

# Supporting Information

Crystal structure refinement and luminescence properties of  
 $\text{Ce}^{3+}$  singly doped and  $\text{Ce}^{3+}/\text{Mn}^{2+}$  co-doped  $\text{KBaY}(\text{BO}_3)_2$  for  
n-UV pumped white-light-emitting diodes

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**Table S1.** The Refined Positions of All Atoms and The Lattice Parameters for A Series of  $\text{KBa}_{0.995-0.5y}\text{Y}_{0.995-0.5y}(\text{BO}_3)_2: 0.01\text{Ce}^{3+}, y\text{Mn}^{2+}$  ( $y=0, 0.01, 0.03, 0.05, \text{ and } 0.07$ )

<b>Sample</b> $\text{KBa}_{0.995}\text{Y}_{0.995}(\text{BO}_3)_2: 0.01\text{Ce}^{3+}$					
<b>Atom</b>	x	y	z	Occupancy	$U_{\text{iso}}$
<b>Y</b>	1.000000(0)	0.000000(0)	1.000000(0)	0.978(0)	0.01784
<b>Ba</b>	1.000000(0)	0.000000(0)	0.7861470(0)	0.479(0)	0.02389
<b>K</b>	1.000000(0)	0.000000(0)	0.7861470(0)	0.519(0)	0.02389
<b>Ce(1)</b>	1.000000(0)	0.000000(0)	0.7861470(0)	0.005(3)	0.02389
<b>Ce(2)</b>	1.000000(0)	0.000000(0)	1.000000(0)	0.005(0)	0.01784
<b>B</b>	0.6667000(0)	0.3333000(0)	0.7436000(0)	1.000(0)	0.02769
<b>O</b>	0.5204000(0)	0.4796000(0)	0.7449300(0)	1.000(0)	0.01512
<b>Sample</b> $\text{KBa}_{0.99}\text{Y}_{0.99}(\text{BO}_3)_2: 0.01\text{Ce}^{3+}, 0.01\text{Mn}^{2+}$					
<b>Atom</b>	x	y	z	Occupancy	$U_{\text{iso}}$
<b>Y</b>	1.000000(0)	0.000000(0)	1.000000(0)	0.959(0)	0.01990
<b>Ba</b>	1.000000(0)	0.000000(0)	0.7862240(0)	0.485(0)	0.02651
<b>K</b>	1.000000(0)	0.000000(0)	0.7862240(0)	0.504(0)	0.02651
<b>Ce(1)</b>	1.000000(0)	0.000000(0)	0.7862240(0)	0.004(0)	0.02651
<b>Ce(2)</b>	1.000000(0)	0.000000(0)	1.000000(0)	0.005(0)	0.01990
<b>Mn(1)</b>	1.000000(0)	0.000000(0)	0.7862240(0)	0.006(0)	0.02651
<b>Mn(2)</b>	1.000000(0)	0.000000(0)	1.000000(0)	0.004(0)	0.01990
<b>B</b>	0.6667000(0)	0.3333000(0)	0.7436000(0)	1.000(0)	0.02335
<b>O</b>	0.5204000(0)	0.4796000(0)	0.7449300(0)	1.000(0)	0.02713
<b>Sample</b> $\text{KBa}_{0.98}\text{Y}_{0.98}(\text{BO}_3)_2: 0.01\text{Ce}^{3+}, 0.03\text{Mn}^{2+}$					
<b>Atom</b>	x	y	z	Occupancy	$U_{\text{iso}}$
<b>Y</b>	1.000000(0)	0.000000(0)	1.000000(0)	0.964(0)	0.01699
<b>Ba</b>	1.000000(0)	0.000000(0)	0.7855820(0)	0.488(0)	0.02575
<b>K</b>	1.000000(0)	0.000000(0)	0.7855820(0)	0.515(0)	0.02575
<b>Ce(1)</b>	1.000000(0)	0.000000(0)	0.7855820(0)	0.003(0)	0.02575
<b>Ce(2)</b>	1.000000(0)	0.000000(0)	1.000000(0)	0.005(2)	0.01699
<b>Mn(1)</b>	1.000000(0)	0.000000(0)	0.7855820(0)	0.016(0)	0.02575
<b>Mn(2)</b>	1.000000(0)	0.000000(0)	1.000000(0)	0.015(0)	0.01699
<b>B</b>	0.6667000(0)	0.3333000(0)	0.7436000(0)	1.000(0)	0.00931

<b>O</b>	0.5204000(0)	0.4796000(0)	0.7449300(0)	1.000(0)	0.01460
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<b>Sample</b>	<b>KBa<sub>0.97</sub>Y<sub>0.97</sub>(BO<sub>3</sub>)<sub>2</sub>: 0.01Ce<sup>3+</sup>, 0.05Mn<sup>2+</sup></b>				
<b>Atom</b>	<b>x</b>	<b>y</b>	<b>z</b>	<b>Occupancy</b>	<b>U<sub>iso</sub></b>
<b>Y</b>	1.0000000(0)	0.0000000(0)	1.0000000(0)	0.896(0)	0.01903
<b>Ba</b>	1.0000000(0)	0.0000000(0)	0.7860870(0)	0.455(0)	0.02936
<b>K</b>	1.0000000(0)	0.0000000(0)	0.7860870(0)	0.522(0)	0.02936
<b>Ce(1)</b>	1.0000000(0)	0.0000000(0)	0.7860870(0)	0.004(0)	0.02936
<b>Ce(2)</b>	1.0000000(0)	0.0000000(0)	1.0000000(0)	0.006(0)	0.01903
<b>Mn(1)</b>	1.0000000(0)	0.0000000(0)	0.7860870(0)	0.025(4)	0.02936
<b>Mn(2)</b>	1.0000000(0)	0.0000000(0)	1.0000000(0)	0.022(0)	0.01903
<b>B</b>	0.6667000(0)	0.3333000(0)	0.7436000(0)	1.000(0)	0.03035
<b>O</b>	0.5204000(0)	0.4796000(0)	0.7449300(0)	1.000(0)	0.02843
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<b>Sample</b>	<b>KBa<sub>0.96</sub>Y<sub>0.96</sub>(BO<sub>3</sub>)<sub>2</sub>: 0.01Ce<sup>3+</sup>, 0.07Mn<sup>2+</sup></b>				
<b>Atom</b>	<b>x</b>	<b>y</b>	<b>z</b>	<b>Occupancy</b>	<b>U<sub>iso</sub></b>
<b>Y</b>	1.0000000(0)	0.0000000(0)	1.0000000(0)	0.953(0)	0.01611
<b>Ba</b>	1.0000000(0)	0.0000000(0)	0.7850790(0)	0.508(0)	0.02927
<b>K</b>	1.0000000(0)	0.0000000(0)	0.7850790(0)	0.505(0)	0.02927
<b>Ce(1)</b>	1.0000000(0)	0.0000000(0)	0.7850790(0)	0.003(0)	0.02927
<b>Ce(2)</b>	1.0000000(0)	0.0000000(0)	1.0000000(0)	0.006(0)	0.01611
<b>Mn(1)</b>	1.0000000(0)	0.0000000(0)	0.7850790(0)	0.038(0)	0.02927
<b>Mn(2)</b>	1.0000000(0)	0.0000000(0)	1.0000000(0)	0.034(0)	0.01611
<b>B</b>	0.6667000(0)	0.3333000(0)	0.7436000(0)	1.000(0)	0.02000
<b>O</b>	0.5204000(0)	0.4796000(0)	0.7449300(0)	1.000(0)	0.01633