

Supporting Information

Effect of Composition Modulation on the Luminescence Properties of Eu^{3+} doped $\text{Li}_{1-x}\text{Ag}_x\text{Lu}(\text{MoO}_4)_2$ Solid-Solution Phosphors

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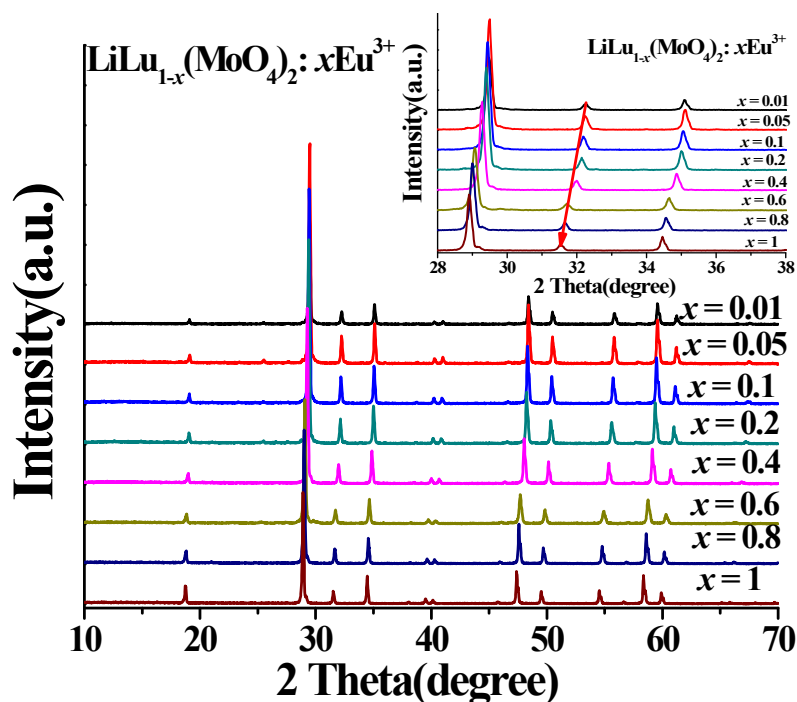
Table S1. Fractional atomic coordinates and isotropic displacement parameters ($\text{\AA}^2$) of $\text{MLn}(\text{MoO}_4)_2$ (M=Ag, Li; Ln = Lu, Eu) samples

	<i>x</i>	<i>y</i>	<i>z</i>	<i>B</i> _{iso}	Occ.
AgLu(MoO ₄) ₂					
Ag	0	0.25	0.625	0.68 (5)	0.5
Lu	0	0.25	0.625	0.68 (5)	0.5
Mo	0	0.25	0.125	0.90 (5)	1
O	0.2430 (10)	0.0946 (8)	0.0435 (4)	1.81 (15)	1
AgEu(MoO ₄) ₂					
Ag	0	0.25	0.625	0.50 (4)	0.5
Eu	0	0.25	0.625	0.50 (4)	0.5
Mo	0	0.25	0.125	0.81 (5)	1
O	0.2415 (6)	0.1017 (4)	0.04352 (18)	1.22 (10)	1
LiLu(MoO ₄) ₂					
Li	0	0.25	0.625	0.70 (8)	0.5
Lu	0	0.25	0.625	0.70 (8)	0.5
Mo	0	0.25	0.125	0.81 (8)	1
O	0.2466 (14)	0.0912 (9)	0.0398 (4)	1.17 (16)	1
LiEu(MoO ₄) ₂					
Li	0	0.25	0.625	0.66 (5)	0.5
Eu	0	0.25	0.625	0.66 (5)	0.5
Mo	0	0.25	0.125	0.69 (4)	1
O	0.2434 (5)	0.0993 (3)	0.04113 (14)	1.00 (8)	1

Table S2. Main bond lengths (Å) of $\text{MLn}(\text{MoO}_4)_2$ ($\text{M}=\text{Ag}, \text{Li}$; $\text{Ln} = \text{Lu}, \text{Eu}$) samples

$\text{AgLu}(\text{MoO}_4)_2$			
$(\text{Ag/Lu})\text{—O}^{\text{i}}$	2.410 (5)	Mo—O	1.757 (5)
$(\text{Ag/Lu})\text{—O}^{\text{ii}}$	2.469 (5)		
$\text{AgEu}(\text{MoO}_4)_2$			
$(\text{Ag/Eu})\text{—O}^{\text{i}}$	2.482 (3)	Mo—O	1.764 (3)
$(\text{Ag/Eu})\text{—O}^{\text{ii}}$	2.499 (2)		
$\text{LiLu}(\text{MoO}_4)_2$			
$(\text{Li/Lu})\text{—O}^{\text{i}}$	2.365 (6)	Mo—O	1.770 (6)
$(\text{Li/Lu})\text{—O}^{\text{ii}}$	2.381 (5)		
$\text{LiEu}(\text{MoO}_4)_2$			
$(\text{Li/Eu})\text{—O}^{\text{i}}$	2.447 (2)	Mo—O	1.767 (2)
$(\text{Li/Eu})\text{—O}^{\text{ii}}$	2.438 (2)		

Symmetry codes: (i) $-x+1/2, -y, z+1/2$; (ii) $-x+1/2, -y+1/2, -z+1/2$.

