

Preparation of Optically Active α -Amino[3]ferrocenophanes – Building Blocks for Chelate Ligands in Asymmetric Catalysis

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We noticed that some mistakes were made in the determination of the $\Delta\epsilon$ and $[\alpha]_D^{20}$ values of compounds **6** and **11** in our article^[1] that need to be corrected. The correct values for **6** (p. 3597) and **11** (p. 3598f.) (including the corrected Figure 8, p. 3596) are as follows:

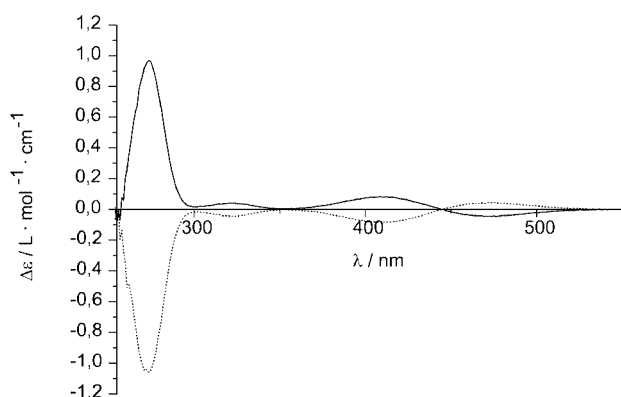


Figure 8. CD spectra of (1*R*,3*R*)-**6** (solid line, 4×10^{-4} g mL⁻¹) and (1*S*,3*S*)-**6** (dotted line, 4×10^{-4} g mL⁻¹) in CH₃CN

(1*S*,3*S*)-**6**: $[\alpha]_D^{20} = -89$ ($c = 2 \times 10^{-3}$ g mL⁻¹, CH₃CN). CD: $\Delta\epsilon$ ($\lambda_{\max}$) = -1.06 (274), -0.04 (326), -0.08 (412), $+0.04$ (476) ($c = 4 \times 10^{-4}$ g mL⁻¹, CH₃CN). (1*R*,3*R*)-**6**: $[\alpha]_D^{20} = +91$ ($c = 2 \times 10^{-3}$ g mL⁻¹, CH₃CN). CD: $\Delta\epsilon$ ($\lambda_{\max}$) = $+0.97$ (274), $+0.04$ (326), $+0.08$ (412), -0.04 (476) ($c = 4 \times 10^{-4}$ g mL⁻¹, CH₃CN).

(1*S*,3*S*)-**11**: $[\alpha]_D^{20} = -109$ ($c = 2 \times 10^{-3}$ g mL⁻¹, CH₃CN). CD: $\Delta\epsilon$ ($\lambda_{\max}$) = -0.23 (275), $+0.02$ (334), -0.10 (416), $+0.03$ (483) ($c = 4 \times 10^{-4}$ g mL⁻¹, CH₃CN). (1*R*,3*R*)-**11**: $[\alpha]_D^{20} = +111$ ($c = 2 \times 10^{-3}$ g mL⁻¹, CH₃CN). CD: $\Delta\epsilon$ ($\lambda_{\max}$) = $+0.23$ (275), -0.02 (334), $+0.10$ (416), -0.02 (483) ($c = 4 \times 10^{-4}$ g mL⁻¹, CH₃CN).

[1] P. Liptau, L. Tebben, G. Kehr, B. Wibbeling, R. Fröhlich, G. Erker, *Eur. J. Inorg. Chem.* **2003**, 3590–3600.

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