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Correction: Synthesis and characterization of novel Li-stuffed garnet-like $\text{Li}_{5+2x}\text{La}_3\text{Ta}_{2-x}\text{Gd}_x\text{O}_{12}$ ($0 \leq x \leq 0.55$): structure–property relationships

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Correction for 'Synthesis and characterization of novel Li-stuffed garnet-like $\text{Li}_{5+2x}\text{La}_3\text{Ta}_{2-x}\text{Gd}_x\text{O}_{12}$ ($0 \leq x \leq 0.55$): structure–property relationships' by Dalia M. Abdel Basset, *et al.*, *Dalton Trans.*, 2017, **46**, 933–946.

There were some errors in Li occupancy in Table 1. The corrected Table 1 is given below. The changes are shown in BLUE.

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Table 1 The powder X-ray Rietveld refinement results for $\text{Li}_{5+2x}\text{La}_3\text{Ta}_{2-x}\text{Gd}_x\text{O}_{12}$ ($0 \leq x \leq 0.55$). The model used for the Li distribution is based on ref. 25, 29, 32 and 33

	Atom	Wyckoff-site	<i>x</i>	<i>y</i>	<i>z</i>	Occupancy	<i>U</i> _{iso} (Å ²)
<i>x</i> = 0 <i>R</i> _p : 9.75, χ^2 : 1.237	Li1	24 <i>d</i>	1/4	7/8	0	0.802	0.025
	Li2	48 <i>g</i>	1/8	0.6826	0.5674	0.139	0.025
	Li3	96 <i>h</i>	0.0927	0.684	0.5795	0.147	0.025
	La	24 <i>c</i>	1/8	0	1/4	1	0.0296(1)
	Ta	16 <i>a</i>	0	0	0	1	0.0268(8)
	O	96 <i>h</i>	0.2879(9)	0.0984(1)	0.2011(2)	1	0.0192(9)
<i>x</i> = 0.15 <i>R</i> _p : 10.2, χ^2 : 3.23	Li1	24 <i>d</i>	1/4	7/8	0	0.718	0.025
	Li2	48 <i>g</i>	1/8	0.6826	0.5674	0.181	0.025
	Li3	96 <i>h</i>	0.0927	0.684	0.5795	0.173	0.025
	La	24 <i>c</i>	1/8	0	1/4	1	0.0208(8)
	Ta	16 <i>a</i>	0	0	0	0.925	0.025
	Gd	16 <i>a</i>	0	0	0	0.075	0.025
	O	96 <i>h</i>	0.2796(5)	0.1056(4)	0.1989(4)	1	0.025
<i>x</i> = 0.25 <i>R</i> _p : 9.97, χ^2 : 1.436	Li1	24 <i>d</i>	1/4	7/8	0	0.681	0.025
	Li2	48 <i>g</i>	1/8	0.6826	0.5674	0.215	0.025
	Li3	96 <i>h</i>	0.0927	0.684	0.5795	0.182	0.025
	La	24 <i>c</i>	1/8	0	1/4	1	0.0221(3)
	Ta	16 <i>a</i>	0	0	0	0.875	0.025
	Gd	16 <i>a</i>	0	0	0	0.125	0.025
	O	96 <i>h</i>	0.2870(1)	0.0949(2)	0.2007(3)	1	0.025
<i>x</i> = 0.35 <i>R</i> _p : 10.07, χ^2 : 1.48	Li1	24 <i>d</i>	1/4	7/8	0	0.678	0.025
	Li2	48 <i>g</i>	1/8	0.6826	0.5674	0.227	0.025
	Li3	96 <i>h</i>	0.0927	0.684	0.5795	0.193	0.025
	La	24 <i>c</i>	1/8	0	1/4	1	0.0197(5)
	Ta	16 <i>a</i>	0	0	0	0.825	0.0114(4)
	Gd	16 <i>a</i>	0	0	0	0.175	0.025
	O	96 <i>h</i>	0.2930(2)	0.0965(2)	0.1968(5)	1	0.025
<i>x</i> = 0.45 <i>R</i> _p : 11.34, χ^2 : 2.304	Li1	24 <i>d</i>	1/4	7/8	0	0.675	0.025
	Li2	48 <i>g</i>	1/8	0.6826	0.5674	0.228	0.025
	Li3	96 <i>h</i>	0.0927	0.684	0.5795	0.210	0.025
	La	24 <i>c</i>	1/8	0	1/4	1	0.0229
	Ta	16 <i>a</i>	0	0	0	0.775	0.0262
	Gd	16 <i>a</i>	0	0	0	0.225	0.025
	O	96 <i>h</i>	0.2851	0.1041(8)	0.2031(2)	1	0.025
<i>x</i> = 0.55 <i>R</i> _p : 13.94, χ^2 : 2.848	Li1	24 <i>d</i>	1/4	7/8	0	0.671	0.025
	Li2	48 <i>g</i>	1/8	0.6826	0.5674	0.229	0.025
	Li3	96 <i>h</i>	0.0927	0.684	0.5795	0.224	0.025
	La	24 <i>c</i>	1/8	0	1/4	1	0.0132(7)
	Ta	16 <i>a</i>	0	0	0	0.725	0.0015
	Gd	16 <i>a</i>	0	0	0	0.275	0.025
	O	96 <i>h</i>	0.2842(6)	0.0908(2)	0.1987(5)	1	0.025

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

