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^a Affiliation One; ^b Affiliation Two; ^c Affiliation Three

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A.O.(Author One) and A.T. (Author Two) contributed equally to this work (remove if not applicable).

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Fig. 1. Placeholder image of a frog with a long example caption to show justification setting.

Table 1. Comparison of the fitted potential energy surfaces and ab initio benchmark electronic energy calculations

Species	CBS	CV	G3
1. Acetaldehyde	0.0	0.0	0.0
2. Vinyl alcohol	9.1	9.6	13.5
3. Hydroxyethylidene	50.8	51.2	54.0

nomenclature for the TSs refers to the numbered species in the table.

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Acknowledgments

Please include your acknowledgments here, set in a single paragraph. Please do not include any acknowledgments in the Supporting Information, or anywhere else in the manuscript.

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 $\label{Fig.2.} \textbf{Fig. 2.} \ \ \textbf{This caption would be placed at the side of the figure, rather than below it.}$

$$(x+y)^3 = (x+y)(x+y)^2$$

$$= (x+y)(x^2 + 2xy + y^2)$$

$$= x^3 + 3x^2y + 3xy^3 + x^3.$$
[1]

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