

## Supporting Information

# **Lu<sub>2</sub>CaMg<sub>2</sub>(Si<sub>1-x</sub>Ge<sub>x</sub>)<sub>3</sub>O<sub>12</sub>: Ce<sup>3+</sup> Solid-solution Phosphors: Bandgap Engineering for Blue-light Activated Afterglow Applicable to AC-LED**

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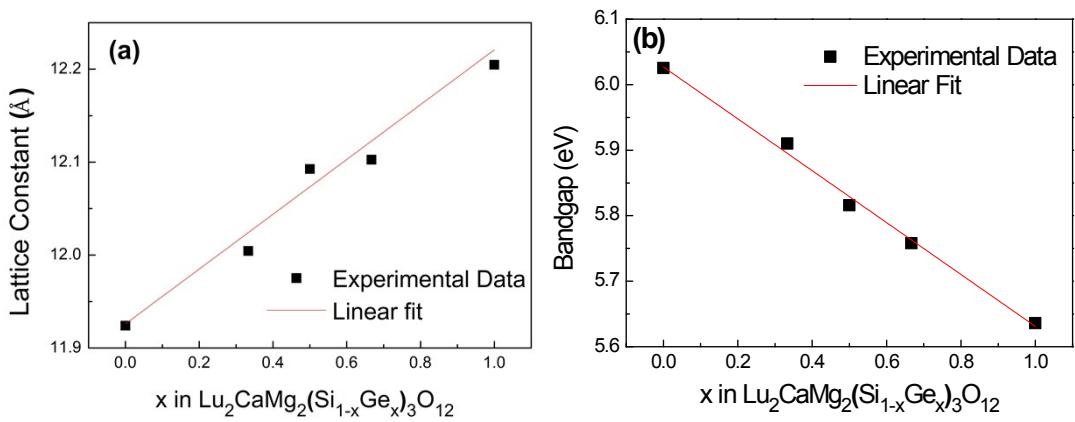
**Table S1.** Refinement results of interatomic distances in Lu<sub>1.99</sub>CaMg<sub>2</sub>Si<sub>1.5</sub>Ge<sub>1.5</sub>O<sub>12</sub> solid-solution phosphors.

Bonds	Distance (nm)	Bonds	Distance (nm)
Lu <sub>1</sub> (Ca <sub>1</sub> )-O <sub>1</sub>	0.2514	Mg <sub>1</sub> -O <sub>1</sub>	0.1979
Lu <sub>1</sub> (Ca <sub>1</sub> )-O <sub>2</sub>	0.2387	Si <sub>1</sub> (Ge <sub>1</sub> )-O <sub>1</sub>	0.1695

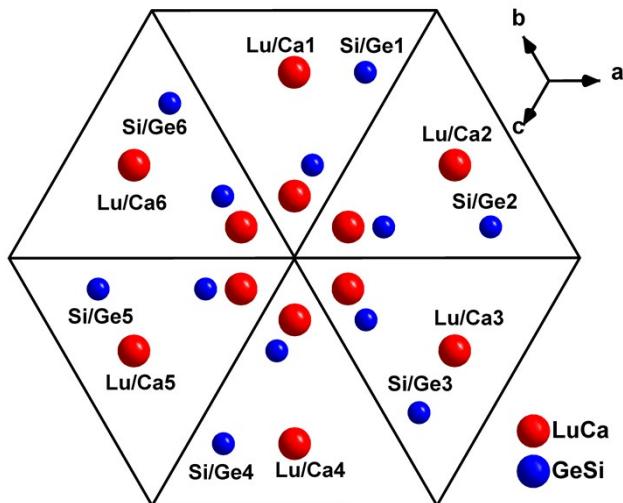
**Table S2.** Estimated thermal quenching parameters of Ce<sup>3+</sup>-doped LCMSGO solid-solution phosphors.

Sample	T <sub>0.5</sub> (K)	E <sub>q</sub> (eV)
Si:Ge=2:1	360	0.207
Si:Ge=1:1	320	0.188
Si:Ge=1:2	290	0.158
LCMGO	240	0.149

