Erratum

Volume 6, 1973, in the article "Ternary Chalcogenide Compounds AB₂X₄: The Crystal Structures of SiPb₂S₄ and SiPb₂Se₄," by J. E. Iglesias and H. Steinfink, pp. 93–98:

The lattice parameters for SiPb₂S₄ given in the paper are related to the previously reported values by Hagenmuller by the transformation $\mathbf{a}_H = -\mathbf{a}$, $\mathbf{b}_H = -\mathbf{b}$, $\mathbf{c}_H = \mathbf{c} + 2\mathbf{a}$, where the subscript refers to Hagenmuller's axes.