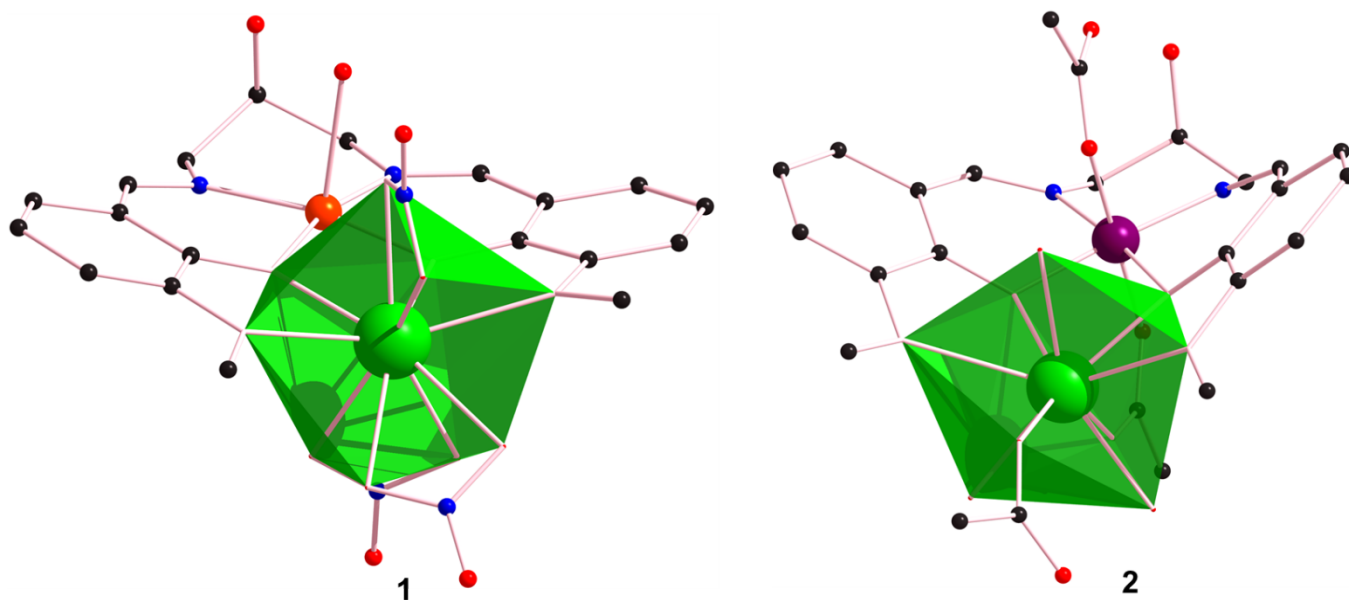


Electronic Supplementary Information (ESI)  
**Cu(II)-Dy(III) and Co(III)-Dy(III) based single molecule magnets with multiple slow magnetic relaxation processes in Cu(II)-Dy(III) complex**

Malay Dolai, Mahammad Ali, Ján Titiš and Roman Boča



**Figure S1.** Distorted bicapped square antiprism geometry and monocapped square antiprism geometry of the complex **1** and **2** respectively.

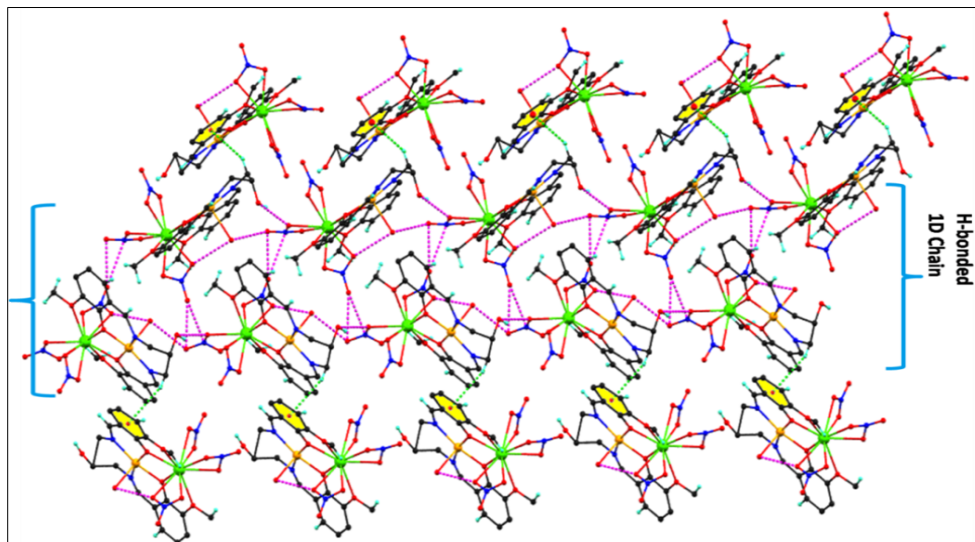
#### **H-bonding, $\pi$ - $\pi$ and C-H $\cdots$ $\pi$ stacking interactions**

The non-covalent interactions e.g. hydrogen bonding,  $\pi$ - $\pi$  and C-H $\cdots$  $\pi$  stacking have been taken into account to clarification of possible supramolecular topologies. The nature of the supramolecular interaction in which the non-coordinated alcoholic -OH group and one of the three coordinated nitrate ion along with the axial aqua ligand bonded to the Cu(II) centre, are involved in the hydrogen bonding to assembled the supramolecular 1D chains along the crystallographic *a* axis (Figure S2) and these 1D chains are moreover extended to 2D network via the inter-locking of C-H $\cdots$  $\pi$  stacking among themselves in the crystallographic *bc* plane in complex **1**. On the other hand, this 2D framework with the combination of H-bonding as well as C-H $\cdots$  $\pi$  stacking extended to form 3D network along crystallographic *a* axis (Figure S3) in complex **1**.

