

## Corrigenda

### The Revised Structure of Bilobanone

By H. IRIE, H. KIMURA, N. OTANI, K. UEDA, and S. UYEO

*Chem. Comm.*, 1967, 678.

On p. 678, l.h.s., line 14: for  $[\alpha]_D + 6.7^\circ$ , read  $+67^\circ$ .  
line 22: for  $[\alpha]_D + 4.6^\circ$ , read  $+46^\circ$ .

### Photofragmentation of Oxirans. Precursors for Phenylcyanocarbene and Phenylmethoxycarbonylcarbene

By P. C. PETRELLIS and G. W. GRIFFIN

*Chem. Comm.*, 1967, 691.

The structure (III) on r.h.s. for read

### The Structure of Jesaconitine

By L. H. KEITH and S. W. PELLETIER

*Chem. Comm.*, 1967, p. 994.

On p. 994, the *upper structure* labelled (III) should be (IV).

### The Structure of Dichlorobis(pentane-2,4-dionato)rhenium(IV)

By C. J. L. LOCK and CHE'NG WAN

*Chem. Comm.*, 1967, 1109.

On p. 1109, l.h.s. line 4, for  $[\text{ReCl}_2(\text{C}_5\text{H}_7\text{O}_2)]$  read  $[\text{ReCl}_2(\text{C}_5\text{H}_7\text{O}_2)_2]_2$ .

### Photochemical Isomerization and Radical Fragmentation of 10-Hydroxymethyl- $\Delta^{1,9}$ -2-octalone

By DAVID I. SCHUSTER and DONALD F. BRIZZOLARA

*Chem. Comm.*, 1967, p. 1158.

On p. 1159, line 4, for 500 hr. read 300 hr.

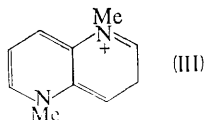
On p. 1159, Table, 2nd column labelled Time, last line, for 300 read 600.

### A Diquaternary Salt of 1,5-Naphthyridine

By L. A. SUMMERS and J. E. DICKESON

*Chem. Comm.*, 1967, 1183.

Formula (III) should be



**Epoxide Cleavage as a Means of Methyl Migration: Model Studies in Cucurbitacin  
Synthesis: Ring A Aromatic 9 $\beta$ -Methyl Steroids**

By J. W. APsIMON and R. R. KING

*Chem. Comm.*, 1967, 1214.

On p. 1215, in formula (III), delete the H attached to C-9, leave blank to represent a methyl group.

**A New General Method for converting  $\alpha\beta$ -Unsaturated Aldehydes into Saturated  
Imino-esters *via*  $\alpha$ -Cyano-amines**

By JASJIT S. WALIA, PARVEEN S. WALIA, LINDA HEINDL, and H. LADER

*Chem. Comm.*, 1967, p. 1290.

On p. 1290, in formula (II) add H to C-2.

On p. 1290, in footnote †, for amino-esters *read* imino-esters.

On p. 1290, in the Table, the heading for columns 2 and 3 should be R and R<sup>1</sup>; and in entry (b) in R<sup>1</sup> column, read PhCH<sub>2</sub> for Ph; the last column heading should read B.p. at 0.2 mm.

On p. 1291, r.h.s., line 2, for (Ib) *read* (If).

**Rotational Isomeration in a Vinylogous Amide**

By RONALD J. PARRY

*Chem. Comm.*, 1967, 1294.

On p. 1295, in formula (Ib) interchange H<sup>5</sup> and H<sup>6</sup>;

On p. 1295, l.h.s., line 6, for H<sub>2</sub> *read* H-2;

On p. 1296, r.h.s., line 30, for H<sub>2</sub> *read* H-2.

**Rearrangement of 2,4-Diphenylthietan Dioxides to 3,5-Diphenyl-1,2-oxathiolan 2-Oxides**

By R. M. DODSON, P. D. HAMMEN, and R. A. DAVIS

*Chem. Comm.*, 1968, 9.

