Additions and Corrections

VOL. 26, 1961

L. H. Zalkow, E. J. Eisenbraun, and J. N. Shoolery: Structures of Carotol and Daucol.

Page 981. N.m.r. peak positions were given in this paper in dimensionless units defined as $10.00 - 10^6(\nu - \nu_{SiMe_4})/\nu_{SiMe_4}$. This definition has come to be generally designated by the symbol τ rather than the symbol δ as used in this paper, the latter usually being recognized as referring to peak positions defined as $10^6(\nu - \nu_{SiMe_4})/\nu_{SiMe_4}$. The data in this paper can be made more consistent with that appearing in other papers either by replacing δ by τ throughout, or by subtracting all chemical shift values from 10.00.

Page 982. The calculated value for the analysis of daucol acetate which appeared on this page should read as follows: "*Anal.* Calcd. for $C_{17}H_{28}O_3$: C, 72.82; H, 10.06; O, 17.12." instead of "*Anal.* Calcd. for $C_{17}H_{26}O_3$: C, 73.34; H, 9.41; O, 17.24."

Samuel C. Temin: Polymers from Bisphenols. Steric Inhibition of Condensation Polymerization.

Page 2519. In Table I, for bisphenol diacetate no. 5, R_2 should be "CH_3" and not "$C(CH_3)_3$.

Page 2519. In Table I, the melting point of bisphenol diacetate no. 3 should be "133–134°C" and not "59–60°C."

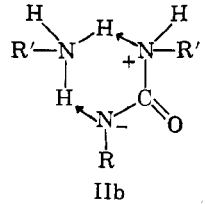
Omar W. Steward and Ogden R. Pierce: Fluoroalkyl and 3-(Fluoroalkoxy)propylpolysiloxanes.

Page 2944. In col. 2, lines 23 to 26, the absorption bands reported as C—F stretching bands probably should not be assigned to any localized vibration; however, they are characteristic of a 3,3,3-trifluoropropyl group attached to silicon. The assignment of a band to the C—Si bond is incorrect.

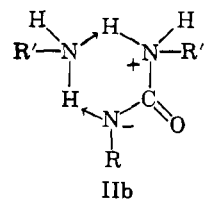
Kenneth Sauer and Marie H. Kasparian: Kinetics of the Reaction of Phenyl Isocyanate with Aniline.

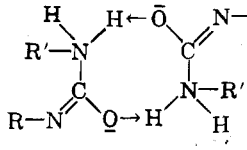
Page 3502. In col. 2 the sentence following Equation 13 should read: A consequence of Equation 13 is the fact that for $K_{-1} > k_2a$ the reaction should be third-order over all and for $k_{-1} < k_2a$ it should be second-order over all.

Page 3503. In col. 2

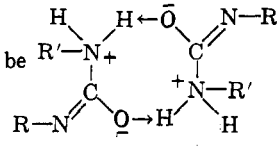


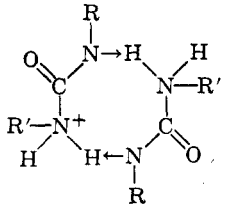
IIb should be IIb



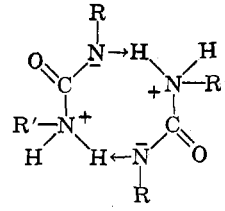


IIIa should be IIIa





IIIb should be IIIb



Konoshin Onodera and Tohru Komano: Debenzyloxycarbonylation of 1,3,4,6-Tetra-*O*-acetyl-2-benzyloxycarbonylamino-2-deoxy-D-hexopyranoses in the Conversion of α,β-Acetoxy to Glycosyl Bromide.

Page 3933. In col. 2, line 42, for "melting at 115°" read "melting at 133°."

Page 3933. In col. 2, line 43 to line 44, for "$C_{22}H_{29}O_{10}N$: C, 56.53; H, 6.21; N, 3.00. Found: C, 56.38; H, 6.27; N, 3.20" read "$C_{22}H_{27}O_{11}N$: C, 54.86; H, 5.65; N, 2.91. Found: C, 54.72; H,5.61; N, 2.93."

Gordon N. Walker and Barbara N. Weaver: Synthesis of Benzo[*f*]quinolines and Ergolines from 5-Phenyl-6-methyl-2-pyridones.

Page 4441. The structure of lysergic acid is depicted incorrectly, and should be altered by the insertion of $—NCH_3$ in the appropriate position in Ring D.

Page 4444. Structure 26 should be altered by insertion of N in ring position 4.

Page 4448. In Experimental, compound "(II. R = C_6H_5)" is named incorrectly and should be 2,4-Diphenyl-3-methylpyrazole.

Robert Barker and Hewitt G. Fletcher, Jr.: 2,3,5-Tri-*O*-benzyl-D-ribosyl and -L-arabinosyl Bromides.

Page 4608. In col. 1, line 34, for "1,5-Anhydro-D-ribitol" read "1,4-Anhydro-D-ribitol."

Elizabeth Dyer and Robert E. Read: New Catalysts for the Conversion of Isocyanates to Carbodiimides.

Page 4678. In the last paragraph, 3rd line, delete "than". That is, the sentence should read: ". . manganese naphthenate was the preferred catalyst due to the formation of fewer by-products with the naphthenate."

H. M. Walborsky, L. Barash, and T. C. Davis: Partial Asymmetric Syntheses: The Diels–Alder Reaction.

Page 4778. On lines 5, 10, 29, wherever dimethyl fumarate appears it should read *dimenthyl* fumarate.

Since the publication of our results we discovered that A. Korolev and V. Mur, *Dokl. Akad. Nauk SSSR*, **59**, 251 (1948) had reported obtaining a partial asymmetric synthesis in the Diels–Alder reaction.

Harold E. Zaugg, Robert W. De Net, and Raymond J. Michaels: Neighboring Group Reactions. II. A Novel Synthesis of Basic Esters of 1-Benzoxacycloalkanecarboxylic Acids.

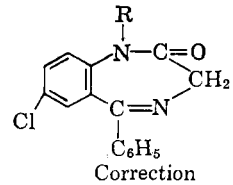
Page 4832. In col. 2, line 5 from end, for "much more nucleophilicity" read "much lower nucleophilicity."

R. K. Summerbell and Sherri R. Forrester: A New Type of Assistance at a Distance.

Page 4837. In Table I, "1-Chloromethyl-4-iodomethyl-*p*-dioxane" should read "2-Chloromethyl-3-iodomethyl-*p*-dioxane".

L. H. Sternbach and E. Reeder: Quinazolines and 1,4-Benzodiazepines. IV. Transformations of 7-Chloro-2-methylamino-5-phenyl-3H-1,4-benzodiazepine 4-Oxide.

Page 4937. Formula Xa,b,d,e should have a single bond between 1 and 2 as shown below.



Correction instead of