

Supporting Information

Title:

Solid-Solution Modulation Strategy in Trivalent Bismuth-Doped Gallate Phosphors for Single Substrate Tunable Emission

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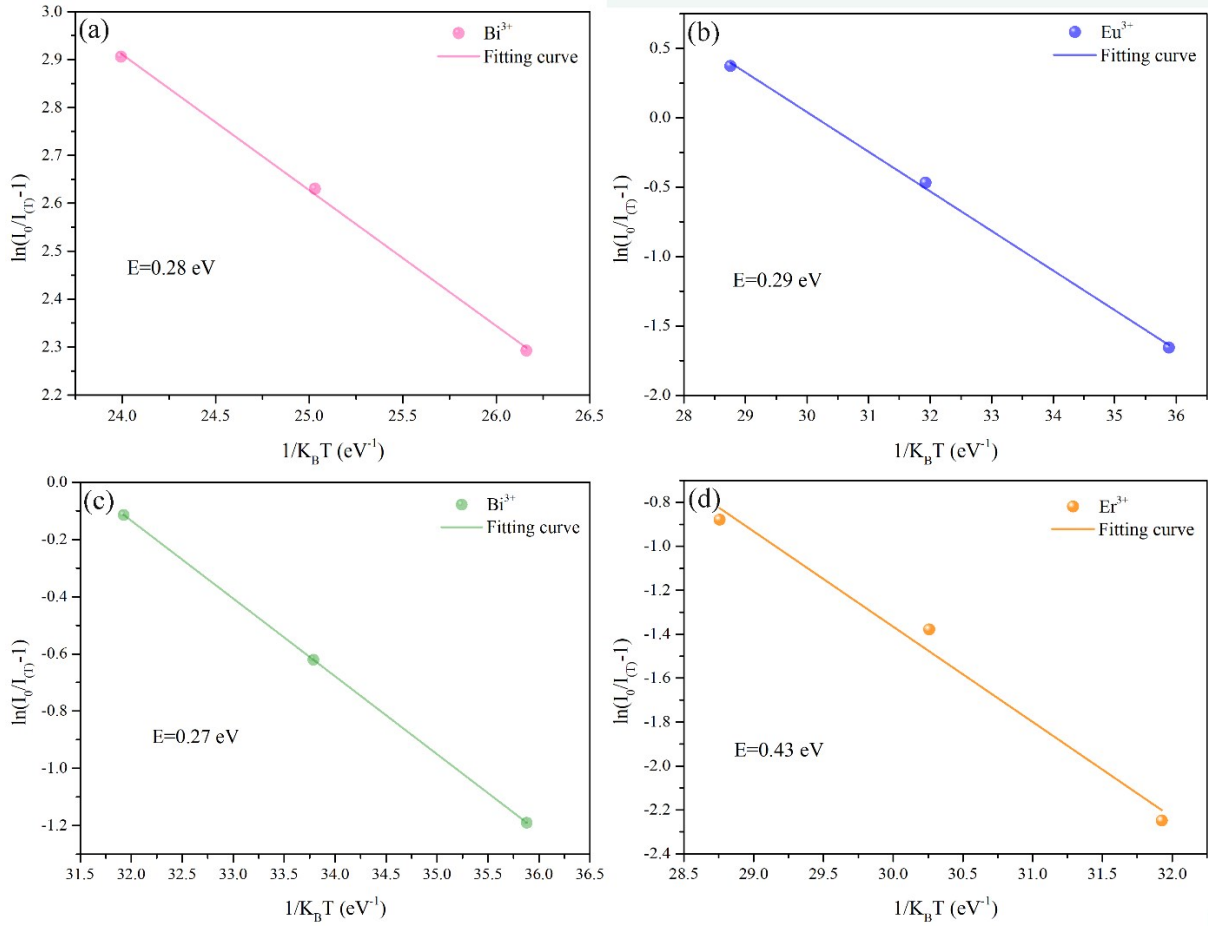


Figure S1. Thermal quenching activation energy fitting diagram of (a) Bi^{3+} , (b) Eu^{3+} in $\text{SLGF}_0:0.01\text{Bi}^{3+}, 0.01\text{Eu}^{3+}$, and (c) Bi^{3+} , (d) Er^{3+} in $\text{SLGF}_{0.6}:0.01\text{Bi}^{3+}, 0.01\text{Er}^{3+}$.

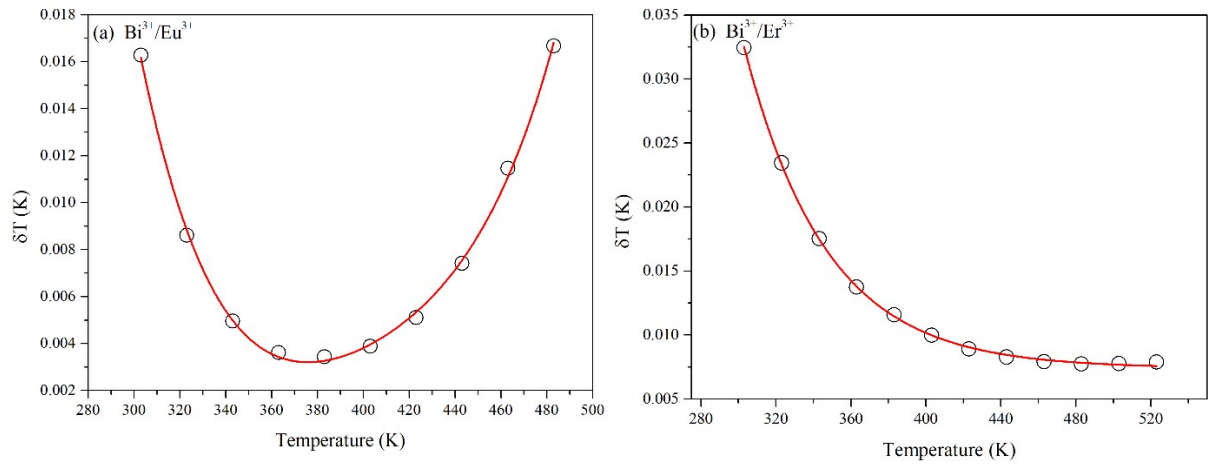


Figure S2. The temperature resolution of (a) $\text{SLGF}_0:0.01\text{Bi}^{3+}, 0.01\text{Eu}^{3+}$ phosphors, (b) $\text{SLGF}_{0.6}:0.01\text{Bi}^{3+}, 0.01\text{Er}^{3+}$ phosphors.

Table S1. Correlated data for Crystallography and Refinement Results for SLGF_{0.2}:0.01Bi³⁺ Samples.

Parameter	Sr ₃ GaO ₄ F standard crystal	SLGF _{0.2} :0.01Bi ³⁺
Space group	Pbcn(60)	I4/mcm
a (Å)	6.7819	6.8982
b (Å)	6.7819	6.8982
c (Å)	11.3662	11.2900
$\alpha=\beta=\gamma$ (deg)	90	90
V (Å ³)	522.78	537.236
Units (Z)	4	4
R _p %	-	9.12
R _{wp} %	-	11.96
χ	-	2.686

Table S2. CIE chromaticity coordinates corresponding to the points in Fig. 3 (d)

x in Sr _{1.985+x} La _{1-x} GaO _{5-x} F _x	CIE x	CIE y
0	0.2091	0.2495
0.05	0.1964	0.2187
0.1	0.1846	0.2012
0.2	0.1726	0.1635
0.4	0.165	0.137
0.6	0.1579	0.1063
0.8	0.1557	0.0925