

Structure and photoluminescence properties of a rare-earth free red-emitting Mn²⁺-activated KMgBO₃

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Table S1 Lattice parameters and fractional atomic site of $\text{KMg}_{0.91}\text{Mn}_{0.09}\text{BO}_3$

lattice		a=6.8494(1) Å		V=321.33(1) Å ³	
atom	site	x	y	z	Occ.
K	4a	0.1332(3)	0.1332(1)	0.1332(2)	1.0000
Mg	4a	0.8543(2)	0.8543(2)	0.8543(1)	0.9153
Mn	4a	0.8543(2)	0.8543(2)	0.8543(1)	0.0847
B	12b	0.3997(2)	0.3997(2)	0.3997(2)	1.0000
O	4a	0.4152(1)	0.2576(1)	0.5373(2)	1.0000

Table S2 Fractional bond length and angle of $\text{KMg}_{0.91}\text{Mn}_{0.09}\text{BO}_3$

Bonds	Distances(Å)	Bonds	Distances(Å)	Bonds	Distances(Å)
K-O ^c	2.8066(2)	Mg(Mn)-O ^b	2.0977(1)	B-O	1.3594(2)
K-O ^g	2.8066(2)	Mg(Mn)-O ^f	2.0977(1)	B-O ^d	1.3594(2)
K-O ^k	2.8066(2)	Mg(Mn)-O ⁱ	2.0977(1)	B-O ^h	1.3594(2)
K-O ^a	2.7764(1)	Mg(Mn)-O ^g	2.1561(2)	O-B-O ^d	119.898(1)
K-O ^e	2.7764(1)	Mg(Mn)-O ^c	2.1561(2)	O-B-O ^h	119.898(1)
K-O ⁱ	2.7764(1)	Mg(Mn)-O ^k	2.1561(2)	O ^d -B-O ^h	119.898(1)

(a) x,y,z; (b) -x+1/2,y,z; (c) x+1/2,-y,z; (d) -x,-y,z; (e) x,y+1/2,z+1/2; (f) -x+1/2,y+1/2,z+1/2; (g) x+1/2, -y+1/2,z+1/2; (h) -x,-y+1/2,z+1/2