On p. 9, r.h.s., line 9 should read as

$$-12.8, J_{AP}=7.4, J_{AX}=2.6, J_{KP}=12.4; J_{KX}=8.4$$

On p. 10, l.h.s., sentence beginning on line 10 should read: 'The dihedral angle between H<sub>A</sub> and H<sub>X</sub> should be between 60° and 120°; thus H<sub>A</sub> and H<sub>X</sub> should be *trans* to each other. This makes H<sub>K</sub> *cis* to H<sub>X</sub> ( $J_{KX} = 8.4$ ) and *trans* to H<sub>P</sub> ( $J_{KP} = 12.4$ ); H<sub>A</sub> is *cis* to H<sub>P</sub> ( $J_{AP} = 7.4$ ).

**Symmetry Selection Rules for Sigmatropic Migrations in C<sub>n</sub>H<sub>n+1</sub> Monocycles**

By A. G. ANASTASSIOU

*Chem. Comm.*, 1968, 15.

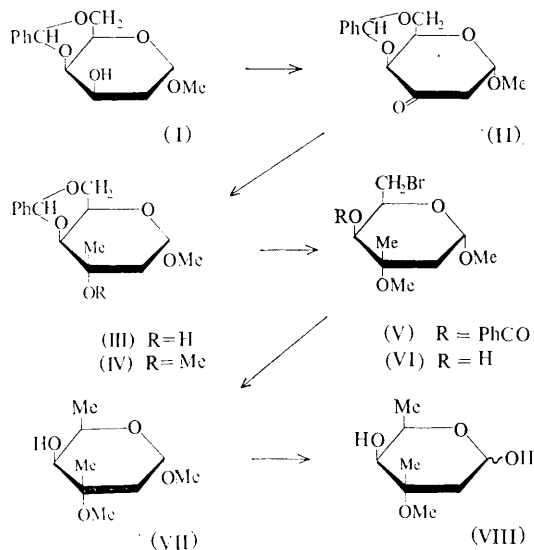
On p. 17, the last entry for Thermal configuration should read  $\psi_1^2\psi_2^2\psi_3^2\psi_4^1$ .

### The Synthesis of D-Arcanose

By G. B. HOWARTH, W. A. SZAERK, and J. K. N. JONES

*Chem. Comm.*, 1968, 62.

On p. 63, the middle two formulae drawings should be interchanged. The complete scheme should read:



### Structure of Deoxycrustecdysone, a Second Crustacean Moulting Hormone

By M. N. GALBRAITH, D. H. S. HORN, E. J. MIDDLETON, and R. J. HACKNEY

*Chem. Comm.*, 1968, 83.

On p. 84, in the upper structural formula interchange  $R^1$  and  $R^2$ .

On p. 84, r.h.s., line 18, for ref. 7 read ref. 4.

### Structure of Liguloxide, Liguloxidol, and Liguloxidol Acetate, Sesquiterpenes of Antipodol Guaian Type

By H. ISHII, T. TOZYO, and H. MINATO

*Chem. Comm.*, 1968, 106.

On p. 106, line 8, for  $[\alpha]_D -52.8^\circ$  read  $(\alpha)_D -58.2^\circ$ .

On p. 106, in formula (I), the double line extending off the first benzene ring should be a single thick line.

**Crystal and Molecular Structures of New Cobalt Carbonyl Clusters**

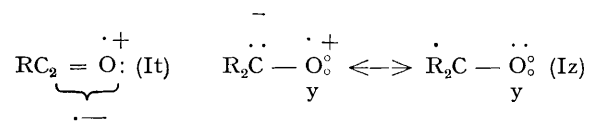
By V. ALBANO, P. CHINI, and V. SCATTURIN

*Chem. Comm.*, 1968, 163.On p. 164, r.h.s., line 7, for a cyclopropenyl methyl group, read a cyclopropenyl (CH)<sub>3</sub> group.**Electronically Excited Species in Organic Photochemistry; A Reply**

By H. E. ZIMMERMAN

*Chem. Comm.*, 1968, 174.

p. 175, Formulae (It) and (Iz) are incorrect and should be

**The Total Synthesis of (±)-4-Demethylaristolone and Related Compounds**

By EDWARD PIERS, WILLIAM DE WAAL, and RONALD W. BRITTON

*Chem. Comm.*, 1968, p. 188.

