ERRATA

J. Organometal. Chem., Vol. 20, No. 1 (November 1969)

Page 118, line 3 should read:

great interest in π-bonding involving elements of the second and higher rows of the

Page 119, compound (X) in Table 1 should read:

$PhMe_2As^+-CH_2CH_2-NHPh\ I^-$ (X)

Page 120, Table 2

The parentheses should be deleted in the 2nd, 3rd and 4th column, for compound (XI).

Acridine[c] should correspond with 250 and 5.3

2nd line up should read: 355[d] 4.2

Page 126, 3rd line up should read:

methyl iodide (42 mg; 1.1 mol) in acetonitrile (2 ml) to stand at room temperature

CORRIGENDA

Re: VIBRATIONAL SPECTRA OF MESITYLENECHROMIUM TRICARBONYL AND MESITYLENEMOLYBDENUM TRICARBONYL; by G. Davidson and E.M. Riley.

J. Organometal. Chem., Vol. 19, No. 1 (September 1969) pages 101–114

The overall symmetry of the $[1,3,5\text{-}C_6H_3(CH_3)_3]M(CO)_3$ molecule will be C_{3v} in both "eclipsed" and "staggered" conformations, not C_1 for the latter as suggested. The "local symmetry" of the two halves of the molecule is unaltered, and the subsequent assignment may be left unchanged.

Re: BIMOLECULAR HOMOLYTIC SUBSTITUTION BY TERT-BUTOXY RADICALS AT METAL ATOMS; by A.G. Davies and B.P. Roberts.

J. Organometal. Chem., Vol. 19, No. 1 (September 1969) page P18

The following words should be delected: "Kochi[3] has observed by ESR the radical $R^\cdot$ formed by displacement by the butylthio radical at phosphorus in a phosphine R_3P."