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Advances in Theoretical Studies of Electron Capture

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Electron Capture (EC)

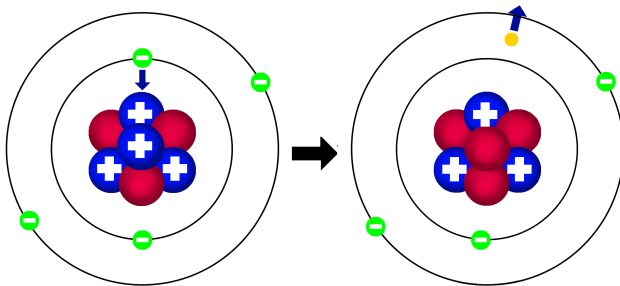


Figure: Electron Capture

Why EC?

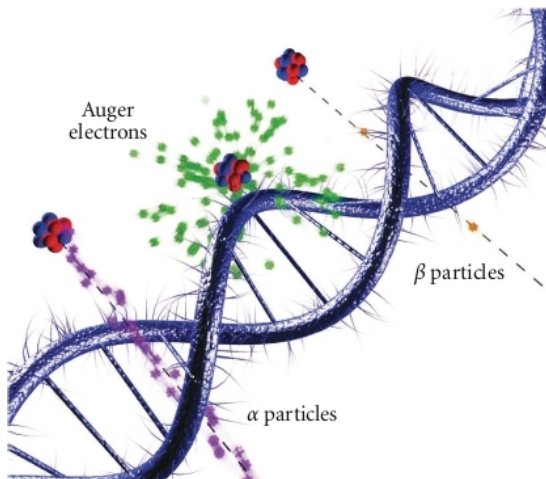


Figure: EC in Nuclear medicine [1]

Why EC?

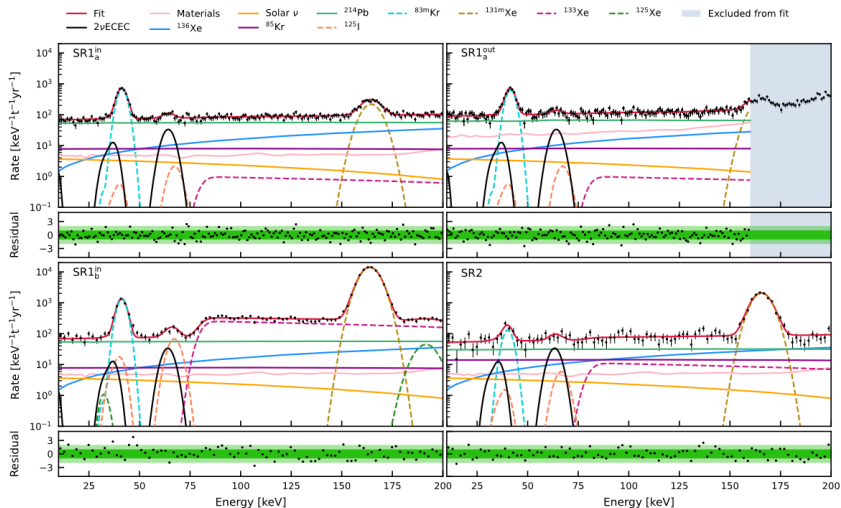


Figure: XENON1T Collaboration experimental results [2]

Why EC?

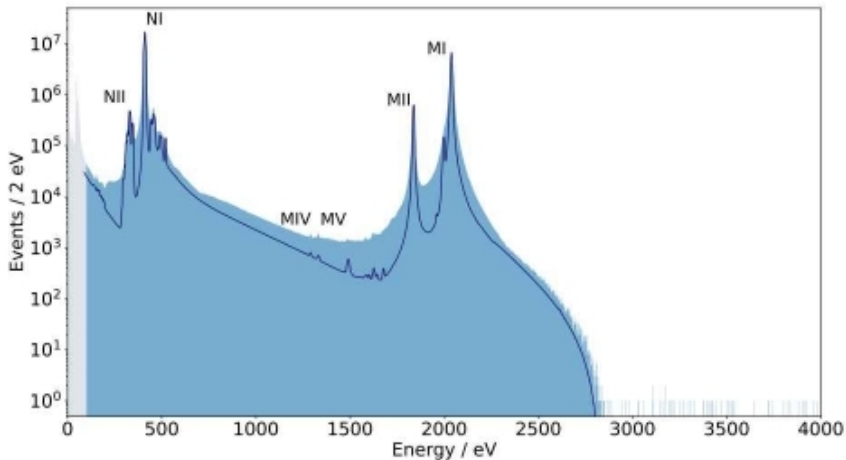


Figure: ECHO-1k experimental results [3]

$$\lambda = \frac{G_{\beta}^2}{2\pi^3} \sum_x N_x C_x F_x S_x \quad (1)$$

$$F_x = \frac{\pi}{2} q_x^2 \beta_x^2 B_x \quad (2)$$

- $x = (n\kappa)$
- $k_x = |\kappa| \leq L$
- N_x - Relative occupation number
- p_x, q_x - Electron, neutrino momentum
- β_x - Coulomb amplitude
- B_x - The exchange and overlap correction factor
- S_x - The shake-up and shake-off correction factor

$$\begin{aligned}
 C_x = & \left[M_L \left(k_x, k_\nu^{(1)} \right) + (\kappa_x/k_x) m_L \left(k_x, k_\nu^{(1)} \right) \right]^2 \\
 & + \left[M_L \left(k_x, k_\nu^{(2)} \right) + (\kappa_x/k_x) m_L \left(k_x, k_\nu^{(2)} \right) \right]^2 \\
 & + \left[M_{L+1} \left(k_x, k_\nu^{(2)} \right) + (\kappa_x/k_x) m_{L+1} \left(k_x, k_\nu^{(2)} \right) \right]^2 \\
 & + \delta_{\Delta J,0} \left[M_0(1,1) + (\kappa_x/k_x) m_0(1,1) \right]^2
 \end{aligned} \tag{3}$$

