## 5. X-ray Energies and Intensities

Tables 7a, 7b, 7c, and 7d list energies and intensities for x-rays with intensities greater than 0.001 per 100 primary vacancies in the K,  $L_1$ ,  $L_2$ , and  $L_3$  atomic shells, respectively. The first column shows the Siegbahn notations for the x-ray transitions (the associations with initial and final atomic-shell vacancies are given in Table 6). The following columns give, for each element, the x-ray energies in keV (boldface) rounded to the nearest eV, and their corresponding intensities directly below. Intensities for the L x-rays are totals from both primary and secondary atomic-shell vacancies.

X-ray energies have been determined from differences between the corresponding atomic-shell binding energies reported by Larkins. Energies of complex x-ray transitions, e.g.,  $L_{\beta_{2,15}}$ , are unweighted averages of those for the single-line components.

X-ray intensities have been determined from the experimental relative emission probabilities of Salem, *et al*,<sup>2</sup> and the atomic yields of Krause.<sup>3</sup> The theoretical emission probabilities of Scofield<sup>4</sup> were occasionally used whenever experimental values were not available.

The relative intensities of x-rays from the same initial atomic shells are independent of the processes creating the shell vacancies. Tables 7a-7d may, therefore, be used to separate experimentally unresolved or complex x-ray intensities from the photon tables of the *Table of Isotopes*. Table 5 shows the initial atomic shells and their associated x-rays, and the procedure below illustrates the separation of an x-ray peak.

 $\begin{array}{|c|c|c|c|c|} \hline \text{Atomic Shell} & \text{Associated x-rays} \\ \hline & K & K_{\alpha_{1}}, K_{\alpha_{2}}, K_{\alpha_{3}}, K_{\beta_{1}}, K_{\beta_{2}}, K_{\beta_{3}}, K_{\beta_{4}}, K_{\beta_{5}}, KO_{2,3}, KP_{2,3} \\ \hline & L_{1} & L_{\beta_{3}}, L_{\beta_{4}}, L_{\gamma_{2}}, L_{\gamma_{3}} \\ \hline & L_{2} & L_{\beta_{1}}, L_{\eta}, L_{\gamma_{1}}, L_{\gamma_{6}} \\ \hline & L_{3} & L_{\alpha_{1}}, L_{\alpha_{2}}, L_{\beta_{2,15}}, L_{\beta_{5}}, L_{\beta_{6}}, L_{I} \\ \hline \end{array}$ 

Table 5

The single-line x-ray intensity of a specific transition *i* from an initial atomic shell *j* is

$$I(ji) = \frac{I}{I^{\circ}} I^{\circ}(ji) \tag{1}$$

where I is the measured (or photon-table) intensity value of a single or complex x-ray transition from atomic-shell j,  $I^0$  is the intensity of the same x-ray transition from Tables 7a-7d, and  $I(ji)^0$  is the intensity of the specific i x-ray transition from atomic-shell j, also from Tables 7a-7d. As an example, the uranium  $K_{\beta_1}$  intensity per 100 disintegrations of  $^{235}$ Np is $^5$ 

$$I(K_{\beta_{1}}) = \frac{I(K_{\alpha_{1}})}{I(K_{\alpha_{1}}^{0})} I(K_{\beta_{1}}^{0}) = \frac{0.957}{45.1} 10.70 = 0.227\%.$$
 (2)

 $I(K_{\alpha_1}^0)$  is from the photons table for <sup>235</sup>Np, and  $I(K_{\alpha_1}^0)$ , and  $I(K_{\beta_1}^0)$  are from Table 7a. Calculations for the L<sub>1</sub> atomic shell may be more complex, because none of the x-ray transitions in the photon tables of reference 5 is associated exclusively with this shell.

<sup>&</sup>lt;sup>1</sup>F.B. Larkins, At. Data and Nucl. Data Tables 20, 313 (1977).

<sup>&</sup>lt;sup>2</sup>S.I. Salem, S.L. Panossian, and R.A. Krause, *Atomic Data and Nucl. Data Tables* 14, 91 (1974).

