

Supporting Information for

Two mono- and dinuclear Eu(III) enantiomeric pairs based on chiral bis-bidentate bridging ligands: synthesis, structures, luminescent and ferroelectric properties

Xi-Li Li,*^a Yu-Liang Gao,^a Xiang-Li Feng,^a You-Xuan Zheng,^b Chun-Lai Chen,^a Jing-Lin Zuo*^b and Shao-Ming Fang*^a

^a Henan Provincial Key Laboratory of Surface and Interface Science, Zhengzhou University of Light Industry, Zhengzhou 450002, P.R. China

E-mail: lixl@zzuli.edu.cn

^b State Key Laboratory of Coordination Chemistry, School of Chemistry and Chemical Engineering, Nanjing University, Nanjing 210093, P.R. China

E-mail: zuojl@nju.edu.cn

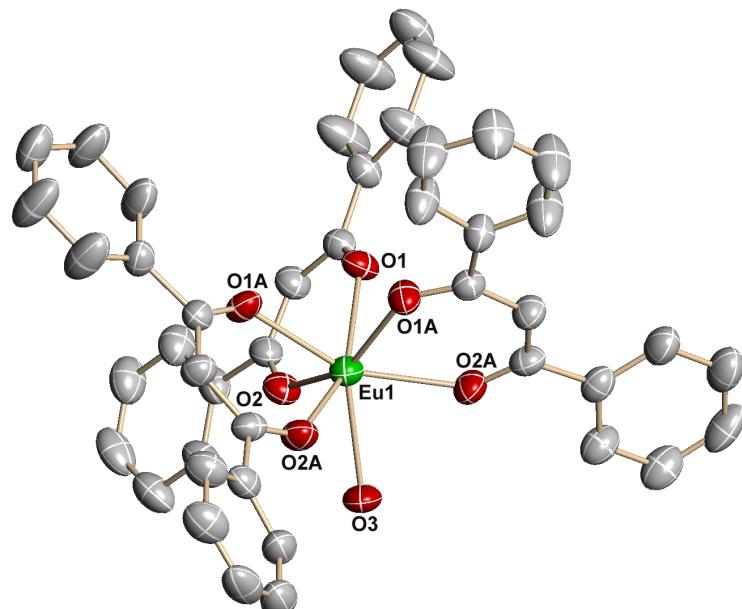


Fig. S1. The structure of Eu(dbm)₃H₂O showing 50% probability displacement ellipsoids of non-hydrogen atoms with partly atom-numbering scheme. H atoms are omitted for clarity.

