75.7, 58.1, 55.8, 55.5, 55.2, 53.6, 38.7, 38.1, 36.4, 34.1, 30.8, 29.7, 24.3; HRMS (M + H)$^+$ calcd for $C_{28}H_{36}N_3O_6$ 510.2604, found 510.2608. **Diasteromer 20b** was prepared from **19b** exactly as described above: HPLC t_r = 8.62 min; ^{1}H NMR (CD_3OD) δ 8.24 (s, 1H), 7.27–7.12 (m, 5H), 7.07 (d, J = 8.1, 2H), 6.74 (d, J = 8.3, 2H), 4.59 (dd, J = 4.6, 9.6, 1H), 4.16 (m, 1H), 4.04 (m, 1H), 3.96 (d, J = 5.9, 1H), 3.22 (m, 1H), 2.98–2.86 (m, 3H), 2.35 (m, 1H), 1.96 (m, 2H), 1.74 (m, 2H), 1.55 (m, 3H), 1.44–1.17 (m, 3H); ^{13}C NMR (CD_3OD) δ 177.3, 169.4, 166.6, 158.2, 139.2, 131.6, 130.3, 129.3, 127.5, 126.3, 116.9, 75.8, 58.6, 58.0, 55.8, 55.1, 53.9, 53.6, 38.8, 37.9, 36.6, 34.1, 33.2, 30.0, 29.2, 24.0; HRMS (M + H)$^+$ calcd for $C_{28}H_{36}N_3O_6$ 510.2604, found 510.2618.

Acknowledgment. We are grateful to the National Institutes of Health (GM-39758) and IGEN, Inc., Rockville, MD, for support. Acknowledgement is made to the donors of the Petroleum Research Fund, administered by the American Chemical Society, for partial support of this research. This research was supported in part by a grant from the Amherst College Faculty Research Award Program. D.E.H. is a National Science Foundation Presidential Young Investigator (MCB-8958239) and a Camille and Henry Dreyfus Teacher-Scholar.

Supporting Information Available: ^{1}H-NMR spectra for compounds **2–9**, **15**, and **17–20b** (15 pages). This material is contained in libraries on microfiche, immediately follows this article in the microfilm version of the journal, and can be ordered from the ACS; see any current masthead page for ordering information.

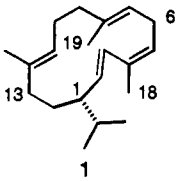
JO950216X

Additions and Corrections

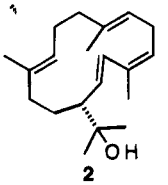
Vol. 60, 1995

Alexey V. Vorobjev,* Makhmut M. Shakirov, Victor A. Raldugin, and Clayton H. Heathcock. Conformational Analysis of the 10- and 13-Hydroxy Derivatives of Cembrene.

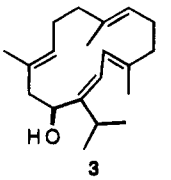
Page 64, column 1. Structures **1–10** should be replaced by the following structures:



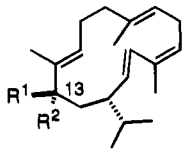
1



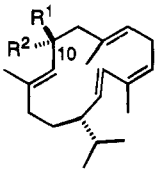
2



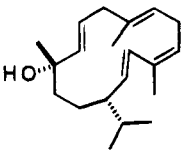
3



4: R^1 = OH, R^2 = H
5: R^1 = H, R^2 = OH
6: R^1 = H, R^2 = OAc



7: R^1 = OH, R^2 = H
8: R^1 = H, R^2 = OH
9: R^1 + R^2 = O



10

JO9540110

David P. Kelly,* Martin G. Banwell, John H. Ryan, James R. Phyland, and Jason R. Quick. ^{13}C–^{1}H Coupling Constants in Carbocations. 8. Application of the ΔJ Equation to Tertiary Dicyclopropylcarbinyl Cations: The Methyl Dicyclopropylcarbinyl, (1α,3β,5β,7α)-2-Methyltricyclo[5.1.0.0^{3,5}]-octan-2-yl, (1α,3α,5α,7α)-2-Methyltricyclo[5.1.0.0^{3,5}]octan-2-yl, and 3-Methyltetracyclo[3.3.1.0^{2,8}.0^{4,6}]nonan-3-yl (Triasteryl) Cations.

Page 1654. The data for compound **20** in Table 1 should read as follows: **20**[h] −110 43.9 (d, 179) 263.3 (s) 43.9 (d, 179) 40.7 (t, 169)[c] 74.2 (d, 171) 21.3 (t, 130) 38.0 (q, 130)

[h]Chemical shifts from internal CD_2Cl_2 taken as 52.8 ppm.

JO9540108

Dieter Seebach,* Robert Dahinden, Roger E. Marti, Albert K. Beck, Dietmar A. Plattner, and Florian N. M. Kühnle. On the Ti-TADDOLate-Catalyzed Diels–Alder Addition of 3-Butenoyl-1,3-oxazolidin-2-one to Cyclopentadiene. General Features of Ti-BINOLate- and Ti-TADDOLate-Mediated Reactions.

Page 1788. The correct receipt date for this manuscript is October 19, 1994.

JO9540099