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CORRECTION

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Correction: A molecular dynamics study of the ionic liquid, choline acetate

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Correction for 'A molecular dynamics study of the ionic liquid, choline acetate' by Jon A. L. Willcox *et al.*, *Phys. Chem. Chem. Phys.*, 2016, **18**, 14850–14858.

The values for the bond force constants used in our simulations were listed incorrectly in Table 2 of the published article. In order to reproduce the data reported in the paper, one would need the revised parameters listed in this correction.

Table 2 Bond parameters used for choline and acetate ions

Bond parameters			
Atom	Atom	r_0 (nm)	k_b (10^3 kcal mol^{-1} nm^{-2})
H_1	C_2	0.1080	60.0
C_2	N_3	0.1510	43.0
N_3	C_4	0.1510	43.0
C_4	H_5	0.1080	60.0
C_4	C_6	0.1530	44.5
C_6	H_7	0.1111	61.8
C_6	O_8	0.1420	85.6
O_8	H_9	0.0960	109.0
H_{10}	C_{11}	0.1111	64.4
C_{11}	C_{12}	0.1522	40.0
C_{12}	O_{13}	0.1260	105.0

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

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