**Fig. S1** XRD patterns of $\text{LiLu}_{1-x}\text{Eu}_x(\text{MoO}_4)_2$ ($x = 0.01, 0.05, 0.1, 0.2, 0.4, 0.6, 0.8, 1.0$), and the inset shows the enlarged patterns in the range of $28\text{--}38^\circ$ highlighting the shift of the peaks.

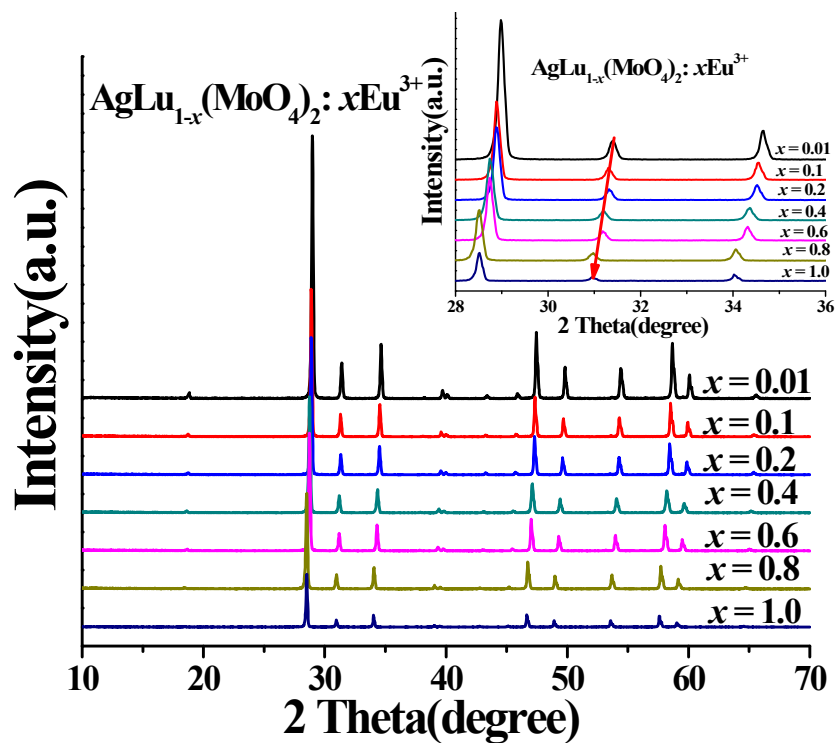


Fig. S2 XRD patterns of $\text{AgLu}_{1-x}\text{Eu}_x(\text{MoO}_4)_2$ ($x = 0.01, 0.1, 0.2, 0.4, 0.6, 0.8, 1.0$), and the inset shows the enlarged patterns in the range of $28-38^\circ$ highlighting the shift of the peaks.

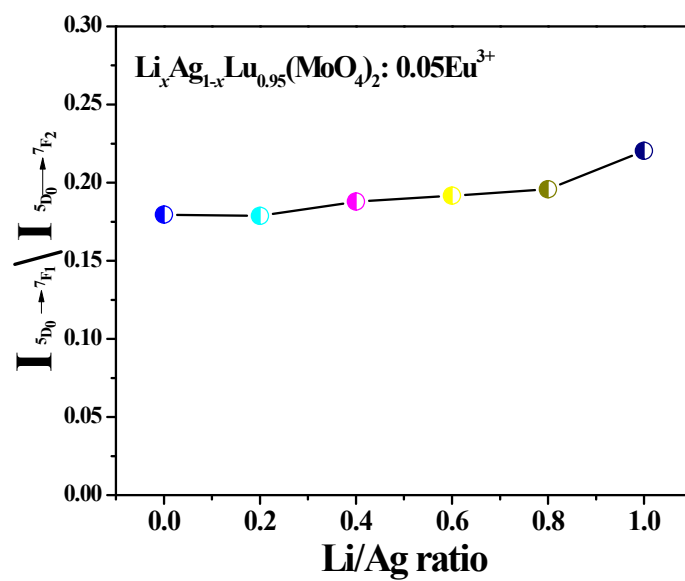


Fig. S3 Li/Ag ratio dependent $I(^5\text{D}_0-^7\text{F}_1)/I(^5\text{D}_0-^7\text{F}_2)$ of $\text{Li}_x\text{Ag}_{1-x}\text{Lu}_{0.95}(\text{MoO}_4)_2 : 0.05\text{Eu}^{3+}$ ($x = 0, 0.2, 0.4, 0.6, 0.8, 1.0$) phosphors measured at room temperature.