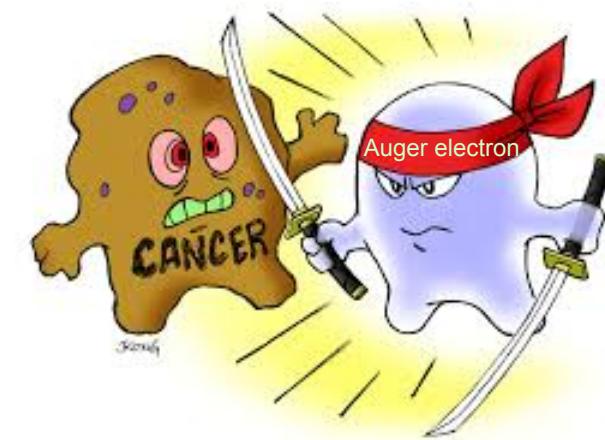




Auger and conversion electron spectroscopy of medical radioisotope ^{125}I

A magic bullet for cancer therapy

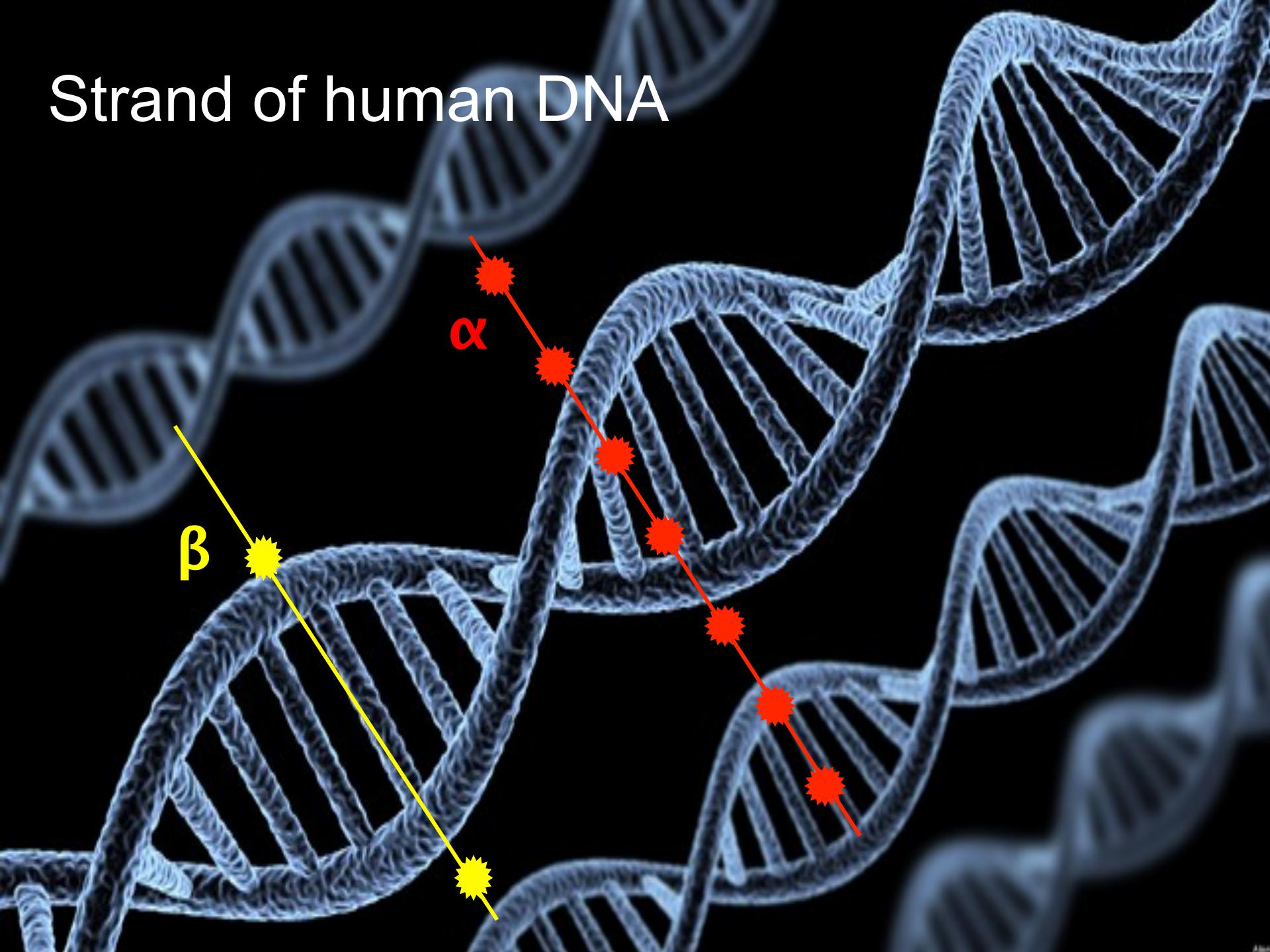


Presented by: **Bryan Tee Pi-Ern**

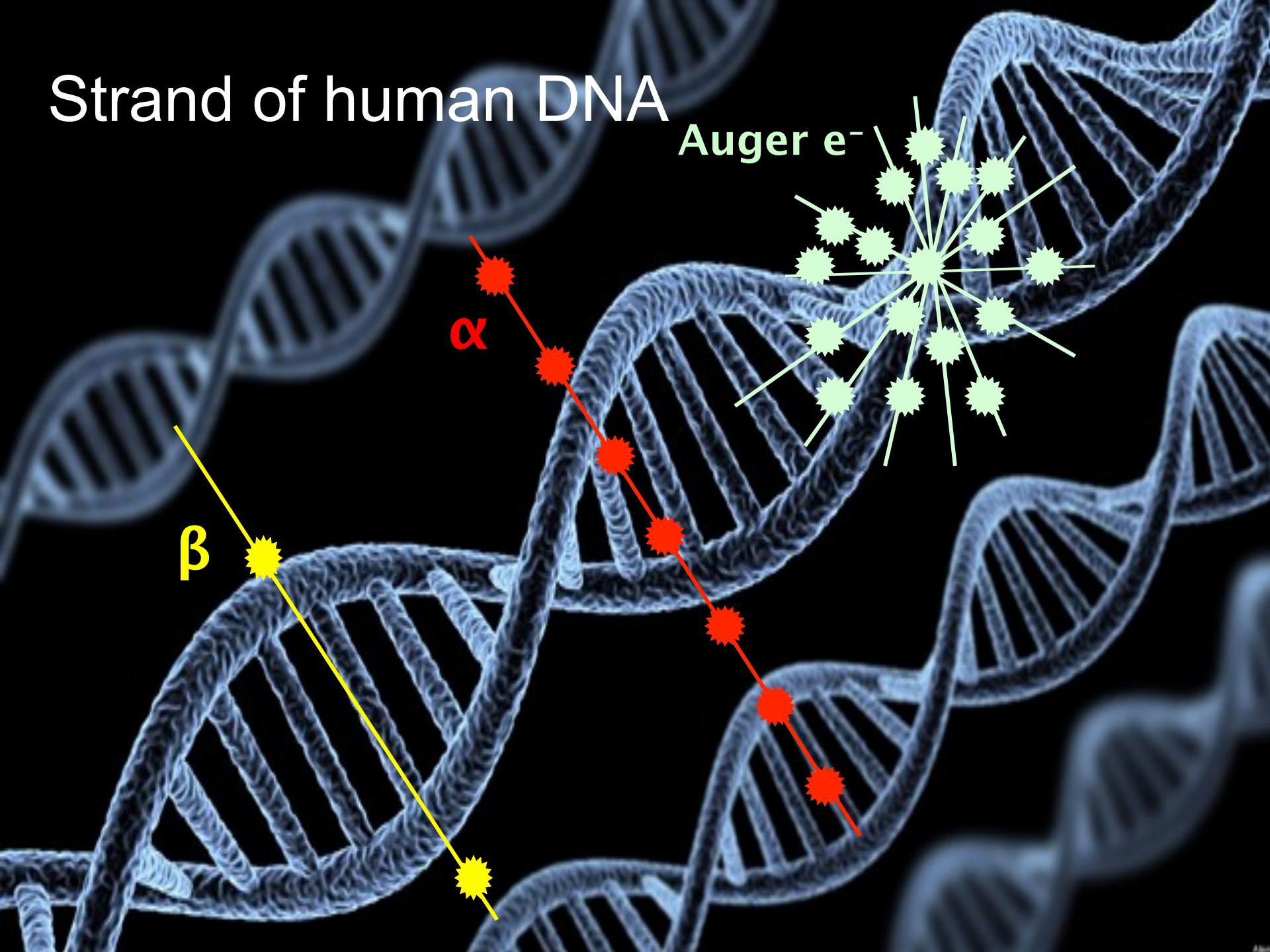
Date: 19th October 2018

Supervised by **Dr Tibor Kibédi, A/Prof Maarten Vos and
Professor Andrew Stuchberry**

