Supporting Information for

Structures, Luminescent and Magnetic Properties of a Series of (3,6)-connected Lanthanide-Organic Frameworks

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Fig. S1 Thermogravimetric analyses (TGA) curves of compounds 1-7.

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Fig. S2 The emission spectrum of the ligand H_2 vspc.



Fig. S3 Luminescence decay curves of 2 (a), 4 (b) and 5 (c).



Fig. S4 The χ_{M}^{-1} versus *T* and the Curie-Weiss linear fit of **3**.

The data over the temperature range of 2-300 K of **3** fit well the Curie-Weiss law: $\chi_{\rm M} = C/(T-\theta)$ (Fig. S4), with Curie constant C = 7.82 cm³·K·mol⁻¹ and Weiss temperature $\theta = -0.69$ K. The negative Weiss temperature $\theta = -0.69$ K for **3** further supports the existence of antiferromagnetic coupling between Gd³⁺.



Fig. S5 Plots of $\chi_{M}' vs.T$ of **5** in a 3Oe ac field at the indicated frequency.



Fig. S6 Plots of natural logarithm of χ''/χ' vs T^{-1} of **5**; the solid lines represent the fitting in the range of 2.0-3.2 K.

Atom	Atom	Length/Å	Atom	Atom	Length/Å				
Compound 1									
Pr1	01	2.508(4)	Pr1	O 9 ¹	2.511(4)				
Pr1	02	2.438(4)	Pr1	O10	2.520(4)				
Pr1	O3 ⁶	2.384(5)	Pr1	O16 ⁵	2.362(5)				
Pr1	$O8^1$	2.635(4)	Pr1	O17 ⁴	2.394(4)				
Compound 2									
Eu1	01	2.453(5)	Eu1	$O5^3$	2.307(5)				
Eu1	02	2.367(6)	Eu1	O11 ⁴	2.463(6)				
Eu1	O3 ²	2.326(5)	Eu1	O16 ¹	2.468(6)				
Eu1	O4	2.329(6)	Eu1	O17 ¹	2.580(6)				
		Comp	oound 3						
Gd1	01	2.443(4)	Gd1	O5 ⁵	2.317(4)				
Gd1	02	2.450(4)	Gd1	O11 ⁶	2.471(4)				
Gd1	03	2.563(4)	Gd1	O16 ⁷	2.310(4)				
Gd1	O4	2.287(4)	Gd1	O17 ⁴	2.354(4)				
Compound 4									
Tb1	01	2.413(4)	Tb1	O9 ²	2.341(4)				
Tb1	O2	2.438(4)	Tb1	O10	2.269(4)				

Table S1. Selected Bond Lengths (Å) for Compounds 1-7

Tb1	O3	2.548(5)	Tb1	O11 ⁶	2.306(4)
Tb1	O8 ⁵	2.288(5)	Tb1	O18 ⁷	2.454(5)
		Comp	ound 5		
Dy1	01	2.283(4)	Dy1	$O9^2$	2.538(4)
Dy1	O2	2.409(5)	Dy1	O10 ²	2.430(4)
Dy1	O3	2.323(5)	Dy1	011	2.255(5)
Dy1	O4 ⁴	2.283(5)	Dy1	O16 ⁵	2.438(5)
		Comp	ound 6		
Ho1	01	2.306(7)	Ho1	O5	2.247(7)
Ho1	O2	2.381(7)	Ho1	O11 ⁴	2.257(7)
Ho1	O3	2.424(8)	Ho1	O16 ³	2.527(8)
Ho1	O4	2.266(7)	Ho1	O17 ³	2.438(7)
		Comp	ound 7		
Er1	01	2.239(8)	Er1	O5	2.521(8)
Er1	O2	2.419(8)	Er1	O6	2.400(7)
Er1	O3	2.258(6)	Er1	O11 ⁶	2.234(8)
Er1	O4	2.369(8)	Er1	O12 ⁵	2.288(6)

 Table S2. Selected Bond Angles (deg) for Compounds 1-7

Atom	Atom	Atom	Angle/°	Atom	Atom	Atom	Angle/°		
Compound 1									
O16 ⁵	Pr1	02	106.77(15)	O3 ⁶	Pr1	O2	89.65(15)		
O16 ⁵	Pr1	O 8 ¹	73.88(15)	O3 ⁶	Pr1	O 8 ¹	69.37(15)		
O16 ⁵	Pr1	O10	74.71(15)	O3 ⁶	Pr1	O10	141.98(16)		
O16 ⁵	Pr1	O3 ⁶	142.77(16)	O3 ⁶	Pr1	O9 ¹	77.69(15)		
O16 ⁵	Pr1	O9 ¹	74.65(14)	01	Pr1	O8 ¹	136.32(15)		
O16 ⁵	Pr1	01	146.70(15)	01	Pr1	O10	72.07(15)		
O16 ⁵	Pr1	O17 ⁴	80.93(14)	01	Pr1	O9 ¹	132.31(14)		
02	Pr1	O8 ¹	124.66(14)	O17 ⁴	Pr1	02	155.89(16)		
02	Pr1	O10	80.92(14)	O17 ⁴	Pr1	O8 ¹	79.29(14)		
02	Pr1	O9 ¹	75.54(14)	O17 ⁴	Pr1	O10	79.20(15)		
02	Pr1	01	70.19(15)	O17 ⁴	Pr1	O9 ¹	128.44(14)		
O 9 ¹	Pr1	O8 ¹	50.65(12)	O17 ⁴	Pr1	01	90.69(14)		
O9 ¹	Pr1	O10	133.54(14)	O10	Pr1	$O8^1$	144.19(15)		
Compound 2									
O5 ³	Eu1	O17 ¹	73.79(19)	O11 ⁴	Eu1	O17 ¹	143.31(19)		

