

## Erratum to: Constitutive Behavior of Mixed Sn-Pb/Sn-3.0Ag-0.5Cu Solder Alloys

J.P. TUCKER,<sup>1</sup> D.K. CHAN,<sup>1</sup> G. SUBBARAYAN,<sup>1,2</sup>  
and C.A. HANDWERKER<sup>1</sup>

1.—Purdue University, West Lafayette, IN 47907, USA. 2.—e-mail: ganeshs@purdue.edu

**Erratum to: Journal of ELECTRONIC MATERIALS**  
**Vol. 41, No. 3, 2012, pp. 596–610**  
**DOI: 10.1007/s11664-011-1812-9**

In Table VI of the original article the values for the  $\zeta$  column were incorrect. The updated Table VI below lists the corrected values for  $\zeta$ . The correct values of the model constants were used for the Anand model predictions in Figs. 11–13 and 16 presented in the original article.

**Table VI. Anand model constants for Sn-3.0Ag-0.5Cu (from Ref. 3) and Sn-2.9Ag-0.5Cu-1.0Pb, Sn-2.6Ag-0.4Cu-5.0Pb, and Sn-1.4Ag-0.2Cu-20.0Pb solder alloys**

	$s_0$ (MPa)	$Q/R$ (K)	$A$ ( $s^{-1}$ )	$\zeta$	$m$	$h_0$ (MPa)	$\hat{s}$ (MPa)	$n$	$a$
Sn-3.0Ag-0.5Cu	2.15	10,000	18.0	0.35	0.15	1500	2.5	0.028	1.7
Sn-2.9Ag-0.5Cu-1.0Pb	0.21	9900	17.0	0.88	0.30	1300	2.5	0.041	1.9
Sn-2.6Ag-0.4Cu-5.0Pb	0.20	9500	10.5	0.90	0.31	850	2.5	0.044	2.0
Sn-1.4Ag-0.2Cu-20.0Pb	0.19	8500	6.0	1.0	0.37	700	2.5	0.050	2.1