Supporting Information

A polymorphic SrGa₂O₄:Eu²⁺ red phosphor for warm illumination and operando visualization of H₂O₂ catalytic reaction

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Figure S1 XRD refinements on the (a) β -SrGa₂O₄:Eu²⁺ synthesized at 1000 °C and (b) γ -SrGa₂O₄:Eu²⁺ synthesized at 1200 °C

Sample	Space group	Cell parameters(Å,°),	$R_p(\%),$
		Cell volume(Å ³)	$R_{wp}(\%)$
γ -SrGa $_2O_4$:Eu $^{2+}$ β -SrGa $_2O_4$:Eu $^{2+}$	P21/n P21/c	<i>a</i> = 8.10924	<u> </u>
		<i>b</i> = 10.75332	
		<i>c</i> =9.04985	5.60
		V=788.872717	3.94
		$\alpha = \gamma = 90$	
		$\beta = 91.5413$ a = 8.38204	
		<i>b</i> = 8.99710	
		c = 10.68064	12.76
		V=803.597830	7.87
		$\alpha = \gamma = 90$	
		$\beta = 93.9076$	

Table S1 The refined structure parameters of β -SrGa₂O₄:Eu²⁺ and γ -SrGa₂O₄:Eu²⁺ red phosphor



Figure S2 (a) PLE and PL spectra of β-SrGa₂O₄:Eu²⁺ phosphor measured at temperature 77 K, the inset gives the enlarge view of the PL spectra. (b) low temperature PL decay curves of β-SrGa₂O₄:Eu²⁺ phosphor by monitoring 580 nm and 680 nm emission



Figure S3 The fitted activation energy for the thermal quenching of β -SrGa₂O₄:Eu²⁺ red phosphor.

The activation energy ΔE_a for the thermal quenching of the Eu²⁺ emission was calculated by the following Arrhenius equation:

$$I_T / I_0 = \frac{1}{1 + A \exp\left(-\frac{\Delta E_a}{kT}\right)}$$

Where I_0 is the initial intensity, I_T is the intensity at given temperature T, A is constant, and k is the Boltzmann constant (8.62×10⁻⁵ eV/K). The thermal activation energy ΔE_a for the β-SrGa₂O₄:Eu²⁺ red phosphor was calculated to be 0.30 eV.



Figure S4 High-resolution Eu 3d XPS spectra of (a) β -SrGa₂O₄:Eu²⁺ phosphor, and (b) the compound collected after immersing β -SrGa₂O₄:Eu²⁺ phosphor in MnO₂@H₂O₂ solution for 3h.