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CORRECTION

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Correction: Effects of mixed-valence states of Eu-doped $FAPbI_3$ perovskite crystals studied by first-principles calculation

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Correction for 'Effects of mixed-valence states of Eu-doped $FAPbI_3$ 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 $MAPb(Eu)I_3$ 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}	η
Eu(III)	0	0.62	−424 857	—	—	—	−28	−63	91	0.38
Eu(II)	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.

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