For (L-1)th unique forbidden transitions:

$$C_x = \frac{(2L-2)!!}{(2L-1)!!} \left({}^A F_{LL-11}^0 \right)^2 R^{2(L-1)} \frac{p_x^{2(k_x-1)} q_x^{2(L-k_x)}}{(2k_x-1)! \{2(L-k_x)+1\}!} \tag{4}$$

For allowed transitions: $C_x = \delta_{\Delta J,0} \left({}^V F_{000}^0 \right)^2 + \left({}^A F_{101}^0 \right)^2$

$$\begin{aligned} k_{\nu}^{(1)} &= L - k_x + 1, \\ k_{\nu}^{(2)} &= L - k_x + 2. \end{aligned} \tag{5}$$

$$\begin{aligned} M_L(k_x, k_{\nu}^{(1)}) &= K_L (p_x R)^{k_x - 1} (q_x R)^{k_{\nu}^{(1)} - 1} \times \\ &\left\{ -[(2L + 1)/L]^{1/2} {}^V F_{LL-11}^0 + (2k_x + 1)^{-1/2} \alpha Z {}^V F_{LL0}^0(k_x, 1, 1, 1) \right. \\ &+ \left[(2k_x + 1)^{-1} (2k_{\nu}^{(1)} + 1)^{-1} q_x R \right] {}^V F_{LL0}^0 - \\ &(2k_x + 1)^{-1} \alpha Z [(L + 1)/L]^{1/2} {}^A F_{LL1}^0(k_x, 1, 1, 1) \\ &\left. - \left[(2k_x + 1)^{-1} W_x R + (2k_{\nu}^{(1)} + 1)^{-1} q_x R \right] [(L + 1)/L]^{1/2} {}^A F_{LL1}^0 \right\} \end{aligned} \tag{6}$$

More decay constant

$$m_L \left(k_x, k_\nu^{(1)} \right) = K_L (p_x R)^{k_x - 1} (q_x R)^{k_\nu^{(1)} - 1} (2k_x + 1)^{-1} R \\ \left\{ {}^V F_{LL0}^0 - [(L+1)/L]^{1/2} A F_{LL1}^0 \right\},$$

$$M_L \left(k_x, k_\nu^{(2)} \right) = -\tilde{K}_L (p_x R)^{k_x - 1} (q_x R)^{k_\nu^{(2)} - 1} \sqrt{\frac{L+1}{(2k_x - 1)(2k_\nu^{(2)} - 1)}} \quad (7)$$

$$\left\{ {}^V F_{LL0}^0 + (k_x - k_\nu^{(2)}) (L+1)^{-1} [(L+1)/L]^{1/2} A F_{LL1}^0 \right\},$$

$$M_{L+1} \left(k_x, k_\nu^{(2)} \right) = -\tilde{K}_L (p_x R)^{k_x - 1} (q_x R)^{k_\nu^{(2)} - 1} A F_{(L+1)L1}^0.$$

$$K_L = (1/2)^{1/2} [(2L)!! / (2L+1)!!]^{1/2} \left[(2k_x - 1)! (2k_\nu^{(1)} - 1)! \right]^{-1/2}, \quad (8)$$

$$\tilde{K}_L = [(2L)!! / (2L+1)!!]^{1/2} \left[(2k_x - 1)! (2k_\nu^{(2)} - 1)! \right]^{-1/2}. \quad (9)$$

Form-factors and nuclear matrix elements

$$\begin{aligned} {}^V F_{KLS}^N(k_x, m, n, \rho) &= (-1)^{K-L} {}^V \mathfrak{M}_{KLS}^N(k_x, m, n, \rho), \\ {}^A F_{KLS}^N(k_x, m, n, \rho) &= (-1)^{K-L} g_A {}^A \mathfrak{M}_{KLS}^N(k_x, m, n, \rho), \end{aligned} \quad (10)$$

$$\begin{aligned} {}^V F_{KLS}^N &= {}^V F_{KLS}^N(k_x, m, n, 0) = (-1)^{K-L} {}^V \mathfrak{M}_{KLS}^N(k_x, m, n, 0), \\ {}^A F_{KLS}^N &= {}^A F_{KLS}^N(k_x, m, n, 0) = (-1)^{K-L} g_A {}^A \mathfrak{M}_{KLS}^N(k_x, m, n, 0), \end{aligned} \quad (11)$$

$$\begin{aligned} (-1)^{J_f - M_f} \begin{pmatrix} J_f & K & J_i \\ -M_f & M & M_i \end{pmatrix} \left\{ {}^V \mathfrak{M}_{KLS}^N(k_x, m, n, \rho) + g_A {}^A \mathfrak{M}_{KLS}^N(k_x, m, n, \rho) \right\} &= \sqrt{\frac{4\pi}{2J_i + 1}} \times \\ \iint \cdots \int \psi_f^\dagger(1, 2, \dots, A; J_f M_f \pi_f) \times \sum_{j=1}^A \left\{ \left(\frac{r}{R} \right)^{L*2N} I(k_x, m, n, \rho; r) (1 + g_A \gamma_5) T_{KLS}^M t^+ \right\}_j \times \\ \psi_i(1, 2, \dots, A; J_i M_i \pi_i) d\tau_1 d\tau_2 \cdots d\tau_A. \end{aligned} \quad (12)$$

Electronic wave function

$$\psi_{n\kappa m}(\vec{r}) = \begin{pmatrix} g_{n\kappa}(r)\Omega_{\kappa,m}(\hat{r}) \\ if_{n\kappa}(r)\Omega_{-\kappa,m}(\hat{r}) \end{pmatrix} \quad (13)$$

$$\begin{aligned} \left(\frac{d}{dr} + \frac{\kappa + 1}{r}\right) g_{n\kappa}(r) &= (E_{n\kappa} - V(r) + m_e) f_{n\kappa}(r), \\ \left(\frac{d}{dr} - \frac{\kappa - 1}{r}\right) f_{n\kappa}(r) &= -(E_{n\kappa} - V(r) - m_e) g_{n\kappa}(r). \end{aligned} \quad (14)$$