<sup>&</sup>lt;sup>3</sup>M.O. Krause, J. Phys. Chem. Ref. Data 8, 307 (1979).

<sup>&</sup>lt;sup>4</sup>J.H. Scofield, Atomic Data and Nucl. Data Tables 14, 121 (1974).

<sup>&</sup>lt;sup>5</sup> E. Browne and R.B. Firestone, Table of Radioactive Isotopes, John Wiley & Sons, Inc. (1986).

**Table 6. Notations for X-ray Transitions** 

Classical designation (Siegbahn notation)	Associated initial - final shell vacancies
$\begin{array}{c} K_{\alpha_{1}} \\ K_{\alpha_{2}} \\ K_{\alpha_{3}} \\ K_{\beta_{1}} \\ K_{\beta_{2}} \\ K_{\beta_{3}} \\ K_{\beta_{5}} \\ K_{\beta_{5}} \\ K_{\beta_{5}} \\ K_{\beta_{2,3}} \\ L_{\alpha_{1}} \\ L_{\alpha_{2}} \\ L_{\beta_{1}} \\ L_{\beta_{5}} \\ L_{\beta_{5}} \\ L_{\beta_{5}} \\ L_{\beta_{5}} \\ L_{\beta_{5}} \\ L_{\gamma_{1}} \\ L_{\gamma_{2}} \\ L_{\gamma_{3}} \\ L_{\gamma_{2}} \\ L_{\gamma_{3}} \\ L_{\gamma_{2}} \\ L_{\gamma_{3}} \\ L_{\gamma_{5}} \\ L_{\gamma_{1}} \\ L_{\gamma_{2}} \\ L_{\gamma_{3}} \\ L_{\gamma_{5}} \\ L_{\gamma_{5}}$	K - L <sub>3</sub> K - L <sub>2</sub> K - L <sub>1</sub> K - M <sub>3</sub> K - N <sub>2</sub> N <sub>3</sub> K - N <sub>4</sub> N <sub>5</sub> K - M <sub>4</sub> M <sub>5</sub> K - O <sub>2</sub> O <sub>3</sub> K - P <sub>2</sub> P <sub>3</sub> L <sub>3</sub> - M <sub>4</sub> L <sub>2</sub> - M <sub>4</sub> L <sub>3</sub> - N <sub>4</sub> N <sub>5</sub> L <sub>1</sub> - M <sub>3</sub> L <sub>1</sub> - M <sub>2</sub> L <sub>3</sub> - O <sub>4</sub> O <sub>5</sub> L <sub>3</sub> - N <sub>4</sub> L <sub>1</sub> - N <sub>2</sub> L <sub>1</sub> - N <sub>3</sub> L <sub>1</sub> - N <sub>2</sub> L <sub>1</sub> - N <sub>3</sub> L <sub>2</sub> - O <sub>4</sub> L <sub>2</sub> - M <sub>4</sub> L <sub>3</sub> - M <sub>4</sub> L <sub>1</sub> - N <sub>2</sub> L <sub>1</sub> - N <sub>3</sub> L <sub>2</sub> - O <sub>4</sub> L <sub>3</sub> - M <sub>1</sub> L <sub>3</sub> - M <sub>1</sub>
Group designation	Associated transitions
$egin{array}{c} K^{'}_{eta_1} \ K^{'}_{eta_2} \ L_{lpha} \ L_{eta} \ L_{\gamma} \end{array}$	$\begin{array}{c} K_{\beta 1} + K_{\beta 3} + K_{\beta 5} \\ K_{\beta 2} + K_{\beta 4} + \dots \\ L_{\alpha_1} + L_{\alpha_2} \\ L_{\beta 1} + L_{\beta 2,15} + L_{\beta 3} + L_{\beta_4} + L_{\beta 5} + L_{\beta 6} \\ L_{\gamma_1} + L_{\gamma_2} + L_{\gamma_3} + L_{\gamma_6} \end{array}$