

## CORRECTION

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Cite this: *New J. Chem.*, 2015,  
39, 1525

DOI: 10.1039/c4nj90052a

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

Bradley J. Butler and Jason B. Harper\*

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.

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  $[\text{bmim}][\text{(CF}_3\text{SO}_2)_2\text{N}]$  in ethanol- $d_6$

$\chi$ $[\text{bmim}][\text{(CF}_3\text{SO}_2)_2\text{N}]$	$\Delta H^\ddagger/\text{kJ mol}^{-1}\text{a}$	$\Delta S^\ddagger/\text{J K}^{-1} \text{mol}^{-1}\text{a}$
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  $[\text{bmim}][\text{(CF}_3\text{SO}_2)_2\text{N}]$ ,  $[\text{bmim}][\text{Cl}]$  and lithium bis(trifluoromethylsulfonyl)imide in ethanol- $d_6$  for the ethanolysis of diethyl chlorophosphate

$\chi_{\text{salt}}$	$\Delta H^\ddagger/\text{kJ mol}^{-1}\text{a}$	$\Delta S^\ddagger/\text{J K}^{-1} \text{mol}^{-1}\text{a}$
0	$62 \pm 2$	$-201 \pm 7$
0.21 $[\text{bmim}][\text{(CF}_3\text{SO}_2)_2\text{N}]$	$46 \pm 2$	$-243 \pm 7$
0.24 $[\text{Li}][\text{(CF}_3\text{SO}_2)_2\text{N}]$	$76 \pm 5$	$-180 \pm 16$
0.22 $[\text{bmim}][\text{Cl}]$	$52 \pm 1$	$-215 \pm 4$

<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.

School of Chemistry, University of New South Wales, UNSW Sydney 2052, Australia. E-mail: [j.harper@unsw.edu.au](mailto:j.harper@unsw.edu.au)

