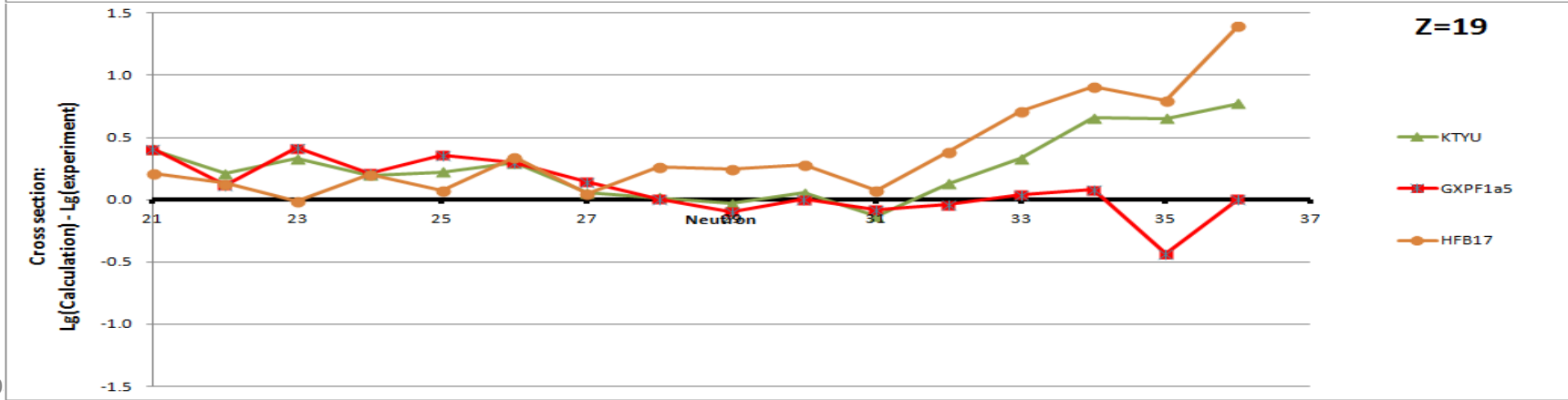
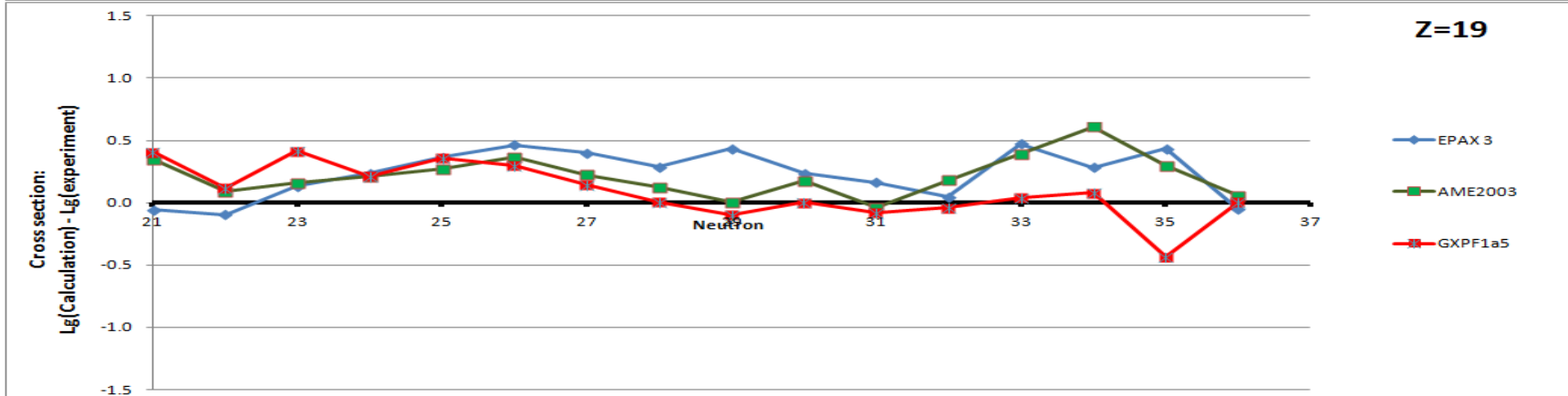
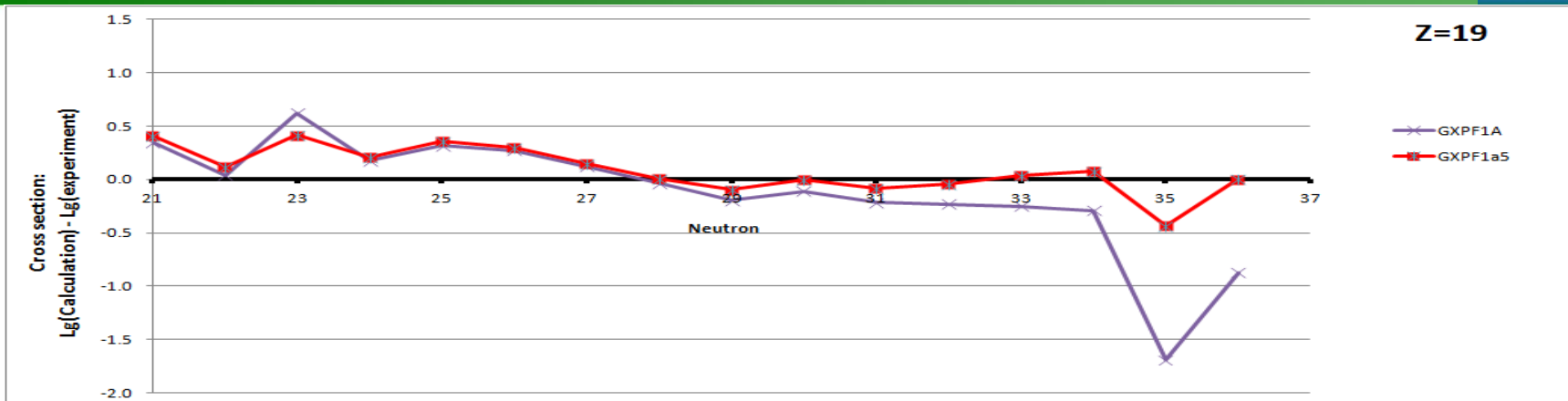
$^{82}\text{Se} + \text{Be} \rightarrow \text{Z}=19$