Strand of human DNA

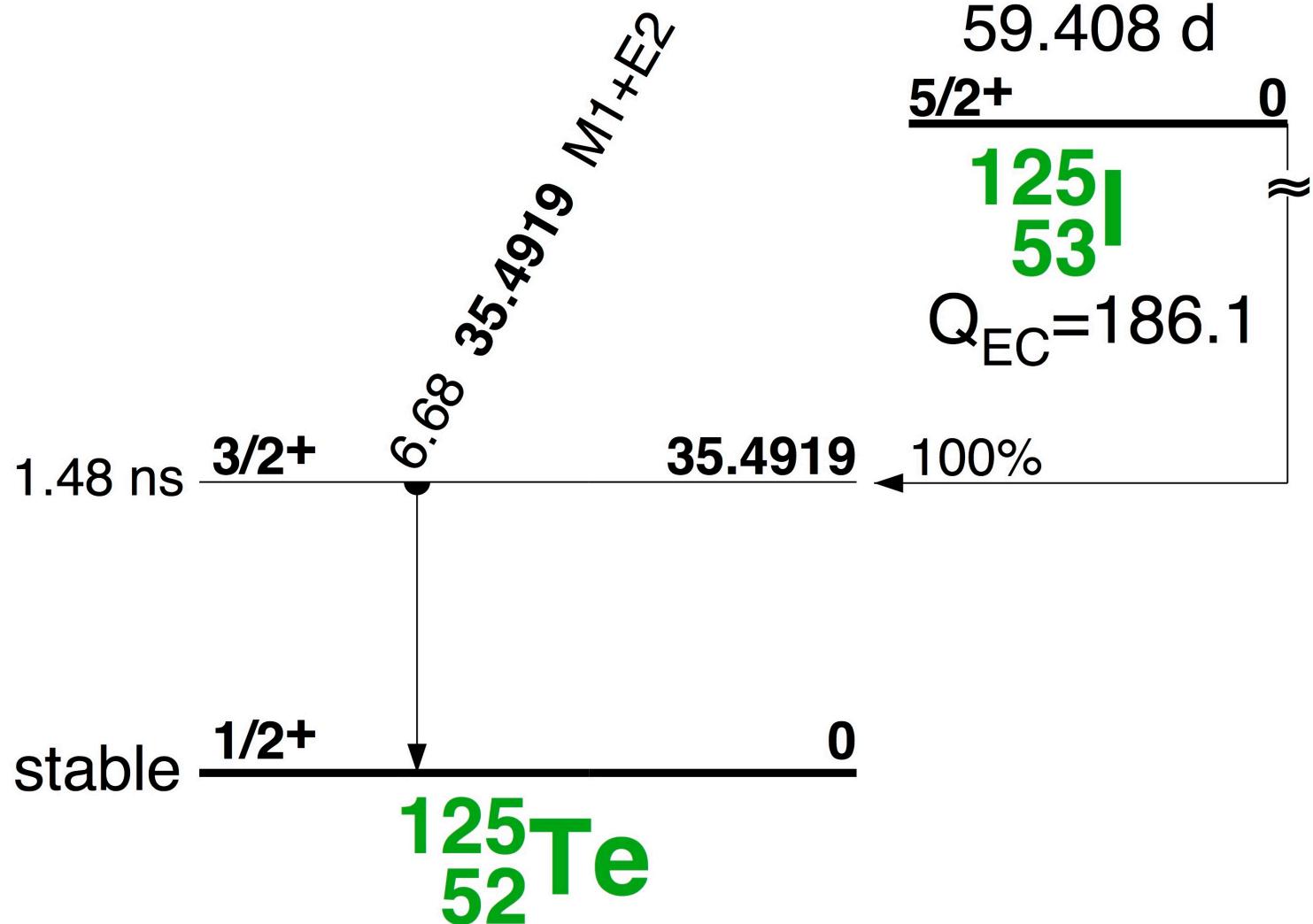


Strand of human DNA





Decay scheme of ^{125}I

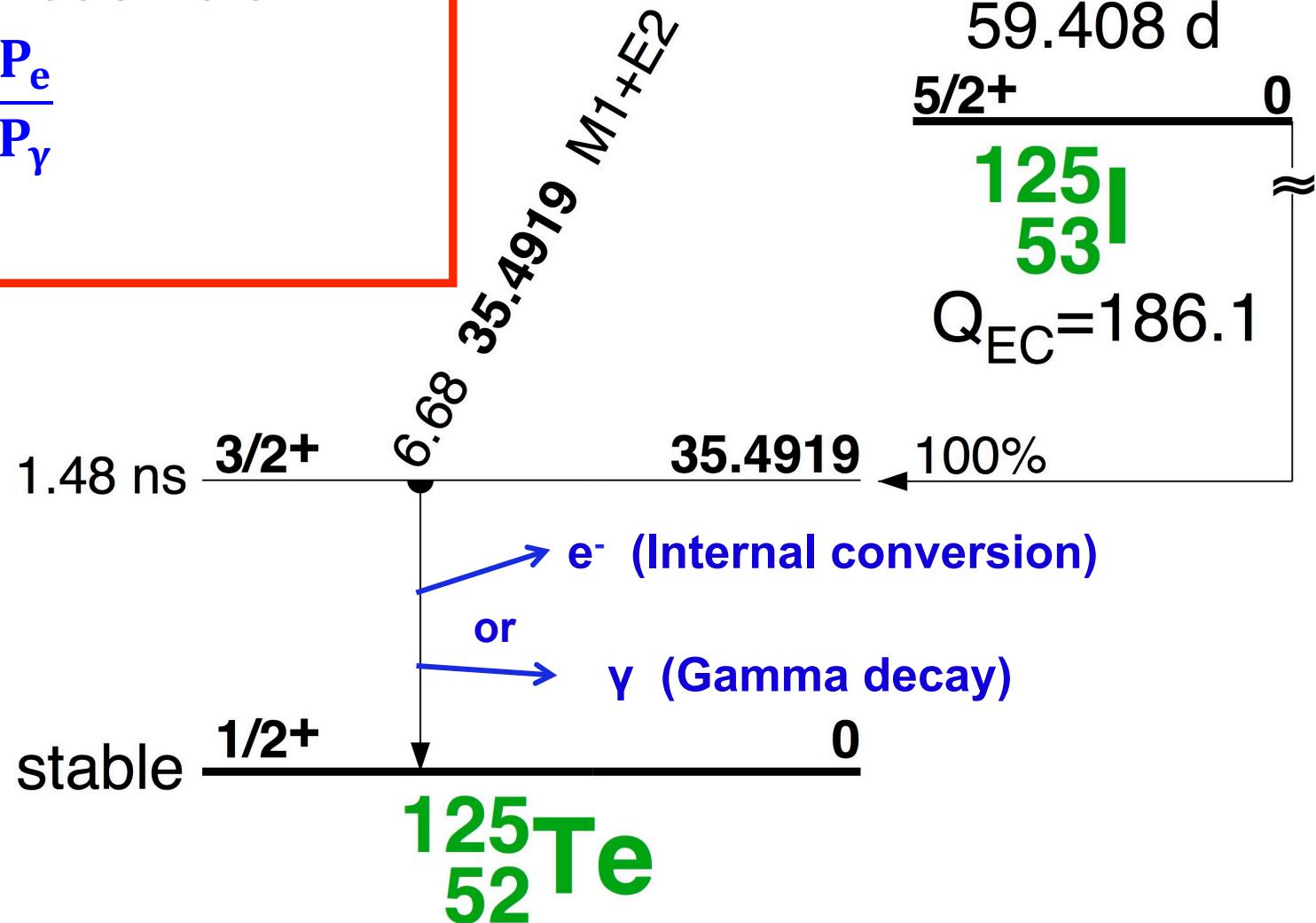




Decay scheme of ^{125}I

Conversion coefficient

$$\alpha = \frac{P_e}{P_\gamma}$$

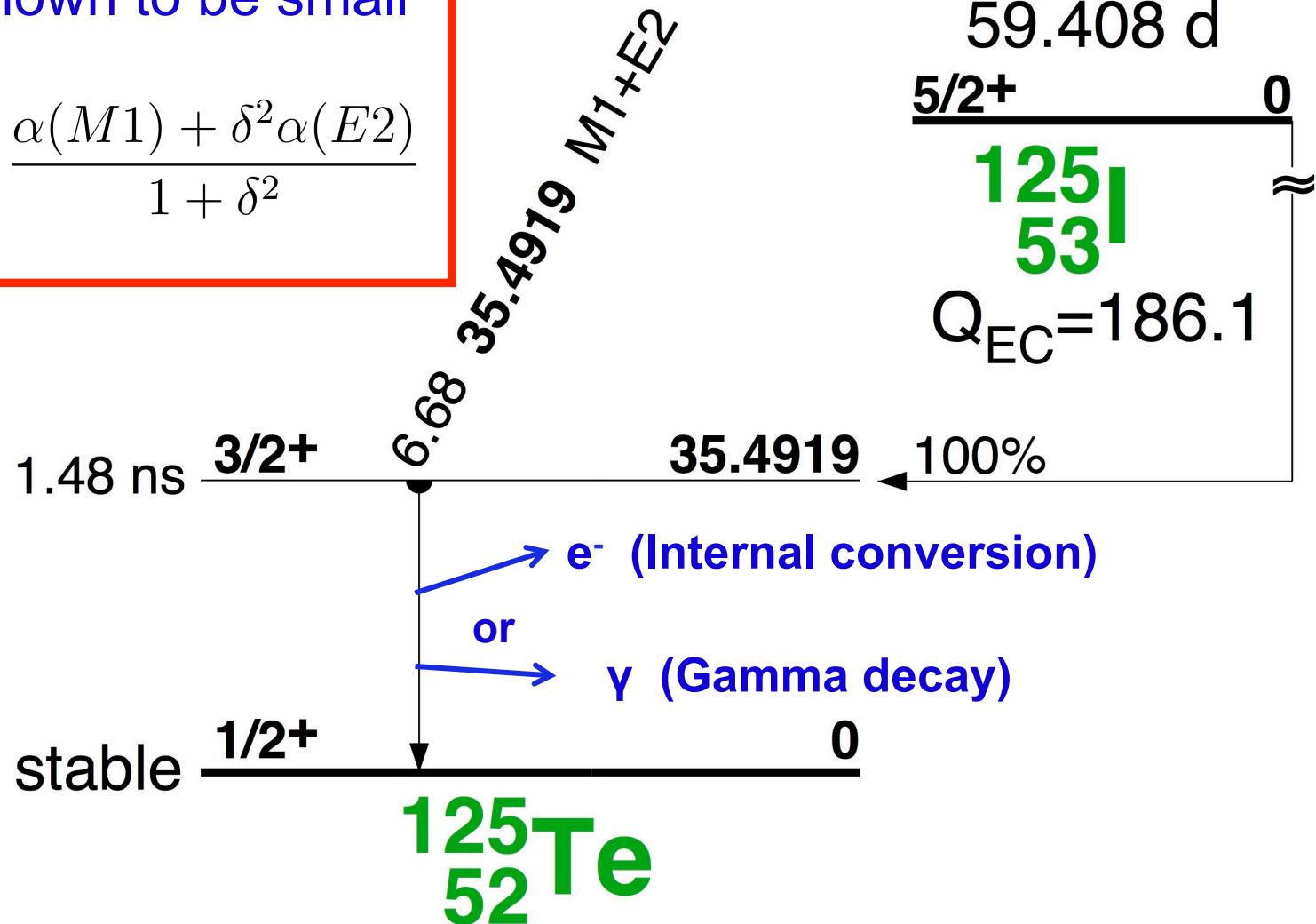




Decay scheme of ^{125}I

Mixing ratio of the M1+E2 transition is known to be small ($\delta \ll 1$).

