



Cite this: *Phys. Chem. Chem. Phys.*, 2022, 24, 28080

DOI: 10.1039/d2cp90215j

rsc.li/pccp

Correction: The structuring effect of the alkyl domains on the polar network of ionic liquid mixtures: a molecular dynamics study

Valerio Mazzilli,^{ab} Yanting Wang^{cd} and Giacomo Saielli^{*ab}

Correction for 'The structuring effect of the alkyl domains on the polar network of ionic liquid mixtures: a molecular dynamics study' by Valerio Mazzilli et al., *Phys. Chem. Chem. Phys.*, 2022, 24, 18783–18792, <https://doi.org/10.1039/D2CP02786K>.

The authors would like to correct an error in the chemical structure in Fig. 1 of the published article. The correct figure is shown below.

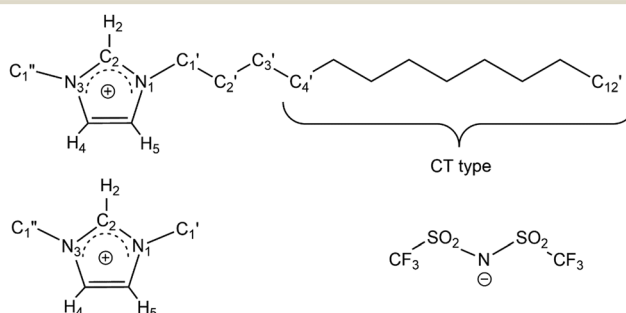


Fig. 1 Structural formula of (top) 1-dodecyl-3-methylimidazolium, [C₁₂C₁im] or C₁₂ in this work; (bottom left) 1,3-dimethylimidazolium, [C₁C₁im] or C₁ in this work; (bottom right) bis(trifluoromethanesulfonyl)imide, [Tf₂N]. The methylene and methyl groups of the alkyl chain of [C₁₂C₁im] from C₄' to the terminal methyl C₁₂' bear no total charge.

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

^a CNR Institute on Membrane Technology, Unit of Padova, Via Marzolo, 1 – 35131 Padova, Italy. E-mail: giacomo.saielli@unipd.it

^b Department of Chemical Sciences, University of Padova, Via Marzolo, 1 – 35131 Padova, Italy

^c CAS Key Laboratory of Theoretical Physics, Institute of Theoretical Physics, Chinese Academy of Sciences, 55 East Zhongguancun Road, P. O. Box 2735, Beijing 100190, China

^d School of Physical Sciences, University of Chinese Academy of Sciences, 19A Yuquan Road, Beijing 100049, China

