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ERRATA

Vol. 1947, page 313, line 38. *For* b. p. 148—150°/22 mm. *read* b. p. 148—150°.

Vol. 1953, page 236, line 7*. *For* m. p. 134—185° *read* 134—135°.

Vol. 1953, page 1900, formula (III). *For* :COPh *read* :CHPh.

Vol. 1954, page 1686, lines 33 and 34. The analytical results *should read* (Found: C, 67·5; H, 5·9; N, 4·1. $C_{20}H_{21}O_5N$ requires C, 67·6; H, 6·0; N, 4·0%).

Vol. 1955, page 383. The arrows between formulæ (X) and (XII) should be reversed.

Vol. 1955, page 1700, line 17*. *For* 2·16 *read* 1·83. *For* 1·98 *read* 1·30.

Vol. 1955, page 2480, summary, line 4. *For* $1k0$, $0kl$, and $1k\bar{l}$ *read* $1k0$, $0k1$, and $1k\bar{1}$.

Vol. 1955, page 2481, line 11. *For* while $0kl$ and $1k\bar{l}$ were *read* while $0k1$ and $1k\bar{1}$ were.

Vol. 1955, page 4023, Table 2. *For* 0·90 *read* 0·090. *For* 0·82 *read* 0·082.

Vol. 1955, page 4024, 2nd line above Table 4. *For* 1·120M-sodium hydroxide *read* 1·120M-potassium hydroxide.

Vol. 1956, page 4, line 30. *For* $C_2H_{188}O_4$ *read* $C_{26}H_{18}O_4$.

Vol. 1956, page 212, line 8. *For* $\log_{10} p$(mm. Hg) = 7·035—1278/(t°C + 219·3°) *read* $\log_{10} p$(mm. Hg) = 7·145—1329/(t°C + 224·0°).

Vol. 1956, page 249, line 3*. *For* C, 60·5 *read* C, 66·5.

Vol. 1956, page 249, line 4*. *For* C, 66·5 *read* C, 60·5.

Vol. 1956, page 361, line 3*. *For* R = H *read* R′ = H.

Vol. 1956, page 531, formula (VI*A*). There should be a single bond between the central oxygen atom and the nearest position on the left-hand benzene ring.

Vol. 1956, page 672. Equation (3) *should read* $K^{-1} = 4\pi N\rho(2q_1)^3 \int_{2q_1/q}^{2q_1/a} e^y y^{-4}\,dy$.

Vol. 1956, page 700, formula (XIII). The thiophen ring should be completed.

Vol. 1956, page 893, formulæ. In the right-hand ring of the 3-isomer NH *should read* NH_2.

Vol. 1956, page 967. Second table *should read*:

$SiHCl_3$ (mmoles)	C_2F_4 (mmoles)	Mole ratio $SiHCl_3 : C_2F_4$	$CHF_2{\cdot}CF_2{\cdot}SiCl_3$ Wt. (g.)	Yield (%)	$H{\cdot}[CF_2{\cdot}CF_2]_m{\cdot}SiCl_3$ (g.)
82·5	68·5	1·2	5·4	34	4·8
110	,,	1·6	7·1	44	4·0
220	,,	3·2	9·8	61	3·1
275	,,	4·0	10·1	63	3·0

* From bottom of main text.

Errata.

Vol. 1956, page 1097, line 7. *For* monomethylhydroxylamine *read* monopropylhydroxylamine.

Vol. 1956, page 1150, line 6. *For* 1-acetoxymethyl-2-oxo*cyclo*pentanecarboxylate *read* 1-acetoxymethyl-2-oxo*cyclo*hexanecarboxylate.

Vol. 1956, page 1302, line 7. *For* $-[H^+]$ *read* $+[OH^-]$ *and for* $+[H^+]$ *read* $-[OH^-]$.

Vol. 1956, page 1554, line 35. *For* 96% *read* 86%.

Vol. 1956, page 1559, last line of text. *For* $A_{1g} \longrightarrow B_{2u}$ *read* $A_{1g} \longrightarrow B_{1u}$.

Vol. 1956, page 1818, line 6*. *For* all but the two cases *read* all but two cases.

Vol. 1956, page 1818, line 1*. *For cyclo*propane *read cyclo*propene.

Vol. 1956, page 1819, line 16. *For* w/v *read* w/w.

Vol. 1956, page 1819, line 15*. *For* C=C *read* C=O.

Vol. 1956, page 1848. On the arrow leading from (XIV) to (XV) *for* Na *read* NaI.

Vol. 1956, page 1970, line 2*. *For* $\psi_2 = \sqrt{\frac{1}{2}}(\phi_1 - \phi_3)$, $\psi_3 = \sqrt{\frac{1}{2}}(\phi_2 - \phi_4)$
read $\psi_2 = \frac{1}{2}(\phi_1 + \phi_2 - \phi_3 - \phi_4)$
$\psi_3 = \frac{1}{2}(\phi_2 + \phi_3 - \phi_4 - \phi_1)$

Vol. 1956, page 1971, figure. The four-membered ring should be rotated through 45° about the z axis in the direction x to y.

Vol. 1956, page 2000, lines 1—2 under formulæ. *For* 2 : 4-diamino-1′-*n*-propyl(2′ : 3′-6 : 7)pteridine *read* 2 : 4-diamino-1′-*n*-propylindolo(2′ : 3′-6 : 7)pteridine.

Vol. 1956, page 2004, Table 2, heading. For *et* read *at*.

Vol. 1956, page 2168. *In* Table 4, *the five entries* 5 : 6 enol *should all read* 1 : 6 enol.

Vol. 1956, page 2210, Table 4, last column. The last three entries should read *l*, *m*, *n*, respectively.

Vol. 1956, page 2701, title. *For* Akhookh *read* Akhnookh.

Vol. 1956, page 3311. *For* Part III.[1] *read* Part III.* *Footnotes should read* : * Part II, *J.*, 1956, 2131.
[1] Albert, Brown, and Wood, *J.*, 1956, 2066.

Vol. 1956, page 3311, summary, lines 4 and 5. *For* 5-amino-4-methylpyrimidines *read* 5-amino-4-methylaminopyrimidines.

Vol. 1956, page 3366, Table 3. *Column headings should read* : Cols. 4—6 E = amount (mmoles) of solute absorbed.* Cols. 7—9 G = wt. (g.) of H_2O absorbed.*

Vol. 1956, 3370, Fig. 8. The labels *A* and *C* should be interchanged.

Vol. 1956, page 3476, figure. The molecular extinction coefficients corresponding to the tie-stubs on the left of the figure should read (from top to bottom) : 50,000, 10,000, 5,000, 1000, 500, 100 (*i.e.* each stub should be numbered).

Vol. 1956, page 3661, Fig. 4. Right-hand scale should read *log* ε (*curves J* and *K*).

* From bottom of main text.

Errata.

Vol. 1956, page 3949, Fig. 2. The ordinate should read $1/\chi_M$.

Vol. 1956, page 3950, Fig. 3. The ordinate should read $1/\chi_M$.

Vol. 1956, page 4597, line 15. *For* m. p. 161—169° *read* m. p. 168—169°.

Vol. 1956, page 4598, line 7*. *For* C, 15·8 *read* C, 45·8.

Vol. 1956, page 4641, Fig. 2(*b*). The captions for the full and broken line should be interchanged.

Vol. 1956, page 4678, line 2*. *For* hydrolysis product *read* product of hydrolysis at room temperature.

* From bottom of main text.