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## Correction: How fast do defects migrate in halide perovskites: insights from on-the-fly machine-learned force fields

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Correction for 'How fast do defects migrate in halide perovskites: insights from on-the-fly machine-learned force fields' by Mike Pols *et al.*, *Chem. Commun.*, 2023, **59**, 4660–4663, <https://doi.org/10.1039/D3CC00953J>.

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The authors regret that there was an error in the units listed for the reported diffusion parameters shown in Table 1, header row, column 3 of the original article.  $D_0$  ( $\times 10^{-6}$  s) should read  $D_0$  ( $\times 10^{-6}$  cm<sup>2</sup> s<sup>-1</sup>). The correct units are shown in the updated table below.

**Table 1** Overview of the diffusion parameters  $D_0$  and  $E_a$  for halide vacancies ( $V_X$ ) and halide interstitials ( $I_X$ ) in CsPbI<sub>3</sub> and CsPbBr<sub>3</sub>

Perovskite	Defect	$D_0$ ( $\times 10^{-6}$ cm <sup>2</sup> s <sup>-1</sup> )	$E_a$ (eV)
CsPbI <sub>3</sub>	$V_I$	$0.7 \pm 0.5$	$0.34 \pm 0.04$
CsPbI <sub>3</sub>	$I_I$	$2.2 \pm 1.0$	$0.20 \pm 0.02$
CsPbBr <sub>3</sub>	$V_{Br}$	$0.1 \pm 0.2$	$0.29 \pm 0.06$
CsPbBr <sub>3</sub>	$I_{Br}$	$2.0 \pm 0.5$	$0.21 \pm 0.01$

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