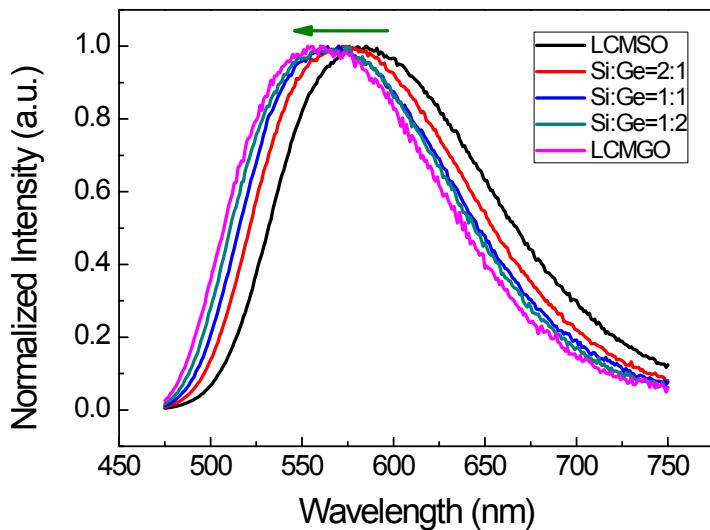
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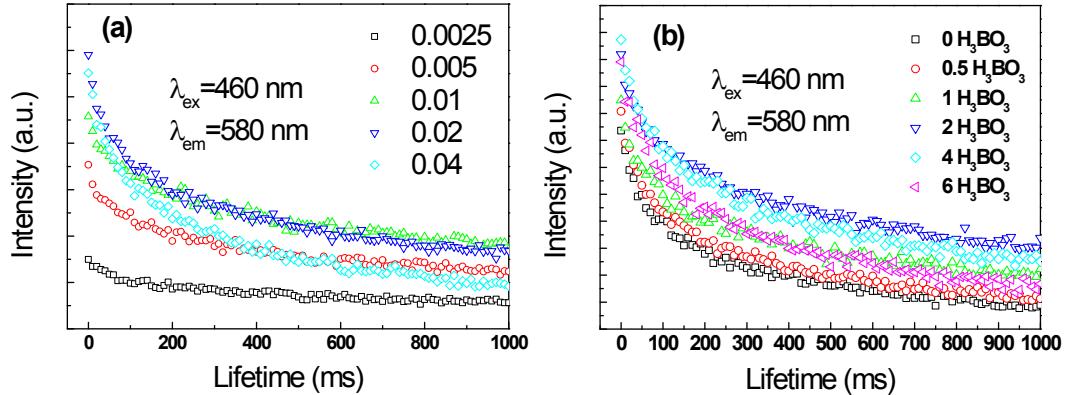
**Figure S1.** Dependence of the (a) lattice constant and (b) bandgap on  $x$  of the compounds  $\text{Lu}_2\text{CaMg}_2(\text{Si}_{1-x}\text{Ge}_x)_3\text{O}_{12}$ .



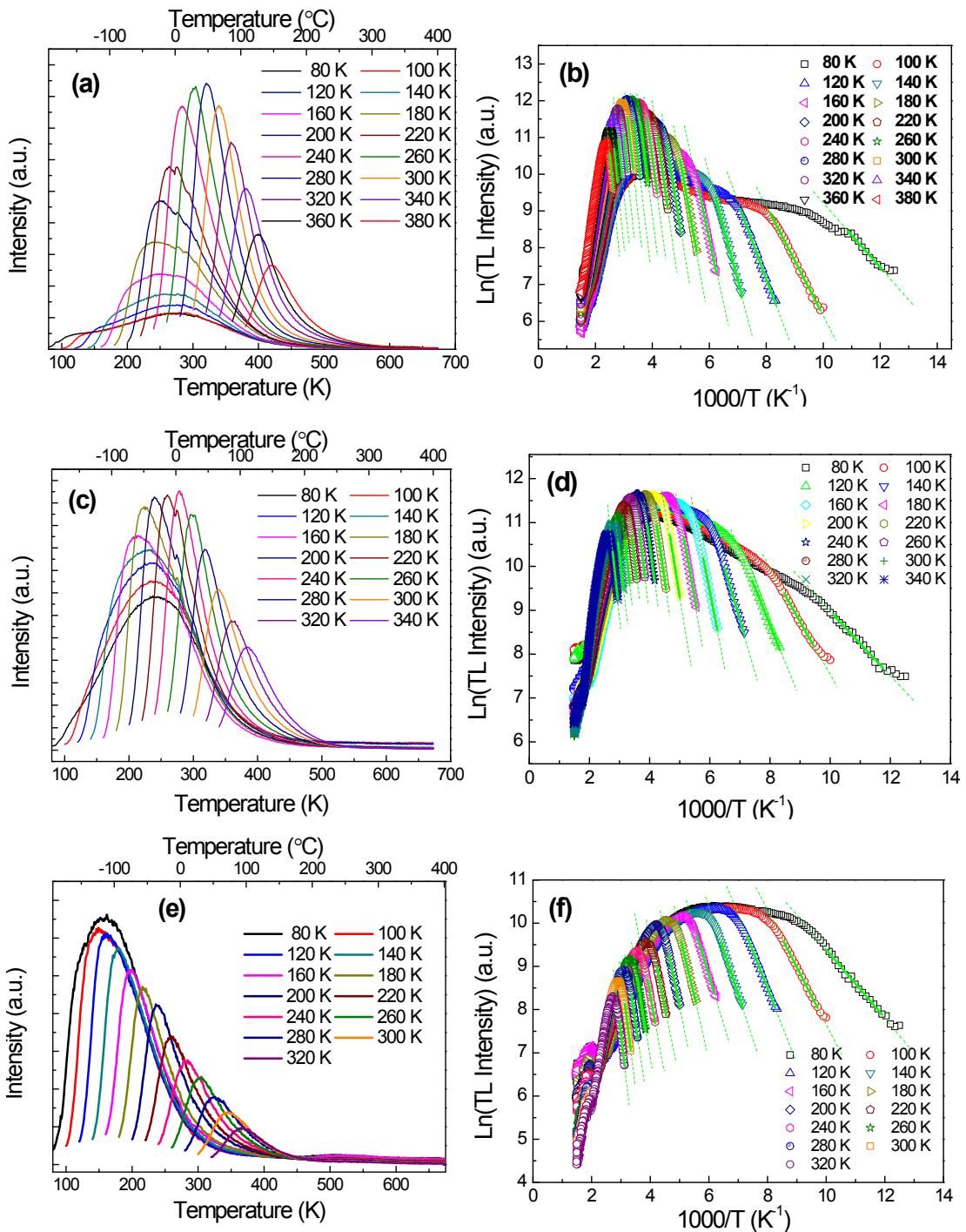
**Figure S2.** Schematic illustration of the method dealing with the Lu/Ca and Ge/Si cation disorders in DFT calculation. In order to remain the symmetry to the most extent, the Si/Ge1, Si/Ge3, and Si/Ge5 sites were set to be occupied by Si, while the Si/Ge2, Si/Ge4, and Si/Ge6 sites by Ge; likewise, the Ca/Lu1 and Ca/Lu4 sites were set to be occupied by Ca, while the Ca/Lu2, Ca/Lu3, Ca/Lu5, and Ca/Lu6 sites by Lu.



