Corrigenda

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Page 78, Fig. 1.A.

The experimental modified molecular intensity points have been plotted with the wrong sign from $s = 6.750$ to 8.375 $Å^{-1}$ inclusive. The experimental as well as the theoretical intensities are *positive* in this range. The difference curve, Fig. 1.B., is correct.

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In Table 3 the correct values of $|J(^{119}Sn-^{13}C)|$ for ring carbons of the two fluoro derivatives should read:

Compound	C(1)	C(2)	C(3)	C(4)	C(5)	C(6)
p-F	467.0	42.0	50.2	11.8	50.2	42.0
m-F	447.4	39.1	62.6	9.0	52.2	32.5

Regression analysis parameters remain unaffected by these changes.

Page 108, The caption to Fig. 3 should read:

Correlation of $|J(Sn-CH_3)|$ with $|J(Sn-C(1))|$ (averaged tin-117,119 values).

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