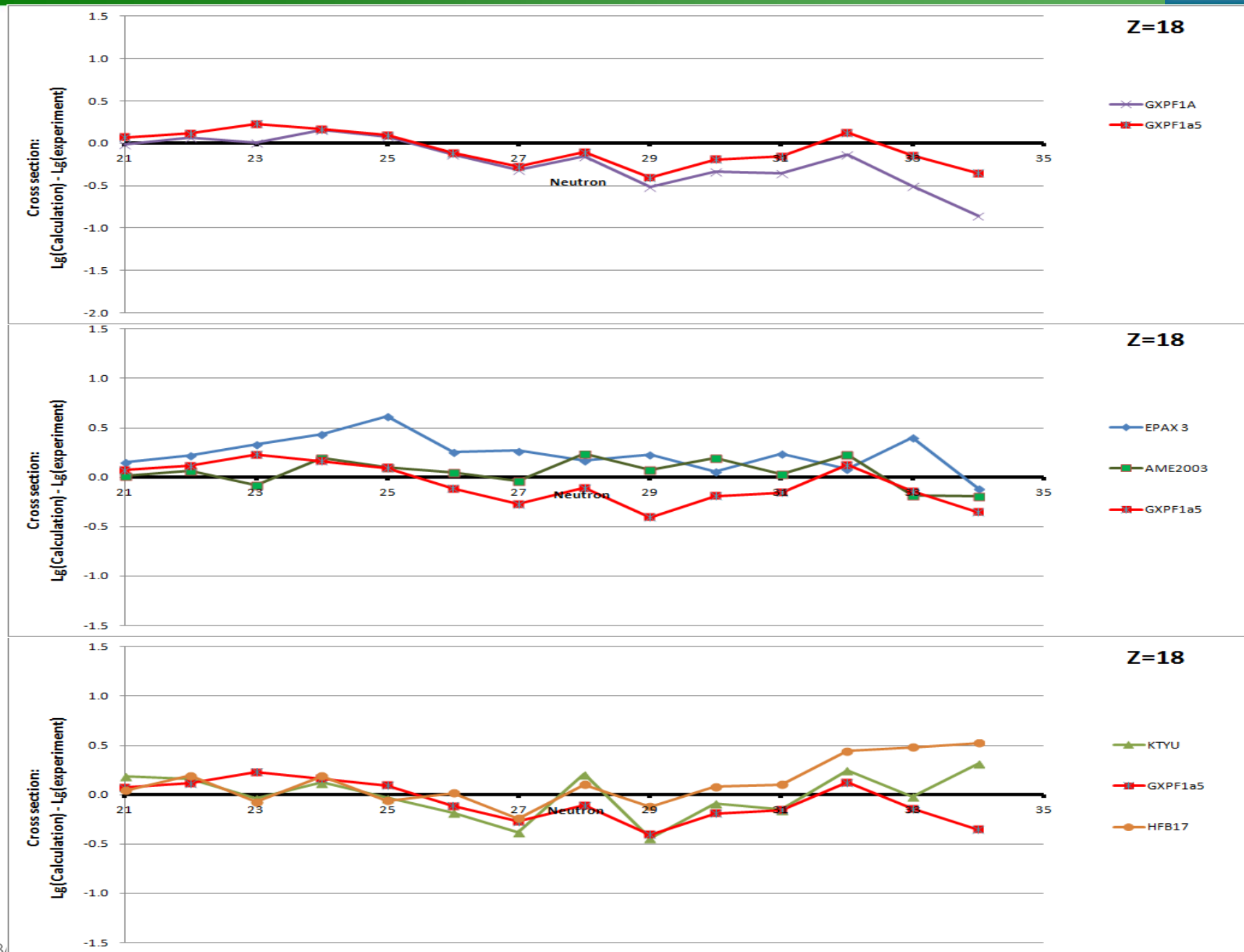
Excit.Energy Method:< 2 >; <E*>:15.0*dA MeV Sigma:9.15; Coef^{Thermalization}=5.00e-22MeV.s DB₁="GXPF1B"
 NP=64; SE:"DB1+Cal0" Density:"auto" GeomCor