$$\alpha(M1 + E2) = \frac{\alpha(M1) + \delta^2 \alpha(E2)}{1 + \delta^2}$$



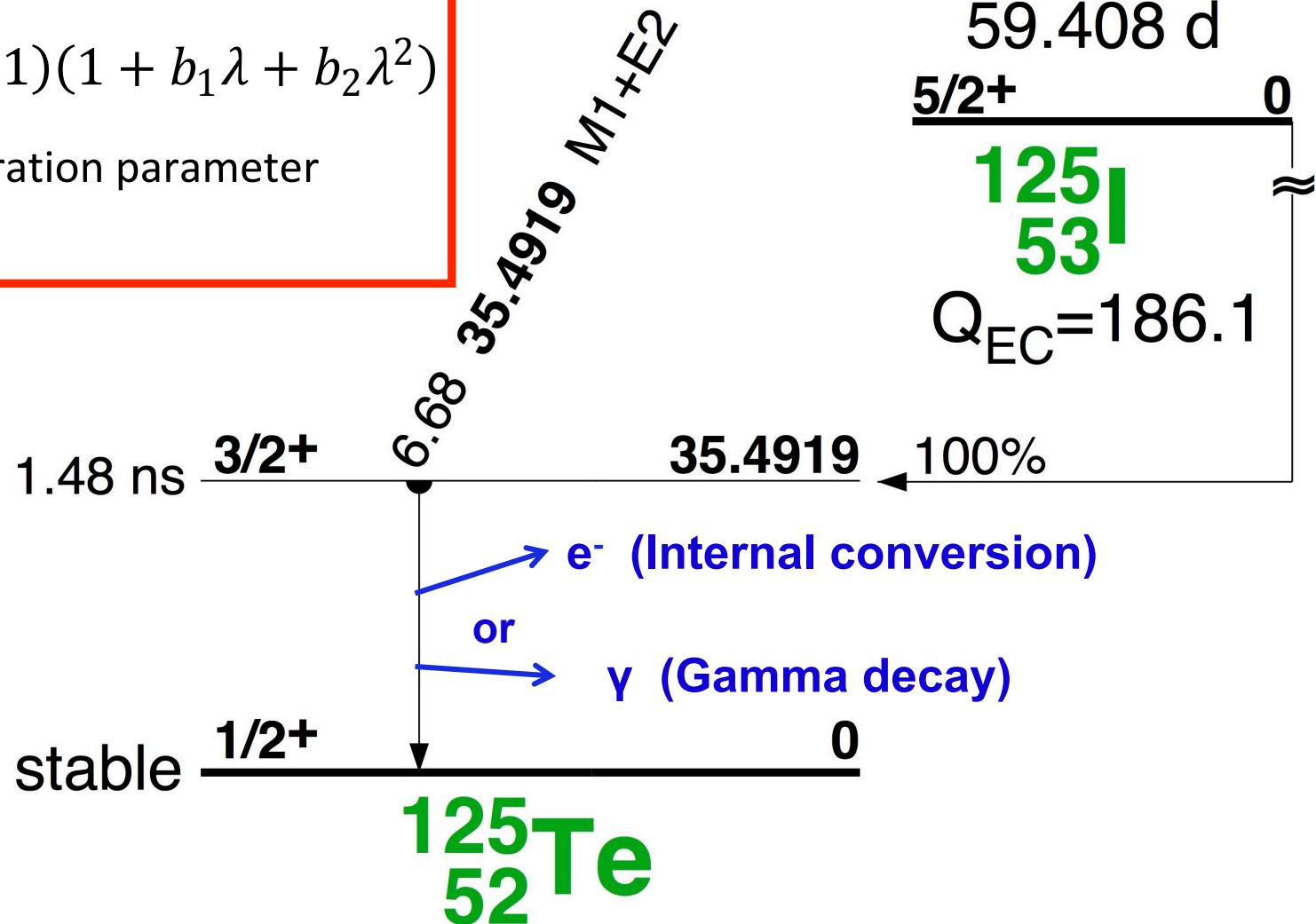


Decay scheme of ^{125}I

Penetration effects

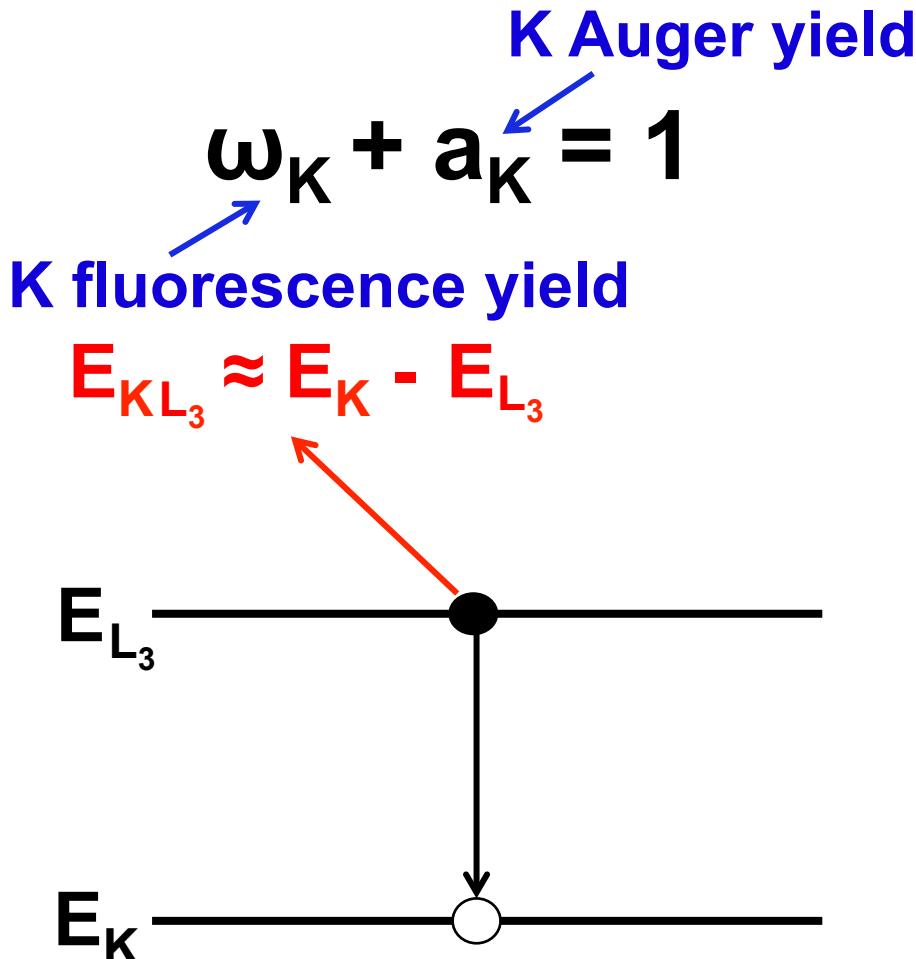
$$\alpha(M1) = \alpha_0(M1)(1 + b_1\lambda + b_2\lambda^2)$$

Where λ = penetration parameter



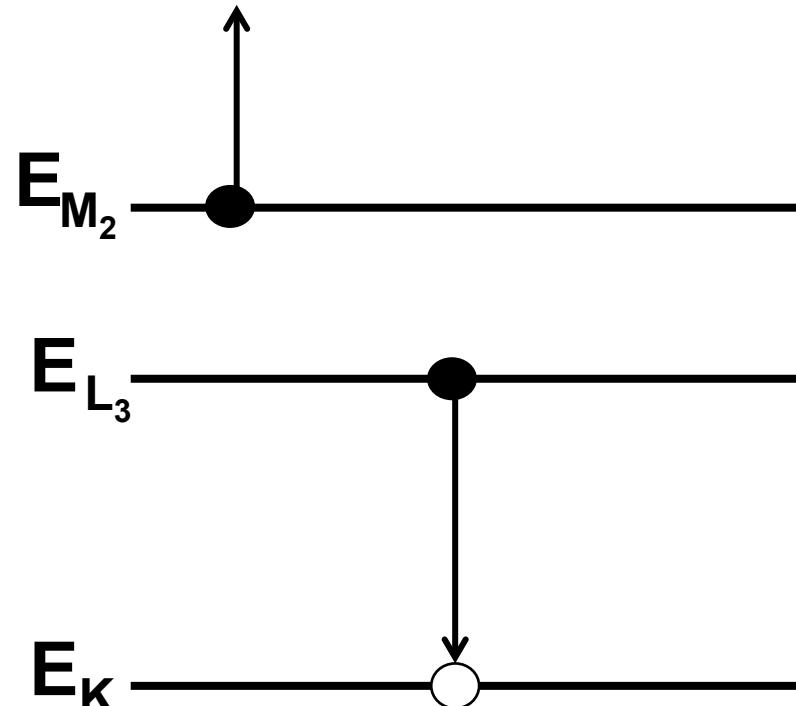


X-ray transition



Auger transition

$$E_{KL_3M_2} \approx E_K - E_{L_3} - E_{M_2}$$

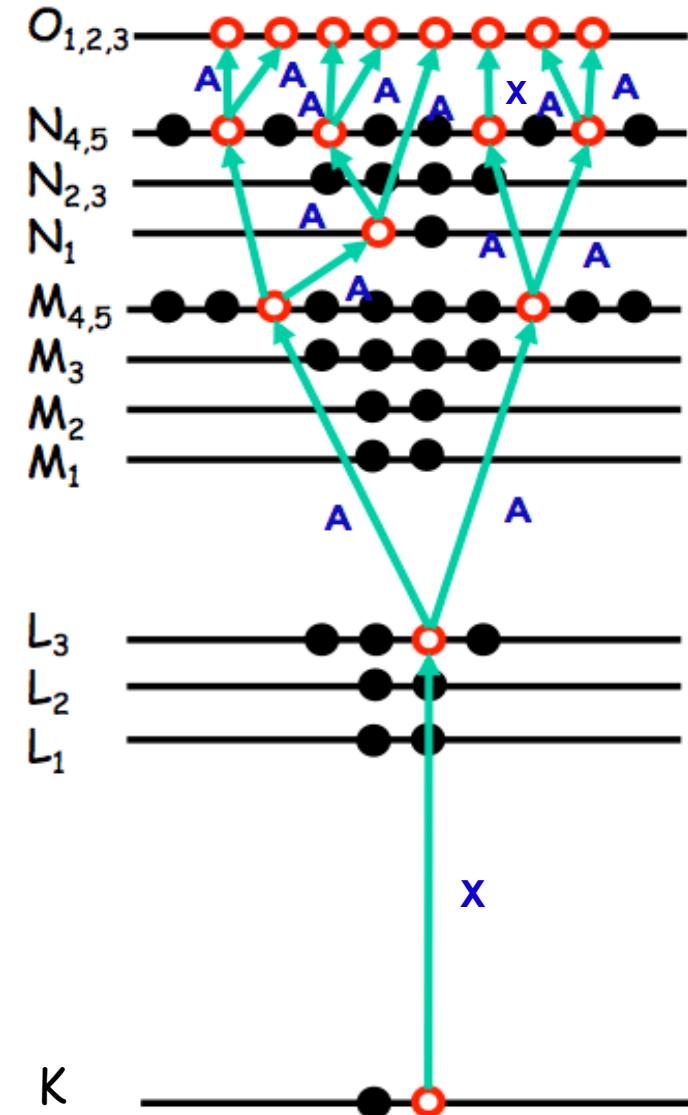


*Atomic notations: K = $1s_{1/2}$, L₃ = $2p_{3/2}$, M₂ = $3p_{1/2}$



Vacancy cascade

- ❖ Resulting in heaps of Auger electrons
- ❖ Energy range: a few eV to 30 keV (for ^{125}I case)

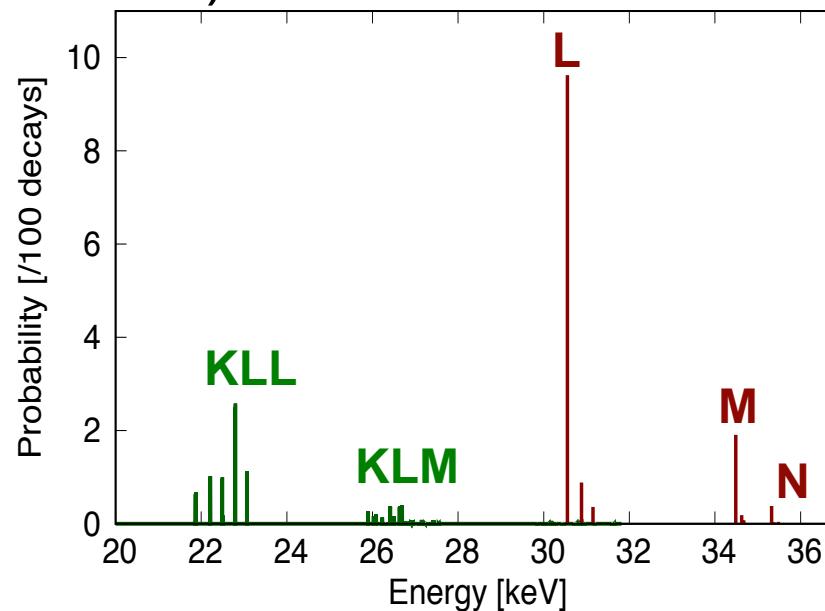


Note: X = X-ray transition, A = Auger transition



Computational model - BriccEmis

- ❖ Calculate the Auger and X-ray spectra using a Monte Carlo approach
- ❖ Transition probabilities from Evaluated Atomic Data Library (EADL) (Perkins 1991)
- ❖ Transition energies are calculated using the relativistic self-consistent-field Dirac Fock method, using RAINIE code (Band 2002)





- ❖ Measure an accurate Auger yield from medical radioisotope ^{125}I

Approach

- I. Determine the nuclear parameters (λ and δ)
- II. Measure the Auger to conversion electrons intensity ratios.
- III. Deduce the absolute intensity of Auger electrons from the conversion coefficients.



