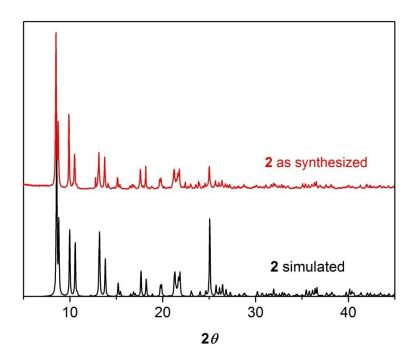
## **Supplementary Information**



**Fig.S1**. Powder X-ray diffraction patterns of **2** (red) compared with the simulation (black) obtained from single crystal data of **2**.

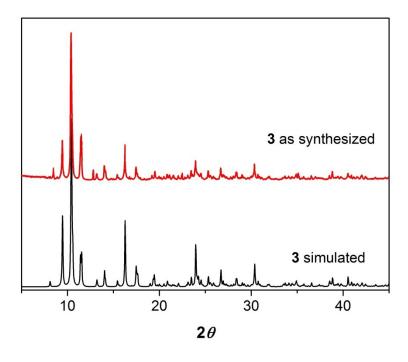


Fig. S2. Powder X-ray diffraction patterns of 3 (red) compared with the simulation (black) obtained from single crystal data of 3.

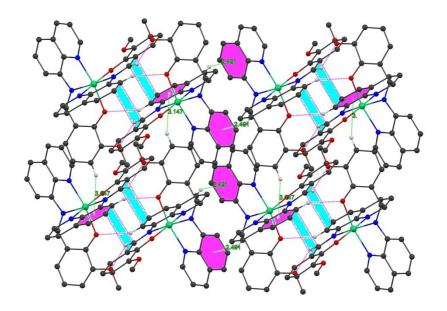


Fig. S3. Packing diagram of complex 1

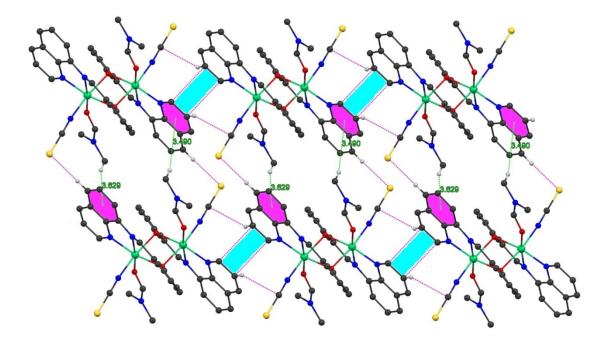


Fig. S4. Packing diagram of complex 2

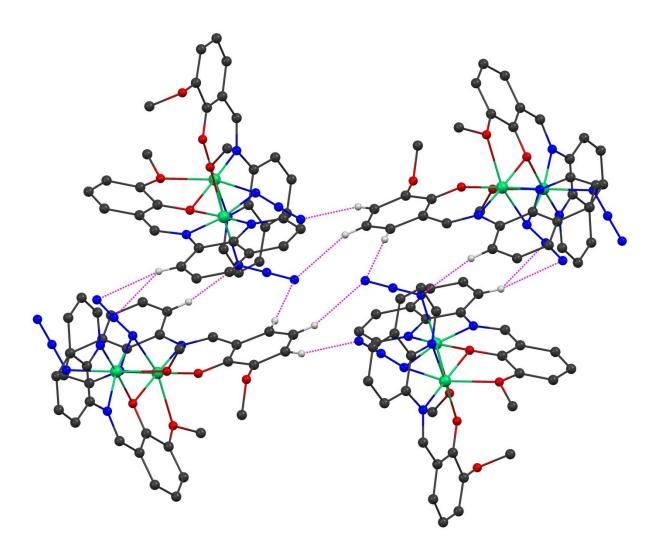
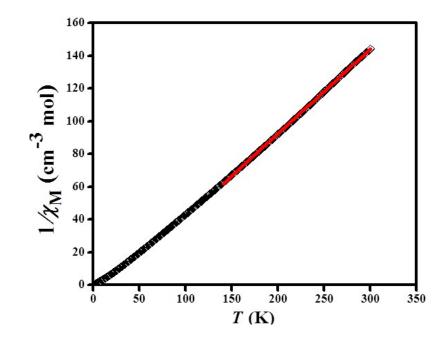
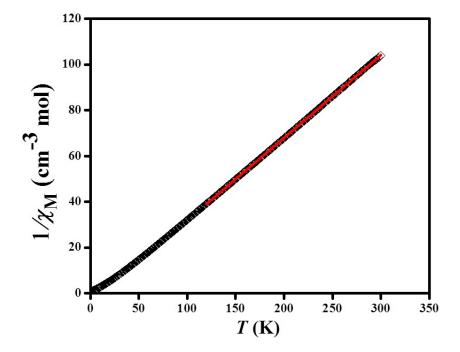


Fig. S5.Packing diagram of complex 3



**Fig. S6**: Reciprocal molar susceptibility as function of temperature for **2**; solid red line indicates Curie-Weiss fitting in the range 140 K–300 K.



**Fig. S7**: Reciprocal molar susceptibility as function of temperature for **3**; solid red line indicates Curie-Weiss fitting in the range 120 K–300 K