

Supporting Information

Large Magnetic Anisotropy of a Decorated Spin-Chain System $K_2Co_3(MoO_4)_3(OH)_2$

Bhakti Patel,^a Feng Ye,^b W. L. N. C. Liyanage,^c Carolyn Buchanan,^c Dusting A. Gilbert,^{c,d} Joseph W. Kolis,^a Liurukara D Sanjeewa^{e,f*}

^a*Department of Chemistry and Center for Optical Materials Science and Engineering Technologies (COMSET), Clemson University, Clemson, SC 29634-0973, USA*

^b*Neutron Scattering Division, Oak Ridge National Laboratory, Oak Ridge, TN 37831, USA*

^c*Department of Physics and Astronomy, University of Tennessee, Knoxville, Tennessee 37996, USA*

^d*Materials Science and Engineering, University of Tennessee, Knoxville, Tennessee 37996, USA*

^e*University of Missouri Research Reactor (MURR), University of Missouri, Columbia, MO 65211, USA*

^f*Department of Chemistry, University of Missouri, Columbia, MO 65211, USA*

*Corresponding Author

Table SI1. X-ray single crystal refinement parameters of $K_2Co_3(MoO_4)_3(OH)_2$.

Figure SI1: Magnetic susceptibilities, $\chi = M/H$, of $K_2Co_3(MoO_4)_3(OH)_2$ as a function of temperature under an applied field of 500 Oe along the *a*-axis.

Figure SI2: Magnetic susceptibilities, $\chi = M/H$, of $K_2Co_3(MoO_4)_3(OH)_2$ as a function of temperature under an applied field of 500 Oe along the *c*-axis.

Table SI1. X-ray single crystal refinement parameters of $\text{K}_2\text{Co}_3(\text{MoO}_4)_3(\text{OH})_2$. The data were refined in the orthorhombic crystal system space group *Pnma* (No. 62) and the unit cell parameters are $a = 17.9394(4)$ Å, $b = 6.0836(2)$ Å, $c = 12.3969(4)$ Å, $V = 1352.95$ Å³ and $Z = 4$.

Atom	<i>x</i>	<i>y</i>	<i>z</i>	<i>U</i> _{eq}	Wyckoff
K(1)	0.36091(6)	0.25000	-0.03775(4)	0.0224(4)	4c
K(2)	0.25395(7)	0.25000	0.66356(5)	0.0342(5)	4c
Co(1)	0.40172(1)	0.25000	0.25767(6)	0.0104(1)	4c
Co(2)	0.55509(5)	0.50205(4)	0.31399(6)	0.0157(4)	8d
Mo(1)	0.44026(5)	0.75000	0.11421(4)	0.0116(5)	4c
Mo(2)	0.60608(4)	0.25000	0.56580(5)	0.0128(1)	4c
Mo(3)	0.20163(6)	0.25000	0.32692(1)	0.0102(5)	4c
O(1)	0.51170(3)	0.25000	0.23014(2)	0.0112(5)	4c
O(2)	0.58890(3)	0.75000	0.40268(3)	0.0255(6)	4c
O(3)	0.52041(6)	0.75000	0.20028(5)	0.0137(5)	4c
O(4)	0.38722(5)	0.51118(3)	0.14285(3)	0.0167(5)	8d
O(5)	0.15512(5)	0.01456(2)	0.27677(3)	0.0142(3)	8d
O(6)	0.56721(4)	0.00932(3)	0.62578(2)	0.0160(3)	8d
O(7)	0.47202(4)	0.75000	-0.01787(1)	0.0222(3)	4c
O(8)	0.58443(6)	0.25000	0.42432(4)	0.0162(3)	4c
O(9)	0.29397(4)	0.25000	0.28221(4)	0.0145(1)	4c
O(10)	0.70082(5)	0.25000	0.58552(5)	0.0145(2)	4c
O(11)	0.19817(3)	0.25000	0.46601(3)	0.0346(1)	4c

Figure SI1: Magnetic susceptibilities, $\chi = M/H$, of $K_2Co_3(MoO_4)_3(OH)_2$ as a function of temperature under an applied field of 500 Oe along the a -axis.

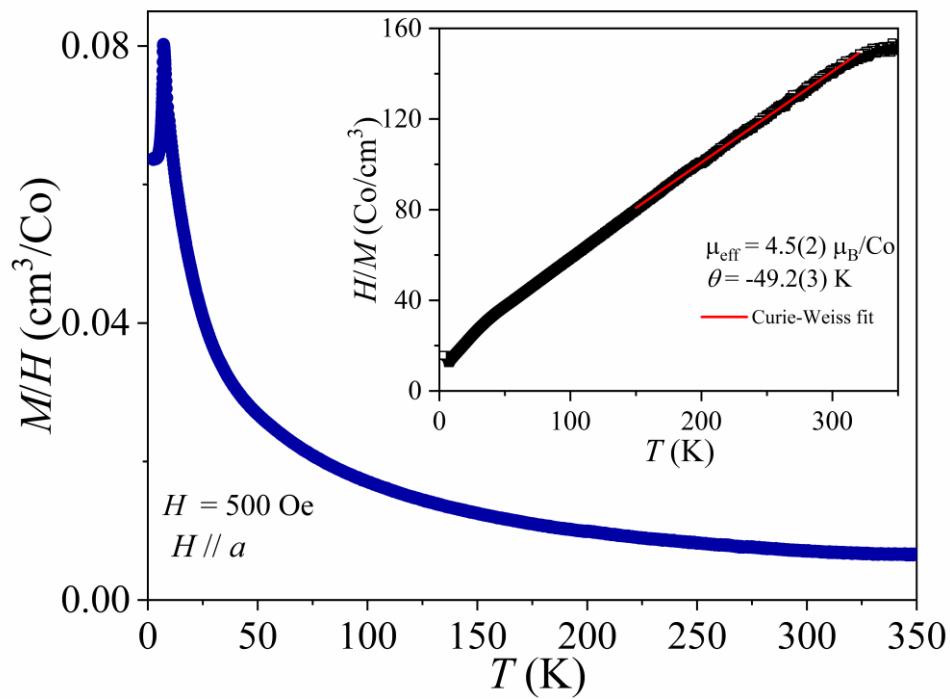


Figure SI2: Magnetic susceptibilities, $\chi = M/H$, of $K_2Co_3(MoO_4)_3(OH)_2$ as a function of temperature under an applied field of 500 Oe along the c -axis.

