ADDITIONS AND CORRECTIONS

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Chia-Ching Wang, Ya-Jen Yang, and Jyh-Chiang Jiang*: DFT Study of NH_x ($x = 1-3$) Adsorption on $RuO_2(110)$ Surfaces.

Page 2816. Because of a mistake of using the wrong electronic energy value for the NH molecule, all the binding energies of the NH molecule on $RuO_2(110)$ surfaces in the article should be reduced by 1.58 eV. Table 1 lists the binding energies of NH molecules after the correction.

TABLE 1: Corrected Binding Energies (eV) of NH on Stoichiometric $RuO_2(110)$ Surfaces ($\theta = 0.5$ or 1 ML) and Oxygen-Rich $RuO_2(110)$ Surface

species	$\theta = 0.5$ ML	$\theta = 1$ ML	oxygen-rich
NH	2.65	2.67	2.71

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