

## CORRECTION

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# Correction: The effect of an ionic liquid on the rate of reaction at a phosphorus centre

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 Correction for 'The effect of an ionic liquid on the rate of reaction at a phosphorus centre' by Bradley J. Butler *et al.*, *New J. Chem.*, 2015, DOI: 10.1039/c4nj01224k.

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In Tables 3 and 5, the units for entropy are incorrectly listed as  $\text{kJ mol}^{-1}$ . The correct unit is  $\text{J K}^{-1} \text{mol}^{-1}$ , and the correct tables are shown below.

**Table 3** Activation parameters for the ethanolysis of diethyl chlorophosphate at different mole fractions of [bmim][(CF<sub>3</sub>SO<sub>2</sub>)<sub>2</sub>N] in ethanol-*d*<sub>6</sub>

$\chi$ [bmim][(CF <sub>3</sub> SO <sub>2</sub> ) <sub>2</sub> N]	$\Delta H^\ddagger/\text{kJ mol}^{-1}{}^a$	$\Delta S^\ddagger/\text{J K}^{-1} \text{mol}^{-1}{}^a$
0	$62 \pm 2$	$-201 \pm 7$
0.31	$47 \pm 2$	$-238 \pm 8$
0.72	$44 \pm 3$	$-253 \pm 11$

<sup>a</sup> Uncertainties quoted are derived from the fit of the linear regression.
**Table 5** Activation parameters in the presence of the salts [bmim][(CF<sub>3</sub>SO<sub>2</sub>)<sub>2</sub>N], [bmim][Cl] and lithium bis(trifluoromethylsulfonyl)imide in ethanol-*d*<sub>6</sub> for the ethanolysis of diethyl chlorophosphate

$\chi_{\text{salt}}$	$\Delta H^\ddagger/\text{kJ mol}^{-1}{}^a$	$\Delta S^\ddagger/\text{J K}^{-1} \text{mol}^{-1}{}^a$
0	$62 \pm 2$	$-201 \pm 7$
0.21 [bmim][(CF <sub>3</sub> SO <sub>2</sub> ) <sub>2</sub> N]	$46 \pm 2$	$-243 \pm 7$
0.24 [Li][(CF <sub>3</sub> SO <sub>2</sub> ) <sub>2</sub> N]	$76 \pm 5$	$-180 \pm 16$
0.22 [bmim][Cl]	$52 \pm 1$	$-215 \pm 4$

<sup>a</sup> Uncertainties quoted are derived from the fit of the linear regression.

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