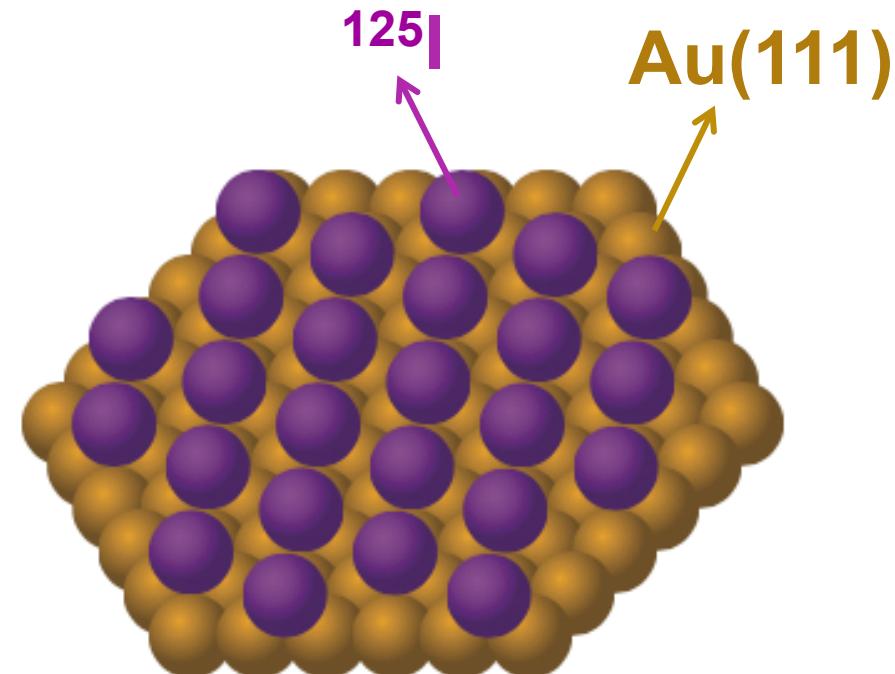
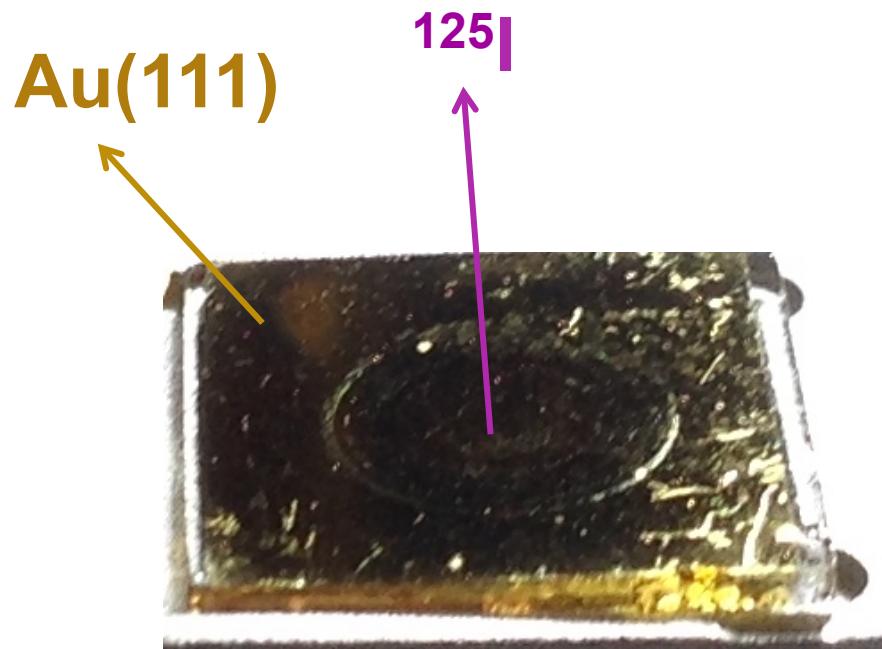
Project Aims

- ❖ Measure an accurate Auger yield from medical radioisotope ^{125}I
- ❖ Test and benchmark the model



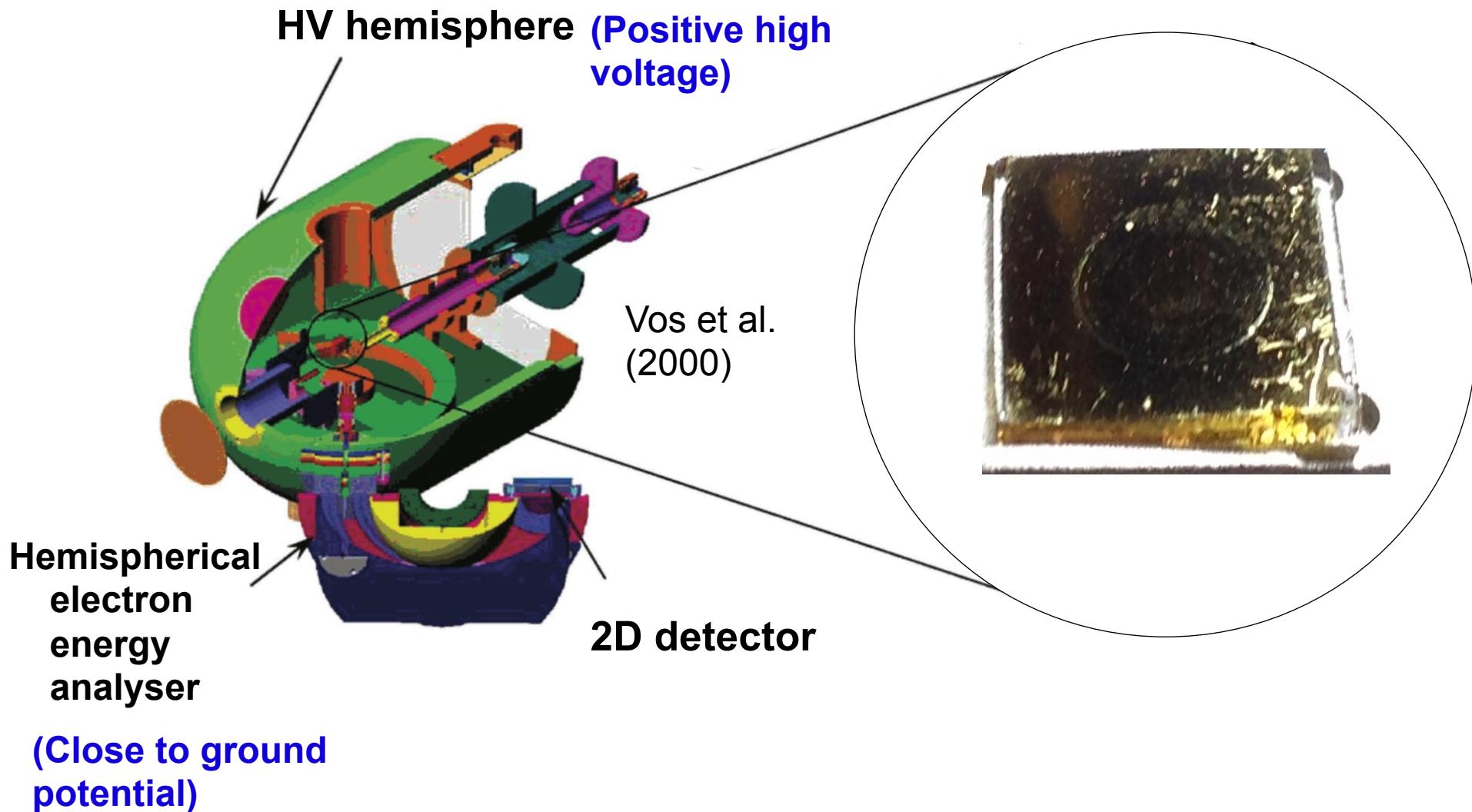
Source preparation

- ❖ Monolayer of ^{125}I on top of a gold substrate





High-energy electrostatic spectrometer

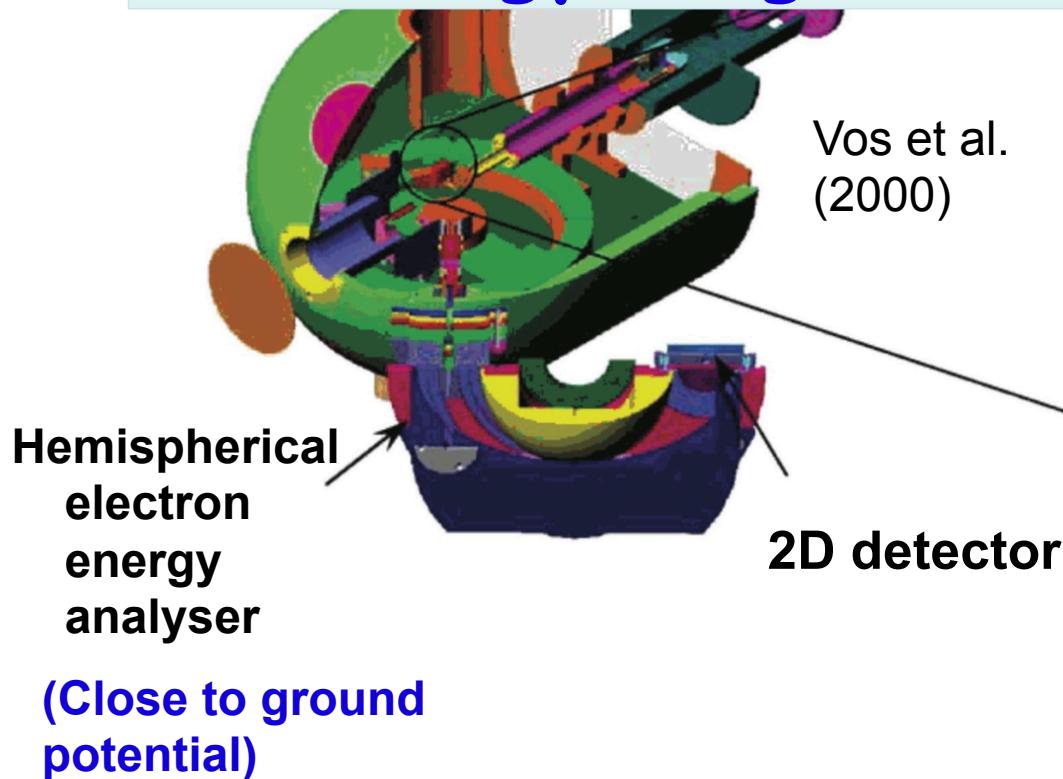




High-energy electrostatic spectrometer

HV hemisphere (Positive high voltage)