On p. 188, r.h.s., last line, for reduction of (VIII) read reduction of (VII).

On p. 189, ref. 7 was omitted and should read

? W. G. Dauben and J. T. Deving, *J. Org. Chem.*, 1966, **31**, 3794.**Unexpected Products obtained in the Thiation of Hypoxanthine Derivatives**

By Z. NEIMAN

*Chem. Comm.*, 1968, p. 200.

On p. 201, in structures (III), (IV), and (VI), for Y = S read Y = SH; in compound (VII), for X = S read X = SH.

**Methyl 2,2,2,4-Tetramethylbicyclo[1,1,0]butane-1-carboxylate**

By MISS C. BURRIDGE and D. P. G. HAMON

*Chem. Comm.*, 1968, 206.

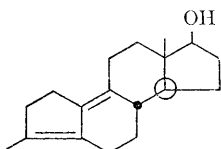
The title compound should read as Methyl 2,2,4,4-Tetramethyl-----

**A Backbone Rearrangement in A, 19-Bisnor-steroids**

By J. BASCOUL and A. CRASTES DE PAULET

*Chem. Comm.*, 1968, 256.

Formula (VI) should be



**Synthetical Approaches to Simple Derivatives of *as*-Indacene**

By R. R. HILL and G. H. MITCHELL

*Chem. Comm.*, 1968, 314.

On p. 314, r.h.s., l. 13, for 4·3 (m, 2H) read 6·3 (m, 2H).

**Synthesis of ( $\pm$ )-Dihydroradicinin**

By K. KATO, Y. HIRATA, and S. YAMAMURA

*Chem. Comm.*, 1968, p. 319.

On p. 319, l.h.s., line 13, for (V), m.p. 48—49° read (V), m.p. 94—95°.

**Nuclear Magnetic Resonance for the Reversible Stereospecific Protonation of Steroidal  $\alpha$ -Diazo-ketones**

By M. AVARO, J. LEVISALLES, and J. M. SOMMER

*Chem. Comm.*, 1968, p. 410.

In the title, add the word 'Evidence' after Resonance.

On p. 410, r.h.s., line 3, for ca. 93 p.p.m. read 9·3 p.p.m.

On p. 411, l.h.s., line 15, for PhC:N:N·CO·Ph read PhCN<sub>2</sub>·COPh.**Synthesis of New Phthalimidines and 3,4-Dihydroisocarbostyrils: New Method involving an Unusual Cyclohydration**

By I. T. BARNISH, C. L. MAO, R. L. GAY, and C. R. HAUSER

*Chem. Comm.*, 1968, 564.

Title, for Cyclohydration read Cyclodehydration.

Delete superscript 1 in title, and in line 1, for <sup>2</sup> read <sup>1</sup>, and in line 2, for <sup>3</sup> read <sup>2</sup>.**Tricarbonyl(*trans*- $\pi$ -pentadienyl)iron Cations**

By NYE A. CLINTON and C. PETER LILLYA

*Chem. Comm.*, 1968, 579.On p. 579, footnote <sup>a</sup> to the Table, for  $k_1 = (2.48 \pm 0.05) \times 10^{-4} \text{ sec.}^{-1}$  read  $k_1 = (2.48 \pm 0.05) \times 10^{-4} \text{ min.}^{-1}$ .**A Quantitative Study of the Quaternisation of Tropanes**

By G. FODOR, J. D. MEDINA, and NAGABHUSHANAM MANDAVA

*Chem. Comm.*, 1968, 581.On p. 581, in formula (I), C-3 hydroxyl group should be axial ( $\alpha$ -position) and the hydrogen should be equatorial ( $\beta$ -hydrogen).On p. 582, Table, Tropane (unsubstituted), 5th entry, for 3 $\beta$ ,6 $\beta$ -dihydroxy, read 3 $\alpha$ ,6 $\beta$ -dihydroxy; 8th entry, should read 3 $\alpha$ -hydroxy-ethiodide  
-N-ethyl nor- ethiodide.

