

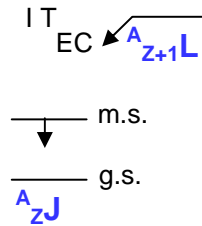
**Branch Codes IND, CUM, M+ and M-**  
(N. Otsuka, 2013-04-15, CP-D/783 Rev.)

The usages of branch codes to indicate contributions of precursors (IND, CUM, M+, M-) are explained in LEXFOR “Independent and Cumulative Data”.

When a nuclide <sup>A</sup>J has an isomer, the cross section derived by observation of the ground state of nuclide <sup>A</sup>gJ (e.g., detection of β delayed gamma from <sup>A</sup>gJ) is

$$\sigma_{\text{obs}}(^A\text{gJ}) = \sigma_{\text{dir}}(^A\text{gJ}) + a \sigma_{\text{dir}}(^A\text{mJ}) + \sum_i b_i \sigma_{\text{pre},i}$$

where “dir” is for the cross section for direct production of the nuclide without decay, and “pre,*i*” is for contribution of the precursor *i* which may decay to <sup>A</sup>gJ. The figure shows a β<sup>+</sup>+EC decay precursor <sup>A</sup>L contributes to the 3rd term. For simplicity we assume there is only one precursor, and



$$\sigma_{\text{obs}}(^A\text{gJ}) = \sigma_{\text{dir}}(^A\text{gJ}) + a \sigma_{\text{dir}}(^A\text{mJ}) + b \sigma_{\text{pre}}$$

The coefficient *a* and *b* may depend on decay data (*T*<sub>1/2</sub>, IT transition probability etc.) as well as measurement condition like cooling time. Combinations of SF4 and SF5 are summarized below for various conditions for *a* and *b*:

	<i>a</i> =0	0< <i>a</i> <1	<i>a</i> =1	<i>a</i> =?	No <i>a</i>
<i>b</i> =0	Z-S-A-G, IND	Z-S-A-G, IND/M+	Z-S-A, IND	Z-S-A-G, (M)	Z-S-A, IND
0< <i>b</i> ≤1	Z-S-A-G, CUM/M-	Z-S-A-G, CUM	Z-S-A, CUM	Z-S-A-G, CUM	Z-S-A, CUM
<i>b</i> =?	Z-S-A-G, (CUM)/M-	Z-S-A-G, (CUM)/M+	Z-S-A, (CUM)	Z-S-A-G, (CUM)	Z-S-A, (CUM)
No <i>b</i>	Z-S-A-G	Z-S-A-G, M+	Z-S-A	Z-S-A-G, (M)	Z-S-A

=0: Contribution is excluded (e.g., subtraction of its contribution, short cooling time, decay-curve analysis, on-line separation etc.)

=1: Full contribution is expected (e.g., 100% branching ratio to <sup>A</sup>gJ, long cooling time, *T*<sub>1/2</sub>(<sup>A</sup>gJ) >> *T*<sub>1/2</sub>(<sup>A</sup>mJ) or *T*<sub>1/2</sub>(<sup>A</sup>L) etc.)

No: <sup>A</sup>J is physical shielded from <sup>A</sup>mJ or <sup>A</sup>L (e.g., stable state, no decay branch to <sup>A</sup>gJ, *T*<sub>1/2</sub>(<sup>A</sup>gJ) << *T*<sub>1/2</sub>(<sup>A</sup>mJ) or *T*<sub>1/2</sub>(<sup>A</sup>L)).

**Proposed corrections to LEXFOR “Independent and Cumulative Data”**

1. Addition of the above table.
2. Corrections to the explanation of (CUM) and (M).  
The current LEXFOR explanation is inconsistent with dictionary 31, and I would believe the parenthesized codes will be more useful if we define (CUM) and (M) as above. The corresponding expansions in dictionary 236 should be corrected so that they are consistent with the expansions in dictionary 31.

<b>Branch Code</b>	<b>Definition</b>
...	...
(CUM)	<i>Uncertain if the formation via radioactive decay and isomeric transition are included.</i>
...	...
(M)	<i>Uncertain if the formation by partial feeding via isomeric transition is included.</i>