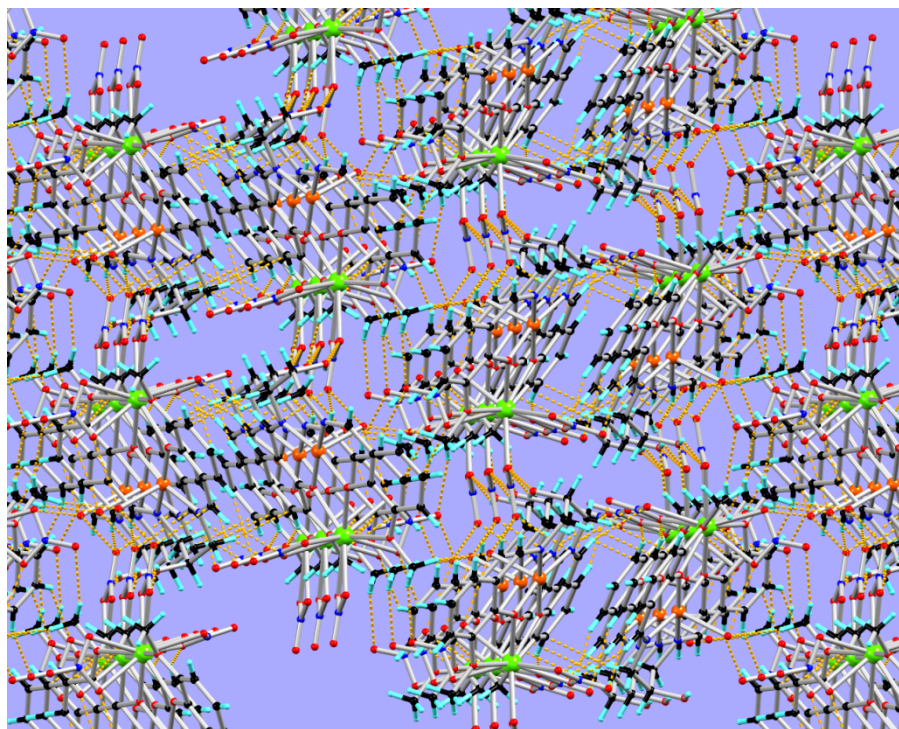
Simultaneously in complex **2**, the non-coordinated alcoholic -OH group and three coordinated aqua ligands along with the axial acetate ligand to the Co(III) centre, are involved in the hydrogen bonding to assembled the supramolecular 1D chains along the crystallographic *c* axis (Figure S4) and these 1D chains are moreover extended to 2D network via the inter-locking of  $\pi$ - $\pi$  stacking among themselves in the crystallographic *ab* plane in complex **1**. On the other hand, this 2D framework with the combination of H-bonding as well as  $\pi$ - $\pi$  stacking extended to form 3D network along crystallographic *a* axis (Figure S5) in complex **1**.

Electronic Supplementary Information (ESI)  
**Cu(II)-Dy(III) and Co(III)-Dy(III) based single molecule magnets with multiple slow magnetic relaxation processes in Cu(II)-Dy(III) complex**

Malay Dolai, Mohammad Ali, Ján Titiš and Roman Boča



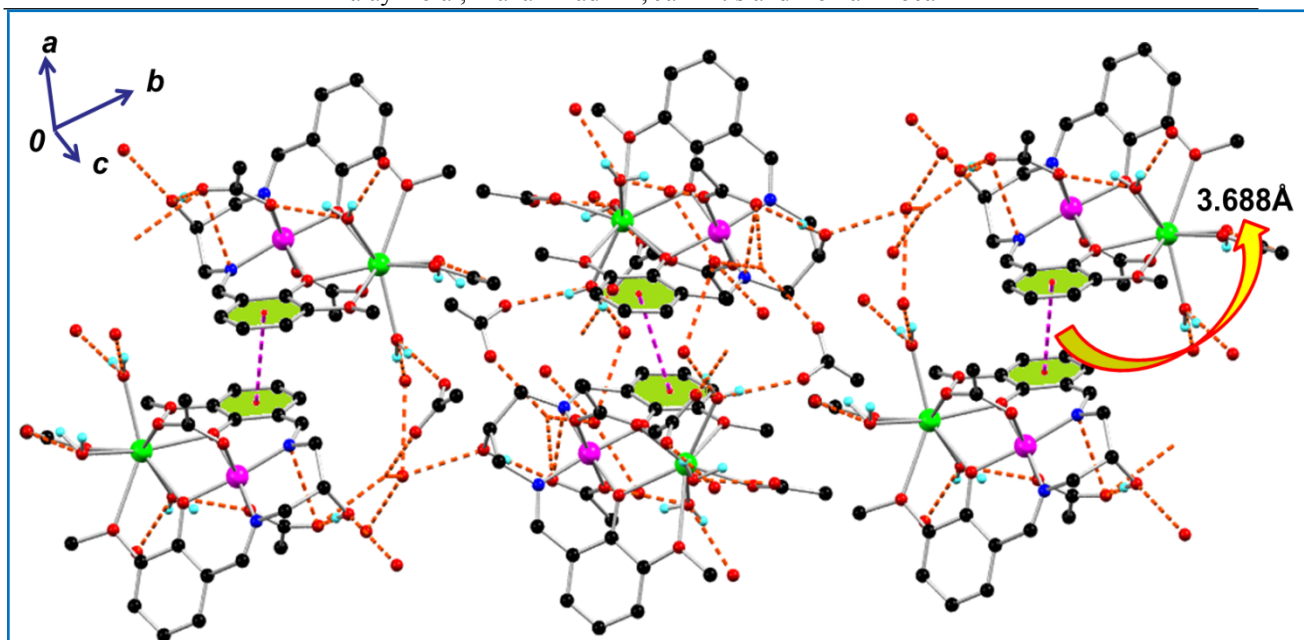
**Figure S2.** The H bonding interaction (magenta color dash line) as well as CH- $\pi$  interaction (green color dash line) leads to formation of 2D supramolecular framework in crystallographic *bc* plane in complex **1**.



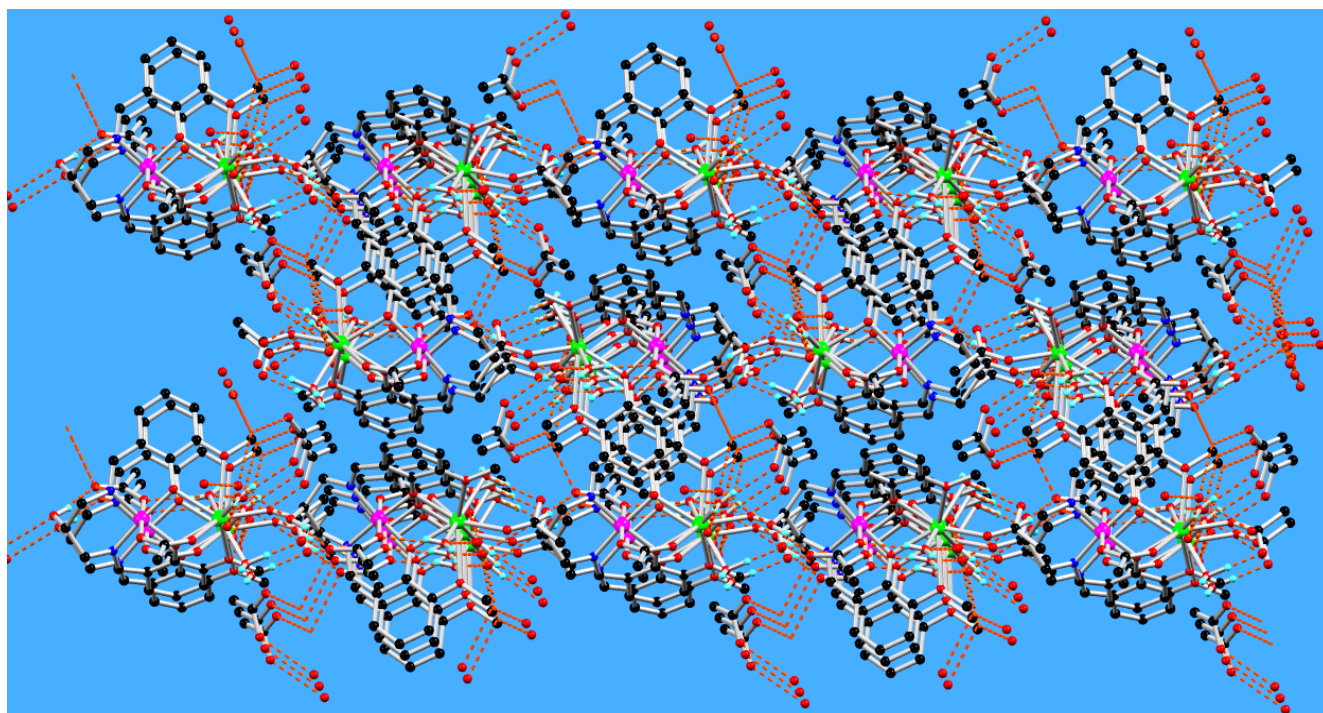
**Figure S3.** The H bonding interaction as well as C-H... $\pi$  interaction leads to formation of 3D supramolecular network along crystallographic *a* axis in complex **1**.

