# **Supporting Information**

Construction of thermally stable Tb<sup>3+</sup>-activated green-emitting phosphors: doping concentration and excitation wavelength dual driving strategy

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**Figure S1.** Rietveld refinement patterns for X-ray diffraction patterns of (a)  $Gd_{0.999}TaO_4:0.1\%Tb^{3+}$  (b)  $Gd_{0.995}TaO_4:0.5\%Tb^{3+}$  (c)  $Gd_{0.98}TaO_4:2\%Tb^{3+}$ .

Figure S2



Figure S2. Raman spectra of  $Gd_{1-x}TaO_4$ :  $xTb^{3+}$  (x = 0.1%, 0.5%, 2%).



**Figure S3.** Gaussian fit to the excitation spectrum of photoluminescence of  $Gd_{0.999}TaO_4:0.1\%Tb^{3+}$ . The wavelength expressed in terms of photon energy (eV) is the horizontal coordinate.  $I(E) = I(\lambda) \times \lambda^2$ .



Figure S4. Diffuse reflectance spectrum of Gd<sub>1-x</sub>TaO<sub>4</sub>:xTb<sup>3+</sup> (x =0.1%, 0.5%, 2%) and computed band gap spectrum fitted with Kubelka-Munk formula.



**Figure S5.** Contour plots of 303-523K variable temperature emission of  $Gd_{1-x}TaO_4:xTb^{3+}$  (x = 0.1%, 0.5%, 2%) phosphor at 232, 264 and 308 nm at different excitation wavelengths (a-c: x = 0.1%, d-f: x = 0.5%, g-i: x = 2%).



**Figure S6.** Excitation line of BaSO<sub>4</sub> and the emission spectrum of Gd<sub>1-x</sub>TaO<sub>4</sub>:xTb<sup>3+</sup> (x =0.1%, 0.5%, 2%) phosphor. (a)  $\lambda_{ex}$ =264nm x=0.1%, (b)  $\lambda_{ex}$ =264nm x=0.5%, (c)  $\lambda_{ex}$ =264nm x=2%, (d)  $\lambda_{ex}$ =308nm x=0.1%, (e)  $\lambda_{ex}$ =308nm x=0.5%, (f)  $\lambda_{ex}$ =308nm x=2%. The data was collected by using an integrating sphere. The inset shows a magnification of the emission spectrum from 450 nm to 700 nm.



**Figure S7 (a-c)**  $Gd_{1-x}TaO_4:xTb^{3+}$  (x = 0.1%, 0.5%, 2%) in variable temperature chromaticity coordinates (x, y), and the insets are the corresponding CIE chromaticity diagrams.



Figure S8 Photoluminescence of Gd<sub>0.999</sub>TaO<sub>4</sub>:0.1%Tb<sup>3+</sup>.



**Figure S9** Variable temperature excitation spectrum of  $Gd_{1-x}TaO_4:xTb^{3+}$  (x = 0.1%, 0.5%) under monitored excitation at 546 nm (a)x = 0.1% (b)x = 0.5%.

Table S1	Data of moles of all reagents used in the synthesis of $Gd_{1-x}TaO_4:xTb^{3+}$ (x =
0.1%, 0.5%,	2%). and masses.

Samples prepared	raw materials	relative molecular mass	mole number/mol	theoretical value/g	Actual weighing value/g
	$Gd_2O_3$	362.4982	0.0050	0.9053	0.9052
$G_{4}$ T <sub>0</sub> O ·0 1% Th <sup>3+</sup>	Ta <sub>2</sub> O <sub>5</sub>	441.8928	0.0050	1.1047	1.1048
$Gu_{0.999}$ 1 $aO_4$ : 0. 1 % 1 $D^{3}$	$Tb_4O_7$	747.6972	0.0050	0.0009	0.0009
	Li <sub>2</sub> CO <sub>3</sub>	73.8909	0.0050	0.0020	0.0022
	$Gd_2O_3$	362.4982	0.0050	0.9017	0.9017
$C_{1}^{-1}$ $T_{2}O_{2}O_{2}O_{2}^{-1}O_{2}^$	$Ta_2O_5$	441.8928	0.0050	1.1047	1.1044
$Gu_{0.995}$ 1 a $O_4$ :0.3% 10°	$Tb_4O_7$	747.6972	0.0050	0.0046	0.0048
	Li <sub>2</sub> CO <sub>3</sub>	73.8909	0.0050	0.0020	0.0022
	$Gd_2O_3$	362.4982	0.0050	0.8881	0.8882
$C_{1}^{-1}$ T <sub>2</sub> O $\cdot 2^{0/2}$ Th <sup>3+</sup>	$Ta_2O_5$	441.8928	0.0050	1.1047	1.1047
$Gu_{0.98}$ 1 a $O_4$ :2% 1 0°	$Tb_4O_7$	747.6972	0.0050	0.0187	0.0185
	Li <sub>2</sub> CO <sub>3</sub>	73.8909	0.0050	0.0020	0.0020