**The Absolute Configuration of Sulphoxides: Mass Spectra of  
3 $\beta$ -Hydroxy-20-thia-17( $\alpha$  and  $\beta$ )-pregn-5-ene Oxides**

By R. M. DODSON, P. J. CAHILL, and V. C. NELSON

*Chem. Comm.*, 1968, 620.

On p. 621, the formula (I,S) should have a =O attached to the r.h.s. of the sulphur atom.

**Polar Solution Behaviour of Selenium Tetrabromide**

By N. KATSAROS and J. W. GEORGE

*Chem. Comm.*, 1968, 662.

p. 662, r.h.s., line 19, for SeBr<sub>4</sub> read SeBr<sub>2</sub>.

**Ruthenium Complexes which catalyse the Dimerization of Acrylonitrile**

By A. MISONO, Y. UCHIDA, M. HIDAI, and I. INOMATA

*Chem. Comm.*, 1968, 704.

On p. 705, r.h.s., line 2 after the Table, for  $\tau$  1.7 and 5.8 read  $\tau$  8.3 and 4.2.

**Synthesis of Unsaturated Sugars containing Vinylic Substituents**

By S. HANESSIAN and N. R. PLESSAS

*Chem. Comm.*, 1968, 706.

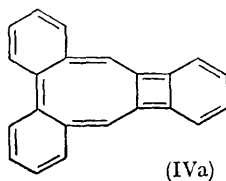
On p. 707, l.h.s., last line, for ( $\tau$  4.18 vinyl H, 6.47 OMe) read ( $\tau$  4.18, 4.67 vinyl H, 6.47 OMe).

**3,4:4,6:9,10-Tribenzobicyclo[6,2,0]decapentanene. A (4n + 4n)  $\pi$ -Electron System**

By P. J. GARRATT and R. H. MITCHELL

*Chem. Comm.*, 1968, 719.

On p. 720, formula (IVa) should be



**Dicarbonyl- $\pi$ -Allylrhodium**

By S. O'BRIEN

*Chem. Comm.*, 1968, 757.

On p. 758, the ordinates on the bottom of both parts of the Figure should read 2100, 2000, 1900, 1800 cm.<sup>-1</sup>.

**Identification of the Sex Pheromone of the False Codling Moth (*Argyroplote leucotreta*)**

By J. S. READ, F. L. WARREN, and P. H. HEWITT

*Chem. Comm.*, 1968, 792.On p. 793, l.h.s., line 8, for 266 ( $M^+$  read 226 ( $M^+$ )).**Mechanism of the Reaction of Methanesulphonylnitrene with Benzene**

By R. A. ABRAMOVITCH and V. UMA

*Chem. Comm.*, 1968, 797.

On p. 798, insert an equilibrium arrow between formulae (III) and (II); in structure (VII), add a positive charge in bottom right corner of the benzene ring.

**Biacetal Photochemistry: Products in Solution**

By WESLEY G. BENTRUDE and K. R. DARNALL

*Chem. Comm.*, 1968, 810.

In the title, for Biacetal read Biacetyl.

On p. 811, Table, Solvent, 4th entry for Cyclohexane read Cyclohexene.

**The Crystal Structure of Dimeric Dichloro-(2,9-dimethyl-1,10-phenanthroline)nickel(II)**

By H. S. PRESTON and C. H. L. KENNARD

*Chem. Comm.*, 1968, 819.On p. 819, l.h.s., 3 lines from bottom, inside the round bracket, for  $x, y, z; \frac{1}{2} = x$  read  $x, y, z; \frac{1}{2} + x$ .**The Crystal Structure of *cis*-Dichloro-*cis*-bis(dimethyl sulphoxide)-*trans*-dimethyltin(IV)**

By N. W. ISAACS, C. H. L. KENNARD, and W. KITCHING

*Chem. Comm.*, 1968, 820.On p. 820, l.h.s., 8 lines from bottom, for  $c = 6 \cdot$  read  $c = 6 \cdot 94 \pm 0 \cdot 01 \text{ \AA}$ .**The Reaction of Lithium Salts with Aluminium Borohydride: Nuclear Magnetic Resonance Evidence for Anionic Aluminium Borohydride Species**

