

Supplementary Information

Highly efficient near-infrared solid solution phosphors with excellent thermal stability and tunable spectra for pc-LED light source toward NIR spectroscopy applications

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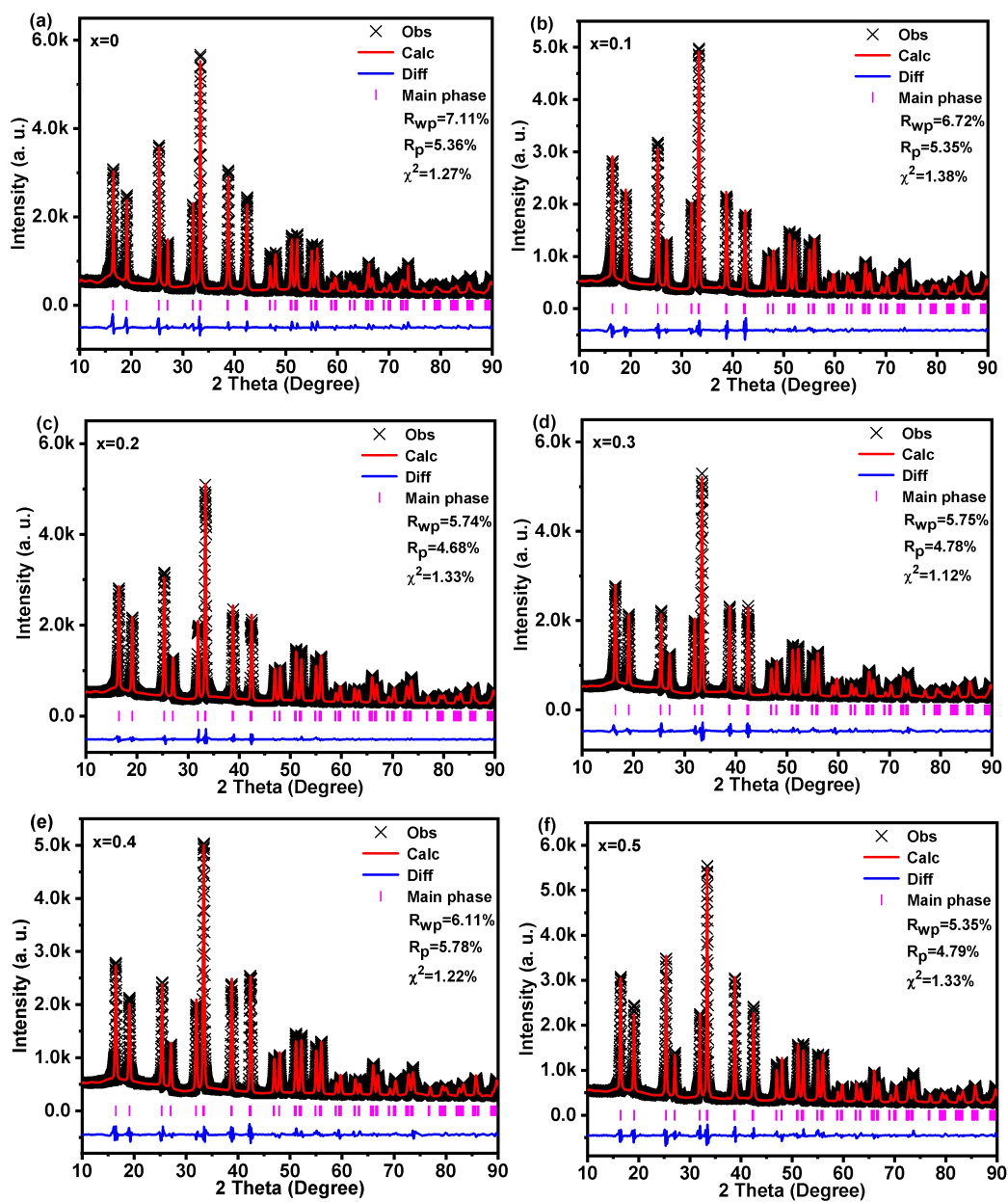


Fig.S1 Rietveld refinement of GLASB:Cr³⁺ (x=0, 0.1, 0.2, 0.3, 0.4 and 0.5).

