

***Errata***

**Ionic Character**

N. C. BAIRD and M. A. WHITEHEAD

Theoret. chim. Acta (Berl.) 2, 259 (1964)

p. 262, Tab. 2: The values of  $E_v$ ,  $b^*$  and  $c^*$  should read:  
for  $Be^-$ : -0.95, +0.71, -0.56 resp.  
for  $B^-$  : -5.32, +4.25, -3.19 resp.

p. 263, Tab. 3: "This Paper" ionic character values for  
NaH, HCl, HBr, HI, FCl, BrCl should be  
25.0, 9.2, 5.5, 4.2, 9.7, 4.8% resp.  
The "Gordy" ionic characters of HBr, HI, BrCl are  
20.5, 15.5, 16.5% resp.

**Group Orbital Electronegativities**

M. A. WHITEHEAD, N. C. BAIRD and M. KAPLANSKY

Theoret. chim. Acta (Berl.) 3, 135 (1965)

p. 140, Eq. (4) should read:

$$c_B = 0.5 (\delta_B - \alpha_B) + 0.5 (n_T^{(1)}) (\varepsilon_B - \beta_B) + 0.5 (n_T^{(1)})^2 (\zeta_B - \gamma_B).$$

p. 142, Tab. 5: The fifth column heading is  $\varrho_{C135}$  rather than  $n_{C135}$ .

Corrected  $E_v$  parameters for C (*tetetete*), C (*tririrr*  $\pi$ ) and the improved data for benzene are listed in Tab. 1 and 6 of N. C. BAIRD and M. A. WHITEHEAD's paper: Theoret. chim. Acta 6, 167 (1966).

We regret misspelling Dr. B. P. DAILEY's name in the references.