Coulomb Amplitudes

$$\begin{Bmatrix} g_{n\kappa}(r) \\ f_{n\kappa}(r) \end{Bmatrix} = \beta_{n\kappa} \frac{(p_{n\kappa} r)^{k-1}}{(2k-1)!!} \sum_{j=0}^{\infty} \begin{Bmatrix} a_j \\ b_j \end{Bmatrix} r^j \quad (15)$$

$$V_{\text{eff}} = \sum_{m=0}^{\infty} v_m r^m \quad (16)$$

$$V(r) \equiv V_{\text{DHFS}}(r) = V_{\text{nuc}}(r) + V_{\text{el}}(r) + V_{\text{ex}}(r) \quad (17)$$

A Fermi distribution for proton density is used to compute the nuclear potential.

$$\rho_p(r) = \frac{\rho_0}{1 + e^{(r-R_n)/a}}, \quad V_{\text{nuc}}(r) = -\alpha \int \frac{\rho_p(r')}{|r-r'|} dr' \quad (18)$$

The electronic potential is computed from the electron density

$$\rho(r) = \sum_{n\kappa} \psi_{n\kappa}^\dagger(r) \psi_{n\kappa}(r), \quad V_{\text{el}}(r) = \alpha \int \frac{\rho(r')}{|r-r'|} dr' \quad (19)$$

The exchange potential is

$$V_{\text{ex}}(r) = \begin{cases} -\frac{3}{2}\alpha \left(\frac{3}{\pi}\right)^{1/3} [\rho(r)]^{1/3} & r < r_{\text{Latter}} \\ -\frac{\alpha(Z-N+1)}{r} - V_{\text{nuc}}(r) - V_{\text{el}}(r) & r \geq r_{\text{Latter}} \end{cases} \quad (20)$$

DHFS Potential

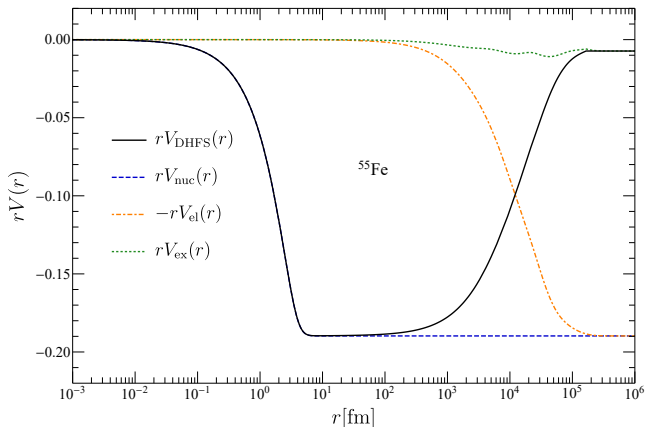


Figure: The DHFS potential (black) for neutral atom ^{55}Fe in ground state electronic configuration multiplied by the radius, along with its components.[4]

Binding energies

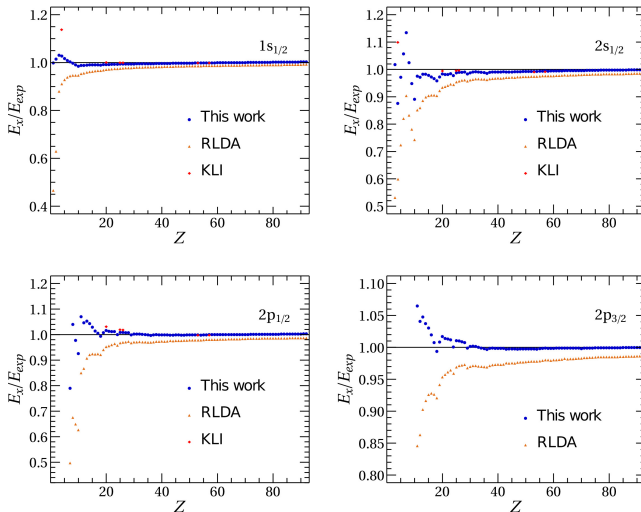


Figure: Comparison of theoretical binding energies with experimental values. [4]

The exchange and overlap correction

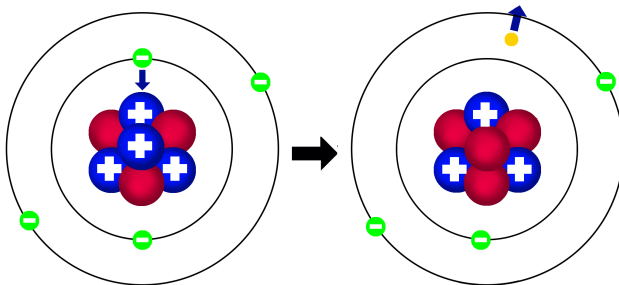


Figure: Electron Capture WITHOUT correction

The exchange and overlap correction

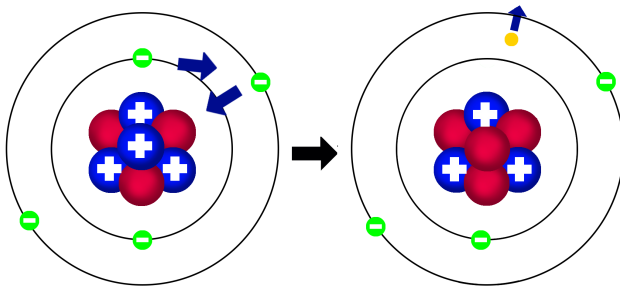


Figure: Electron Capture WITH correction

The exchange and overlap correction factor

The exchange and overlap correction takes into accounts the exchange between the captured electron and the measured position of the hole, and the imperfect overlap between initial and final wave function of the spectator electrons.

