Proposal: 5-31-24		415			<b>Council:</b> 4/2015	
Title:	Relaxo	Relaxor ferromagnetism in nickel perovskites				
Research area: Chemistry						
This proposal is a new proposal						
Main proposer:		Peter BATTLE				
Experimental team:		Peter BATTLE Chun Mann CHIN Emily HUNTER				
Local contacts:		Emmanuelle SUARD				
Samples: La3Ni2TaO9 SrLa2Ni2WO9 SrLa2Ni2TeO9 Dy2CoGe4O12						
Instrument			Requested days	Allocated days	From	То
D2B			3	3	04/12/2015	07/12/2015
D1B			2	2	02/12/2015	04/12/2015
Abstract						

## ADSIFACI:

We have previously shown that La3Ni2SbO9 behaves as a relaxor ferromagnet below ~100 K. There are many other A3B2MO9 oxides that do not show this behaviour. We want to understand what features of the crystal structure play a role in determining whether relaxor behaviour is observed, with the eventual aim of producing a compound with a higher transition temperature. We also want to determine what other magnetic behaviours are observed when an A3B2MO9 compound is not a relaxor. We propose to study three A3B2MO9 oxides all with B = Ni and with M = Te, W, or Ta. Our magnetometry experiments suggest that one is a relaxor but the other two are not. We hope to be able to understand this by determining the details of the crystal structures, including the degree of cation ordering. In the case of the two non-relaxors we hope to determine what kind of magnetic ordering, if any, they do show below the transition temperatures that are clearly present in their magnetic susceptibilities.

The following is the abstract of a paper that has been published as *Journal of Solid State Chemistry* **243** 304 – 311 (2016).

## Structural Chemistry and Magnetic Properties of the Perovskite SrLa<sub>2</sub>Ni<sub>2</sub>TeO<sub>9</sub>

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## Abstract

A polycrystalline sample of SrLa<sub>2</sub>Ni<sub>2</sub>TeO<sub>9</sub> has been synthesized using a standard ceramic method and characterized by neutron diffraction, magnetometry and electron microscopy.

The compound adopts a monoclinic, perovskite–like structure with space group  $P2_1/n$  and unit cell parameters a = 5.6008(1), b = 5.5872(1), c = 7.9018(2) Å,  $\beta = 90.021(6)$  ° at room temperature. The two crystallographically–distinct B sites are occupied by Ni<sup>2+</sup> and Te<sup>6+</sup> in ratios of 83:17 and 50:50.

Both ac and dc magnetometry suggest that the compound is a spin glass below 35 K but the neutron diffraction data show that some regions of the sample are antiferromagnetic. Electron microscopy revealed twinning on a nanoscale and local variations in composition. These defects are thought to be responsible for the presence of two distinct types of antiferromagnetic ordering.