

Errata**Erratum: Bull. Chem. Soc. Jpn., 70, 957 (1997)****Interfacial Mechanism in the Extraction Kinetics of Ni(II)
with 2-(5-Bromo-2-pyridylazo)-5-diethylaminophenol and
Molecular Dynamics Simulation of Interfacial Reactivity of the Ligand****Hitoshi Watarai,* Masashi Gotoh, and Narushi Gotoh†**

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On Page 963, Figure 9 was printed incorrectly. The corrected figure is given below.

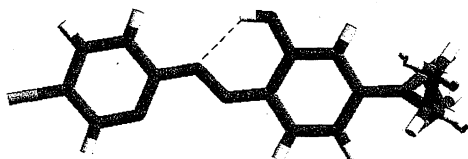
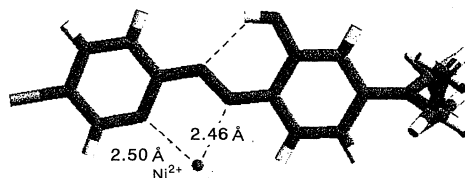
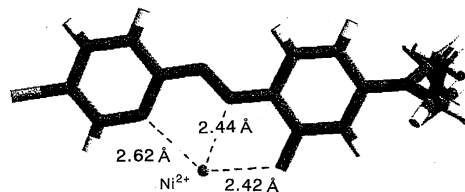
E = 48.30 kcal/mol**E = 48.34 kcal/mol**↓ **ΔE = -1.34 kcal/mol****E = 47.00 kcal/mol**

Fig. 9. Molecular Mechanics calculation of the ring-closure step in the complexation of Ni(II) with 5-Br-PADAP in water.