:"On" Tunlg:"auto" FisBar=#1 Bar^{Fac}=1.00 Modes=1010 1010 010

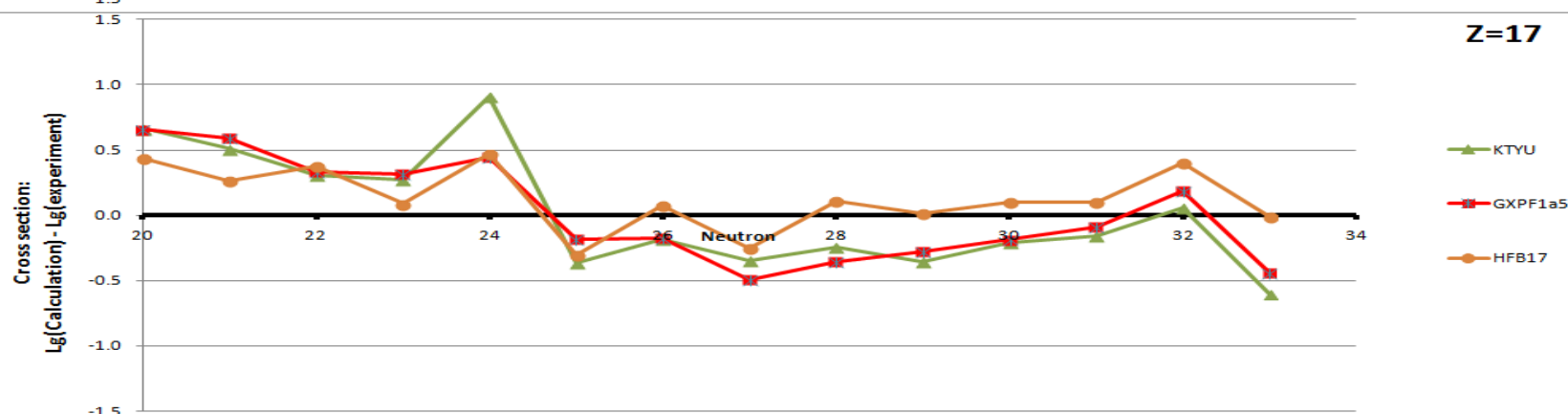
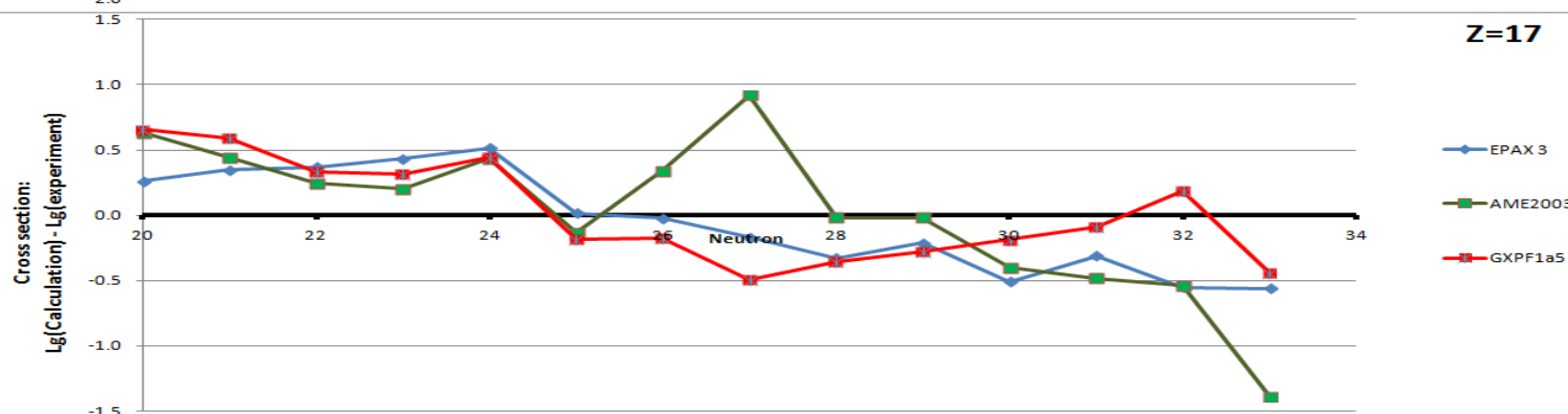
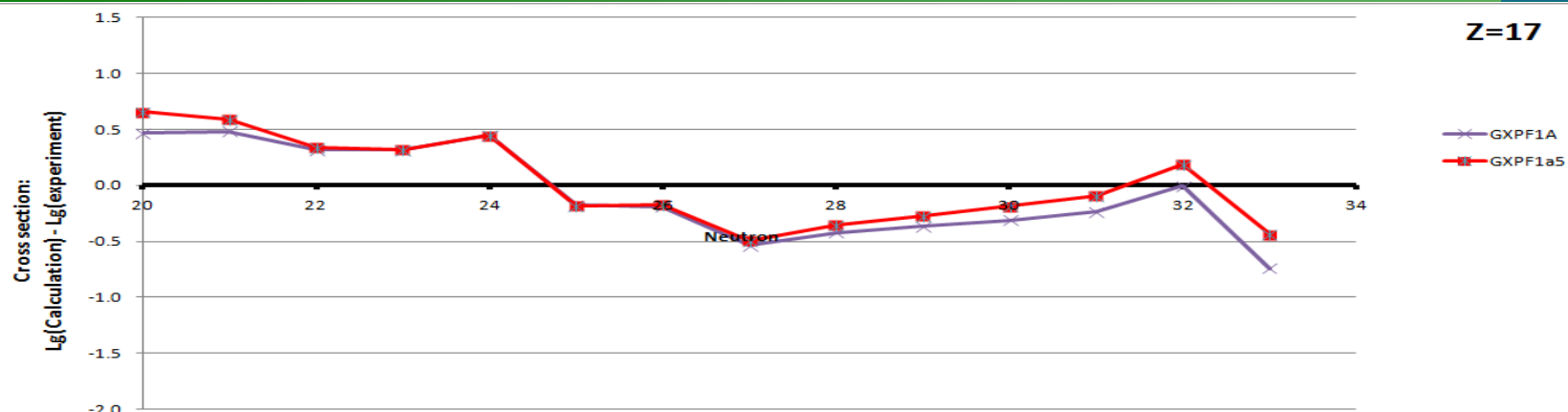