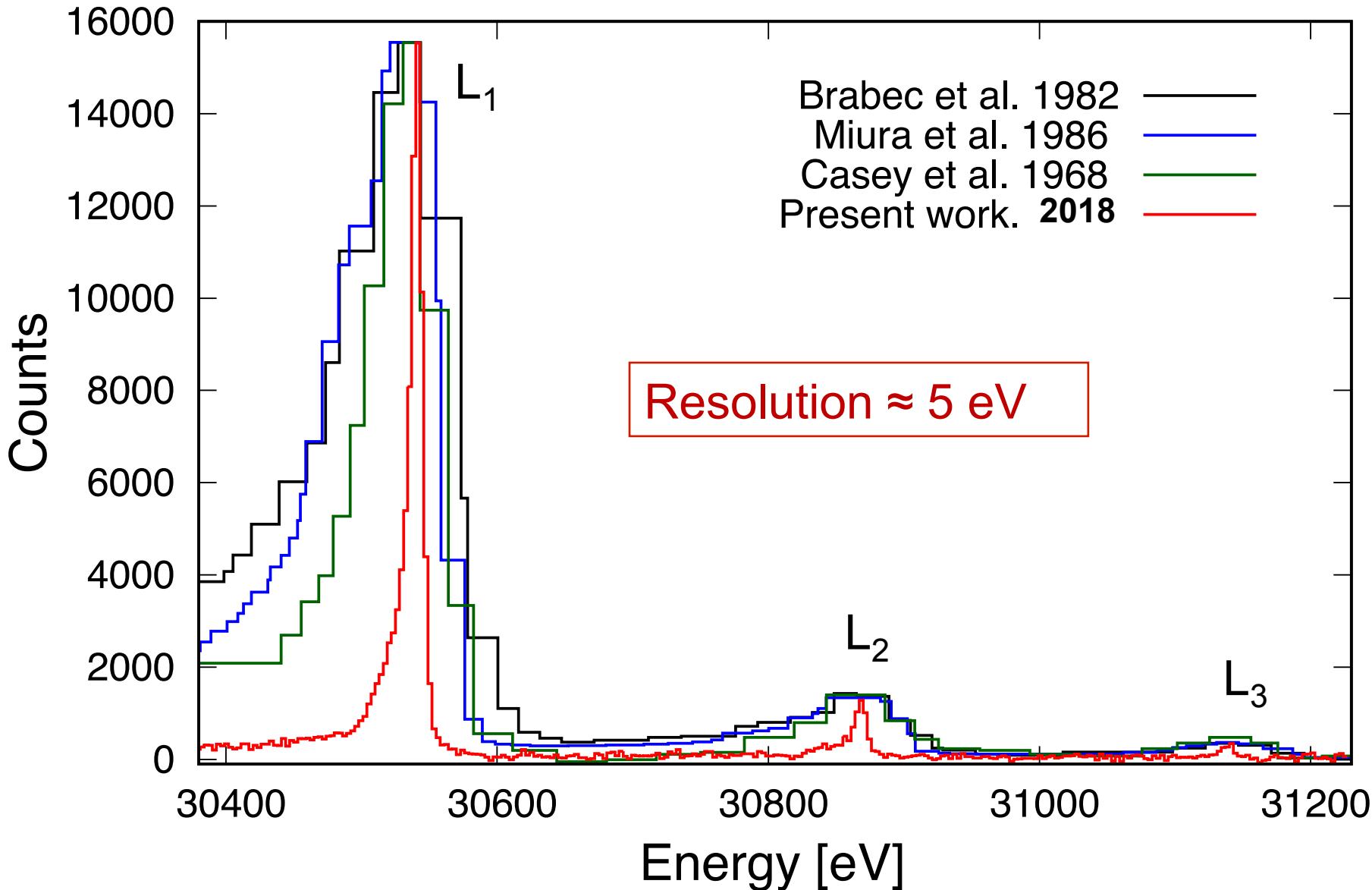
Energy range: 2 keV to 40 keV



2D detector

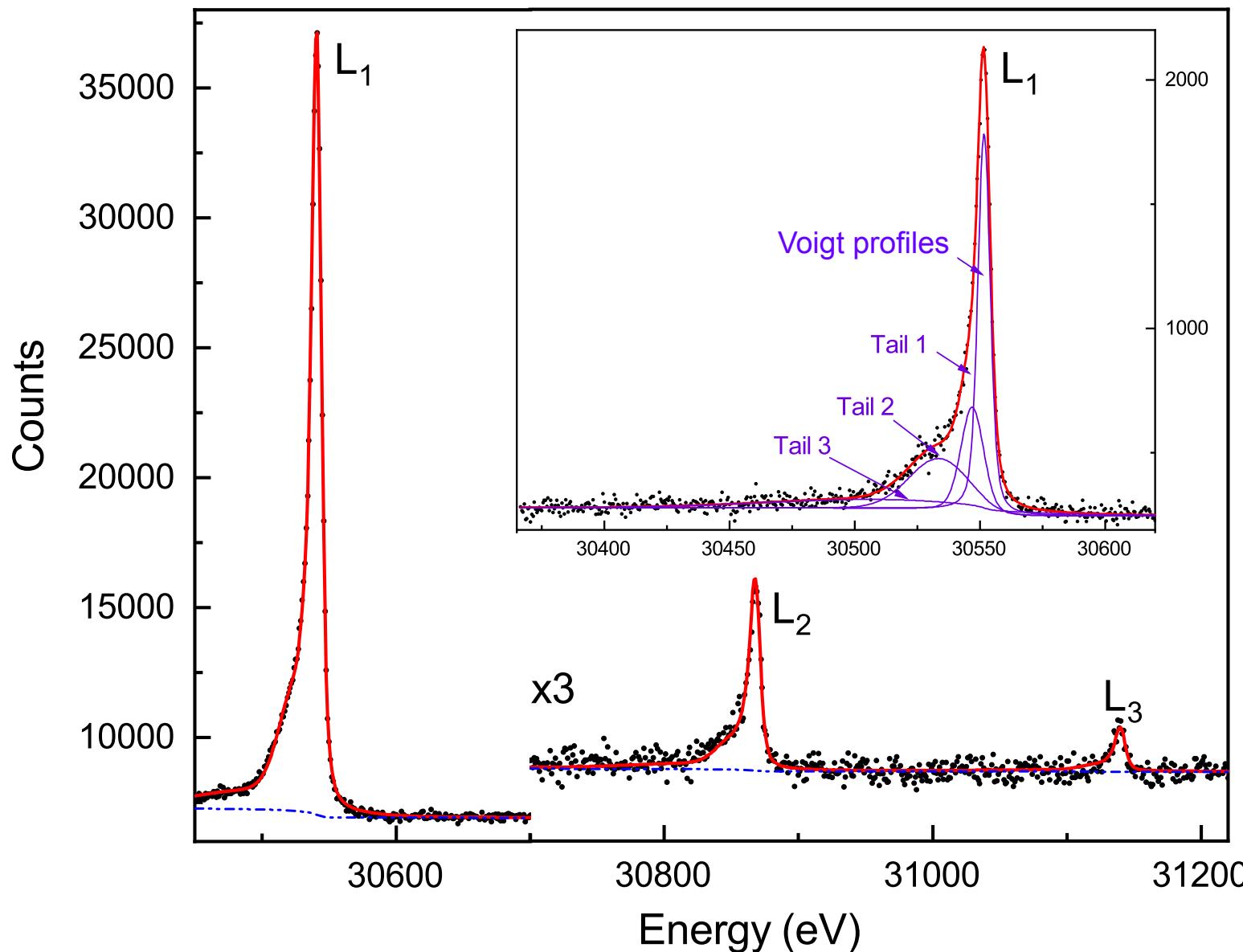


Conversion electron measurements





Conversion electron line shapes





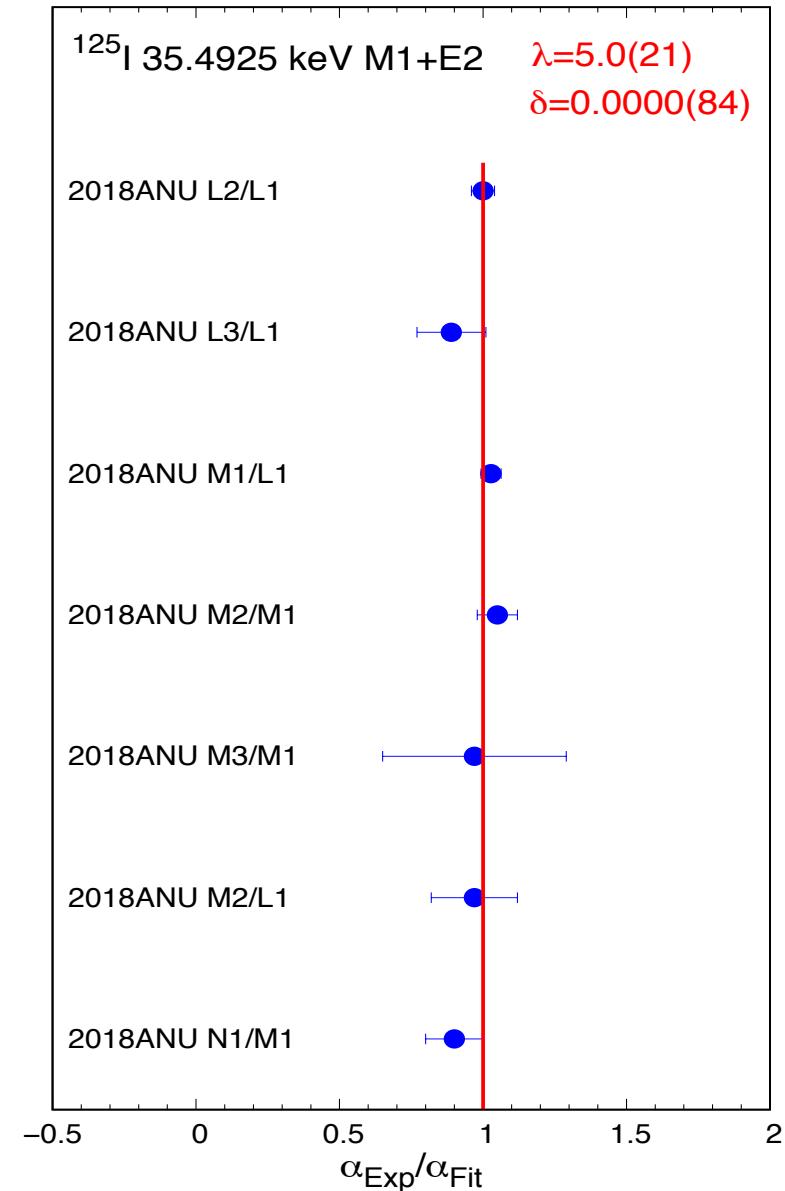
Nuclear parameters determination

ANU data only

- ❖ Chi-square fitting method
- ❖ Reduced $\chi^2 = 0.63$
- ❖ $\lambda = 5.0(21)$, $\delta = 0.0000(84)$