$$B_{n\kappa} = \left| \frac{b_{n\kappa}}{\beta_{n\kappa}} \right|^2 \quad (21)$$

$$b_{n\kappa} = \left[\prod_{m,\mu} \langle (m, \mu)' | (m, \mu) \rangle^{n_{m\mu}} \right] \langle (n, \kappa)' | (n, \kappa) \rangle^{-\frac{1}{2|\kappa|}} \\ \times \left[\beta_{n\kappa} - \sum_{m \neq n} \beta_{m\kappa} \frac{\langle (m, \kappa)' | (n, \kappa) \rangle}{\langle (m, \kappa)' | (m, \kappa) \rangle} \right] \quad (22)$$

$$B_{1-1} \approx 0.92 - 0.99; B_{n\kappa} \approx 0.8 - 2.1$$

Shake-up and shake-off corrections

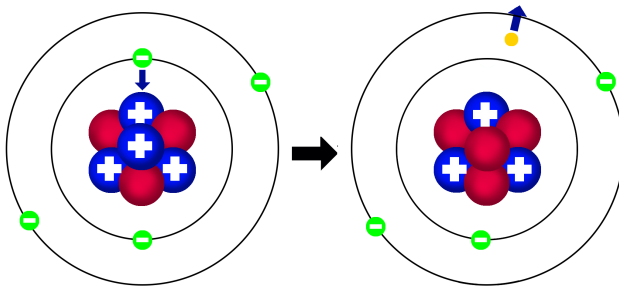


Figure: Electron Capture WITHOUT corrections

Shake-up and shake-off corrections

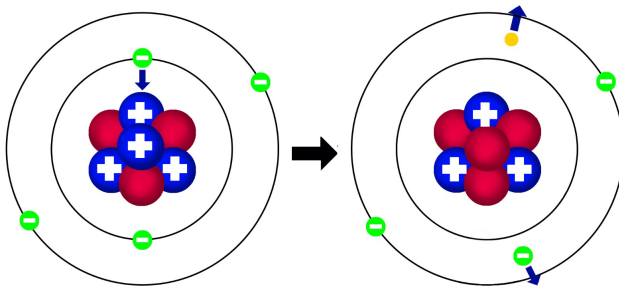


Figure: Electron Capture WITH shake-up and shake-off corrections

Shake-up and shake-off corrections

The shake-up and shake-off corrections takes into account the probability that an spectator electron is promoted to a upper vacant shell or is ejected in continuous.

$$S_{n\kappa} = 1 + \sum_{m,\mu} P_{m\mu} \quad (23)$$

$$P_{m\mu} = 1 - |\langle (m, \mu)' | (m, \mu) \rangle|^{2n_{m\mu}} - \sum_{l \neq m} n'_{l\mu} n_{m\mu} |\langle (l, \mu)' | (m, \mu) \rangle|^2 \quad (24)$$

$$S_{1-1} \approx 1 - 1.05; S_{n\kappa} \approx 1 - 1.21$$

Energy conservation

Approximate formula:

$$M_i + m_e - |E_{n\kappa}| = M_f + R_\gamma + q_{n\kappa} \quad (25)$$

Refined formula:

$$\mathcal{M}_{gs}(A, Z) = \mathcal{M}_{n\kappa}(A, Z - 1) + R_\gamma + q_{n\kappa} \quad (26)$$

$$\mathcal{M}_{gs}(A, Z) - \mathcal{M}_{gs}(A, Z - 1) = \mathcal{M}_{n\kappa}(A, Z - 1) - \mathcal{M}_{gs}(A, Z - 1) + R_\gamma + q_{n\kappa} \quad (27)$$

$$\mathcal{M}_{gs}(A, Z) - \mathcal{M}_{gs}(A, Z - 1) = |B_{gs}(Z - 1)| - |B_{n\kappa}(Z - 1)| + R_\gamma + q_{n\kappa} \quad (28)$$

$$R_{n\kappa} = |B_{gs}(Z - 1)| - |B_{n\kappa}(Z - 1)| - \text{Released energy after EC} \quad (29)$$

$$\lambda_{n\kappa} \propto q_{n\kappa}^{2(L-k+1)} \text{ (for unique transitions)} \quad (30)$$

Framework published in [4]

Atomic relaxation energy

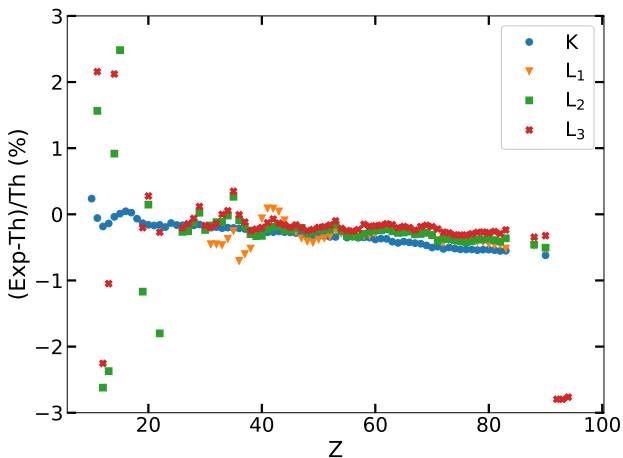


Figure: Comparison between theoretical and experimental atomic relaxation energy. [4]