O5 ³	Eu1	O3 ²	142.9(2)	O11 ⁴	Eu1	O16 ¹	133.08(19)	
O5 ³	Eu1	O16 ¹	75.28(19)	O4	Eu1	O17 ¹	78.9(2)	
O5 ³	Eu1	O11 ⁴	74.1(2)	O4	Eu1	O16 ¹	129.7(2)	
O5 ³	Eu1	O4	82.0(2)	O4	Eu1	O11 ⁴	79.5(2)	
O5 ³	Eu1	O2	105.2(2)	O4	Eu1	02	154.9(2)	
O5 ³	Eu1	01	145.5(2)	O4	Eu1	01	88.8(2)	
O3 ²	Eu1	O17 ¹	69.7(2)	02	Eu1	O17 ¹	126.1(2)	
O3 ²	Eu1	O16 ¹	77.2(2)	02	Eu1	O16 ¹	75.2(2)	
O3 ²	Eu1	O11 ⁴	142.6(2)	02	Eu1	O11 ⁴	79.6(2)	
O3 ²	Eu1	O4	97.2(2)	02	Eu1	01	71.6(2)	
O3 ²	Eu1	O2	91.1(2)	01	Eu1	O17 ¹	136.87(19)	
O3 ²	Eu1	01	71.2(2)	01	Eu1	O16 ¹	132.91(19)	
O16 ¹	Eu1	O17 ¹	51.9(2)	01	Eu1	O11 ⁴	71.50(19)	
			Comp	ound 3				
O5 ⁵	Gd1	01	87.65(15)	O16 ⁷	Gd1	O5 ⁵	96.44(14)	
O5 ⁵	Gd1	O2	130.00(14)	O16 ⁷	Gd1	01	71.33(16)	
O5 ⁵	Gd1	O11 ⁶	79.25(15)	O16 ⁷	Gd1	02	77.63(15)	
O5 ⁵	Gd1	03	79.12(14)	O16 ⁷	Gd1	O11 ⁶	142.22(17)	
O5 ⁵	Gd1	O17 ⁴	154.31(16)	O16 ⁷	Gd1	03	69.60(15)	
01	Gd1	O2	133.65(15)	O16 ⁷	Gd1	O17 ⁴	92.49(14)	
01	Gd1	O11 ⁶	71.00(16)	O4	Gd1	O5 ⁵	82.58(14)	
01	Gd1	03	136.78(14)	O4	Gd1	01	144.66(16)	
O2	Gd1	O11 ⁶	133.29(13)	O4	Gd1	02	75.44(15)	
O2	Gd1	03	51.99(12)	O4	Gd1	O11 ⁶	73.82(16)	
O17 ⁴	Gd1	01	72.47(14)	O4	Gd1	03	74.35(15)	
O17 ⁴	Gd1	O2	75.46(14)	O4	Gd1	O16 ⁷	143.40(17)	
O17 ⁴	Gd1	O11 ⁶	78.96(14)	O4	Gd1	O17 ⁴	104.12(14)	
O17 ⁴	Gd1	03	126.54(14)	O11 ⁶	Gd1	03	143.37(15)	
Compound 4								
01	Tb1	O2	133.21(15)	O9 ²	Tb1	03	126.31(16)	
01	Tb1	O3	136.76(15)	O9 ²	Tb1	O18 ⁵	79.40(16)	
01	Tb1	O18 ⁵	71.30(16)	O10	Tb1	01	145.18(17)	
02	Tb1	C1	26.23(16)	O10	Tb1	02	75.35(16)	
O2	Tb1	O3	52.25(14)	O10	Tb1	03	73.92(16)	
O2	Tb1	O18 ⁵	133.09(14)	O10	Tb1	O8 ⁶	142.83(18)	
O8 ⁶	Tb1	01	71.47(17)	O10	Tb1	O9 ²	104.67(16)	
O8 ⁶	Tb1	O2	77.00(16)	O10	Tb1	O11 ⁷	82.52(15)	