Electronic Supplementary Information (ESI)  
**Cu(II)-Dy(III) and Co(III)-Dy(III) based single molecule magnets with multiple slow magnetic relaxation processes in Cu(II)-Dy(III) complex**

Malay Dolai, Mohammad Ali, Ján Titiš and Roman Boča



**Figure S4.** The H bonding interactions (brown color dash line) as well as  $\pi$ - $\pi$  interaction (magenta color dash line) leads to formation of 2D supramolecular framework in crystallographic  $ab$  plane in complex 2.



**Figure S5.** The H bonding interaction as well as  $\pi$ - $\pi$  interaction leads to the formation of 3D supramolecular network along crystallographic  $a$  axis in complex 2.



**Cu(II)-Dy(III) and Co(III)-Dy(III) based single molecule magnets with multiple slow magnetic relaxation processes in Cu(II)-Dy(III) complex**

Malay Dolai, Mohammad Ali, Ján Titiš and Roman Boča

**Table S1.** Selected bond lengths (Å) and bond angles (°) of complex **1**

Dy1 -O1	2.502(5)	O1 -Dy1-O2	64.85(16)
Dy1 -O2	2.345(5)	O1 -Dy1-O3	127.15(17)
Dy1 -O3	2.343(5)	O1 -Dy1-O4	147.06(17)
Dy1 -O4	2.494(5)	O1 -Dy1 -O7	77.03(19)
Dy1 -O7	2.458(5)	O1 -Dy1 -O8	71.86(18)
Dy1 -O8	2.452(5)	O1 -Dy1 -O10	132.2(3)
Dy1 -O10	2.461(9)	O1 -Dy1 -O11	84.0(3)
Dy1 -O11	2.478(10)	O1 -Dy1 -O13	71.5(2)
Dy1 -O13	2.491(7)	O1 -Dy1 -O14	108.9(2)
Dy1 -O14	2.599(8)	O2 -Dy1 -O3	63.90(17)
Cu1 -O2	1.963(5)	O2 -Dy1 -O4	125.03(16)
Cu1 -O3	1.946(5)	O2 -Dy1 -O7	72.65(18)
Cu1 -O5	2.329(7)	O2 -Dy1 -O8	114.92(17)
Cu1 -N1	1.999(6)	O2 -Dy1 -O10	107.4(3)
Cu1 -N2	1.980(6)	O2 -Dy1 -O11	70.5(3)
		O2 -Dy1 -O13	120.7(2)
		O2 -Dy1 -O14	168.4(2)
		O3 -Dy1 -O4	64.95(16)
		O3 -Dy1 -O7	76.47(18)
		O3 -Dy1 -O8	120.49(17)
		O3 -Dy1 -O10	76.1(3)
		O3 -Dy1 -O11	90.7(3)
		O3 -Dy1 -O13	150.5(2)
		O3 -Dy1 -O14	123.8(2)
		O4 -Dy1 -O7	77.26(19)
		O4 -Dy1 -O8	76.08(18)
		O4 -Dy1 -O10	78.1(3)
		O4 -Dy1 -O11	128.6(3)
		O4 -Dy1 -O13	113.6(2)
		O4-Dy1 -O14	65.8(2)
		O7 -Dy1 -O8	51.49(17)
		O7-Dy1 -O10	148.9(3)
		O7 -Dy1 -O11	142.9(3)
		O7-Dy1 -O13	132.9(2)
		O7 -Dy1 -O14	116.5(3)
		O8 -Dy1 -O10	137.7(3)
		O8 -Dy1 -O11	148.1(3)
		O8 -Dy1 -O13	85.4(2)
		O8 -Dy1 -O14	70.0(3)
		O10 -Dy1 -O11	51.5(4)
		O10 -Dy1 -O13	75.0(3)
		O10 -Dy1 -O14	68.9(3)
		O11 -Dy1 -O13	66.9(3)
		O11 -Dy1 -O14	99.7(3)
		O13 -Dy1 -O14	48.0(3)
		O2 -Cu1 -N1	167.9(2)
		O2 -Cu1 -N2	91.4(2)
		O2 -Cu1 -O3	78.8(2)
		O2 -Cu1 -O5	93.8(2)
		O3 -Cu1 -O5	91.4(3)
		N1 -Cu1 -N2	97.3(3)
		O3 -Cu1 -N1	91.6(2)
		O3 -Cu1 -N2	168.4(2)
		O5 -Cu1 -N1	93.7(3)
		O5 -Cu1 -N2	95.5(3)
		Dy1 -O2 -Cu1	108.3(2)
		Dy1 -O3 -Cu1	109.0(2)

**Cu(II)-Dy(III) and Co(III)-Dy(III) based single molecule magnets with multiple slow magnetic relaxation processes in Cu(II)-Dy(III) complex**

Malay Dolai, Mahammad Ali, Ján Titiš and Roman Boča

**Table S2.** Hydrogen bonding interactions in **1** (Å, °).

D–H•••A	Distances (Å)			Angles (°)
	D–H	H•••A	D•••A	D–H•••A
O6 -H6 •••O13 <sup>i</sup>	0.82	2.19	2.855(12)	138
C16 -H5 •••O11 <sup>ii</sup>	0.93	2.54	3.327(13)	143
C19 -H7A •••O6 <sup>iii</sup>	0.96	2.59	3.380(14)	140
C19 -H7B •••O8	0.96	2.53	3.187(11)	126
C19 -H7B •••O14	0.96	2.44	2.783(12)	101
C5 -H15•••O9 <sup>iv</sup>	0.93	2.46	3.343(11)	159
C1 -H19A •••O5 <sup>v</sup>	0.96	2.55	3.415(11)	151
C1 -H19B •••O13	0.96	2.47	2.996(12)	114

Symmetry code: (i) = x,y,-1+z; (ii) = 1+x,y,z; (iii) = 1+x,y,1+z; (iv) = x,1/2-y,1/2+z; (v) = -1+x,1/2-y,-1/2+z.