Energy conservation

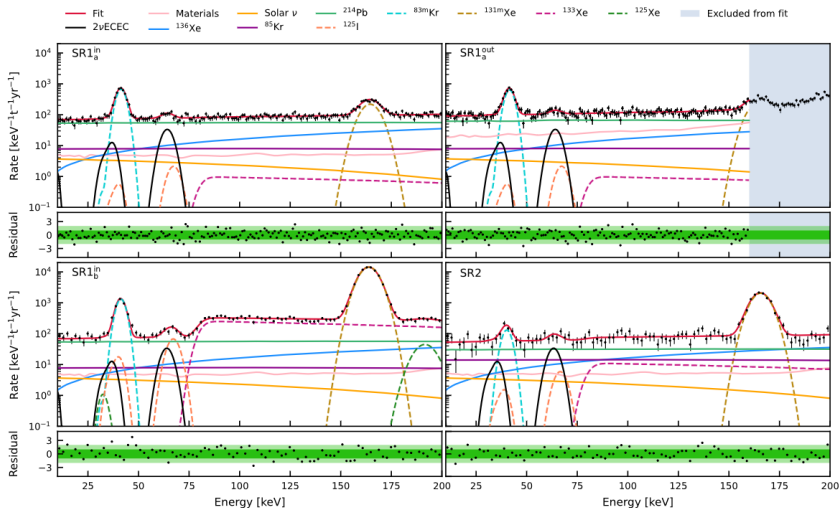


Figure: XENON1T Collaboration experimental results [2]

Comparison between formulas - K shell

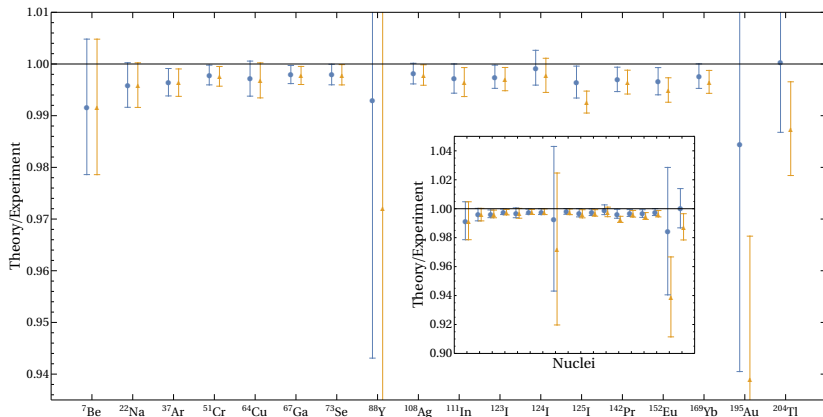


Figure: The ratio between the theoretical and experimental values for relative capture probability from the *K* shell considering both the approximate (orange triangle) and the refined (blue circle) formulas for energy conservation. [4]

Comparison between formulas - L shell

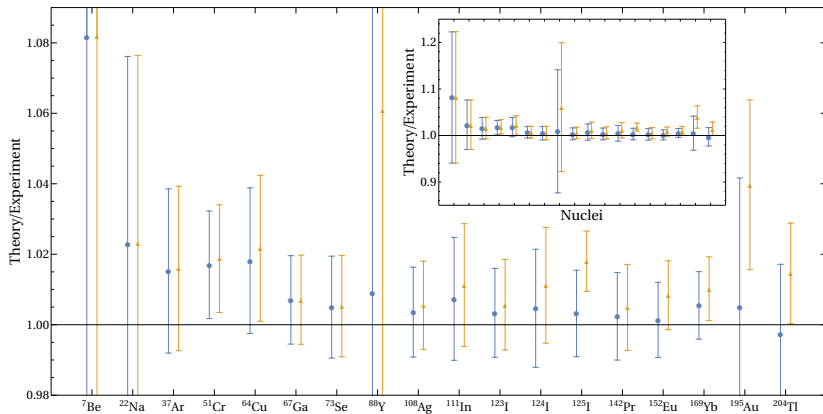


Figure: The ratio between the theoretical and experimental values for relative capture probability from the L shell considering both the approximate (orange triangle) and the refined (blue circle) formulas for energy conservation. [4]

EC ratios

Isotope	Q (keV)	R_γ (keV)	Type	Quantity	OW	RD
^7Be	861.89(7)	0	Allowed	λ_K/λ	0.90047(2)	0.908(12)
				λ_L/λ	0.09952(2)	0.092(12)
	861.89(7)	477.612(3)	Allowed	λ_K/λ	0.90046(5)	0.908(12)
				λ_L/λ	0.09954(5)	0.092(12)
^{64}Cu	1674.62(21)	1345.777(23)	Allowed	λ_K/λ	0.8815(2)	0.884(3)
				λ_L/λ	0.1008(2)	0.099(2)
				λ_M/λ	0.01714(3)	0.0162(5)
^{124}I	3159.6(19)	2335.03(1)	2nd UF	λ_K/λ	0.82040(273)	0.82099(43)
				λ_L/λ	0.14024(233)	0.13959(19)
				λ_M/λ	0.03099(54)	0.03135(15)
	3159.6(19)	2483.362(13)	1st UF	λ_K/λ	0.83148(215)	0.83184(41)
				λ_L/λ	0.13175(183)	0.13127(18)
				λ_M/λ	0.02891(43)	0.02928(14)
^{195}Au	226.8(10)	98.880(2)	1st UF	λ_K/λ	0.445(19)	0.452(6)
				λ_L/λ	0.400(14)	0.398(4)
				λ_{M+}/λ	0.1552(53)	0.1499(18)

Table: Comparison of our theoretical predictions and measured experimental data for electron capture ratios. This table is taken from [4].

L-shell dominance

Isotope	$Q(\text{keV})$ [5]	$R_{\gamma}(\text{keV})$ [6]		λ_K/λ	λ_L/λ	λ_M/λ	λ_{M+}/λ
^{127}Xe	662.33(546)	618.31(13)	OW	0.303(179)	0.520(125)	0.137(39)	0.177(50)
^{127}Xe	662.33(546)	618.31(13)	RD	0.310(80)	0.523(44)	0.137(12)	-
^{207}Bi	2397.41(266)	2339.92(0.8)	OW	0	0.638(8)	0.260(5)	0.362(7)
^{207}Bi	2397.41(266)	2339.92(0.8)	RD	0	0.651(6)	-	0.349(6)
^{236}Np	933.51(5043)	848.3(10)	OW	0	0.651(133)	0.245(79)	0.349(128)
^{236}Np	933.51(5043)	848.3(10)	RD	0	0.6(-)	-	0.4(-)

Table: Comparison of theoretical predictions and experimental results of electron capture probabilities for selected isotopes and transitions. This table is taken from [7].

