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CORRIGENDA

The assignment of the ^{15}N NMR signals of N_5^+ in the Communication by **K. O. Christe et al.** in Issue 13/14, 1999, pp. 2004–2010 were inadvertently reversed. The signal at $\delta = -237.3$ (calculated: -235) corresponds to the central N atom N3, and that at $\delta = -100.4$ (calculated: -95) to N1 (see for example Table 1 in this contribution).

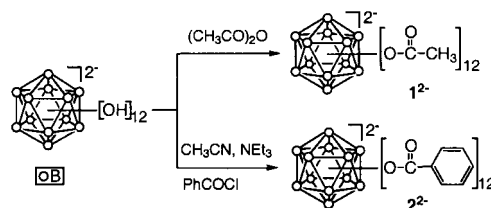
The ^{13}C NMR data given in the Communication by **R. Langer et al.** in Issue 4, 2001, pp. 726–728 for dinitro[2.2]paracyclophane **2b** were for a mixture of isomers. Data for the pseudo-*para* isomer as well as revised data for the nitro[2.2]paracyclophane **2a** and the amino[2.2]paracyclophane **3a** are given here.

2a: $\delta = 34.36, 34.69, 34.91, 35.93, 129.45, 129.86, 132.31, 133.07, 133.12, 136.37, 137.27, 137.73, 139.26, 139.68, 142.02, 149.17$.

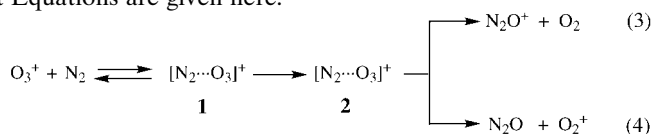
2b: $\delta = 34.17, 34.60, 129.13, 133.87, 136.46, 137.36, 142.00, 149.71$.

3a: $\delta = 32.23, 33.03, 34.95, 35.37, 122.36, 122.99, 124.60, 126.82, 131.48, 132.43, 133.23, 135.44, 138.91, 138.97, 141.06, 144.72$.

In the Communication by **M. F. Hawthorne et al.** in Issue 9, 2001, pp. 1661–1664 there was an error in the equation at the top of the frontispiece: it should read R_2O not $(\text{R}^2\text{CO})_2\text{O}$, and the text at the bottom the page should read “Globular molecules with an icosahedral core—12(12)-closomers—are obtained by total esterification and etherification of $[\text{B}_{12}(\text{OH})_{12}]^{2-}$. For details as well as for the transformation of an anionic closomer ether into the neutral *hypercloso* analogue see the following two communications”. An error in Scheme 1 was also inadvertently introduced; the correct scheme is given here.



In the Communication by **G. de Petris et al.** in Issue 10, 2001, pp. 1938–1941 complexes **1** and **2** in Equations (3) and (4) were erroneously written. The correct Equations are given here.



In the minireview by **P. Braunstein** and **N. M. Boag** in Issue 13, 2001, pp. 2427–2433 the isolobal signs were inadvertently missing from the second sentence of the abstract, the second sentence should read:

The C, Si, and P donor atoms of these ligands are sp^3 -hybridized and the ligands are related to each other by the isolobal analogy: $(\text{CR}_3)^- \rightarrow (\text{SiR}_3)^- \rightarrow \text{PR}_3$.

We would like to apologize for this error.

In the communication by **Y. Gu et al.** in Issue 12, 2001, pp. 2320–2322 the protein surface coverages listed in Table 1 were incorrect. Since that publication, the authors identified and corrected a systematic error in the method used to determine surface coverages. The corrected Table is given here. The corrected data show that a spacer arm of three ethylene glycol units (in Gal-3) is sufficient for rgp120 recognition of a galatosyl receptor presented at the surface of a planar DOPC membrane. Longer spacer arm lengths (in Gal-4 and -5) result in less efficient recognition. The authors apologize for this error.

Table 1. Summary of rgp120 binding behavior at planar DOPC membranes, either pure or doped with 5 % (mol mol $^{-1}$) of Gal-3, -4, or -5 in the outer leaflet.^[a]

	DOPC	Gal-3	Gal-4	Gal-5
K_a ($\times 10^6$)	–	5.4 ± 1.8	3.8 ± 0.57	2.5 ± 0.57
cooperativity coefficient (ω)	–	1.44 ± 0.20	1.50 ± 0.28	1.59 ± 0.11
surface coverage [mol cm $^{-2} \times 10^{-13}$] ^[b]	1.6 ± 0.23	3.6 ± 0.57	2.2 ± 0.19	2.2 ± 0.13

[a] All values listed are the mean and standard deviation of two experimental trials. [b] Surface coverages were determined at a dissolved rgp120 concentration of 208 nM, using a modification of the method described by Haldy et al.^[26]