

Electronic supplementary information (ESI)

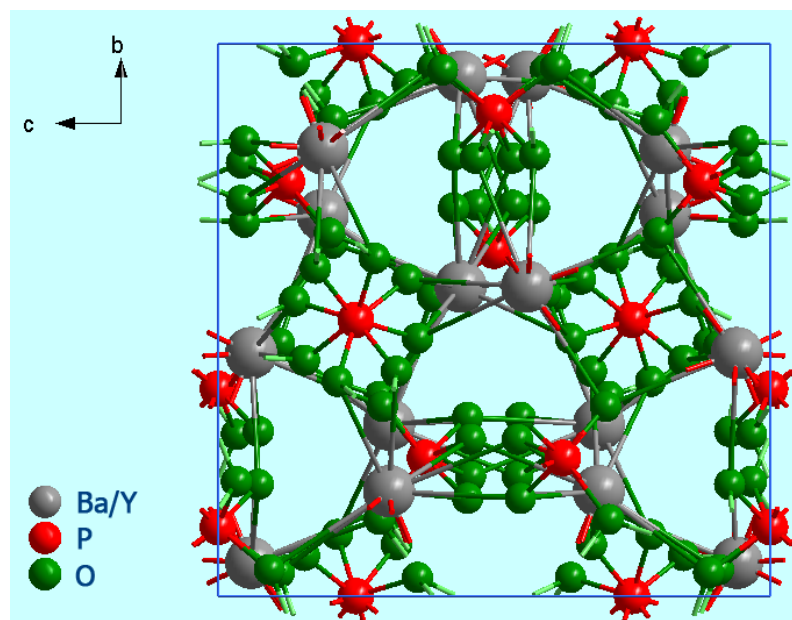


Figure S1. The crystal structure of $\text{Ba}_3\text{Y}(\text{PO}_4)_3$ viewing along $[100]$ direction.

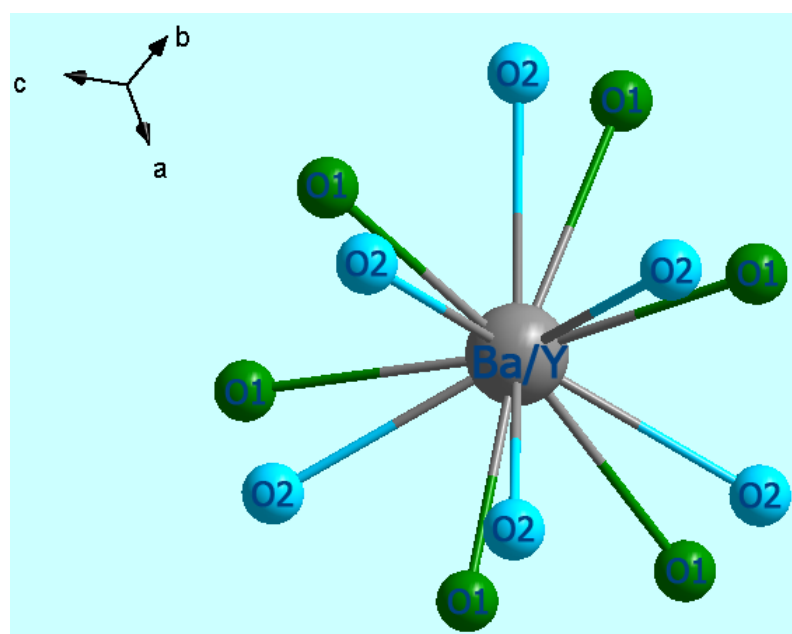


Figure S2. The coordination environment of Ba/Y viewed along $[111]$ direction.

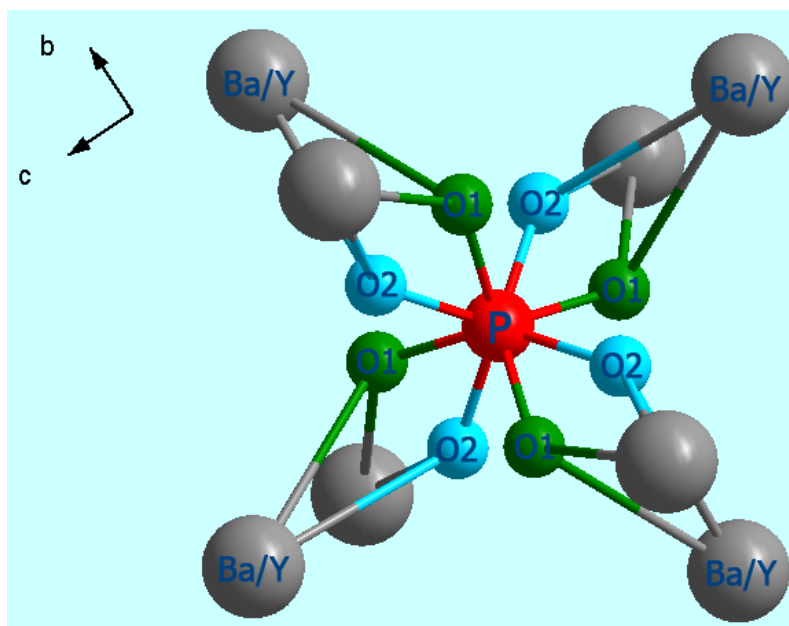


Figure S3. The coordination environment of P viewed along [100] direction.

Description of the Crystal Structure: In $\text{Ba}_3\text{Y}(\text{PO}_4)_3$ crystal, Ba^{2+} and Y^{3+} are disordered in a single crystallographic site, whereas the oxygen atoms are disordered over two orientations (O1 and O2 with 35% and 65% occupancy factors, respectively), giving rise to rotational disorder of the phosphate anion.

Table S1. Comparison of integral area of the four Gaussian components for $\text{Ba}_3\text{Ln}(\text{PO}_4)_3:\text{Eu}^{2+}$

$\text{Ba}_3\text{Ln}(\text{PO}_4)_3:\text{Eu}^{2+}$	Integral area of the four Gaussian components								
	A		B		C		D		SUM
Ln									
Lu	0.052	↑	0.775	↑	0.038		0.135	↓	1
Y	0.050		0.767		0.036		0.147		1
Gd	0.041		0.695		0.037		0.227		1