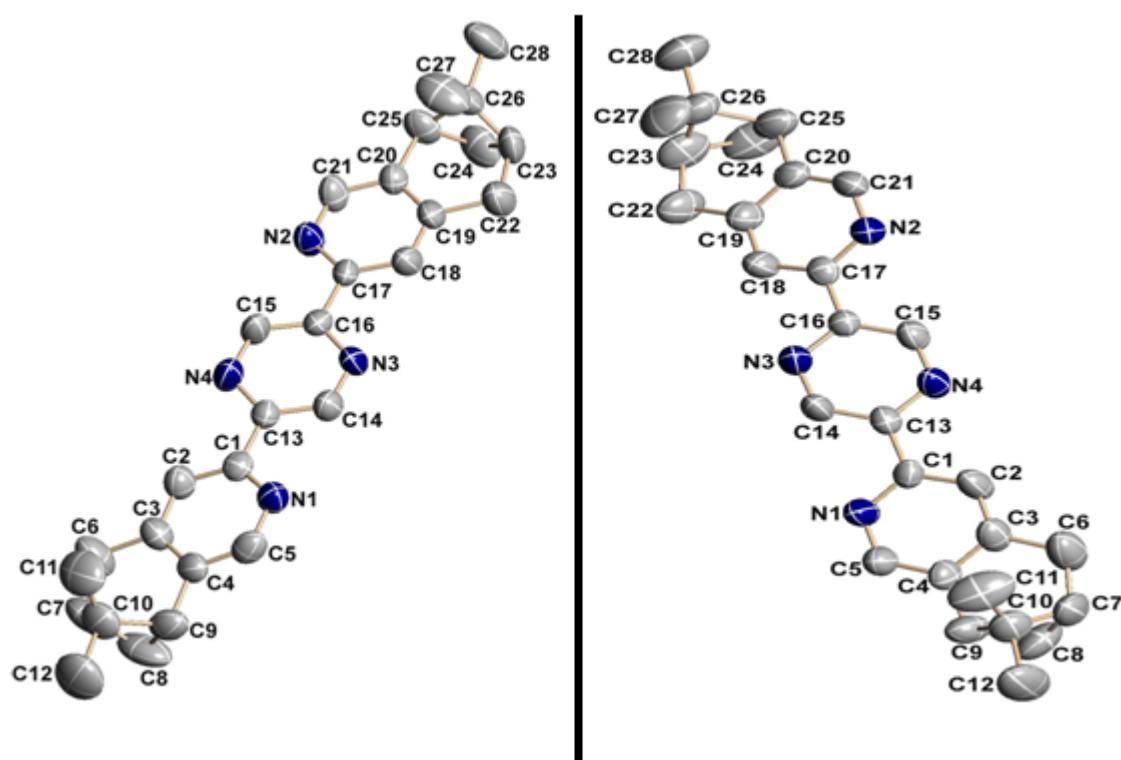


Fig. S2. The structures of \mathbf{L}_S (left) and \mathbf{L}_R (right), showing 50% probability displacement ellipsoids of non-hydrogen atoms. H atoms are omitted for clarity.

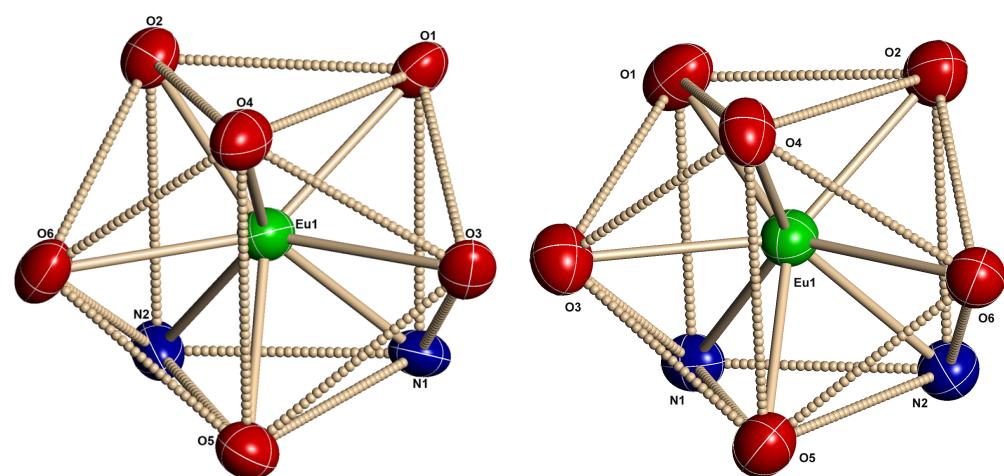


Fig. S3. Coordination geometries of $\mathbf{R}\text{-}1$ (left) and $\mathbf{S}\text{-}1$ (right) around the Eu(III) ions.

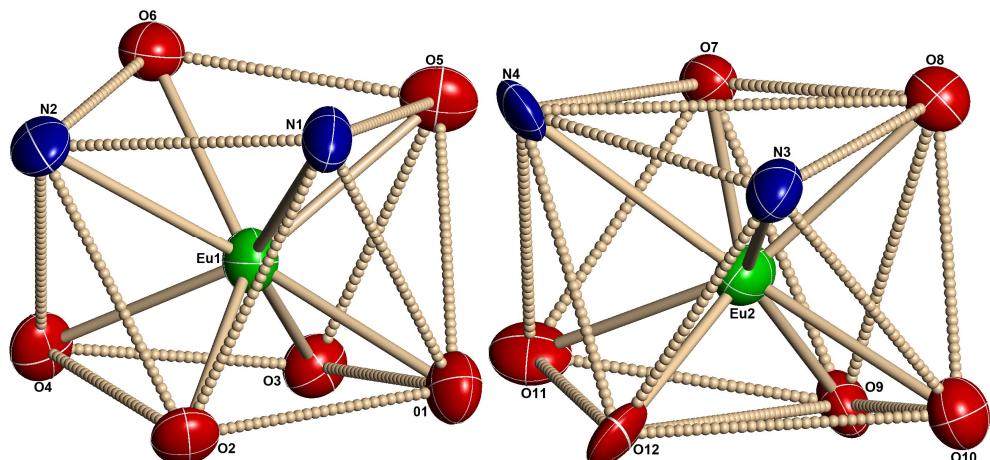


Fig. S4. Coordination geometries of **R-2** around the Eu(III) ions.