By M. EHEMANN, H. NÖTH, N. DAVIES, and M. G. H. WALLBRIDGE

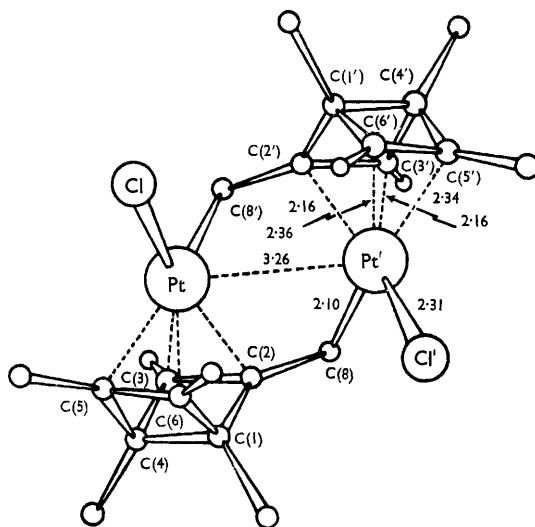
*Chem. Comm.*, 1968, 862.On p. 862, l.h.s., l. 5; replace the 3rd sentence with "Wiberg and Schrauzer<sup>2</sup> first treated  $B_2H_6$  with  $LiAlH_4$  in tetrahydrofuran . . . decompose to solvated hydrides of the type  $AlH_n(BH_4)_{3-n} \cdot R_2O$  ( $R_2O = Et_2O$ , tetrahydrofuran) and  $LiBH_4$ , and in ether to  $(LiBH_4)_2AlH_3$  or  $LiAlBH_7$ , also".

### A Dehydro-hexamethyl Dewar Benzene Complex of Platinum(II)

By R. MASON, G. B. ROBERTSON, P. O. WHIMP, B. L. SHAW, and G. SHAW

*Chem. Comm.*, 1968, 868.

This Figure was unfortunately omitted from the Communication.



### Synthesis of Hashish Cannabinoids by Terpenic Cyclisation

By L. CROMBIE and R. PONSFORD

*Chem. Comm.*, 1968, 894.

On p. 894, the double bond in the chroman ring of (X) should be deleted and inserted in the chroman ring of (IX).

On p. 894, footnote †, line 2, for B. Yagintsky read B. Yagnitinsky.

### The Nomenclature of the Sinensals

By R. TERANISHI, A. F. THOMAS, P. SCHUDEL, and G. BÜCHI

*Chem. Comm.*, 1968, 928.

In formula (II), the methyl group on the right hand end should be omitted; the double bond should be linked with a methylene group.

### A Novel Decarboxylation of an $\alpha$ -Keto-ester: Methyl Phenylpyruvate

By A. E. OPARA and G. READ

*Chem. Comm.*, 1968, 937.

On p. 937, l.h.s. lines 4 and 5. For dimethyl 2-benzylsuccinate read dimethyl 2-benzyl-3-phenylsuccinate.



**Charge-transfer Interactions of the Highest Valency Halides, Oxyhalides, and Oxides with Aromatic Hydrocarbons and Fluorocarbons: Ball-Plane Interactions**

By P. R. HAMMOND and R. R. LAKE

*Chem. Comm.*, 1968, 987.On p. 988, l.h.s., line 8. For  $\text{ClO}_2\text{F}$  read  $\text{ClO}_3\text{F}$ .**Selective Nucleophilic Attack by Peroxy-acid Anions on Nucleic Acid Components**

By L. R. SUBBARAMAN, JIJI SUBBARAMAN, and E. J. BEHRMAN

*Chem. Comm.*, 1968, 1024.

Authors: for JIJI SUBBARAMAN read JIJIE SUBBARAMAN.

Table: in heading to second column, for  $(\text{mole}^{-1} \text{min.}^{-1})$  read  $(\text{litre mole}^{-1} \text{min.}^{-1})$   
second column, four lines from bottom, for 8.3 read 83.