EC in nuclear medicine

Isotope	$Q(\text{keV})[5]$	$R_\gamma(\text{keV})[6]$		λ_K/λ	λ_L/λ	$\lambda_M/\lambda \times 10^2$	$\lambda_N/\lambda \times 10^3$	$\lambda_O/\lambda \times 10^3$
^{67}Ga	1001.2(11)	0	This work	0.8818(3)	0.0995(3)	1.715(5)	1.504(5)	0
			RD	0.8836(15)	0.0989(12)	1.640(40)	-	-
^{67}Ga	1001.2(11)	93.312(5)	This work	0.8816(3)	0.0997(3)	1.718(6)	1.507(5)	0
			RD	0.8834(15)	0.0991(12)	1.640(40)	-	-
^{67}Ga	1001.2(11)	184.579(6)	This work	0.8814(4)	0.1000(3)	1.722(6)	1.510(6)	0
			RD	0.8832(15)	0.0993(12)	1.640(40)	-	-
^{67}Ga	1001.2(11)	393.531(7)	This work	0.8806(5)	0.1005(4)	1.734(8)	1.521(8)	0
			RD	0.8824(15)	0.0999(12)	1.650(40)	-	-
^{111}In	860(3)	416.72(3)	This work	0.8494(24)	0.11922(206)	2.520(46)	5.532(101)	0.609(11)
			RD	0.8518(2)	0.11835(13)	-	-	-
^{123}I	1228(3)	158.994(22)	This work	0.8513(10)	0.1166(8)	2.516(2)	5.913(45)	0.942(7)
			RD	0.8533(14)	0.1163(10)	2.48(5)	-	-
^{123}I	1228(3)	440.00(4)	This work	0.8489(13)	0.1185(11)	2.563(26)	6.027(61)	0.960(10)
			RD	0.8510(14)	0.1181(10)	2.53(5)	-	-
^{123}I	1228(3)	489.78(5)	This work	0.8482(14)	0.1190(12)	2.580(28)	6.057(66)	0.965(11)
			RD	0.8503(14)	0.1186(10)	2.54(5)	-	-
^{123}I	1228(3)	505.35(4)	This work	0.8480(14)	0.1191(13)	2.580(28)	6.067(68)	0.966(11)
			RD	0.8501(14)	0.1187(10)	2.54(5)	-	-
^{123}I	1228(3)	687.97(3)	This work	0.8444(20)	0.1219(17)	2.649(40)	6.233(94)	0.993(15)
			RD	0.8464(14)	0.1216(10)	2.62(5)	-	-
^{123}I	1228(3)	783.62(3)	This work	0.8412(25)	0.1243(21)	2.709(49)	6.380(117)	1.017(19)
			RD	0.8436(14)	0.1237(10)	2.67(5)	-	-

Table: Comparison theoretical and experimental values of relative capture probabilities for nuclei of interest in nuclear medicine. This table is taken from [4]

EC as a tool for neutrino mass determination

We consider $m_\nu \neq 0$ eV and take into account Γ_x the Breit-Wigner resonances centered at R_x for allowed and UF transitions:

$$\rho(E) = \frac{G_\beta^2}{(2\pi)^2} \sum_x N_x C_x \beta_x^2 B_x S_x q_x E_{\nu x} \frac{\Gamma_x/(2\pi)}{(E - R_x)^2 + \Gamma_x^2/4} \quad (31)$$

with

$$R_x = |B_{gs}(Z - 1)| - |B_x(Z - 1)| \quad (32)$$

Normalized distributions of released energy for ^{95}Tc

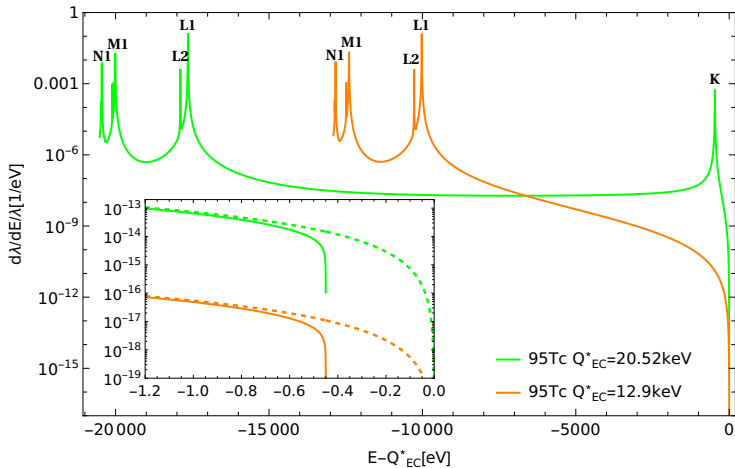


Figure: Normalized distributions of released energy in the EC decay of ^{95}Tc in the transitions to the two excited states of ^{95}Mo [8]

Comparison of ^{95}Tc with ^{163}Ho

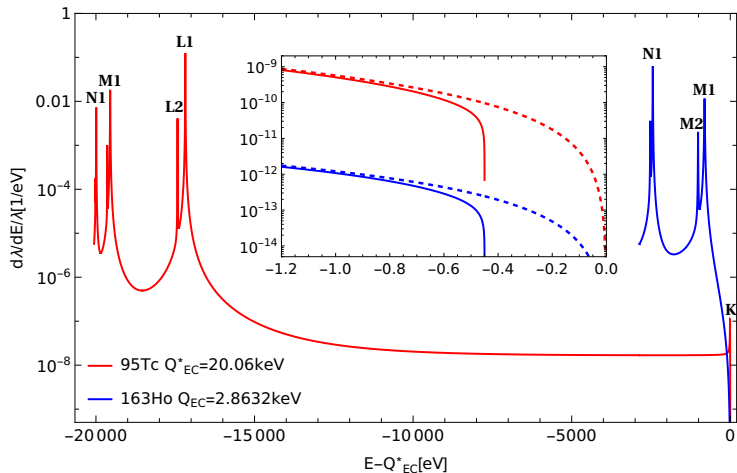


Figure: Normalized distributions of released energy in the EC decay of ^{95}Tc with Q_{EC}^* within 1σ range; and of ^{163}Ho [8]

