Chemical Science

ROYAL SOCIETY OF CHEMISTRY

CORRECTION

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Cite this: *Chem. Sci.*, 2025, **16**, 10063

Correction: Uncovering diverse reactivity of NHCs with diazoalkane: C–H activation, C=C bond formation, and access to N-heterocyclic methylenehydrazine

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DOI: 10.1039/d5sc90098k

rsc.li/chemical-science

Correction for 'Uncovering diverse reactivity of NHCs with diazoalkane: C–H activation, C=C bond formation, and access to N-heterocyclic methylenehydrazine' by Kajal Balayan *et al.*, *Chem. Sci.*, 2024, **15**, 18387–18394, https://doi.org/10.1039/D4SC05740F.

The authors regret that there is an error in the previously published version of Fig. 12. The structure was inadvertently labelled as COOEt, whereas it should be COO^{t}Bu. However, all calculations were performed using COO^{t}Bu, so the reported energy values remain unchanged. The corrected scheme is shown below. This correction does not affect any of the conclusions presented in the work.

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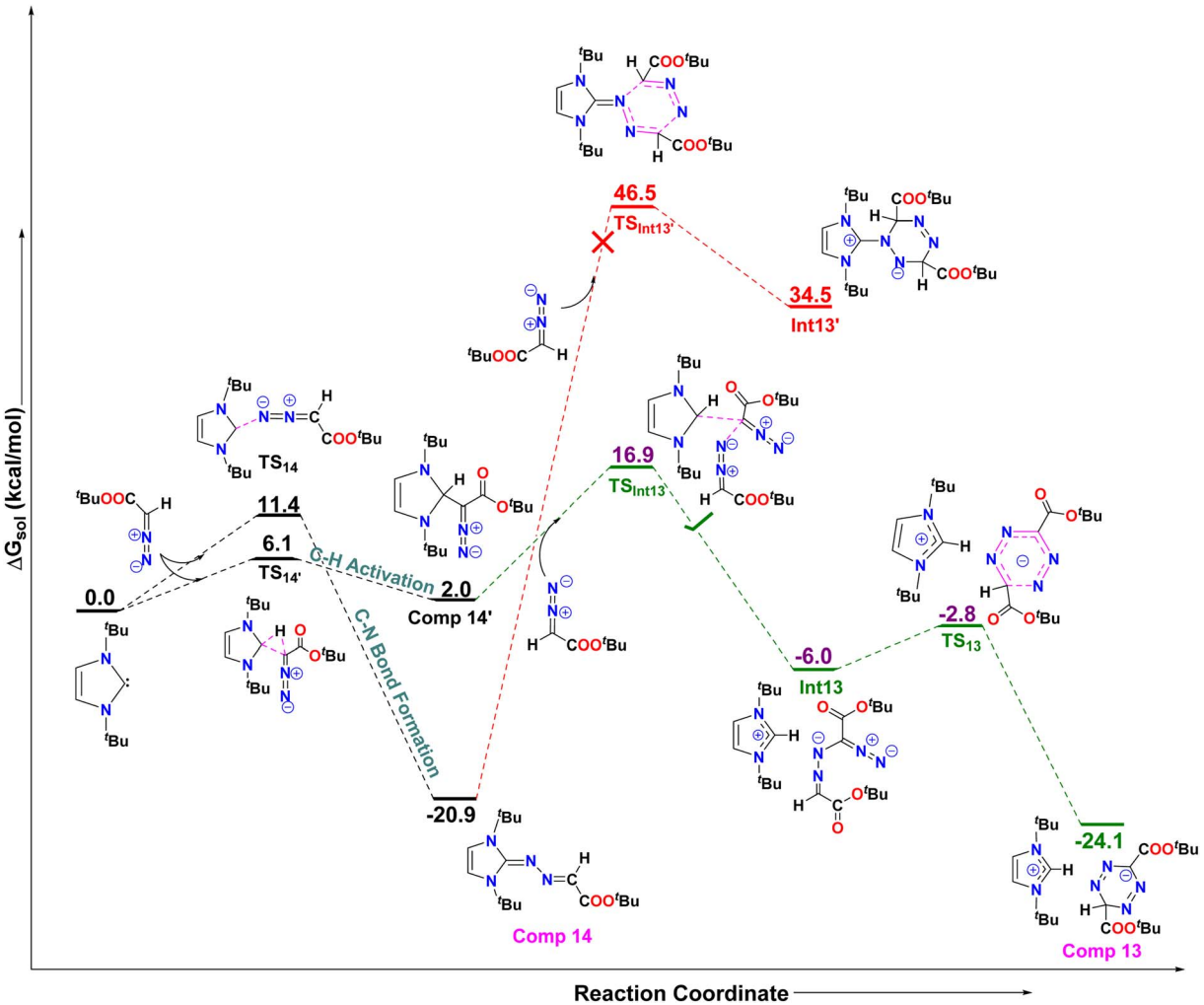

Fig. 12 The free energy profiles for the formation of compound **13** and compound **14** have been shown here. All values are in kcal mol^{-1}. Level of theory: PBE0-D3/def2TZVP//PBE-D3/defTZVP with solvent toluene.

The Royal Society of Chemistry apologises for these errors and any consequent inconvenience to authors and readers.