**Table S3.**  $\pi$ - $\pi$  interactions and C–H••• $\pi$  interaction in complex **1**<sup>a,b</sup>

Ring(i)...ring(j)	Distance of centroid(i) from ring(j) (Å)	Dihedral angle (i, j) (°)	Distance between the (i, j) ring centroids (Å)
R(1)...R(2) <sup>i</sup>	4.414(5)	22.9(4)	3.664(4)
R(2)...R(1) <sup>ii</sup>	4.414(5)	22.9(4)	3.664(4)
C–H...ring(j)	H...R distance (Å)	C–H...R angle (°)	C...R distance (Å)
C11-H9A...R(2) <sup>ii</sup>	2.95	141	3.751(9)

<sup>a</sup> Symmetry code: (i) = -1+x,y,z; (ii) = 1+x,y,z; (iii) = 1-x,-y,-z.

<sup>b</sup> R(i)/R(j) denotes the ith/jth rings:

R(1) = C(2) C(3) C(4) C(5) C(6) C(7);

R(2) = C(13) C(14) C(15) C(16) C(17) C(18).

**Cu(II)-Dy(III) and Co(III)-Dy(III) based single molecule magnets with multiple slow magnetic relaxation processes in Cu(II)-Dy(III) complex**

Malay Dolai, Mohammad Ali, Ján Titiš and Roman Boča

**Table S4.** Selected bond lengths (Å) and bond angles (°) of complex **2**

Dy1 -O1	2.633(2)	O1 -Dy1-O2	63.04(6)
Dy1 -O2	2.3545(19)	O1 -Dy1-O3	121.12(7)
Dy1 -O3	2.373(2)	O1 -Dy1-O4	129.02(10)
Dy1 -O4	2.596(3)	O1 -Dy1 -O9	117.37(9)
Dy1 -O9	2.346(2)	O1 -Dy1 -O10	77.20(8)
Dy1 -O10	2.262(2)	O1 -Dy1 -O12	68.97(10)
Dy1 -O12	2.385(3)	O1 -Dy1 -O13	141.45(9)
Dy1 -O13	2.409(3)	O1 -Dy1 -O14	65.76(10)
Dy1 -O14	2.430(2)	O2 -Dy1 -O3	66.12(7)
Co1 -O2	1.9251(19)	O2 -Dy1 -O4	122.72(7)
Co1 -O3	1.914(2)	O2 -Dy1 -O9	72.39(7)
Co1 -O7	1.919(2)	O2 -Dy1 -O10	139.69(8)
Co1 -O8	1.886(2)	O2 -Dy1 -O12	91.48(8)
Co1 -N1	1.923(3)	O2 -Dy1 -O13	141.19(9)
Co1 -N2	1.917(2)	O2 -Dy1 -O14	73.79(8)
		O3 -Dy1 -O4	62.75(7)
		O3 -Dy1 -O9	70.66(7)
		O3 -Dy1 -O10	138.95(8)
		O3 -Dy1 -O12	139.67(9)
		O3 -Dy1 -O13	97.21(8)
		O3 -Dy1 -O14	73.36(8)
		O4 -Dy1 -O9	111.39(8)
		O4 -Dy1 -O10	77.18(8)
		O4 -Dy1 -O12	145.29(9)
		O4 -Dy1 -O13	70.21(10)
		O4 -Dy1 -O14	68.53(9)
		O9 -Dy1 -O10	137.38(8)
		O9 -Dy1 -O12	70.70(9)
		O9 -Dy1 -O13	68.97(9)
		O9-Dy1 -O14	138.02(9)
		O10 -Dy1 -O12	79.73(9)
		O10-Dy1 -O13	76.04(9)
		O10 -Dy1 -O14	84.52(9)
		O12-Dy1 -O13	79.32(10)
		O12 -Dy1 -O14	134.29(10)
		O13 -Dy1 -O14	137.28(10)
		O2 -Co1 -N1	90.64(9)
		O2 -Co1 -N2	174.57(9)
		O2 -Co1 -O3	84.41(8)
		O2 -Co1 -O7	86.53(8)
		O2 -Co1 -O8	92.78(9)
		O3 -Co1 -O7	86.30(9)
		O3 -Co1 -O8	91.97(9)
		O3 -Co1 -N1	174.85(10)
		O3 -Co1 -N2	90.17(10)
		O7 -Co1 -N1	94.85(9)
		O7 -Co1 -N2	92.98(9)
		O7 -Co1 -O8	178.19(9)
		O8 -Co1 -N1	86.83(10)
		O8 -Co1 -N2	87.55(10)
		N1 -Co1 -N2	94.79(11)
		Dy1 -O2 -Co1	104.25(8)
		Dy1 -O3 -Co1	103.91(9)

Electronic Supplementary Information (ESI)  
**Cu(II)-Dy(III) and Co(III)-Dy(III) based single molecule magnets with multiple slow magnetic relaxation processes in Cu(II)-Dy(III) complex**

Malay Dolai, Mohammad Ali, Ján Titiš and Roman Boča

**Table S5.** Hydrogen bonding interactions in **2** (Å, deg).

D–H•••A	Distances (Å)			Angles (°)
	D–H	H•••A	D•••A	D–H•••A
O5 - H12•••O6	0.84	1.8200	2.652(4)	171
O12 - H57•••O15	0.75(4)	2.09(4)	2.837(7)	172(5)
O12 - H58•••O3S	0.94(5)	1.78(5)	2.695(6)	166(4)
O13 - H64•••O11	1.05(7)	1.76(7)	2.617(4)	136(6)
O14 - H66 •••O2S <sup>i</sup>	0.83(5)	1.94(5)	2.685(5)	150(4)
O14 - H70 ••• O7	0.82(6)	2.08(6)	2.835(4)	153(7)
C1 - H7B ••• O10	0.98	2.42	3.098(4)	125
C9 - H9B •••O8	0.99	2.57	2.963(4)	103
C11 - H12A •••O8	0.99	2.55	2.956(4)	104
C3 - H16 •••O11 <sup>ii</sup>	0.95	2.51	3.448(4)	171
C19 - H18B •••O10	0.98	2.54	3.161(5)	121
C27 - H27A •••O3S <sup>ii</sup>	0.98	2.45	3.265(8)	141
C27 - H27A •••O13 <sup>ii</sup>	0.98	2.56	3.092(7)	114

Symmetry code: (i) = -1+x,y,z; (ii) = x,3/2-y,1/2+z.

**Table S6.**  $\pi$ - $\pi$  interactions and C–H••• $\pi$  interaction in complex **2**<sup>c,d</sup>

Ring(i)...ring(j)	Distance of centroid(i) from ring(j) (Å)	Dihedral angle (i, j) (°)	Distance between the (i, j) ring centroids (Å)
R(1)...R(2) <sup>i</sup>	3.6878(19)	21.81	3.4239(13)
R(2)...R(1) <sup>ii</sup>	3.9088(19)	27.72	3.4603(13)

<sup>c</sup> Symmetry code: (i) = 1-x, 1-y, 1-z; (ii) = -x, 1-y, -z.

<sup>d</sup> R(i)/R(j) denotes the ith/jth rings:

R(1) = C(2) C(3) C(4) C(5) C(6) C(7);

R(2) = C(13) C(14) C(15) C(16) C(17) C(18).