Atomic shell	<i>Experiment</i>	
	Present work	Literature
100/(1+ <i>Tot</i>)		6.68(14) [12] 6.55(13) [13]
<i>Tot</i>		12.95(28) [15] ^a 14.25(64) [8]
<i>K</i> /(1 + <i>Tot</i>)		0.80(5) [16] 0.804(10) [17]
<i>L</i> /(1 + <i>Tot</i>)		0.11(2) [16]
<i>M</i> /(1 + <i>Tot</i>)		0.020(4) [16]
<i>K</i>		11.78(18) ^a [15] 11.90(31) [8]
<i>L</i>		1.4(1) [18]
<i>K/L</i>		12.3(25) [10]
<i>L/M</i>		5.21(26) [9]
<i>M/N</i>		4.87(20) [9]
<i>L</i> ₁ : <i>L</i> ₂ : <i>L</i> ₃	1:0.085(2):0.019(2)	1:0.089(4):0.024(2) [7] 1:0.106(22):0.041(2) [10] 1:0.082(4):0.019(3) [8] 1:0.095(2):0.023(5) [9]
<i>L</i> ₁ : <i>M</i> ₁	1: 0.204(7)	-
<i>M</i> ₁ : <i>M</i> ₂ : <i>M</i> ₃	1:0.094(6):0.022(7)	1:0.092(5):0.044(3) [8] 1:0.101(5):0.030(5) [9]
<i>L</i> ₁ : <i>M</i> ₂ <i>M</i> ₁ : <i>N</i> ₁	1:0.0173(26) 1:0.179(20)	- 1:0.214(6) [9]

^a Corrected ω_K to 0.875



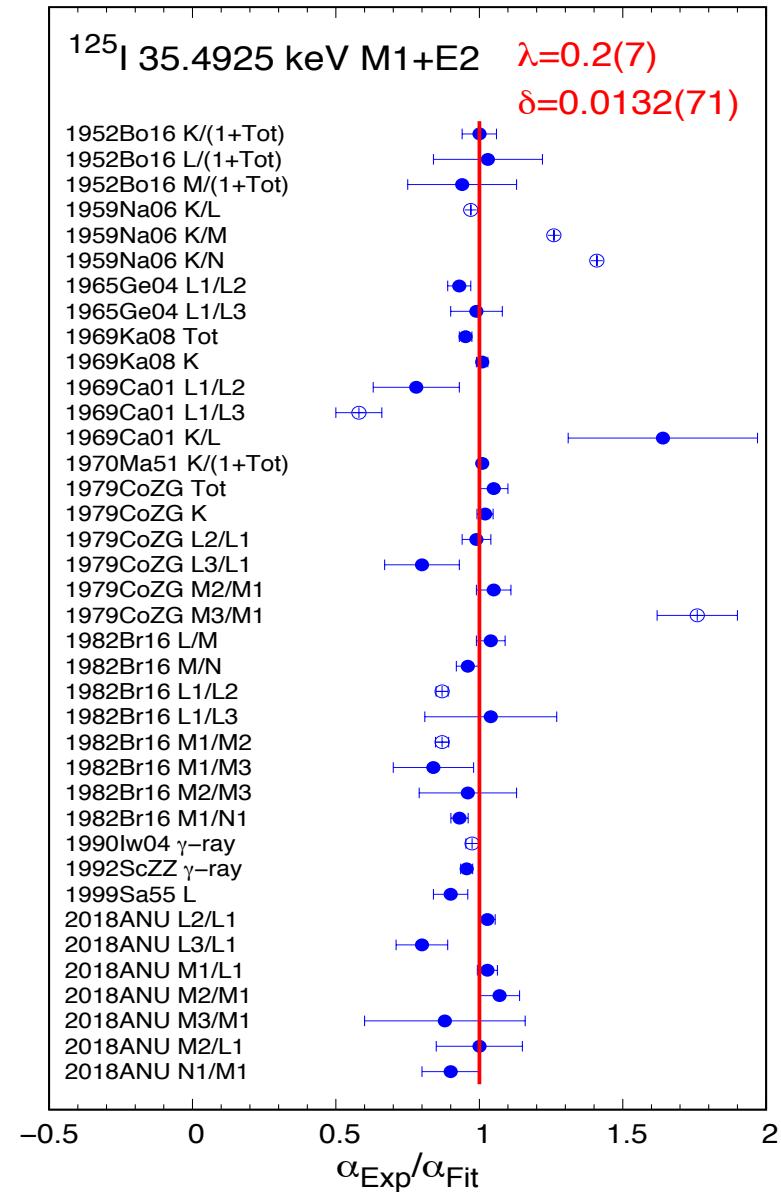
Nuclear parameters determination

All data

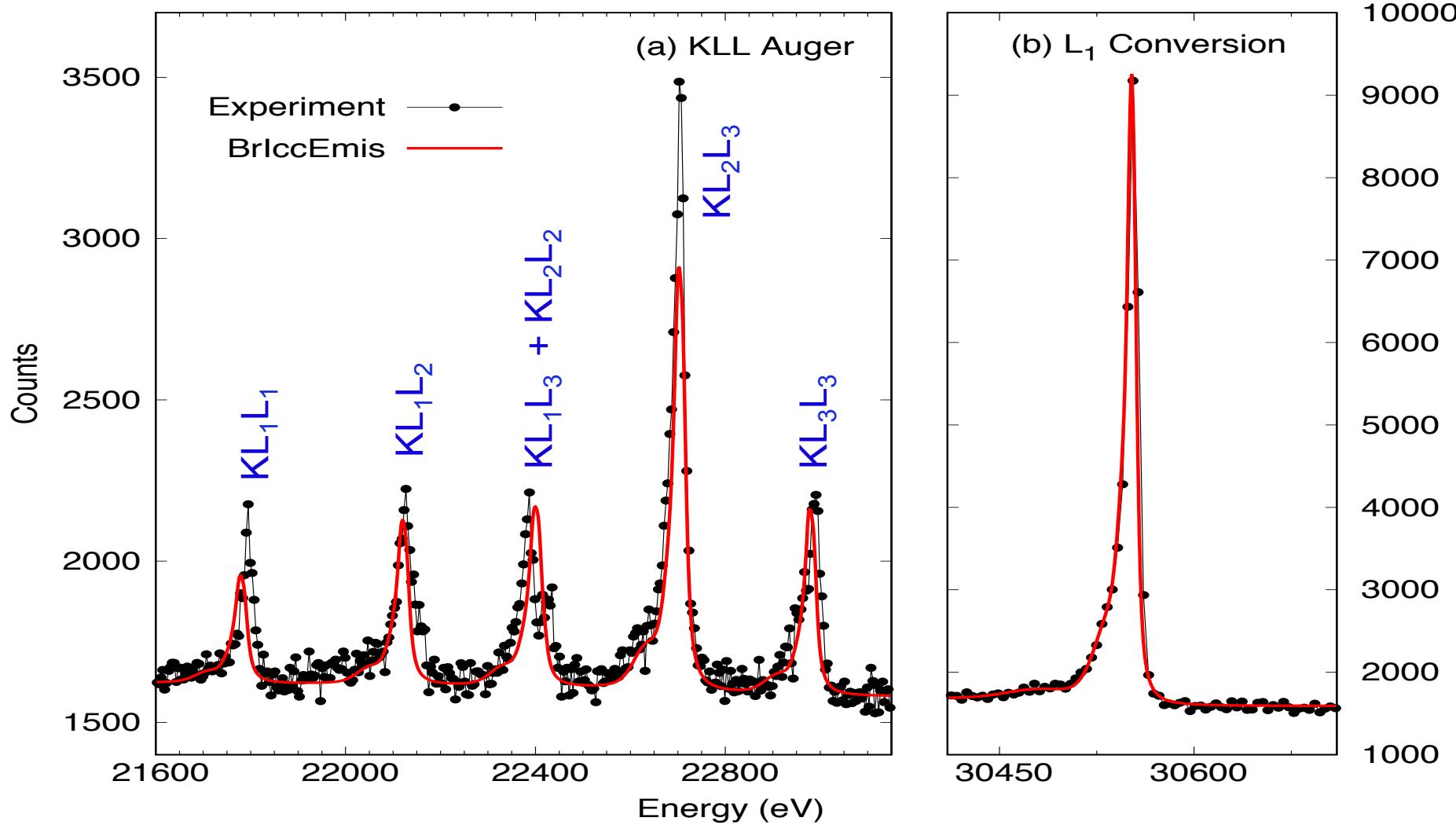
- ❖ Chi-square fitting method
- ❖ Reduced $\chi^2 = 1.55$
- ❖ $\lambda = 0.2(7)$, $\delta = 0.0132(71)$

Atomic shell	<i>Experiment</i>	
	Present work	Literature
100/(1+ <i>Tot</i>)		6.68(14) [12] 6.55(13) [13]
<i>Tot</i>		12.95(28) [15] ^a 14.25(64) [8]
<i>K</i> /(1 + <i>Tot</i>)		0.80(5) [16] 0.804(10) [17]
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<i>L</i> ₁ : <i>L</i> ₂ : <i>L</i> ₃	1:0.085(2):0.019(2)	1:0.089(4):0.024(2) [7] 1:0.106(22):0.041(2) [10] 1:0.082(4):0.019(3) [8] 1:0.095(2):0.023(5) [9]
<i>L</i> ₁ : <i>M</i> ₁	1: 0.204(7)	-
<i>M</i> ₁ : <i>M</i> ₂ : <i>M</i> ₃	1:0.094(6):0.022(7)	1:0.092(5):0.044(3) [8] 1:0.101(5):0.030(5) [9]
<i>L</i> ₁ : <i>M</i> ₂	1:0.0173(26)	-
<i>M</i> ₁ : <i>N</i> ₁	1:0.179(20)	1:0.214(6) [9]

