



Contribution ID: 256

Type: Oral

Spin dynamics of V-based molecular magnets with integer spin values

Tuesday, 30 August 2022 09:40 (20 minutes)

In the present work, we investigate the spin dynamics of one-dimensional spin-integer molecular nanomagnets $((\text{CH}_3)_2\text{NH}_2)\text{V}_7\text{MF}_8(\text{O}_2\text{CtBu})_{162}\text{C}_7\text{H}_8$, with $\text{M}=\text{Ni}/\text{Mn}$, in short V_7M [1,2,3], by means of magnetization, susceptibility and MuSR measurements. These heterometallic nanomagnets contain seven vanadium ions ($s=1$) and one Ni^{2+} ($s=1$) or Mn^{2+} ($s=5/2$) ion, arranged in the form of regular rings. The theoretical studies of rings with a finite number of integer spins indicate a gapped ground state and a significant deviation from the Landé rule, valid for semi-integer spins [4,5]. On the other hand, the infinite spin-integer chain exhibits a topological Haldane gap between the ground state and the first excited state [6]. As confirmed by experimental data, the ground state of V_7Ni and V_7Mn is expected to be antiferromagnetic, similarly to the molecular nanomagnet V_7Zn [1,2,7], and the exchange coupling constants among the nearest neighbour magnetic ions are estimated to be of the order of a few Kelvin degrees. Susceptibility and magnetization measurements at low temperatures display anisotropy effects when an external magnetic field is applied. The muon longitudinal relaxation rate λ vs temperature, at magnetic fields $\mu_0 H \geq 500$ G, in the range $1.5 \leq T \leq 100\text{K}$, follows a heuristic Bloembergen-Purcell-Pound model [8]. No effect related to a topological gap is evinced.

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Session Classification: Oral contributions

Track Classification: Molecular chemistry and chemical physics