Electronic Supplementary Information (ESI)  
**Cu(II)-Dy(III) and Co(III)-Dy(III) based single molecule magnets with multiple slow magnetic relaxation processes in Cu(II)-Dy(III) complex**

Malay Dolai, Mohammad Ali, Ján Titiš and Roman Boča

---

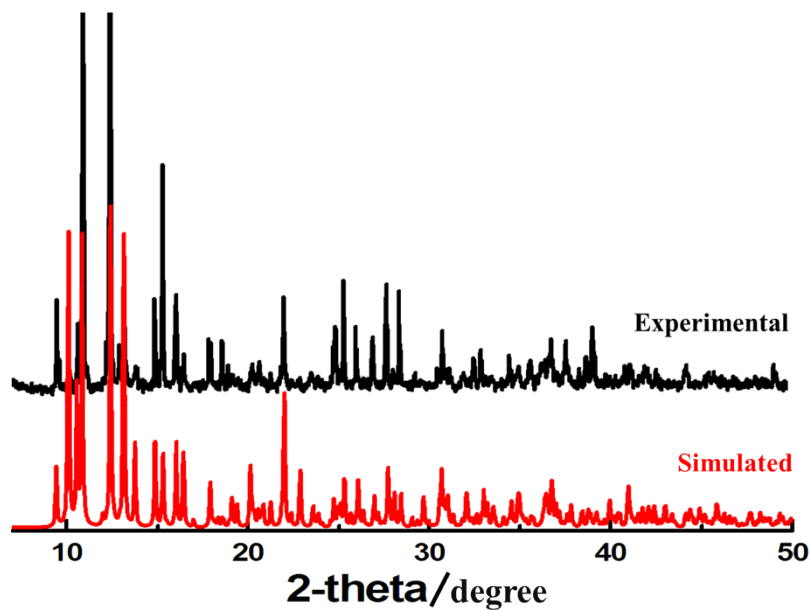


Figure S6(a). The powder diffraction patterns confirming purity of 1.

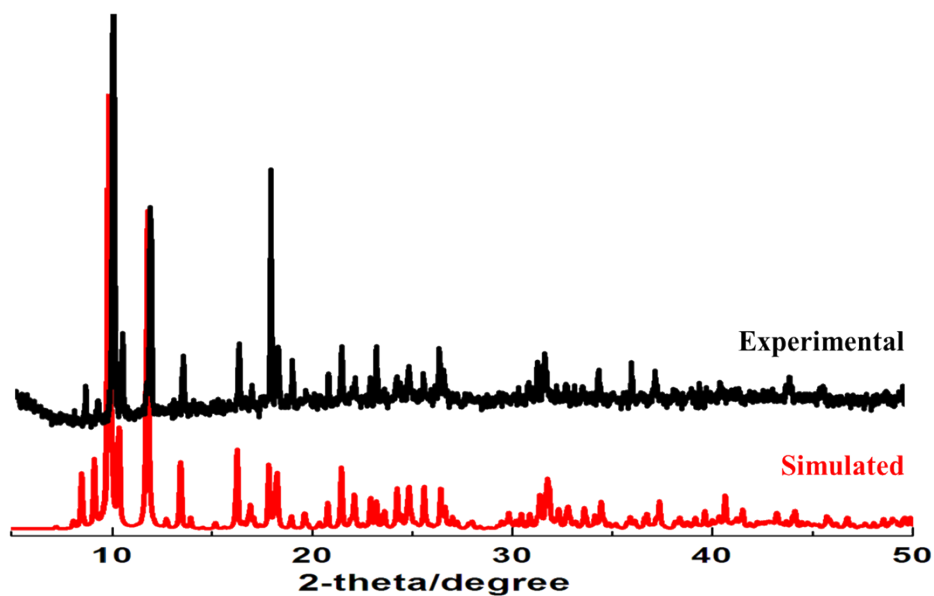
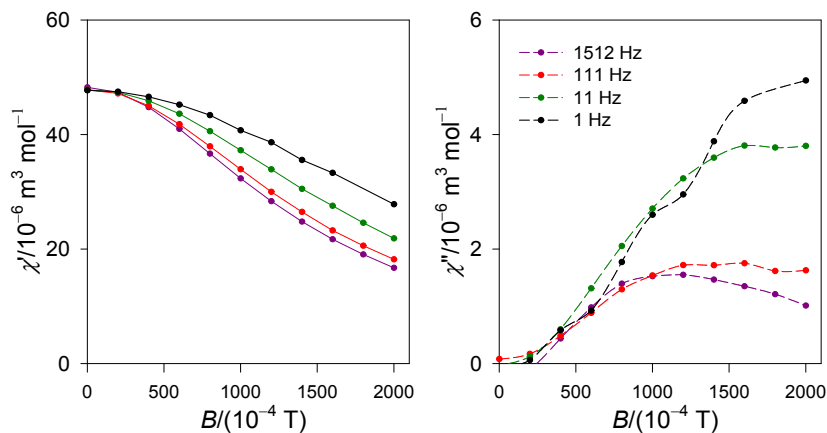


Figure S6(b). The powder diffraction patterns confirming purity of 2.

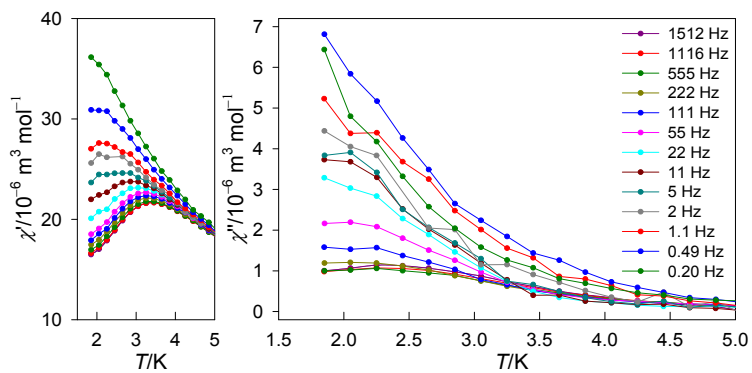


Electronic Supplementary Information (ESI)  
**Cu(II)-Dy(III) and Co(III)-Dy(III) based single molecule magnets with multiple slow magnetic relaxation processes in Cu(II)-Dy(III) complex**

Malay Dolai, Mohammad Ali, Ján Titiš and Roman Boča



**Figure S7.** The in-phase  $\chi'$  and out-of-phase  $\chi''$  molar susceptibility (SI units) for **1** in dependence upon external magnetic field at  $T = 2.0$  K. Lines serve as a guide for eyes.



**Figure S8.** The AC susceptibility data for **1** at  $B_{DC} = 0.2$  T. Left – temperature dependence of the in-phase molar susceptibility; right – temperature dependence of the out-of-phase molar susceptibility. Lines serve as a guide for eyes.