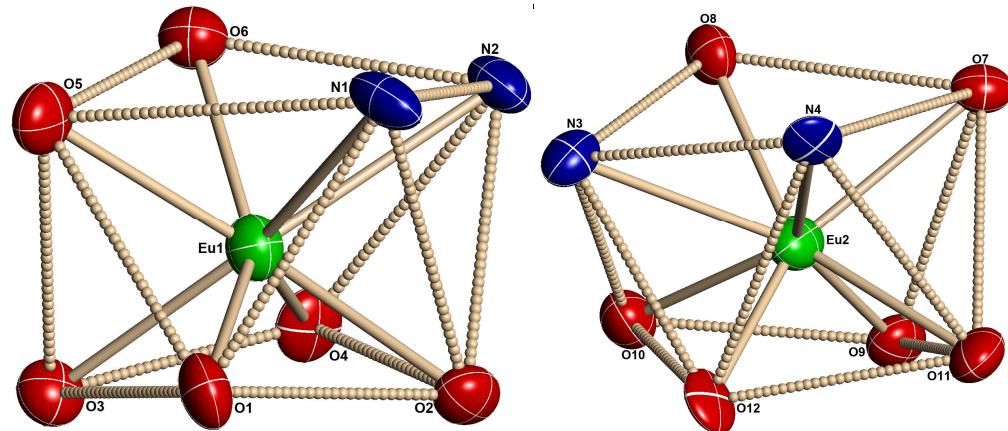


Fig. S5. Coordination geometries of **S-2** around the Eu(III) ions.

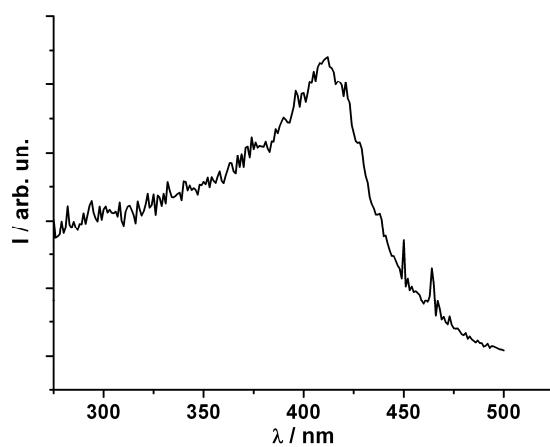


Fig. S6. The excitation spectrum for **R-1** and **R-2** in solid state.

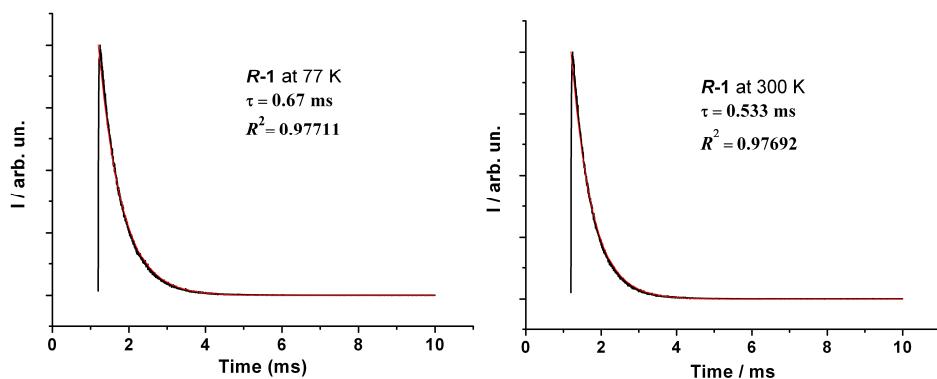


Fig. S7. Photoluminescence lifetime decay measurement of **R-1** in solid state at 77 and 300 K.

