Corrigendum

The First Crystal Structure Characterization of an Alkali Metal Monothiocarboxylate, $(PhCOSLi{\cdot}TMEDA)_2$ (TMEDA = tetramethylethylenediamine): a Chair-shaped Eight-membered $(COSLi)_2$ Ring System

David R. Armstrong, Arthur J. Banister, William Clegg, and Wendy R. Gill

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Atomic co-ordinates, bond lengths and angles, and thermal parameters have been deposited at the Cambridge Crystallographic Data Centre. See Notice to Authors, Issue No. 1, 1986.

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