Electronic Supplementary Information (ESI)  
**Cu(II)-Dy(III) and Co(III)-Dy(III) based single molecule magnets with multiple slow magnetic relaxation processes in Cu(II)-Dy(III) complex**

Malay Dolai, Mohammad Ali, Ján Titiš and Roman Boča

**Table S7.** Parameters of the extended Debye model for **1** using data taken at  $B_{DC} = 0.2$  T. <sup>a</sup>

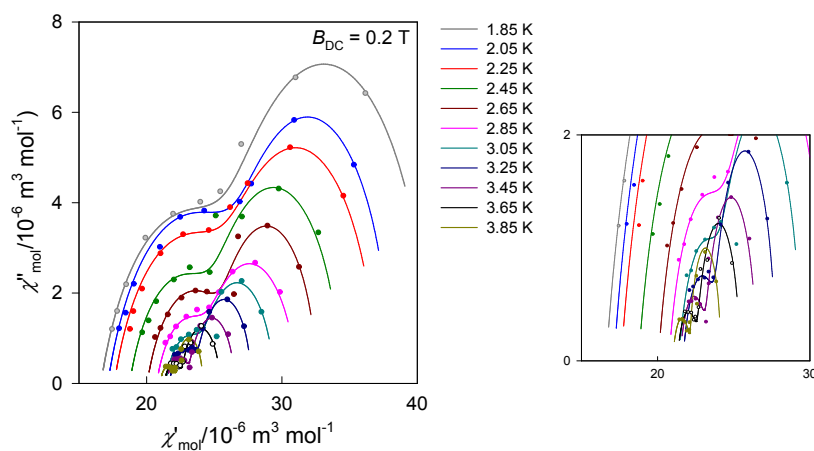
$T/K$	Low-frequency (LF)				Intermediate frequency (IF)			$R(\chi')$ /%	$R(\chi'')$ /%
	$\chi_s$	$\chi_{T1}$	$\alpha_1$	$\tau_1 / s$	$\chi_{T2}$	$\alpha_2$	$\tau_2 / 10^{-3} s$		
1.85b	16.7(4)	28.9(24)	0.00(8)	0.528(38)	40.7(8)	0.33(7)	17.9(52)	1.3	2.0
2.05	17.1(3)	27.2(16)	0.00(6)	0.440(28)	38.1(5)	0.29(5)	15.4(30)	0.91	3.0
2.25	17.6(3)	27.8(17)	0.07(7)	0.383(31)	37.1(5)	0.30(7)	11.3(27)	0.92	2.5
2.45b	19.0(3)	28.1(11)	0.09(5)	0.349(23)	34.7(4)	0.20(7)	7.69(158)	0.56	2.9
2.65b	19.9(2)	27.8(9)	0.12(5)	0.289(18)	32.9(3)	0.28(6)	5.98(111)	0.42	3.3
2.85	20.7(2)	27.3(10)	0.18(6)	0.293(26)	31.2(3)	0.30(10)	4.91(121)	0.45	3.8
3.05	21.4(2)	26.5(7)	0.13(6)	0.272(21)	29.5(2)	0.30(10)	4.34(108)	0.32	4.8
3.25	21.7(2)	25.5(5)	0.05(4)	0.298(16)	27.7(1)	0.30(10)	3.16(60)	0.32	4.6
3.45	21.5(5)	24.8(8)	0.10(8)	0.309(32)	26.5(2)	0.29(22)	1.38(70)	0.50	6.7
3.65	21.3(3)	24.1(4)	0.10(5)	0.308(19)	25.4(12)	0.30(15)	1.26(47)	0.17	7.8
3.85	21.0(5)	22.9(7)	0.00(9)	0.329(37)	24.1(17)	0.30(27)	1.15(93)	0.52	4.5
4.05	20.6(3)	22.1(3)	0.02(5)	0.336(23)	23.1(2)	0.30(16)	1.04(52)	0.24	3.8

<sup>a</sup> SI unit for the molar magnetic susceptibility [ $10^{-6} \text{ m}^3 \text{ mol}^{-1}$ ].  $R$  – discrepancy factor of the fit for dispersion  $\chi'$  and absorption  $\chi''$ , respectively. Two components of the magnetic susceptibility: a) in phase

$$\chi'(\omega) = \chi_s + (\chi_{T1} - \chi_s) \frac{1 + (\omega\tau_1)^{1-\alpha_1} \sin(\pi\alpha_1/2)}{1 + 2(\omega\tau_1)^{1-\alpha_1} \sin(\pi\alpha_1/2) + (\omega\tau_1)^{2-2\alpha_1}} + (\chi_{T2} - \chi_{T1}) \frac{1 + (\omega\tau_2)^{1-\alpha_2} \sin(\pi\alpha_2/2)}{1 + 2(\omega\tau_2)^{1-\alpha_2} \sin(\pi\alpha_2/2) + (\omega\tau_2)^{2-2\alpha_2}} \quad (\text{S1})$$

b) out of phase

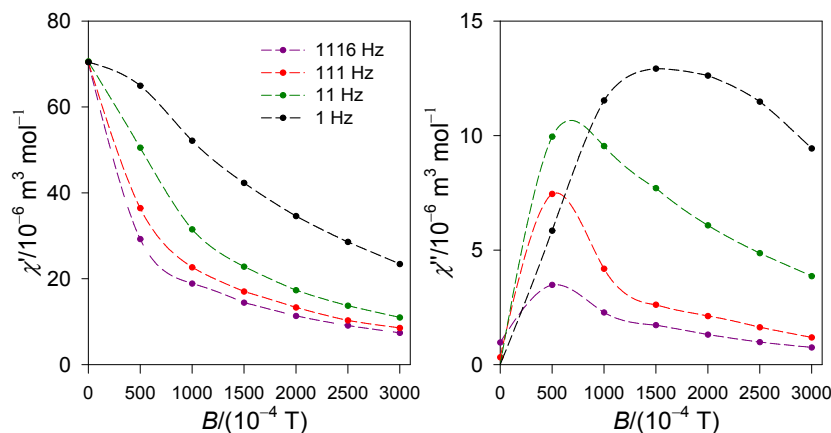
$$\chi''(\omega) = (\chi_{T1} - \chi_s) \frac{(\omega\tau_1)^{1-\alpha_1} \cos(\pi\alpha_1/2)}{1 + 2(\omega\tau_1)^{1-\alpha_1} \sin(\pi\alpha_1/2) + (\omega\tau_1)^{2-2\alpha_1}} + (\chi_{T2} - \chi_{T1}) \frac{(\omega\tau_2)^{1-\alpha_2} \cos(\pi\alpha_2/2)}{1 + 2(\omega\tau_2)^{1-\alpha_2} \sin(\pi\alpha_2/2) + (\omega\tau_2)^{2-2\alpha_2}} \quad (\text{S2})$$



**Figure S9.** Zoomed Argand plot with experimental data and fitted lines for **1**.

Electronic Supplementary Information (ESI)  
**Cu(II)-Dy(III) and Co(III)-Dy(III) based single molecule magnets with multiple slow magnetic relaxation processes in Cu(II)-Dy(III) complex**

Malay Dolai, Mohammad Ali, Ján Titiš and Roman Boča



**Figure S10.** The in-phase  $\chi'$  and out-of-phase  $\chi''$  molar susceptibility (SI units) for **2** in dependence upon external magnetic field at  $T = 2.0$  K. Lines serve as a guide for eyes.

