

Figure S1. IR curve of **Eu1** and **Eu2**.

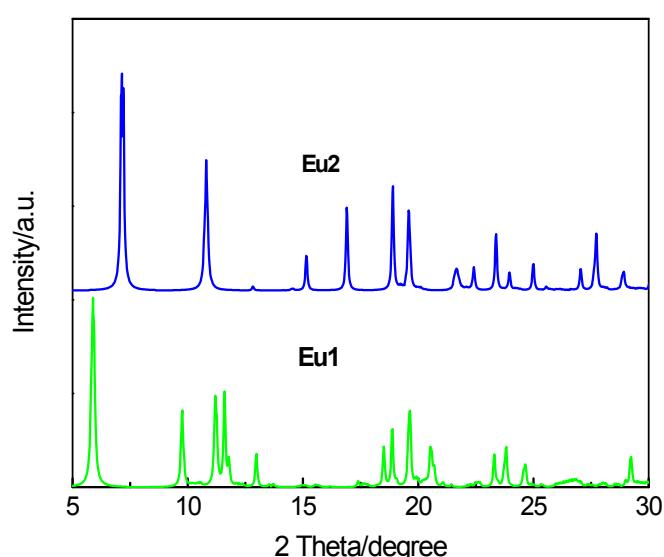


Figure S2. The XRD patterns of **Eu1** and **Eu2**.

Table S1 crystallographic data and structural refinement for **Eu1** and **Eu2**

	<b>Eu1</b>	<b>Eu2</b>
formula	EuC <sub>28</sub> H <sub>19</sub> O <sub>9</sub>	Eu <sub>2</sub> C <sub>28</sub> H <sub>23</sub> NO <sub>15</sub>
Formula weight	651.39	917.39
Temperature	293(2) K	293(2) K
Crystal system	Triclinic	Mono clinic
space group	P-1	<i>P2<sub>1</sub>/c</i>
a/nm	0.92634(19)	1.9239(4)
b/nm	0.94204(19)	1.6373(3)
c/nm	1.5549(3)	0.93961(19)
α	83.92(3)	90
β	75.00(3)	102.68(3)
γ	74.13(3)	90
Volume/nm <sup>3</sup>	1259.8(4)	2887.6(10)
Z	2	4
D <sub>c</sub> / Mg.m <sup>-3</sup>	1.717	2.110
μ/ mm <sup>-1</sup>	2.544	4.383
F(000)	644	1776
θ range/ <sup>o</sup>	2.99~27.48	3.01~27.48
GOF	1.097	1.197
Final R indices/ I>2σ(I)	R <sub>1</sub> = 0.0386 wR <sub>2</sub> = 0.0981	R <sub>1</sub> = 0.0436 wR <sub>2</sub> = 0.0932
R indices (all data)	R <sub>1</sub> = 0.0448 wR <sub>2</sub> = 0.1061	R <sub>1</sub> = 0.0561 wR <sub>2</sub> = 0.0968

$$^{\text{a}}R_1 = \sum |F_o| - |F_c| / \sum |F_o|$$

$$^{\text{b}}wR_2 = \{\sum [w(F_o^2 - F_c^2)^2] / \sum [w(F_o^2)]^2\}^{1/2}$$