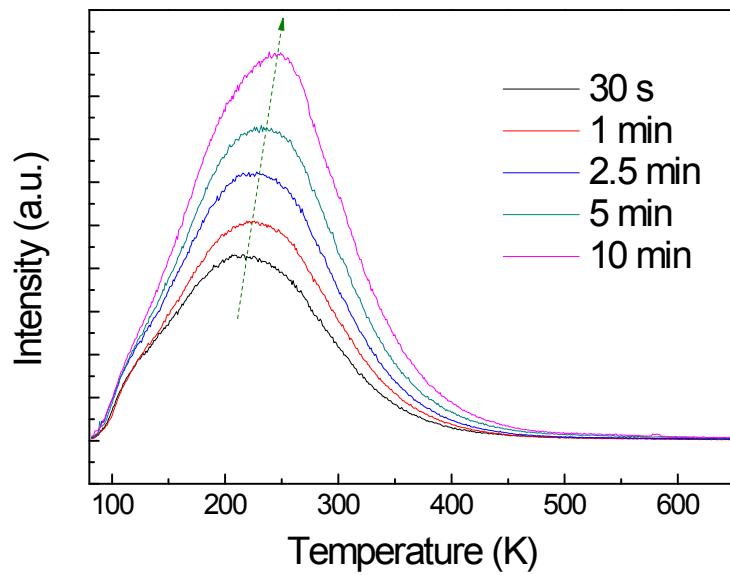
**Figure S3.** Normalized emission spectra of the  $\text{Lu}_2\text{CaMg}_2(\text{Si}_{1-x}\text{Ge}_x)_3\text{O}_{12}:0.01\text{Ce}^{3+}$  ( $x=0-1$ ) samples with various Si:Ge ratio upon 460 nm excitation.



**Figure S4.** PersL decay curves for (a) the  $\text{Lu}_2\text{CaMg}_2(\text{Si}_{1-x}\text{Ge}_x)_3\text{O}_{12}:x\text{Ce}^{3+}$  phosphors ( $x=0.005-0.04$ ) and (b) the  $\text{Lu}_2\text{CaMg}_2\text{Si}_{1.5}\text{Ge}_{1.5}\text{O}_{12}:0.01\text{Ce}^{3+}$  phosphors with various the  $\text{H}_3\text{BO}_3$  amount (0-6 wt%).



**Figure S5.** TL curves by pre-heating/cooling the (a) Si:Ge=1:1, (c) Si:Ge=1:2, and (e) LCMGO samples at different T<sub>exc</sub> (excitation wavelength: 460 nm, heating rate: 1K/s). Corresponding initial rise analyses on the TL curves of (b) Si:Ge=1:1, (d) Si:Ge=1:2, and (f) LCMGO samples as a function of T<sub>exc</sub>.



**Figure S6.** Excitation time dependent TL glow curves of the  $\text{Lu}_2\text{CaMg}_2\text{Si}_1\text{Ge}_2\text{O}_{12}:0.01\text{Ce}^{3+}$  at a heating rate of 1 K/s ( $\lambda_{\text{ex}}=460$  nm,  $\lambda_{\text{em}}=580$  nm).