Acta Crystallographica Section C
Crystal Structure Communications
ISSN 0108-2701

A three-dimensional tin(II) phosphonatobenzenesulfonate with Sn_4O_{12} clusters. Corrigendum

Palanikumar Maniam and Norbert Stock*

Institut für Anorganische Chemie, Christian-Albrechts-Universität zu Kiel, Max-Eyth-Strasse 2, 24118 Kiel, Germany
Correspondence e-mail: stock@ac.uni-kiel.de

Received 15 April 2011
Accepted 27 April 2011
Online 5 May 2011

An error in the scattering factors in the paper by Maniam & Stock [*Acta Cryst.* (2011), C**67**, m73–m76] is corrected.

In the paper by Maniam & Stock (2011), the S and P atoms were inadvertently refined using the scattering factors of the other element. A corrected CIF and structure factors are now available. Minor changes in the geometric parameters reported in the paper have occurred as a consequence, but all of the conclusions about the structure remain unchanged.

References

Maniam, P. & Stock, N. (2011). *Acta Cryst.* C**67**, m73–m76.

  doi:10.1107/S0108270111016076