## **Electronic Supplementary Information**

## {Ni<sup>II</sup><sub>8</sub>Ln<sup>III</sup><sub>6</sub>} (Ln=Gd, Dy) rod-like nano-sized heteronuclear coordination cluster with a double carbonate bridge skeleton and remarkable MCE behaviour

Eliana Guarda<sup>a</sup>, Katharina Bader<sup>b</sup>, Joris van Slageren<sup>b</sup>, Pablo Alborés<sup>a</sup>\*

<sup>a</sup> Departamento de Química Inorgánica, Analítica y Química Física/ INQUIMAE (CONICET), Facultad de Ciencias Exactas y Naturales Universidad de Buenos Aires, Pabellón 2, Ciudad Universitaria, C1428EHA Buenos Aires, Argentina.; <u>Tel:+54(0)1145763380</u>; E-mail: albores@.qi.fcen.uba.ar
<sup>b</sup> Institut für Physikalische Chemie, Universität Stuttgart, Pfaffenwaldring 55, D-70569,

Stuttgart, Germany.



**Figure ESI1.** Space filling representation of  $Ni_8Dy_6$  complexes showing their rod-like overall shape.



Figure ESI2. Molecular representation of the central double  $\mu_5$ -carbonate bridged octanuclear {Ni<sub>4</sub>Dy<sub>4</sub>} moiety. Carbonate bridges are green highlighted.



Figure ESI3. Molecular representation of the opened cubane  $Ni_3Dy$  units located at both sides of the central octanuclear { $Ni_4Dy_4$ } moiety.





**Figure ESI4.** Polyhedra representation of coordination environment of all Ni(II) sites (top) and all Dy(III) sites (bottom).



Figure ESI5. Intra-molecular H-bond interactions present in complex Ni<sub>8</sub>Dy<sub>6</sub>.



**Figure ESI6.** Unit cell crystal packing of Ni<sub>8</sub>Dy<sub>6</sub>. View along monoclinic *b*-axis.



Figure ESI7. Spin ladder energies arising from Ni<sub>4</sub> model fitting (see text).



**Figure ESI8.** Magnetic entropy change field dependence at different fixed temperatures of complexes 1-3, as evaluated from magnetization data.



**Figure ESI9.** PXRD of complexes **1-3** (top to bottom) at room temperature. Black: single crystal simulated PXRD.