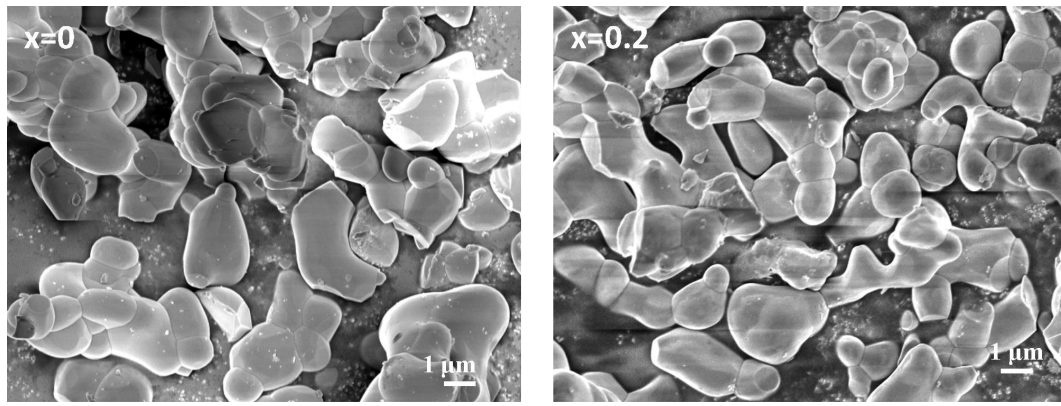


Fig. S2 SEM images of GLASB:Cr³⁺ (left, $x = 0$; and right, $x = 0.2$) with $\times 5$ kx magnifications.

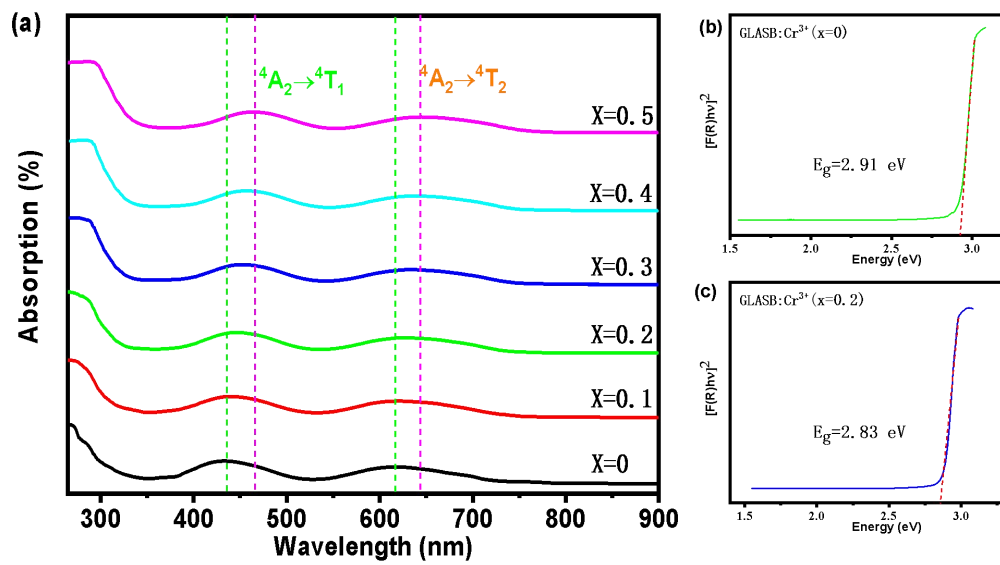


Fig. S3 (a) UV-vis-NIR absorption spectra of GLASB:Cr³⁺ (x=0-0.5). (b, c) The fitting of the corresponding band gap energies of GLASB:Cr³⁺ (x=0 and 0.2) phosphors.

