

## CORRECTION

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## Correction: Effects of mixed-valence states of Eu-doped $\text{FAPbI}_3$ perovskite crystals studied by first-principles calculation

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Correction for 'Effects of mixed-valence states of Eu-doped  $\text{FAPbI}_3$  perovskite crystals studied by first-principles calculation' by Atsushi Suzuki *et al.*, *Mater. Adv.*, 2021, DOI: 10.1039/D0MA00994F.

The authors regret that there was an error in the way Table 2 was displayed in the original manuscript, which could potentially have led to misunderstanding. The correct version of the table is given here.

Table 1 Magnetic parameters of the  $\text{MAPb}(\text{Eu})_3$  perovskite cluster model

Chemical shift	g-Tensor			EFG ( $\times 10^{-4}$ )						
	Spin	Charge (e)	Eu (ppm)	$g_{xx}$	$g_{yy}$	$g_{zz}$	$V_{xx}$	$V_{yy}$	$V_{zz}$	$\eta$
Eu(m)	0	0.62	-424 857	—	—	—	-28	-63	91	0.38
Eu(n)	7.08	0.68	-424 868	2.27070	2.27615	2.30975	-3	-5	9	0.20

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

