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Erratum

A CASSCF and CCI study of the formation of the $Ni_2(C_2H_4)$ complex

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In the quoted paper, the binding energies between Ni_2 and C_2H_4 were erroneously given as 24.6 kcal/mol at the CCI level of approximation and 28.1 kcal/mol with cluster corrections included. The error was due to an incorrect calculation of the CI wavefunction at the asymptotic limit. The correct binding energies are 15.8 and 20.9 kcal/mol respectively.

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