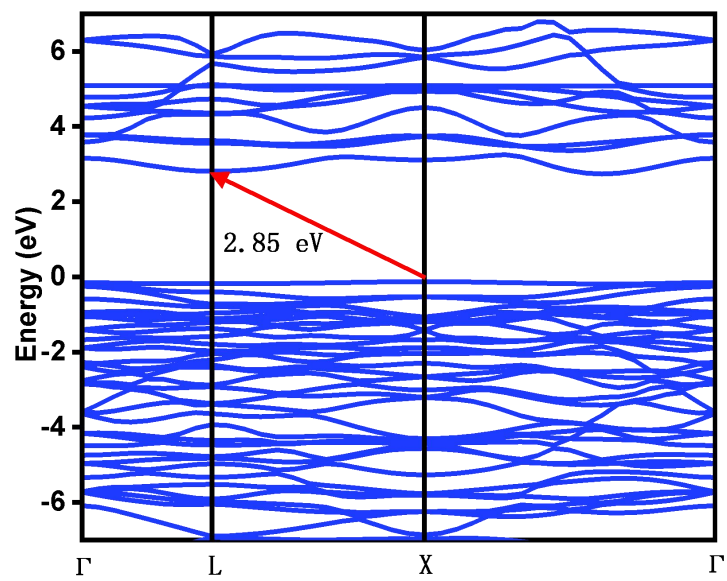


Fig. S4 Calculated electronic band structure

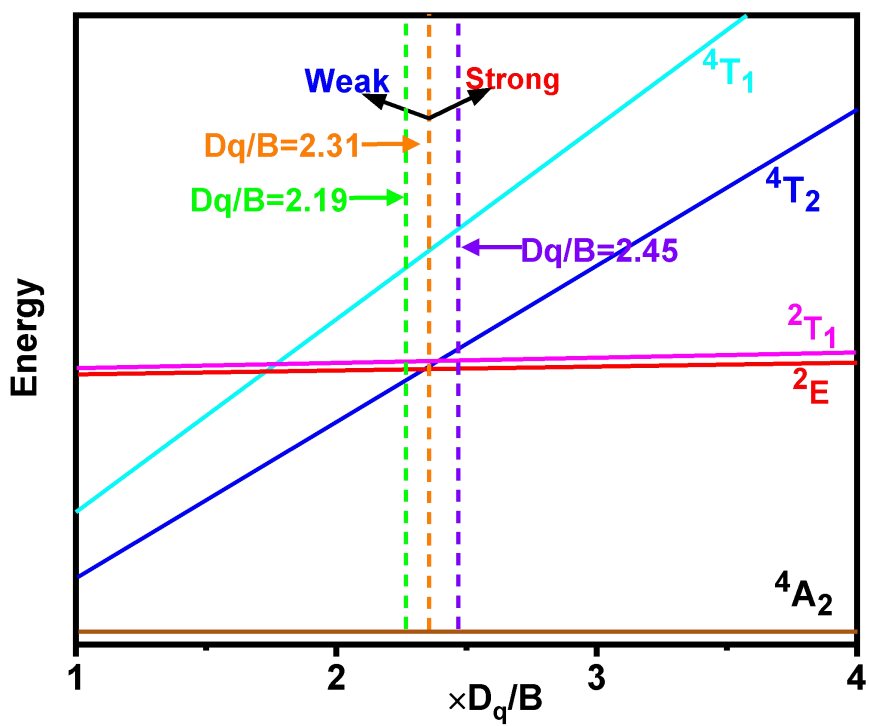


Fig. S5 Tanabe-Sugano ($3d^3$) diagram.

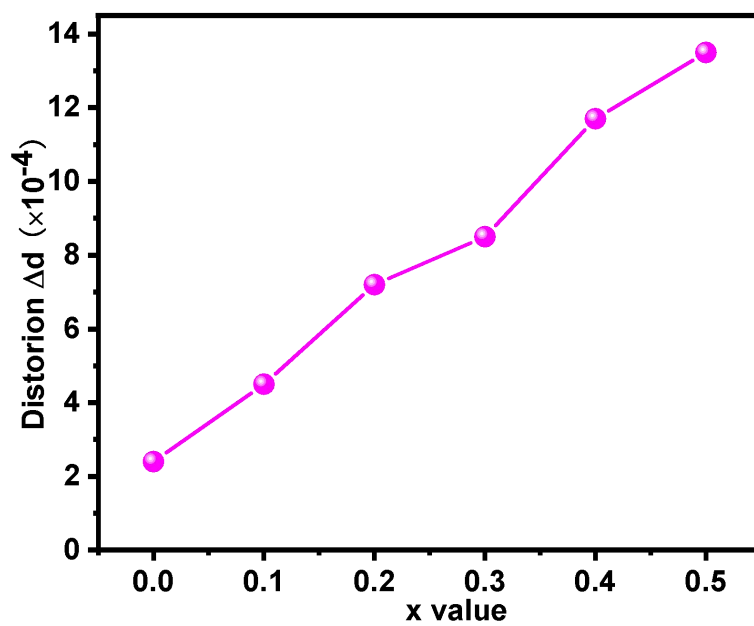


Fig. S6 The linear relationship between the octahedral distortion (Δd) and dissolved x content

Table S1. Refined Crystallographic Parameters for GLASB:Cr³⁺.

