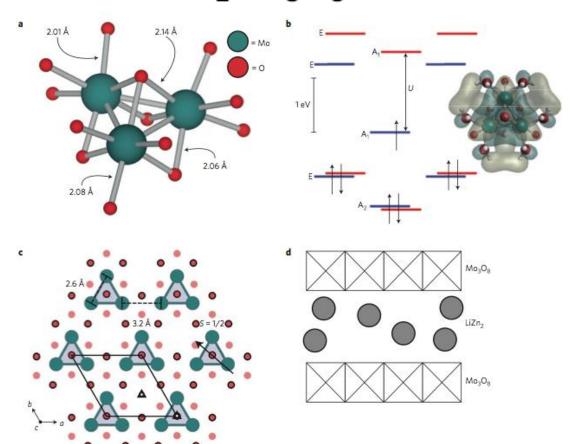
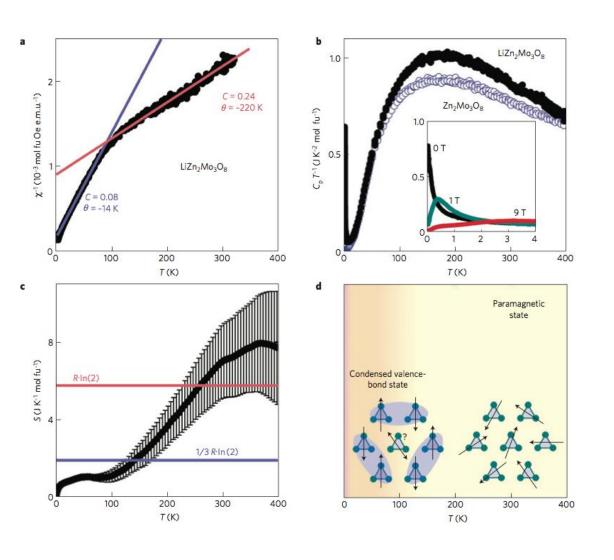
## LiZn<sub>2</sub>Mo<sub>3</sub>O<sub>8</sub>

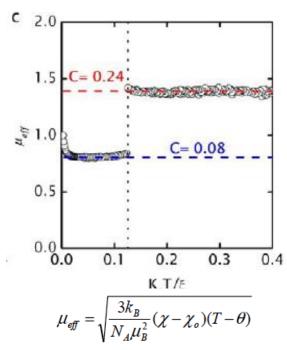
Sheckelton et al, Nature Materials 11, 493 (2012)



Molecular orbital picture of Mo triangle gives S=1/2 per triangle.

These triangles form a triangular lattice.



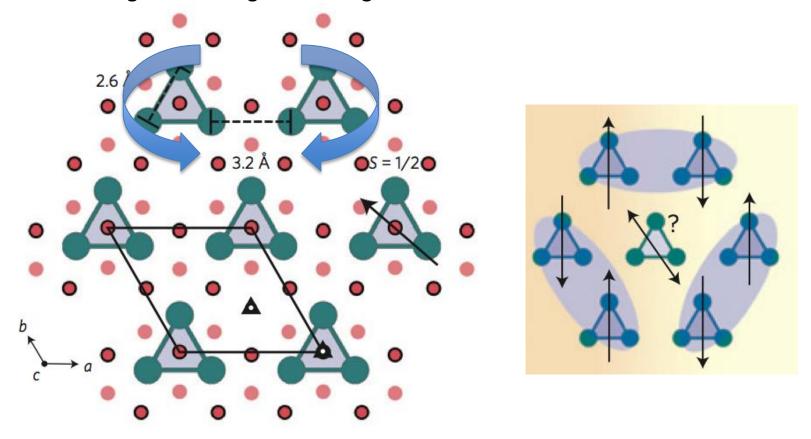


From Curie constant, g factor is 1.6, vs 2 for free spin ½. This may be due to spin orbit coupling.

$$g_m = g_e (1 - A \frac{\xi}{\Delta E})$$

for Mo is  $\xi = 0.068$  eV, similar to the value for Cu,  $\xi = 0.100$  eV

Ongoing work with Rebecca Flint. Rotation of triangles shorten bonds between triangles on hexagon and lengthen bonds to central site.



Spin ½ on hexagonal lattice forms a gapped spin liquid which co-exists with isolated moments.

J1-J2 model?

Detect rotation experimentally?