**Table S8.** Parameters of the extended Debye model for **2** at  $B_{DC} = 0.2$  T.

$T/K$	$\chi_s$	$\chi_r$	$\alpha$	$\tau/s$	$R(\chi')/\%$	$R(\chi'')/\%$
1.9	10.9(3)	66.6(12)	0.44(1)	0.268(19)	2.3	6.7
2.3	11.3(2)	58.1(8)	0.44(1)	0.158(9)	1.9	6.9
2.7	11.5(3)	50.6(7)	0.45(1)	0.055(6)	2.1	7.7
3.1	11.2(3)	45.0(7)	0.47(2)	0.066(5)	2.2	8.4
3.5	10.7(3)	40.3(5)	0.48(2)	0.041(3)	2.2	8.7
3.9	10.2(2)	37.6(3)	0.48(1)	0.029(1)	0.8	7.4
4.3	9.9(2)	33.8(2)	0.46(1)	0.0159(7)	1.3	7.7
4.7	10.3(2)	30.4(2)	0.40(1)	0.0107(5)	1.4	9.3
5.1	10.6(2)	27.9(1)	0.33(1)	0.0080(3)	1.3	10.1
5.5	10.5(1)	26.1(1)	0.29(1)	0.0064(2)	0.97	8.5
5.9	10.1(1)	24.4(1)	0.27(1)	0.0047(1)	1.00	9.2
6.7	9.61(8)	21.6(1)	0.16(1)	0.0026(1)	0.72	6.2
7.1	9.27(6)	20.6(1)	0.13(1)	0.00199(3)	0.52	4.6
7.5	8.92(6)	19.6(1)	0.12(1)	0.00149(2)	0.51	4.0
8.7	8.21(6)	17.1(1)	0.07(1)	0.00066(1)	0.49	3.1
9.9	7.68(5)	15.2(1)	0.06(1)	0.00028(0)	0.13	3.1

<sup>a</sup> SI unit for the molar magnetic susceptibility [ $10^{-6} \text{ m}^3 \text{ mol}^{-1}$ ].  $R$  – discrepancy factor of the fit for dispersion  $\chi'$  and absorption  $\chi''$ , respectively.

**Cu(II)-Dy(III) and Co(III)-Dy(III) based single molecule magnets with multiple slow magnetic relaxation processes in Cu(II)-Dy(III) complex**

Malay Dolai, Mohammad Ali, Ján Titiš and Roman Boča

**Table S9.** Calculated energy levels by DFT (UHF) truncated after 200 levels

		Spin up		Spin down
NO	OCC	E(eV)	OCC	E(eV)
0.000	1.000	-57143.380	1.000	-57143.380
1.000	1.000	-9095.699	1.000	-9095.684
2.000	1.000	-8992.068	1.000	-8992.063
3.000	1.000	-8064.062	1.000	-8064.052
4.000	1.000	-8064.012	1.000	-8064.003
5.000	1.000	-8063.986	1.000	-8063.977
6.000	1.000	-2026.624	1.000	-2024.447
7.000	1.000	-1712.892	1.000	-1710.566
8.000	1.000	-1712.588	1.000	-1710.351
9.000	1.000	-1712.431	1.000	-1710.236
10.000	1.000	-1310.176	1.000	-1308.055
11.000	1.000	-1310.002	1.000	-1307.905
12.000	1.000	-1309.670	1.000	-1307.857
13.000	1.000	-1309.610	1.000	-1307.787
14.000	1.000	-1309.565	1.000	-1307.772
15.000	1.000	-1083.957	1.000	-1083.390
16.000	1.000	-935.877	1.000	-935.233
17.000	1.000	-935.811	1.000	-935.202
18.000	1.000	-935.135	1.000	-934.977
19.000	1.000	-523.167	1.000	-523.168
20.000	1.000	-522.896	1.000	-522.897
21.000	1.000	-522.197	1.000	-522.199
22.000	1.000	-522.116	1.000	-522.117
23.000	1.000	-522.066	1.000	-521.998
24.000	1.000	-521.756	1.000	-521.756
25.000	1.000	-521.732	1.000	-521.732
26.000	1.000	-521.656	1.000	-521.656
27.000	1.000	-521.605	1.000	-521.536
28.000	1.000	-521.336	1.000	-521.336
29.000	1.000	-521.082	1.000	-521.081
30.000	1.000	-520.656	1.000	-520.658
31.000	1.000	-520.636	1.000	-520.634
32.000	1.000	-520.566	1.000	-520.565
33.000	1.000	-518.436	1.000	-518.435
34.000	1.000	-412.758	1.000	-407.477
35.000	1.000	-397.166	1.000	-397.166
36.000	1.000	-396.097	1.000	-396.097
37.000	1.000	-396.022	1.000	-396.022
38.000	1.000	-392.277	1.000	-392.205
39.000	1.000	-391.964	1.000	-391.880
40.000	1.000	-309.222	1.000	-303.600
41.000	1.000	-308.740	1.000	-303.339
42.000	1.000	-308.486	1.000	-303.193
43.000	1.000	-279.814	1.000	-279.817
44.000	1.000	-279.793	1.000	-279.800
45.000	1.000	-279.711	1.000	-279.711
46.000	1.000	-279.693	1.000	-279.697
47.000	1.000	-279.580	1.000	-279.575
48.000	1.000	-279.476	1.000	-279.477
49.000	1.000	-279.440	1.000	-279.435

Electronic Supplementary Information (ESI)

**Cu(II)-Dy(III) and Co(III)-Dy(III) based single molecule magnets with multiple slow magnetic relaxation processes in Cu(II)-Dy(III) complex**

Malay Dolai, Mahammad Ali, Ján Titiš and Roman Boča

50.000	1.000	-279.082	1.000	-279.081
51.000	1.000	-278.457	1.000	-278.455
52.000	1.000	-278.241	1.000	-278.237
53.000	1.000	-278.210	1.000	-278.208
54.000	1.000	-278.065	1.000	-278.065
55.000	1.000	-277.798	1.000	-277.799
56.000	1.000	-277.586	1.000	-277.586
57.000	1.000	-277.475	1.000	-277.476
58.000	1.000	-277.459	1.000	-277.459
59.000	1.000	-277.395	1.000	-277.395
60.000	1.000	-277.366	1.000	-277.366
61.000	1.000	-277.175	1.000	-277.174
62.000	1.000	-168.173	1.000	-161.859
63.000	1.000	-167.811	1.000	-161.366
64.000	1.000	-167.024	1.000	-161.285
65.000	1.000	-166.862	1.000	-161.152
66.000	1.000	-166.815	1.000	-161.131
67.000	1.000	-124.325	1.000	-123.134
68.000	1.000	-81.061	1.000	-79.346
69.000	1.000	-80.936	1.000	-79.328
70.000	1.000	-79.543	1.000	-79.135
71.000	1.000	-54.979	1.000	-52.667
72.000	1.000	-37.935	1.000	-37.931
73.000	1.000	-35.066	1.000	-35.061
74.000	1.000	-33.693	1.000	-33.687
75.000	1.000	-31.406	1.000	-31.399
76.000	1.000	-30.630	1.000	-30.629
77.000	1.000	-30.043	1.000	-29.769
78.000	1.000	-29.852	1.000	-29.304
79.000	1.000	-29.720	1.000	-29.190
80.000	1.000	-29.309	1.000	-28.998
81.000	1.000	-29.055	1.000	-28.884
82.000	1.000	-28.886	1.000	-28.492
83.000	1.000	-28.741	1.000	-28.195
84.000	1.000	-28.480	1.000	-27.890
85.000	1.000	-28.305	1.000	-27.594
86.000	1.000	-27.965	1.000	-27.561
87.000	1.000	-27.636	1.000	-27.478
88.000	1.000	-27.553	1.000	-27.142
89.000	1.000	-27.062	1.000	-26.954
90.000	1.000	-25.719	1.000	-25.634
91.000	1.000	-25.369	1.000	-25.330
92.000	1.000	-24.462	1.000	-24.444
93.000	1.000	-24.239	1.000	-24.225
94.000	1.000	-23.097	1.000	-23.084
95.000	1.000	-21.970	1.000	-21.963
96.000	1.000	-21.654	1.000	-21.644
97.000	1.000	-21.475	1.000	-21.461
98.000	1.000	-21.409	1.000	-21.386
99.000	1.000	-20.838	1.000	-20.821
100.000	1.000	-20.534	1.000	-20.529
101.000	1.000	-19.986	1.000	-19.982
102.000	1.000	-19.397	1.000	-19.380
103.000	1.000	-18.845	1.000	-18.790
104.000	1.000	-18.448	1.000	-18.413
105.000	1.000	-17.892	1.000	-17.863
106.000	1.000	-17.663	1.000	-17.641



