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CORRECTION

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Correction: High-level *ab initio* characterization of the OH + CH_3NH_2 reaction

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Correction for 'High-level *ab initio* characterization of the OH + CH_3NH_2 reaction' by Balázs Gruber *et al.*, *Phys. Chem. Chem. Phys.*, 2024, **26**, 28543–28553, https://doi.org/10.1039/d4cp01470g.

After publication of our paper[1] on the benchmark *ab initio* stationary-point characterization of the OH + CH_3NH_2 reaction, a relevant reference[2] came to our attention, which studied the same system with combined experimental and theoretical methods. The focus of the two studies was different, as ref. 1 provided a high-level *ab initio* description of the title reaction revealing new pathways leading to H atom as well as CH_3 and NH_2 radical formation, whereas ref. 2 reported a detailed kinetics investigation for the major H-abstraction channels. In the case of the H-abstraction processes, ref. 1 confirms and shows the uncertainty of the lower-level energy predictions of ref. 2. The two studies support and complement each other; thus, the interested reader may consult with both ref. 1 and 2 to get a more complete picture on the kinetics and mechanisms of the title reaction.

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

References

1 B. Gruber and G. Czakó, *Phys. Chem. Chem. Phys.*, 2024, **26**, 28543.
2 D. González, A. Lema-Saavedra, S. Espinosa, E. Martínez-Núñez, A. Fernández-Ramos, A. Canosa, B. Ballesteros and E. Jiménez, *Phys. Chem. Chem. Phys.*, 2022, **24**, 23593.

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