molecular-orbital calculation (Segal, 1966; Pople & Beveridge, 1970), which gives +0·087 and −0·166 e for the charges on the C and N respectively. A more extensive experimental and theoretical comparison of the crystal energies of these two compounds would be a worthwhile undertaking.

We thank the National Science Foundation, the University of Minnesota Computer Center, and the University of Minnesota Graduate School for their support of this work.

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SHORT COMMUNICATIONS

Contributions intended for publication under this heading should be expressly so marked; they should not exceed about 1000 *words; they should be forwarded in the usual way to the appropriate Co-editor; they will be published as speedily as possible.*

Acta Cryst. (1977). B**33**, 1303

The crystal structure of 6b,10b-dihydrobenzo[*j*]cyclobut[*a*]acenaphthylene, $C_{18}H_{12}$. Corrigenda.

By A. C. HAZELL and R. G. HAZELL, *Department of Inorganic Chemistry, Aarhus University, DK*-8000 *Aarhus C, Denmark*

(*Received* 15 *February* 1977)

A printer's error in the paper by Hazell & Hazell [*Acta Cryst.* (1977), B**33**, 360–365] is corrected: In Table 2 the fractional atomic coordinates (reported as multiplied by 10E) are multiplied by 10^4. Also, in Table 3 the units of **T** should be $Å^2 \times 10^{-4}$ and those of **L**, $(°)^2$.

All the relevant information is given in the abstract.

Acta Cryst. (1977). B**33**, 1303

Bis(bis-η-cyclopentadienylmolybdenumdi-μ-oxo)phosphorus hexafluorophosphate: erratum.

By KEITH PROUT, M. CLAIRE COULDWELL and ROGER A. FORDER, *Chemical Crystallography Laboratory, Oxford University, 9 Parks Road, Oxford, OX*1 3*PD, England*

(*Received* 3 *March* 1977)

The abstract of Prout, Couldwell & Forder [*Acta Cryst.* (1977), B**33**, 218–221] contains a number of errors. The beginning of the abstract should read: **Abstract.** $[(C_5H_5)_2MoO_2PO_2Mo(C_5H_5)_2]PF_6$, $C_{20}H_{20}O_4F_6P_2Mo_2$, $M_r = 692{\cdot}2$. Monoclinic, $C2/c$, $a = 17{\cdot}079$ (9), $b = 23{\cdot}619$ (13), $c = 12{\cdot}399$ (7) Å, $\beta = 117{\cdot}02$ (2)°, $U = 4455{\cdot}7$ Å^3, $D_c = 2{\cdot}06$ g cm^{-3} for $Z = 8$. Mo $K\alpha$ radiation, $\lambda = 0{\cdot}71069$ Å, $\mu = 11$ cm^{-1}. The structure of the cation...

All information is given in the abstract.