Electronic Supplementary Information (ESI)

**Cu(II)-Dy(III) and Co(III)-Dy(III) based single molecule magnets with multiple slow magnetic relaxation processes in Cu(II)-Dy(III) complex**

Malay Dolai, Mahammad Ali, Ján Titiš and Roman Boča

---

107.000	1.000	-17.438	1.000	-17.406
108.000	1.000	-17.134	1.000	-17.104
109.000	1.000	-16.766	1.000	-16.762
110.000	1.000	-16.630	1.000	-16.607
111.000	1.000	-16.274	1.000	-16.257
112.000	1.000	-16.061	1.000	-16.025
113.000	1.000	-15.851	1.000	-15.818
114.000	1.000	-15.734	1.000	-15.713
115.000	1.000	-15.540	1.000	-15.517
116.000	1.000	-15.368	1.000	-15.336
117.000	1.000	-15.256	1.000	-15.240
118.000	1.000	-15.168	1.000	-15.110
119.000	1.000	-15.096	1.000	-15.078
120.000	1.000	-14.957	1.000	-14.910
121.000	1.000	-14.805	1.000	-14.771
122.000	1.000	-14.656	1.000	-14.633
123.000	1.000	-14.575	1.000	-14.551
124.000	1.000	-14.504	1.000	-14.459
125.000	1.000	-14.451	1.000	-14.421
126.000	1.000	-14.343	1.000	-14.320
127.000	1.000	-14.144	1.000	-14.122
128.000	1.000	-14.087	1.000	-14.031
129.000	1.000	-14.046	1.000	-13.960
130.000	1.000	-13.969	1.000	-13.947
131.000	1.000	-13.849	1.000	-13.829
132.000	1.000	-13.832	1.000	-13.809
133.000	1.000	-13.707	1.000	-13.688
134.000	1.000	-13.613	1.000	-13.553
135.000	1.000	-13.555	1.000	-13.418
136.000	1.000	-13.446	1.000	-13.291
137.000	1.000	-13.298	1.000	-13.238
138.000	1.000	-13.238	1.000	-13.054
139.000	1.000	-12.872	1.000	-12.841
140.000	1.000	-12.786	1.000	-12.728
141.000	1.000	-12.662	1.000	-12.593
142.000	1.000	-12.604	1.000	-12.558
143.000	1.000	-12.571	1.000	-12.524
144.000	1.000	-12.264	1.000	-12.185
145.000	1.000	-12.202	1.000	-11.807
146.000	1.000	-12.165	1.000	-11.759
147.000	1.000	-12.134	1.000	-11.617
148.000	1.000	-12.055	1.000	-11.531
149.000	1.000	-11.939	1.000	-11.439
150.000	1.000	-11.869	1.000	-11.289
151.000	1.000	-11.685	1.000	-11.015
152.000	1.000	-11.660	1.000	-10.696
153.000	1.000	-11.599	1.000	-10.533
154.000	1.000	-11.505	1.000	-10.517
155.000	1.000	-11.398	1.000	-10.448
156.000	1.000	-11.382	1.000	-10.341
157.000	1.000	-11.214	1.000	-10.138
158.000	1.000	-11.094	1.000	-10.064
159.000	1.000	-10.939	1.000	-9.900
160.000	1.000	-10.798	1.000	-9.870
161.000	1.000	-10.715	1.000	-9.846
162.000	1.000	-10.559	1.000	-9.645
163.000	1.000	-10.409	1.000	-9.531

Electronic Supplementary Information (ESI)

**Cu(II)-Dy(III) and Co(III)-Dy(III) based single molecule magnets with multiple slow magnetic relaxation processes in Cu(II)-Dy(III) complex**

Malay Dolai, Mahammad Ali, Ján Titiš and Roman Boča

---

164.000	1.000	-10.342	1.000	-9.388
165.000	1.000	-10.261	1.000	-9.287
166.000	1.000	-10.133	1.000	-9.077
167.000	1.000	-9.987	1.000	-8.872
168.000	1.000	-9.920	1.000	-8.818
169.000	1.000	-9.786	1.000	-8.704
170.000	1.000	-9.665	1.000	-8.673
171.000	1.000	-9.607	1.000	-8.505
172.000	1.000	-9.406	1.000	-8.495
173.000	1.000	-9.209	1.000	-8.343
174.000	1.000	-9.046	1.000	-8.324
175.000	1.000	-8.824	1.000	-8.231
176.000	1.000	-8.764	1.000	-8.068
177.000	1.000	-8.684	1.000	-7.976
178.000	1.000	-8.645	1.000	-7.847
179.000	1.000	-8.552	1.000	-7.742
180.000	1.000	-8.461	1.000	-7.563
181.000	1.000	-8.293	1.000	-7.486
182.000	1.000	-8.118	1.000	-7.413
183.000	1.000	-7.966	1.000	-7.208
184.000	1.000	-7.768	1.000	-7.099
185.000	1.000	-7.604	1.000	-6.705
186.000	1.000	-7.494	1.000	-6.553
187.000	1.000	-7.442	1.000	-6.282
188.000	1.000	-7.379	1.000	-5.541
189.000	1.000	-7.252	1.000	-4.767
190.000	1.000	-7.104	0.000	-3.160
191.000	1.000	-6.737	0.000	-2.748
192.000	1.000	-6.559	0.000	-2.058
193.000	1.000	-6.263	0.000	-1.925
194.000	1.000	-5.547	0.000	-1.815
195.000	1.000	-4.771	0.000	-1.458
196.000	0.000	-2.822	0.000	-1.362
197.000	0.000	-2.115	0.000	-1.083
198.000	0.000	-1.457	0.000	-0.687
199.000	0.000	-0.696	0.000	-0.655
200.000	0.000	-0.475	0.000	-0.439