O8 ⁶	Tb1	03	69.58(17)	O10	Tb1	O18 ⁵	74.05(16)			
$O8^6$	Tb1	O9 ²	91.52(16)	O11 ⁷	Tb1	01	87.86(16)			
$O8^6$	Tb1	O11 ⁷	97.09(16)	O11 ⁷	Tb1	02	130.23(16)			
$O8^6$	Tb1	O18 ⁵	142.70(17)	O11 ⁷	Tb1	O3	79.06(16)			
$O8^6$	Tb1	01	71.47(17)	O11 ⁷	Tb1	O9 ²	154.57(18)			
O9 ²	Tb1	01	72.25(16)	O11 ⁷	Tb1	O18 ⁵	79.29(16)			
O9 ²	Tb1	02	75.01(16)	O18 ⁵	Tb1	O3	143.21(15)			
	Compound 5									
O16 ⁵	Dy1	O9 ²	O16 ⁵	01	Dy1	O16 ⁵	79.31(16)			
011	Dy1	O16 ⁵	73.86(17)	01	Dy1	O10 ²	130.37(15)			
011	Dy1	O10 ²	75.16(16)	01	Dy1	O3	154.33(17)			
011	Dy1	03	103.97(17)	01	Dy1	O9 ²	78.79(15)			
011	Dy1	O9 ²	73.91(17)	01	Dy1	$O4^4$	96.65(17)			
011	Dy1	01	83.01(17)	01	Dy1	02	87.94(16)			
011	Dy1	$O4^4$	142.71(18)	$O4^4$	Dy1	O16 ⁵	142.97(17)			
011	Dy1	02	145.78(17)	$O4^4$	Dy1	O10 ²	77.12(17)			
O10 ²	Dy1	O16 ⁵	132.93(16)	$O4^4$	Dy1	O3	92.34(17)			
O10 ²	Dy1	O9 ²	52.61(14)	$O4^4$	Dy1	O9 ²	69.52(17)			
O3	Dy1	O16 ⁵	79.06(16)	$O4^4$	Dy1	02	71.03(17)			
O3	Dy1	O10 ²	75.09(16)	02	Dy1	O16 ⁵	72.04(17)			
O3	Dy1	O9 ²	126.85(16)	02	Dy1	O10 ²	132.72(16)			
O3	Dy1	02	72.33(17)	02	Dy1	O9 ²	136.34(17)			
			Comp	oound 6						
01	Ho1	02	73.0(3)	05	Ho1	01	103.6(3)			
01	Ho1	03	78.5(3)	05	Ho1	O2	145.1(3)			
01	Ho1	O16 ³	126.9(3)	05	Ho1	O3	73.7(3)			
01	Ho1	O17 ³	74.8(3)	05	Ho1	O4	82.7(3)			
O2	Ho1	03	71.6(3)	05	Ho1	O11 ⁴	143.5(3)			
O2	Ho1	O16 ³	136.4(2)	05	Ho1	O16 ³	74.1(3)			
O2	Ho1	O17 ³	132.9(3)	05	Ho1	O17 ³	75.8(3)			
O3	Ho1	O16 ³	143.1(3)	O11 ⁴	Ho1	01	92.5(3)			
O3	Ho1	O17 ³	132.9(2)	O11 ⁴	Ho1	O2	70.8(3)			
O4	Ho1	01	154.5(3)	O11 ⁴	Ho1	O3	142.4(3)			
O4	Ho1	02	87.6(3)	O11 ⁴	Ho1	O4	96.7(3)			
O4	Ho1	03	79.7(3)	O11 ⁴	Ho1	O16 ³	70.0(3)			
04	Ho1	O16 ³	78.7(3)	O11 ⁴	Ho1	O17 ³	77.2(3)			
O4	Ho1	O17 ³	130.5(3)	O17 ³	Ho1	O16 ³	52.7(2)			

Compound 7								
O1	Er1	02	74.0(3)	O4	Er1	O6	133.5(3)	
01	Er1	03	83.2(3)	06	Er1	02	133.1(3)	
01	Er1	O4	144.9(3)	06	Er1	O5	52.6(2)	
01	Er1	05	73.9(3)	O11 ⁶	Er1	O1	142.8(3)	
01	Er1	06	75.6(3)	O11 ⁶	Er1	02	142.8(3)	
01	Er1	O12 ⁵	103.2(3)	O11 ⁶	Er1	O3	96.5(3)	
O2	Er1	05	143.1(3)	O11 ⁶	Er1	O4	71.7(3)	
O3	Er1	O2	79.8(3)	O11 ⁶	Er1	05	69.6(3)	
O3	Er1	O4	86.8(3)	O11 ⁶	Er1	O6	76.7(3)	
O3	Er1	05	78.8(3)	O11 ⁶	Er1	O12 ⁵	93.0(3)	
O3	Er1	06	130.5(3)	O12 ⁵	Er1	02	78.4(3)	
O3	Er1	O12 ⁵	154.5(3)	O12 ⁵	Er1	O4	73.8(3)	
O4	Er1	02	71.1(3)	O12 ⁵	Er1	05	126.6(3)	
O4	Er1	05	136.7(3)	O12 ⁵	Er1	O6	74.7(3)	