Comparison of ^{97}Tc with ^{163}Ho

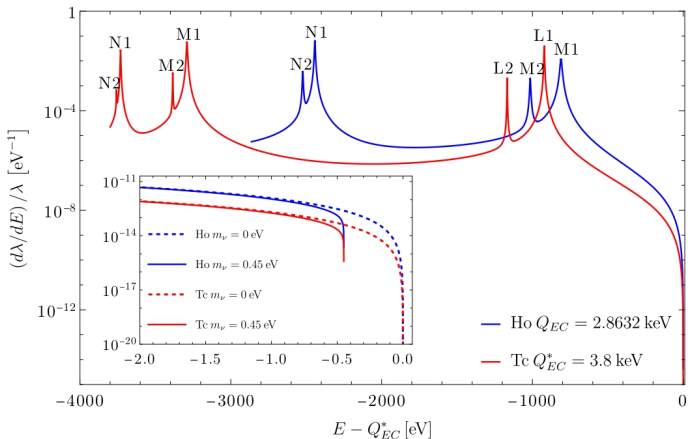


Figure: Normalized distributions of released energy in the EC decay of ^{97}Tc , assuming allowed transition type in comparison with ^{163}Ho . [9]

Statistical approach for ^{97}Tc

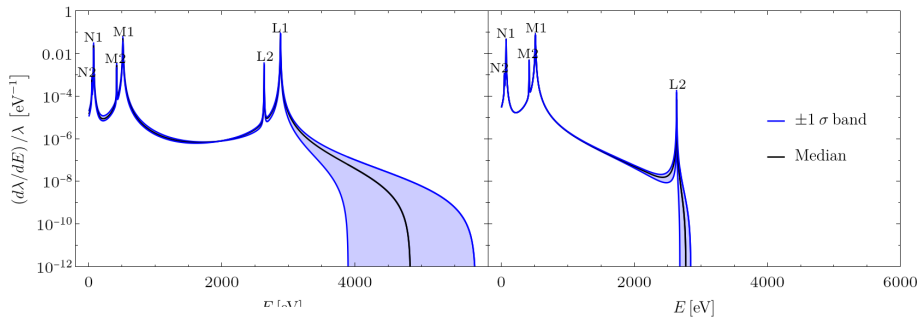


Figure: Normalized distributions of released energy as a function of E in the statistical approach for ^{97}Tc with Q_{EC}^* of 4.8(10) keV.[9]

Comparison of ^{113}Sn with ^{163}Ho

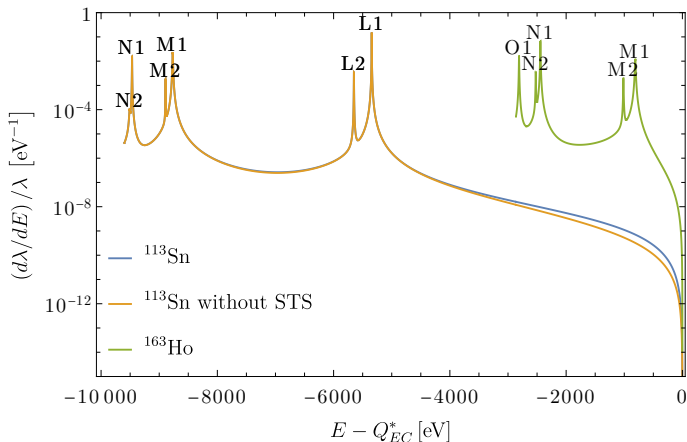


Figure: Normalized distributions of released energy in the EC decay of ^{113}Sn , as functions of $E - Q_{EC}^*$, assuming allowed transition type in comparison with ^{163}Ho .

Comparison of ^{113}Sn with ^{163}Ho - ZOOM

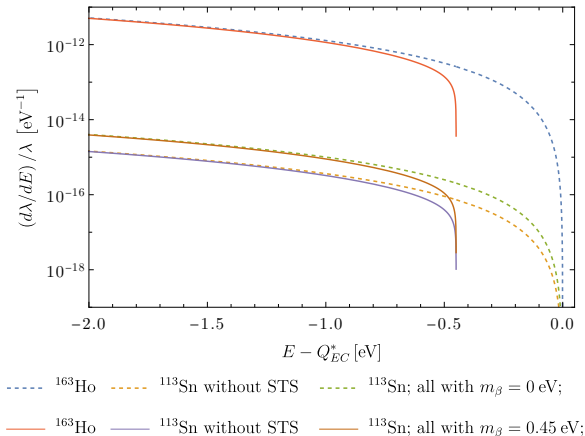


Figure: Zoom of the endpoint for the normalized distributions of released energy in the EC decay of ^{113}Sn in comparison with ^{163}Ho .

To be published...

Comparison of ^{148}Eu with ^{163}Ho

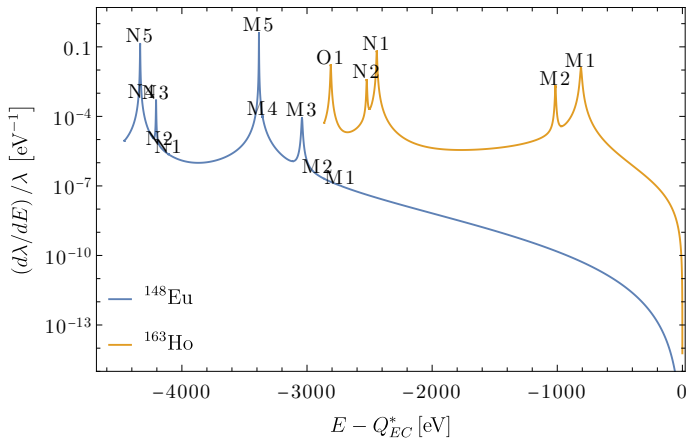


Figure: Normalized distributions of released energy in the EC decay of ^{148}Eu - 3rd UF, as functions of $E - Q_{EC}^*$, in comparison with ^{163}Ho .