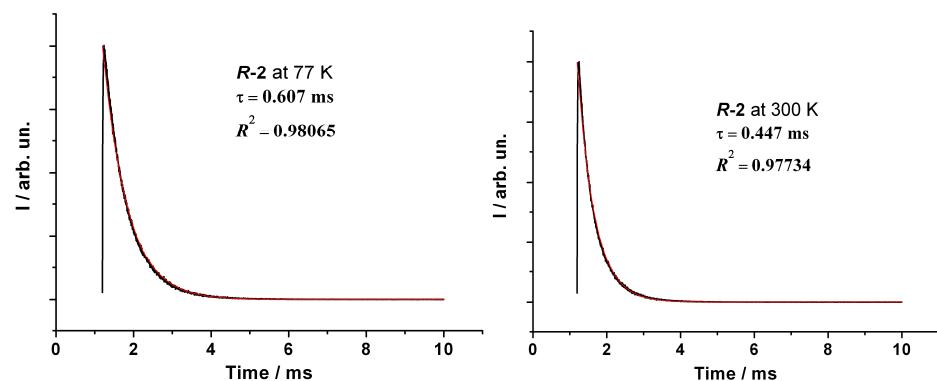


Fig. S8. Photoluminescence lifetime decay measurement of **R-2** in solid state at 77 and 300 K.

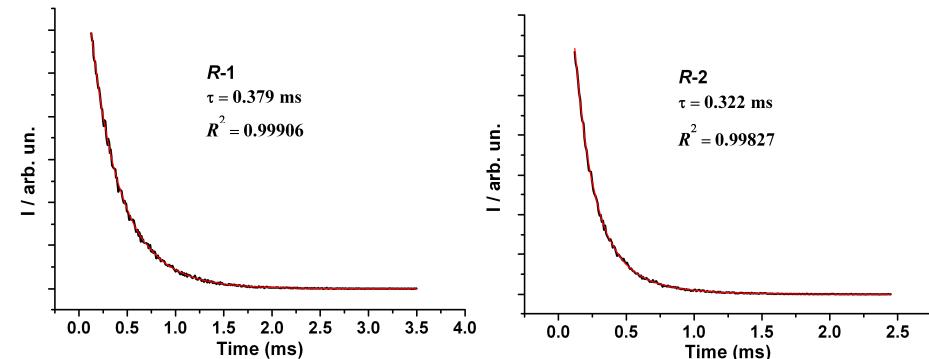


Fig. S9. Photoluminescence lifetime decay measurement of **R-1** and **R-2** in CH_2Cl_2 at room temperature.

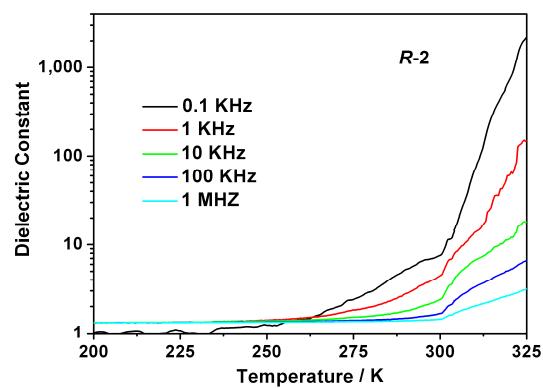


Fig. S10. The temperature dependence of the dielectric constant of **R-2** at various frequencies.

Table S1. X-ray crystallographic data for complexes L_R, L_S and Eu(dbm)₃H₂O

	L _R	L _S	Eu(dbm) ₃ H ₂ O
Chemical formula	C ₂₈ H ₃₀ N4	C ₂₈ H ₃₀ N4	C ₄₅ H ₃₅ O ₇ Eu
Formula weight	422.56	422.56	837.67
Crystal system	Monoclinic	Monoclinic	Trigonal
Space group	P2 ₁	P2 ₁	R3
a /Å	7.0760(5)	7.0752(4)	22.551(3)
b /Å	8.8805(7)	8.8818(6)	22.551(3)
c /Å	19.0843(13)	19.0839(14)	6.3892(15)
α /°	90	90	90
β /°	91.697(7)	91.796(6)	90
γ /°	90	90	120
V/Å ³	1198.70(15)	1198.65(14)	2813.8(9)
Z	2	2	3
D/g cm ⁻³	1.171	1.171	1.483
μ/mm ⁻¹	0.07	0.07	1.724
GOF	1.033	1.022	1.008
R _I ^a /wR ₂ ^b	0.0482/0.0813	0.0531/0.1149	0.0514/0.1072

^a R₁ = $\sum |Fo| - |Fc| / \sum |Fo|$. ^b wR₂ = $[\sum w(Fo^2 - Fc^2)^2 / \sum w(Fo^2)]^{1/2}$