Parameters	x=0	x=0.1	x=0.2	x=0.3	x=0.4	x=0.5
Space group	Hexagonal					
$\alpha/\beta/\gamma$	90°/90°/120					
a/b/c (Å)	7.229/	7.232/	7.249/	7.267/	7.315/	7.332/
	9.313/	9.324/	9.328/	9.351/	9.386/	9.392/
	16.187	16.209	16.222	16.243	16.302	16.336
Cell volume (Å ³)	1089	1091.1	1091.8	1093.9	1095.3	1095.9
R _{wp}	7.11%	6.72%	5.74%	5.75%	6.11%	5.35%
R _p	5.36%	5.35%	4.68%	4.78%	5.78%	4.79%
χ^2	1.27%	1.38%	1.33%	1.12%	1.22%	1.33%

Table S2. Selected (Al, Sc, Cr)-O bond lengths (Å) of GLASB:Cr³⁺.

bond lengths	X=0	X=0.2
(Al, Sc, Cr)-O1	1.8672 (5)	1.8678(2)
(Al, Sc, Cr)-O2	1.9012 (4)	1.9115(3)
(Al, Sc, Cr)-O3	1.9081 (1)	1.9085(1)
(Al, Sc, Cr)-O4	1.9080 (4)	1.9083(3)
(Al, Sc, Cr)-O5	1.9009 (2)	1.9112(4)
(Al, Sc, Cr)-O6	1.8688 (3)	1.8791(5)
Average	1.8923	1.8978

Table S3. Performance parameters for GLASB:Cr³⁺ (x=0-0.5).

Compound	Peak	FWHM	$\Delta d (\times 10^{-4})$	λ_{oct}	σ_{oct}^2
x=0.0	747	128	2.31	1.0235	5.08
x=0.1	789	167	4.38	1.0265	5.17
x=0.2	806	196	7.08	1.0338	5.39
x=0.3	843	182	8.22	1.0440	5.51
x=0.4	858	179	11.57	1.0537	5.60
x=0.5	866	177	13.15	1.0654	5.76

Table S4. Luminescent parameters of Cr³⁺ in different host materials and photoelectric properties of the fabricated NIR pc-LED

Phosphor	λ_{max} (nm)	FWHM (nm)	IQE (%)	TQ at 150 °C (%)	NIR output power/efficiency of pc-LED	Ref.
GLASB:Cr ³⁺	806	196	93.6	105.7	46 mW/37% @100 mA	This work
Ca ₃ Sc ₂ Si ₃ O ₁₂ :Cr ³⁺	740	140	91	100	650-1000	[1]
Ca ₂ LuHf ₂ Al ₃ O ₁₂ :Cr ³⁺	785	145	-	-	3.1 mW/14% @120 mA	[2]
Ca ₃ Ga ₂ Ge ₃ O ₁₂ :Cr ³⁺	803	267	-	90%	27.1 mW/16.3% @100 mA	[3]
ScBO ₃ :Cr ³⁺	800	120	65	51	26 mW/7% @120 mA	[4]
La ₃ Ga ₅ GeO ₁₄ :Cr ³⁺	750	330	-	-	18.2 mW/- @350 mA	[5]
Ca ₂ LuZr ₂ Al ₃ O ₁₂ :Cr ³⁺	752	117	69	60	2.44 mW/4.1% @120 mA	[6]
NaInP ₂ O ₇ :Cr ³⁺	856	133	71.3	47	39.1 mW/10.9% @120 mA	[7]
NaScSi ₂ O ₆ :Cr ³⁺	840	140	-	-	-	[8]
LiScP ₂ O ₇ :Cr ³⁺	880	170	38	-	19 mW@100 mA	[9]
InBO ₃ :Cr ³⁺	820	138	46	-	37.50 mW@120 mA	[10]
LiInP ₂ O ₇ :Cr ³⁺	860	165	20	-	6.42 mW@100 mA	[11]
NaScGe ₂ O ₆ :Cr ³⁺	895	162	40.2	-	12.07 mW@350 mA	[12]
LiGaP ₂ O ₇ :Cr ³⁺	846	170	48	-	28.10 mW@120 mA	[13]

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