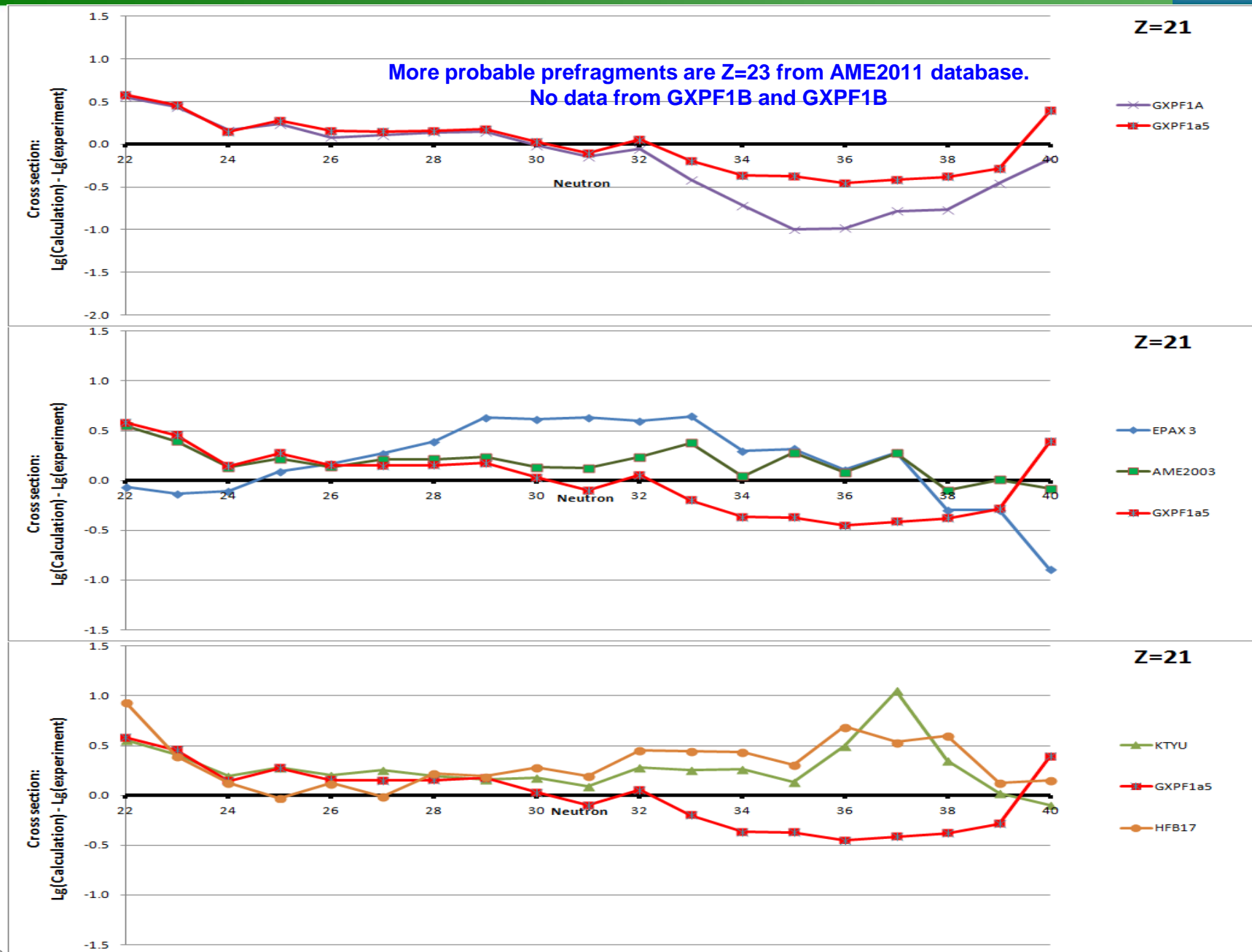












Cross sections (Projectile Fragmentation)

$^{76}\text{Ge} + \text{Be} \rightarrow Z=20$

Excit.Energy Method:< 2 >; <E*>:15.0*dA MeV Sigma:8.60; Coef^{Thermalization}=5.00e-22^{MeV.s} DB₁="GXPF1B"
 NP=64; SE:"DB1+Cal0" Density:"auto" GeomCor:"On" Tunlg:"auto" FisBar=#1 Bar^{Fac}=1.00 Modes=1010 1010 010