Work in progress

Half-life time predictions

Table: Theoretical predictions for partial half-life time for the EC decay of ^{95}Tc . [8]

Q_{EC}^* (keV)	interaction	M_{GT}	Total (10^3yr)	K (10^7yr)	L1 (10^4yr)	L2 (10^5yr)	M1 (10^4yr)
20.52(61)	jj45pna	-0.00696	877	20.0	117	708	640
	jj45pnb	-0.016533	155	3.55	20.8	125	81.6
	glekpn	0.0070667	850	19.4	114	587	446
12.9(10)	jj45pna	-0.02412	173	-	24.1	143	79.7
	jj45pnb	-0.0198667	255	-	35.6	210	117
	glekpn	0.2296	1.9	-	0.266	1.57	0.880

Q_{EC}^* (keV)	Final J	Total (yr)	L1 (yr)	L2 (yr)	L3 (10^9yr)	M1 (yr)	M2 (yr)
9.60	3/2	33.56	59.60	2327.62	-	105.54	4158.28
9.60	1/2	238.08	422.75	16510.2	-	748.59	29495.3
14.97	5/2	4.04×10^9	7.15×10^{11}	2.79×10^{13}	5.85	1.12×10^{12}	4.40×10^{13}

Table: Computed half-lives for the electron capture decay of ^{113}Sn to the excited states of ^{113}In . The shell-model interaction used is jj45pnb. To be published.

Half-life time predictions

Table: Theoretical predictions for ^{97}Tc assuming L1 energetically available [9]

Type	Total	L1	L2	M1	M2	N1	N2	O1	
	(10^4)	(10^4)	(10^7)	(10^4)	(10^6)	(10^5)	(10^7)	(10^6)	
allowed	$19^{+17}_{-0.7}$	46^{+120}_{-26}	$2.2^{+5.0}_{-1.2}$	47^{+32}_{-17}	26^{+18}_{-10}	20^{+12}_{-7}	13^{+8}_{-4}	42^{+25}_{-15}	
2nd	Total	L1	L2	L3	M1	M2	M3	M4	M5
	(10^{16})	(10^{23})	(10^{24})	(10^{18})	(10^{21})	(10^{23})	(10^{18})	(10^{20})	(10^{16})
UF	$11^{+6.1}_{-3.4}$	$1.0^{+50}_{-0.92}$	$3.1^{+84}_{-2.7}$	$5.3^{+38}_{-4.0}$	$4.4^{+15}_{-3.1}$	$2.3^{+7.2}_{-1.6}$	$2.0^{+3.1}_{-1.1}$	$11^{+17}_{-6.2}$	$11^{+6.2}_{-3.4}$
4th	Total	L1	L2	L3	M1	M2	M3	M4	M5
	(10^{21})	(10^{28})	(10^{30})	(10^{24})	(10^{26})	(10^{28})	(10^{23})	(10^{25})	(10^{21})
NUF	$10^{+10}_{-3.6}$	$5.7^{+1066}_{-5.5}$	$1.4^{+116}_{-1.3}$	$2.6^{+58}_{-2.3}$	$5.0^{+31}_{-4.0}$	$2.5^{+1.4}_{-2.0}$	$2.6^{+8.4}_{-1.8}$	$14^{+42}_{-1.0}$	$7.2^{+11}_{-3.9}$

Table: Theoretical predictions for ^{97}Tc assuming L1 energetically unavailable [9]

Type	Total	L2	M1	M2	N1	N2	O1	
	(10^4)	(10^9)	(10^5)	(10^6)	(10^5)	(10^7)	(10^7)	
allowed	93^{+7}_{-6}	$5.7^{+40}_{-3.4}$	$16.9^{+5}_{-3.4}$	91^{+27}_{-18}	60^{+18}_{-12}	38^{+11}_{-8}	$12.7^{+4}_{-2.5}$	
2nd	Total	L2	L3	M1	M2	M3	M4	M5
	(10^{16})	(10^{31})	(10^{22})	(10^{21})	(10^{23})	(10^{18})	(10^{21})	(10^{16})
UF	$33^{+2.5}_{-1.9}$	$4.9^{+2326}_{-4.5}$	$4.1^{+21}_{-2.7}$	215^{+59}_{-38}	98^{+26}_{-17}	$24^{+4.0}_{-2.8}$	$12^{+1.9}_{-1.4}$	$35^{+2.6}_{-2.0}$
4th	Total	L2	L3	M1	M2	M3	M4	M5
	(10^{21})	(10^{39})	(10^{30})	(10^{28})	(10^{29})	(10^{23})	(10^{26})	(10^{21})
NUF	$74^{+11}_{-8.3}$	$6.1^{+25320}_{-5.9}$	$1.9^{+26}_{-1.5}$	$9^{+3.4}_{-2.1}$	$38^{+14}_{-8.7}$	110^{+29}_{-19}	$51^{+12}_{-8.4}$	$78^{+12}_{-8.8}$

Conclusion and outlook

- Relative capture probability, half-life times and other intermediate quantities were calculated
- We considered the exchange and overlap corrections and the shake-up and shake-off corrections
- We improved the way the energy conservation was taken into account
- We implemented the error propagation through pseudo-experiments
- We plotted the distribution of released energy for both ^{95}Tc , ^{97}Tc , and ^{113}Sn in comparison with ^{163}Ho



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