Electronic Supplementary Information

Table S1. Rietveld refined crystallographic parameters of $Na_3CsMg_7(PO_4)_6:xEu^{2+}(x = xEu^{2+})$

Samples	x = 0	x = 0.06	x = 0.10	<i>x</i> = 0.14	<i>x</i> = 0.18	<i>x</i> = 0.20	<i>x</i> = 0.25				
Crystal	Monoclinic										
system											
Space	<i>C12/c1</i> (15)										
group											
a (Å)	12.762506	12.762642	12.763464	12.763513	12.763892	12.765026	12.765918				
	(187)	(144)	(181)	(143)	(172)	(160)	(148)				
b (Å)	10.720152	10.721246	10.722427	10.723418	10.723445	10.724007	10.725558				
	(105)	(107)	(110)	(140)	(135)	(119)	(145)				
c (Å)	15.545872	15.546492	15.547940	15.549800	15.550003	15.551211	15.556384				
	(194)	(191)	(200)	(236)	(216)	(246)	(249)				
$\alpha = \gamma(^{\circ})$	90	90	90	90	90	90	90				
β (°)	113	113	113	113	113	113	113				
$V(Å^3)$	1957.613	1958.598	1958.757	1958.804	1958.840	1959.178	1959.789				
2θ-interval	6-75°	6-75°	6-75°	6-75°	6-75°	6-75°	6-75°				
Ζ	16	16	16	16	16	16	16				
$R_{\rm wp}$ (%)	6.71 %	6.02 %	6.42 %	6.57 %	6.38 %	6.19 %	6.88 %				
<i>R</i> _p (%)	4.45 %	4.06 %	4.23 %	4.17 %	4.11 %	4.06 %	4.15 %				
χ^2	1.309	1.059	1.205	1.264	1.181	1.118	1.406				

0, 0.06, 0.10, 0.14, 0.18, 0.20 and 0.25) samples.

Table S2. Rietveld refined crystallographic parameters of Na₃CsMg_{7-y}(PO₄)₆: yMn²⁺ (y = 0, 0.2, 0.4, 0.6, 0.8, 1.0) and Na₃CsMg_{6.6}(PO₄)₆: $0.18Eu^{2+}/0.4Mn^{2+}$ series samples.

Samples	<i>y</i> = 0	<i>y</i> = 0.2	<i>y</i> = 0.4	<i>y</i> = 0.6	<i>y</i> = 0.8	<i>y</i> = 1.0	x = 0.18,				
Crystal				Monoclinic			y - 0.4				
system											
Space group	<i>C12/c1</i> (15)										
a (Å)	12.762506	12.770974	12.774812	12.784986	12.790604	12.804872	12.773269				
	(187)	(147)	(149)	(142)	(157)	(138)	(161)				
b (Å)	10.720152	10.725223	10.725857	10.733162	10.734987	10.739982	10.729808				
	(105)	(113)	(109)	(102)	(113)	(99)	(117)				
<i>c</i> (Å)	15.545872	15.556590	15.558581	15.576183	15.584209	15.602550	15.568131				
	(194)	(203)	(207)	(193)	(212)	(183)	(131)				
$\alpha = \gamma(^{\circ})$	90	90	90	90	90	90	90				
β (°)	113	113	113	113	113	113	113				
V (Å ³)	1957.613	1961.356	1962.185	1967.493	1969.792	1975.278	1963.588				
2θ-interval	6-75°	6-75°	6-75°	6-75°	6-75°	6-75°	6-75°				
Ζ	16	16	16	16	16	16	16				
$R_{\rm wp}$ (%)	6.71 %	6.49 %	7.07 %	6.66 %	6.17 %	5.12 %	5.46 %				
$R_{\rm p}$ (%)	4.45 %	4.40 %	4.86 %	4.60 %	4.16 %	3.41 %	3.79 %				
χ^2	1.309	1.225	1.501	1.329	1.185	0.825	0.8206				



Figure S1. (a) XRD patterns of the Na₃CsMg₇(PO₄)₆:*x*Eu²⁺(*x* = 0, 0.06, 0.10, 0.14, 0.18, 0.20 and 0.25); (b) XRD patterns of the the Na₃CsMg_{7-y}(PO₄)₆: *y*Mn²⁺ (y = 0, 0.2, 0.4, 0.6, 0.8, 1.0) and Na₃CsMg_{6.6}(PO₄)₆: 0.18Eu²⁺, 0.4Mn²⁺ series samples.



Figure S2. (a), (b) (c) (d) (e) (f) and (g) Rietveld refinements of the XRD files for Na₃CsMg₇(PO₄)₆: $xEu^{2+}(x = 0, 0.06, 0.10, 0.14, 0.18, 0.20 \text{ and } 0.25)$, respectively; (h) The unit cell volume V for Na₃CsMg₇(PO₄)₆: $xEu^{2+}(x = 0, 0.06, 0.10, 0.14, 0.18, 0.20 \text{ and } 0.25)$ samples.



Figure S3. (a), (b) (c) (d) and (e) Rietveld refinements of the XRD files for Na₃CsMg₇. $_{y}(PO_{4})_{6}$: yMn^{2+} (y = 0, 0.2, 0.4, 0.6, 0.8, 1.0), respectively; (f) The unit cell volume V for a₃CsMg_{7-y}(PO₄)₆: yMn^{2+} (y = 0, 0.2, 0.4, 0.6, 0.8, 1.0) samples.



Figure S4. SEM images of $Na_3CsMg_7(PO_4)_6:0.18Eu^{2+},0.4Mn^{2+}$ at different magnifications (a) and elemental mappings of (b) Na, (c) Cs, (d) Mg, (e) P, (f) O, (g) Eu and (h) Mn.



Figure S5. Fitting line of log(I/x) versus log(x) in Na₃CsMg_{7-y}(PO₄)₆: yMn^{2+} (y = 0,

0.2, 0.4, 0.6, 0.8, 1.0) phosphors.