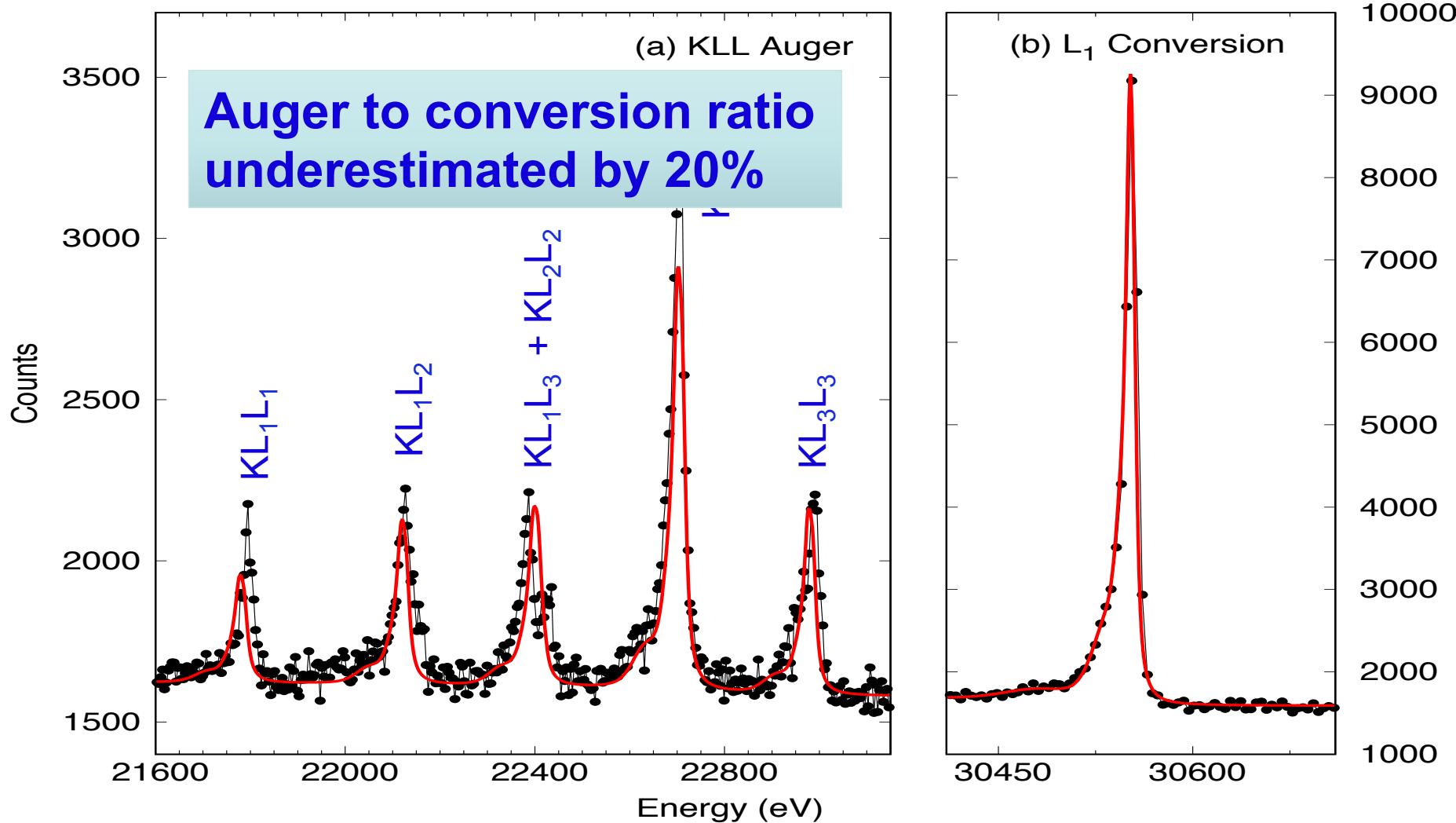
^a Corrected ω_K to 0.875



KLL Auger electron measurements

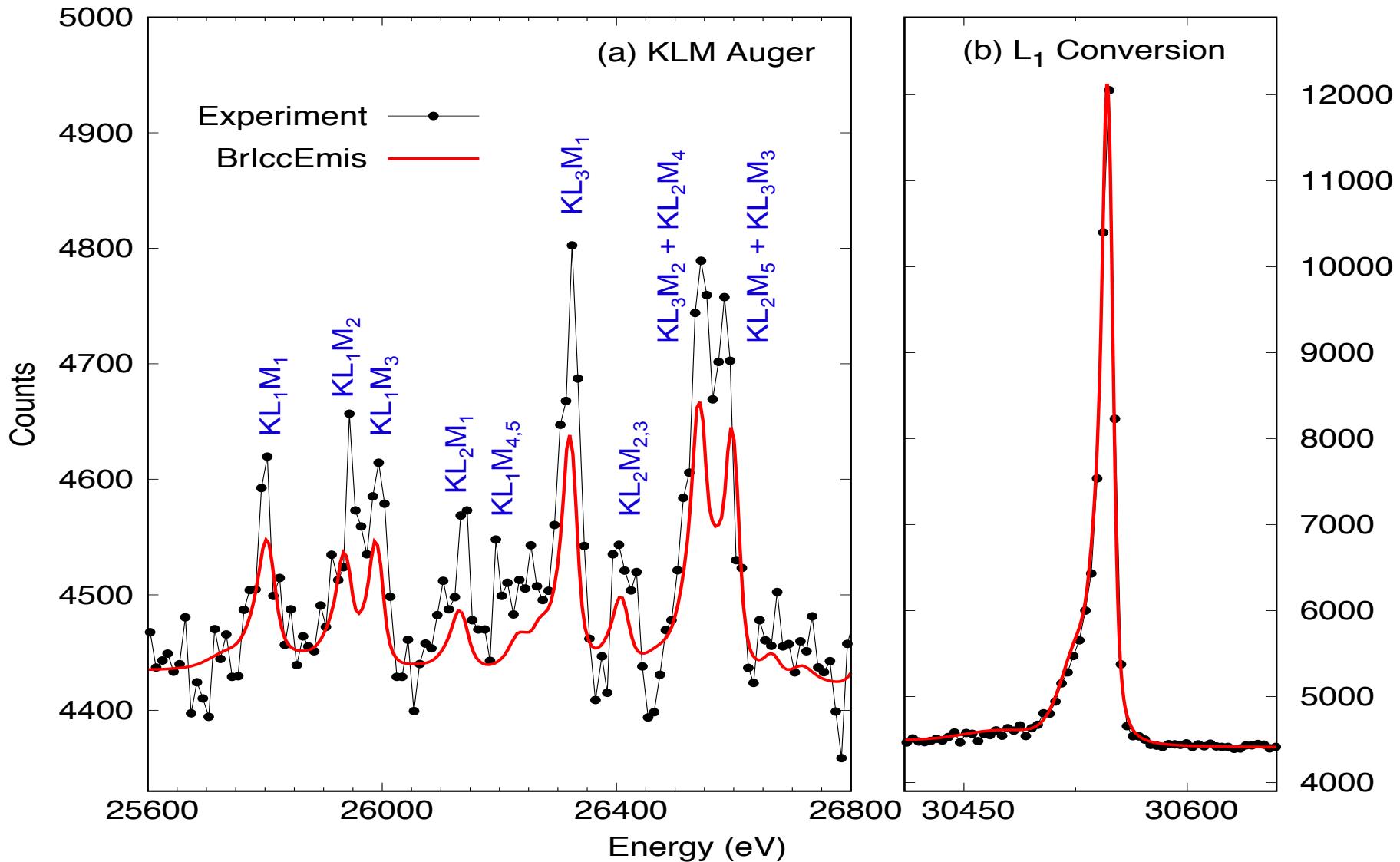


KLL Auger electron measurements



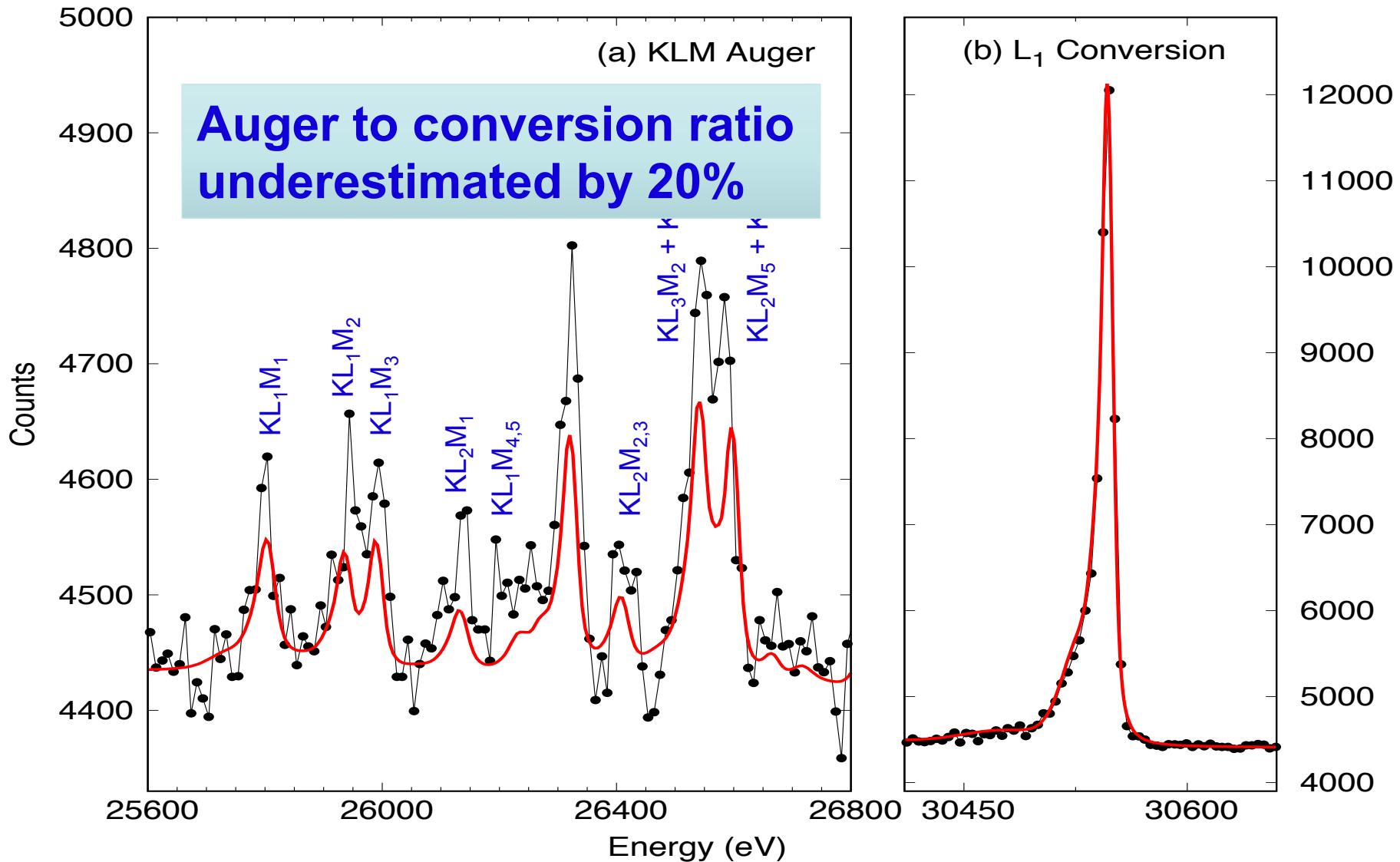


KLM Auger electron measurements





KLM Auger electron measurements





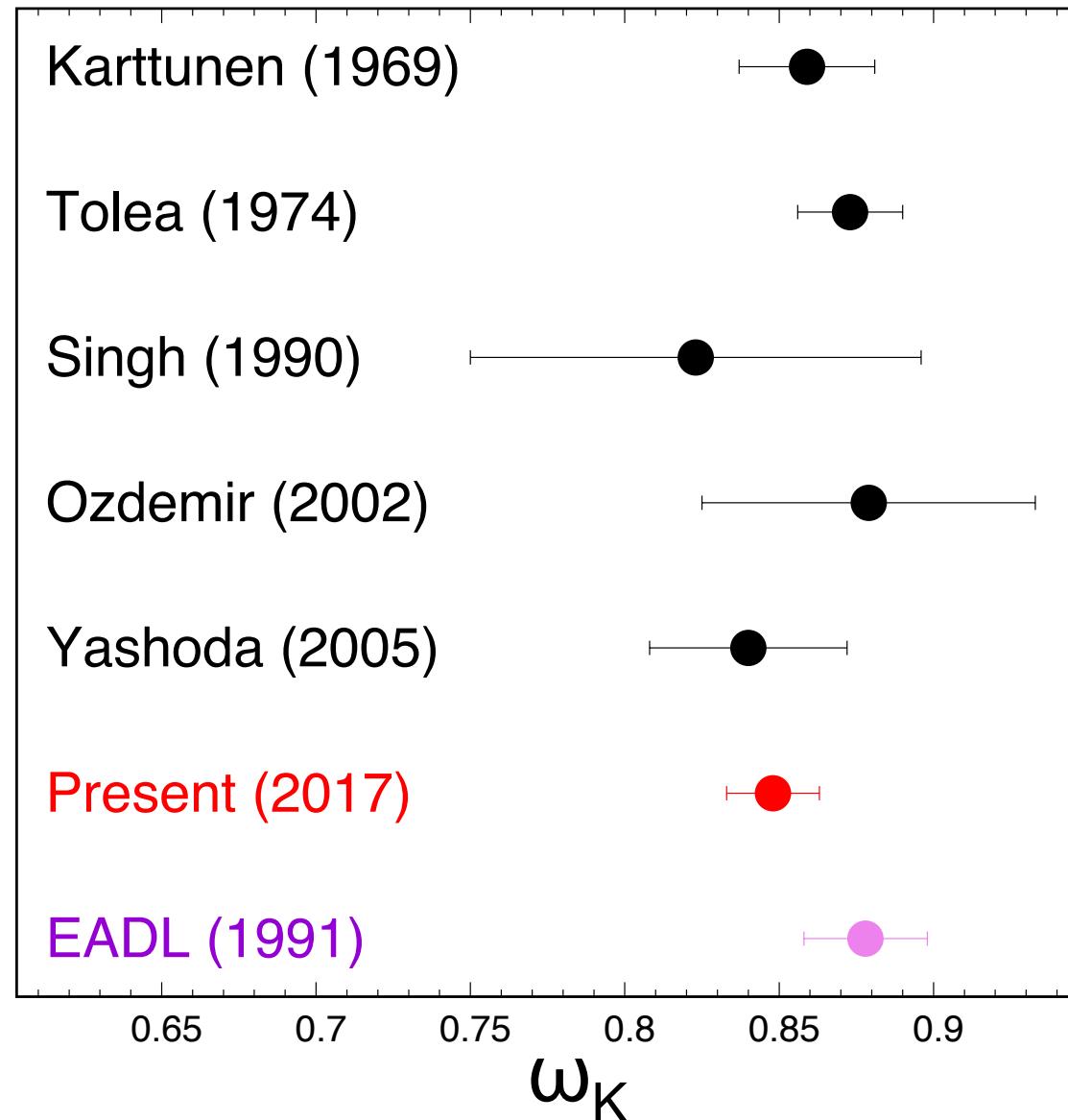
K fluorescence yield determination

$$\omega_K + a_K = 1$$

↑ ↑
0.87 0.13
5% **30%**

0.85 0.15
2% **10%**

0.88 0.12
3% **20%**





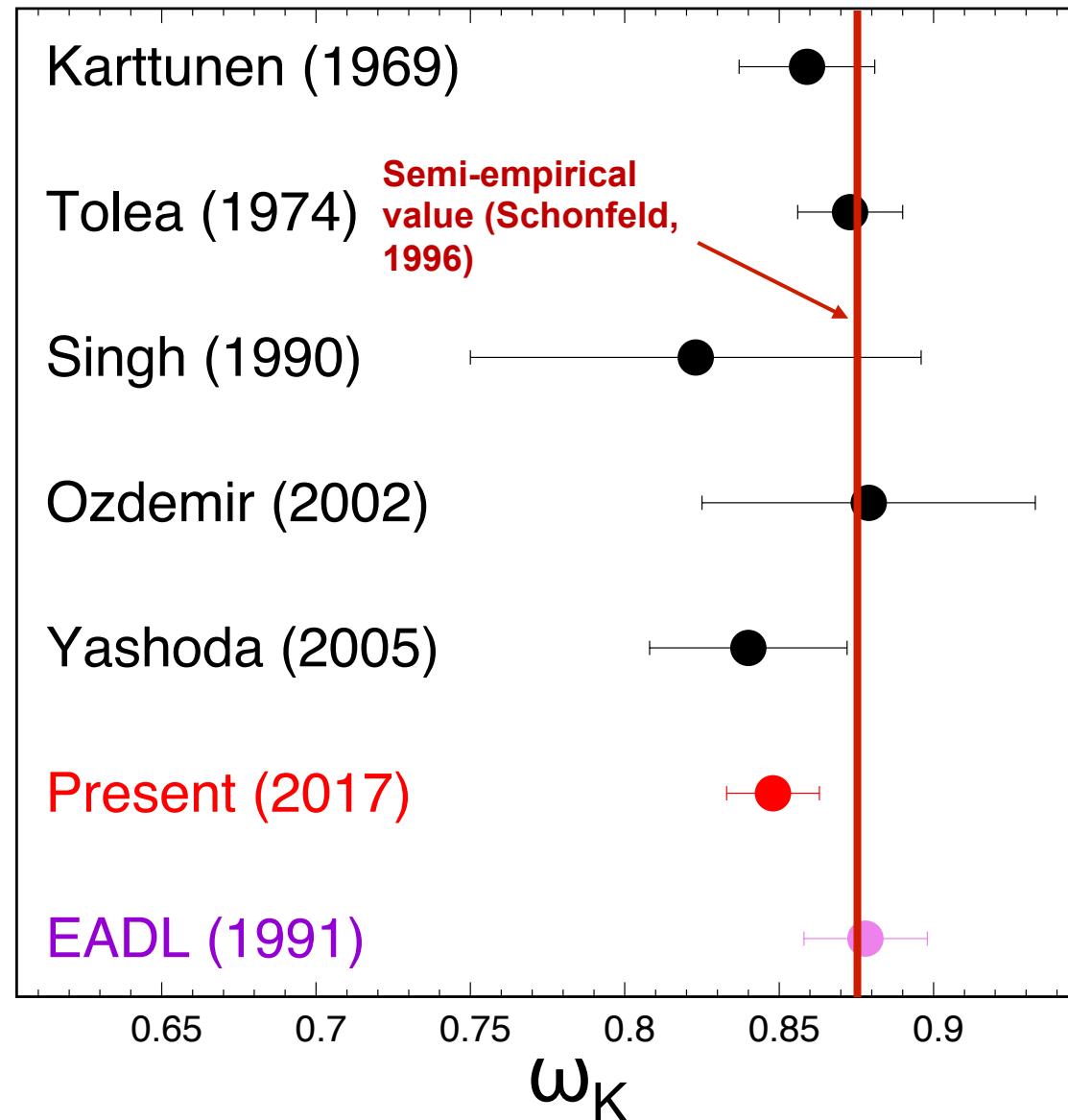
K fluorescence yield determination

$$\omega_K + a_K = 1$$

↑ ↑
0.87 0.13
5% 30%

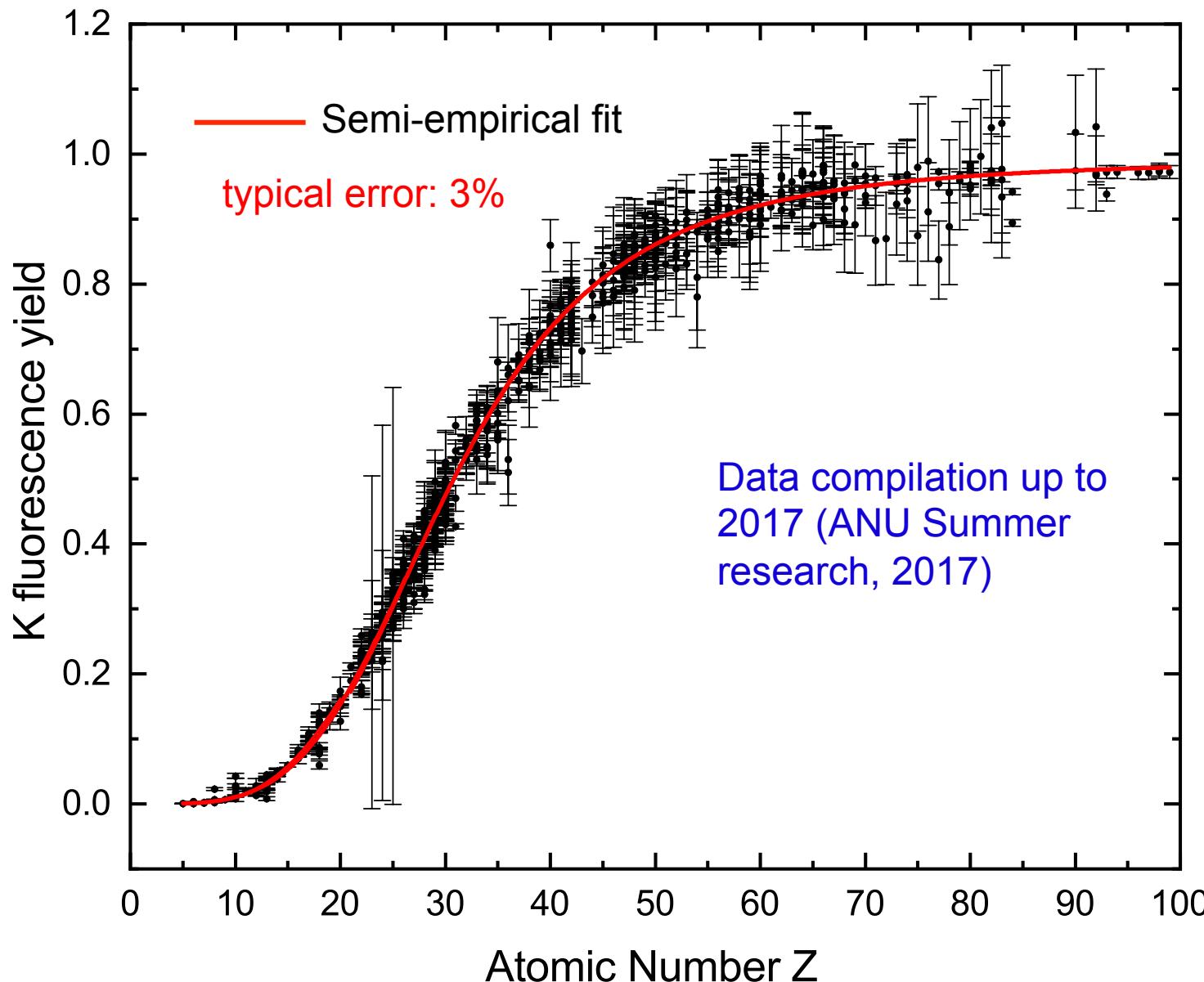
0.85 0.15
2% 10%

0.88 0.12
3% 20%





K fluorescence yield global fit curve





What to do next?

- ❖ Quantify the Auger electrons with energy < 1 keV
- ❖ Effects of electron shake-off following internal conversion and Auger transition (the tails)
- ❖ Atomic structure effect: What is the atomic field after electron capture
- ❖ Potential medical isotopes to study: ^{80m}Br , ^{99m}Tc , ^{99}Mo , ^{119}Sc , ^{153}Sm , ^{177}Lu , ^{193m}Pt , ^{195m}Pt , ^{201}TI , ^{80m}Br



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3. Prof. Andrew Stuchbery

- Collaborators

1. M. Alotiby
2. B.Q Lee
3. ANSTO (Australian Nuclear Science and Technology Organisation)

Thank you

