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Annotationes et Errata

Porphyryns. IV. Extended Hückel Calculations on Transition Metal Complexes

Theoret. chim. Acta (Berl.) **4**, 44 (1966)

By

M. ZERNER and M. GOUTERMAN

Eq. (4), page 46, should read

$$H_{pq} = \langle \chi_p | H_{eff} | \chi_q \rangle = \frac{1}{2} (H_{pp} + H_{qq}) S_{pq} [\alpha + (1 - \alpha) \delta_{pq}].$$

Valency Structures for N₂O₄

Theoret. chim. Acta (Berl.) **2**, 437—452 (1964)

By

R. D. HARCOURT

p. 444 Line 17 down: Read VII for II.

p. 445 Table 1: Headings should be i, ii, iii and iv.

p. 447 Line 2 up: Read an orthogonal for orthogonal an.

p. 449 Line 15 down: Read \mathcal{O} for \mathcal{P}

p. 451 Line 10 down: Read $\left(\frac{s_2 + s_4}{\sqrt{2}}\right)^2$ for $\left(\frac{s_2 + s_4}{2}\right)^2$

Valency Structures for Oxalate and Dithionite Anions and Dimers of Nitrosoalkanes

Theoret. chim. Acta (Berl.) **3**, 194—201 (1965)

By

R. D. HARCOURT

p. 195 Table 1: first row: 0.32 for 0.32.

p. 196 Table 2: The C₂O₄²⁻ weights have been incorrectly reported — they obtain for $\mu = 0.5$ instead of 0.6. For $\mu = 0.6$, the weights reported as 0.32, 0.16 and 0.02 should read 0.28, 0.19 and 0.03. The weights 0.24, —0.06 and 0.32 become 0.19, —0.07 and 0.38. These alterations do not affect the conclusions of the paper.

p. 199 Line 23 down: Read

$$\left| \varphi_4^\alpha \left(\frac{\varphi_4 + \varphi_3}{\sqrt{2}} \right)^\beta \varphi_3^\alpha \varphi_2^\beta \left(\frac{\varphi_1 + \varphi_2}{\sqrt{2}} \right)^\alpha \varphi_1^\beta \right|$$

for

$$\left| \varphi_4^\alpha \left(\frac{\varphi_4 + \varphi_3}{\sqrt{2}} \right)^\beta \varphi_3^\alpha \varphi_3^\beta \left(\frac{1}{\sqrt{2}} \right)^\alpha \varphi_1^\beta \right|$$

p. 200 Line 2 of footnote: Read σ_4- for σ_2- .