Parameter	$Gd_{0.999}TaO_4:0.1\%Tb^{3+}$	$Gd_{0.995}TaO_4{:}0.5\% Tb^{3+}$	$Gd_{0.98}TaO_4:2\%Tb^{3+}$
Space group	I 2/a	I 2/a	I 2/a
Structure	Monoclinic	Monoclinic	Monoclinic
a (Å)	5.3541	5.3478	5.3508
b (Å)	11.0199	11.0108	11.0083
c (Å)	5.1620	5.1579	5.1633
$\alpha = \gamma (deg)$	90.000	90.000	90.000
β(deg)	96.480	96.506	96.646
Unit cell volume $(Å^3)$	302.623	301.759	302.090
$R_p(\%)$	3.25	3.03	2.91
$R_{wp}$ (%)	4.49	4.10	3.81
$\chi^2$	1.828	1.888	1.574

**Table S2** The relevant Rietveld refinement parameters and crystallographic data.

Sample (Bond length(Å))	$Gd_{0.999}TaO_4:$ 0.1%Tb <sup>3+</sup>	$Gd_{0.995}TaO_4$ : 0.5%Tb <sup>3+</sup>	Gd <sub>0.98</sub> TaO <sub>4</sub> : 2%Tb <sup>3+</sup>
Ta-Tb <sup>3+</sup>	3.5029	3.5013	3.4972
Ta-Tb <sup>3+</sup>	3.9855	3.9457	3.9279
Ta-Tb <sup>3+</sup>	3.9230	3.9457	3.9943
Ta-Tb <sup>3+</sup>	3.9629	3.9223	3.9943
Ta-Tb <sup>3+</sup>	3.6648	3.6737	3.6320
Average	3.8078	3.7977	3.8091

**Table S3** The bond length in GSAS Refined Phosphors.

Excitation wavelength	${ m Gd}_{0.999}{ m TaO}_4$ : 0.1%Tb <sup>3+</sup> (ms)	Gd <sub>0.995</sub> TaO <sub>4</sub> : 0.5%Tb <sup>3+</sup> (ms)	Gd <sub>0.98</sub> TaO <sub>4</sub> : 2%Tb <sup>3+</sup> (ms)
232 nm	1.301	1.307	1.249
264 nm	2.425	2.330	1.165
308 nm	1.419	2.345	1.157

**Table S4** The data of the lifetime of  $Gd_{1-x}TaO_4:xTb^{3+}$  (x = 0.1%, 0.5%, 2%) under different excitations.

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Sample	$Gd_{0.999}TaO_4$ : 0.1%Tb <sup>3+</sup>	$Gd_{0.995}TaO_4:$ 0.5%Tb <sup>3+</sup>	$Gd_{0.98}TaO_4:$ 2%Tb <sup>3+</sup>
I <sub>abs</sub>	60.35%	60.71%	55.71%
IQE	12.33%	27.89%	95.29%
EQE	7.44%	16.93%	53.09%

**Table S5** The data of  $I_{abs}$ , IQE and EQE on  $Gd_{1-x}TaO_4:xTb^{3+}$  (x = 0.1%, 0.5%, 2%). ( $\lambda_{ex}=264 \text{ nm}$ )

Sample	$Gd_{0.999}TaO_4:$ 0.1%Tb <sup>3+</sup>	$Gd_{0.995}TaO_4:$ 0.5%Tb <sup>3+</sup>	$Gd_{0.98}TaO_4:$ 2%Tb <sup>3+</sup>
I <sub>abs</sub>	25.60%	32.15%	27.60%
IQE	3.45%	5.61%	10.07%
EQE	0.88%	1.80%	2.78%

**Table S6** The data of  $I_{abs}$ , IQE and EQE on  $Gd_{1-x}TaO_4$ : $xTb^{3+}$  (x = 0.1%, 0.5%, 2%). ( $\lambda_{ex}$ =308nm)