



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

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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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