The following addition should be made to Table III.

| Prod-uct model	f_P	Reactant model				
		R_1	R_2	R_3	R_4	R_5
		1.1708	1.1741	1.1808	1.1841	1.2121
		f_R				
P6	1.1734	0.9978	1.0006	1.0091	1.0091	1.0330
P7	1.1831	*0.9896*	0.9924	0.9980	1.0008	1.0245
P8	1.1982	*0.9771*	*0.9799*	**0.9855**	0.9882	1.0116
P9	1.2082	0.9690	0.9718	*0.9773*	*0.9800*	1.0032

In addition, the entry for P4–R3 should be italicized.—R. E. WESTON, JR.

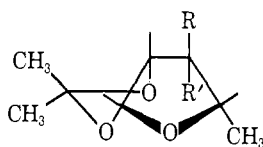**Patricia S. Traylor and F. H. Westheimer:** Mechanisms in the Hydrolysis of Phosphorodiamidic Chlorides.

Page 559. The heading of the first column of Table V should read: in H_2O^a.—F. H. WESTHEIMER.

John R. Dyer, W. E. McGonigal, and K. C. Rice: Streptomycin. II. Streptose.
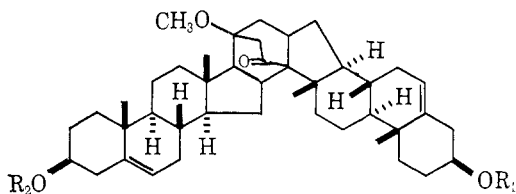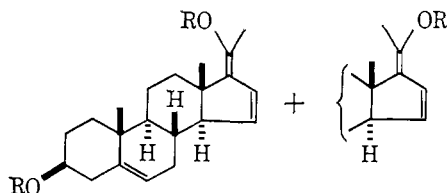
Page 654.
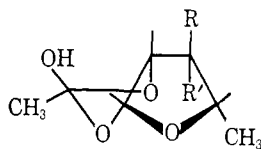
should be

JOHN R. DYER.

S. A. Latt, H. T. Cheung, and E. R. Blout: Energy Transfer. A System with Relatively Fixed Donor–Acceptor Separation.
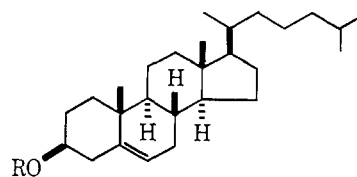
Page 997. Structures XIb, XIIb, XIII, XV, XVI, and XVII should be as shown below.

XIb, R_1 = p-methoxyphenylacetyl; R_2 = 1-naphthoyl
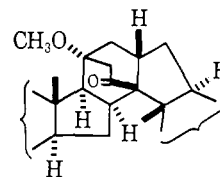XIIb, R_1 = 1-naphthoyl; R_2 = anthracene-9-carbonyl

XIII, R = p-methoxyphenylacetyl

XV, R = p-methoxyphenylacetyl
XVI, R = 1-naphthoyl
XVII, R = anthracene-9-carbonyl

Page 1002. Structure XVIII should be as shown below.

XVIII

E. R. BLOUT.

Ernest Wenkert and Börje Wickberg. General Methods of Synthesis of Indole Alkaloids. IV. A Synthesis of dl-Eburnamonine.

Page 1583. In column 1, line 10, alkalids should read alkaloids. In footnote 24 anhydroxymethylene should read a hydroxymethylene. In Chart III the indole ring of the formula under XVII should possess a radical cation sign of plus, dot instead of the minus sign shown, while the same ring in the formula under XVIII should have a plus instead of minus sign.—ERNEST WENKERT.

**Myron L. Bender and James K. Stoops:** Titration of the Active Sites of Acetylcholinesterase.

Page 1622. Equation 2 should read

$$N = \frac{F(A_1 - A_2) + A_3 - A_4 - A_2}{4.0 \times 10^8}$$

In the last paragraph, second sentence, *directly* proportional should be substituted for *inversely* proportional.—MYRON L. BENDER.

Anthony C. Waiss, Jr., and Joseph Corse: Photooxidative Cyclization of Quercetin Pentamethyl Ether.

Page 2068. Structure V should be
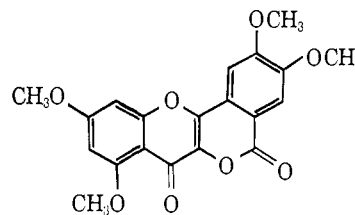
ANTHONY C. WAISS, JR.

Neville Finch, C. W. Gemenden, Iva Hsiu-Chu Hsu, Ann Kerr, G. A. Sim, and W. I. Taylor: Oxidative Transformations of Indole Alkaloids. III. Pseudoindoxyls from Yohimbinoid Alkaloids and Their Conversion to "Invert" Alkaloids.

Page 2232 ff. In the recorded data for the optical rotatory dispersion curves $[\alpha]$ read $[\phi]$.

Page 2232. In column 2, line 9, $[\alpha]_{286}$ −2850 (inflection), $[\alpha]_{272}$ −2150 (inflection) should read $[\phi]_{286-272}$ −2850 to −2150 (inflection).

Page 2235. In column 2, lines 16 and 17, −6450 and −47,000 should read +6450 and +47,000.—WILLIAM I. TAYLOR.

Dwight R. Robinson and William P. Jencks: The Effect of Concentrated Salt Solutions on the Activity Coefficient of Acetyltetraglycine Ethyl Ester.

Page 2478. In Table VIII, the third ion listed should be Cl_3-CCOO⁻, not $(CH_3)_3$CCOO⁻.—WILLIAM P. JENCKS.

B. J. Herold, A. F. Neiva Correia, and J. dos Santos Veiga: Electron Paramagnetic Resonance Studies on Chelation of Alkali Cations by the o-Dimesitoylbenzene Radical Anion.