ERRATUM

Self-consistent, nonorthogonal group function approximation: an ab initio approach for modelling interacting fragments and environmental effects, E.L. Mehler, J. Math. Chem. 10(1992)57.

On page 84, figure 4 is missing. This figure is as follows:

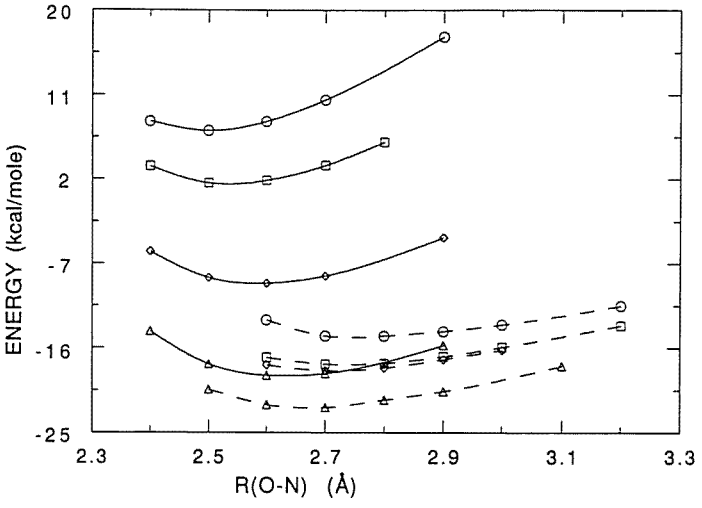

Fig. 4. Interaction potentials for hydrated ammonia–formic acid complex. Ionized pair, solid curves; neutral pair, dashed curves. O AB; □ 2SAB; ◇ AB2S; Δ 2SAB2S; S = H_2O, A = HCOOH($HCOO^-$); B = NH_3(NH_4^+).