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CORRECTION

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Correction: Salt parameterization can drastically affect the results from classical atomistic simulations of water desalination by MoS_2 nanopores

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Correction for 'Salt parameterization can drastically affect the results from classical atomistic simulations of water desalination by MoS_2 nanopores' by João P. K. Abal *et al.*, *Phys. Chem. Chem. Phys.*, 2020, **22**, 11053–11061, DOI: 10.1039/d0cp00484g.

Our manuscript contained a NaCl-ε charge parameter error in Table 1. The corrected Table 1 is shown below, the references are the same as in the original article but are labelled 1–6 here:

Table 1 The Lennard-Jones parameters and atoms charges employed in the simulations

	σ_{LJ} [Å]	ε_{LJ} [kcal mol^{-1}]	Charge (e)
Na-ε [1]	2.52	0.0346	0.885
Cl-ε [1]	3.85	0.3824	−0.885
Na-J [2]	2.18	0.1684	1.0
Cl-J [2]	4.92	0.0117	−1.0
O-Tip4p/ε [3]	3.165	0.1848	−1.054
H-Tip4p/ε [3]	0.0	0.0	0.5270
O-Tip4p/2005 [4]	3.1589	0.1852	−1.1128
H-Tip4p/2005 [4]	0.0	0.0	0.5564
Mo [5]	4.20	0.0135	0.6
S [5]	3.13	0.4612	−0.3
C [6]	3.40	0.0860	0.0

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

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