On p. 1025, r.h.s., line 5, after uracil insert "formic,".

References: in ref. 1, line 2, after *Biochem. Acta*, insert "1967".**Ozonolysis of Hydroxymethylene Ketones: the Baeyer-Villiger Reaction as a Source of Anhydride Formation**

By DOMINIC YOUNG and S. W. PELLETIER

*Chem. Comm.*, 1968, 1055.

On p. 1056, Table, last column, the values should read: 56, 4, 4, 14, 13.

**Alkaline Hydrolysis of 2,2,3,4,4-Pentamethyl-1-phenylphosphetanium Bromide**

By SHELDON E. CREMER

*Chem. Comm.*, 1968, 1132.

The compound named in the title should read 1,2,2,3,4,4-Hexamethyl-1-phenylphosphetanium Bromide.

**Mechanism for the Photohydration of Pyrimidines**

By SHIH YI WANG and JOHN C. NNADI

*Chem. Comm.*, 1968, 1160.On p. 1161, diagram, add a positive charge to the quaternary nitrogen in structural formula ( $\text{D}_1$ ).**Complexes of Platinum(II) with Carborane [1,2- or 1,7-dicarba-closo-dodecaborane(12)] and Tertiary Phosphines: Loss of Hydrogen from the Ligands**

By S. BRESADOLA, P. RIGO, and A. TURCO

*Chem. Comm.* 1968, 1205.

On p. 1206, l.h.s., line 7 from bottom, for 2 mol. of iodine read 1 mol. of iodine.

**Rearrangement of 1,2,3-Triphenylcyclopropene to 1,2-Diphenylindene catalyzed by Di- $\mu$ -chloro-dichlorobis(ethylene)diplatinum(II)**

By JOHN A. WALKER and MILTON ORCHIN

*Chem. Comm.*, 1968, 1239.On p. 1239, r.h.s., line 2, for  $\tau$  6.1 read  $\tau$  5.1.**Hindered Internal Rotation in 5-(*NN*-Dimethylamino)-1,3,4-oxathiazole 3,3-Dioxide**

By H. J. JAKOBSEN and A. SENNING

*Chem. Comm.*, 1968, 1245.On p. 1245, the second footnote, line 1, for 5-(*NN*-Dimethylamino)-1,3,4-oxathiazole read 5-(*NN*-Dimethylamino)-1,3,4-oxathiazole.On p. 1246, r.h.s., line 2, for RCNSMe<sub>2</sub> read RCSNMe<sub>2</sub>.**The Reduction of Nitroarenes with Sodium Dihydrobis(2-methoxyethoxy)aluminate**

By JOHN F. CORBETT

*Chem. Comm.*, 1968, 1257.

On p. 1258, Table, third column (headed "Product"), lines 15—20 inclusive, for initial capital "D" read "B", throughout.

**Synthesis of Cyano-sugars: A Route to Branched-chain Sugars**

By B. E. DAVIDSON, R. D. GUTHRIE, and A. T. MCPHAIL

*Chem. Comm.*, 1968, 1273.

Authors: for B. E. Davidson read B. E. Davison.

**The Reaction of Alcohols with Carbon Tetrachloride and Phosphorus Trisdimethylamide**

By I. M. DOWNIE, J. B. LEE, and M. F. S. MATOUGH

*Chem. Comm.*, 1968, 1350.

In the title, for Phosphorus read Phosphorous.

On p. 1351, r.h.s., the first displayed formula, the plus sign relates to the left hand phosphorus atom.

**The Reaction of Pentacarbonyliron with Phenyl- and Methyl-lithium; a Crystal Structure Analysis of a Two-atom Bridged Complex**

By E. O. FISCHER, V. KIENER, D. ST. P. BUNBURY, E. FRANK, P. F. LINDLEY, and O. S. MILLS

*Chem. Comm.*, 1968, 1378.

On p. 1378, l.h.s., line 14, for trimethylammonium tetrafluoroborate read trimethyloxonium tetrafluoroborate.

On p. 1380, reference 10, for R. E. Schaer read R. E. Sacher.