

# Additions and Corrections

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**Pascal Maire, Frank Breher, Hartmut Schönberg, and Hansjörg Grützmacher\*:** Chiral Rhodium(I) and Iridium(I) Amino–Olefin Complexes:  $pK_a$ , N–H Bond Dissociation Energy, and Catalytic Transfer Hydrogenation.

Pages 3207–3218. Due to an erroneous conversion of energy units, the equation used to estimate the NH bond dissociation energy (BDE) is wrong. The correct formula is  $BDE = 5.74pK_a + 96.5E^\circ_{Fc^+/Fc} + C$  ( $C = 306.9 \text{ kJ mol}^{-1}$  for dmso). Therewith, the NH BDE in (*R,R*)-**6** is  $363(5) \text{ kJ mol}^{-1}$ , which is about the same as previously observed for other Rh(I) amine complexes **H** and **J**.

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