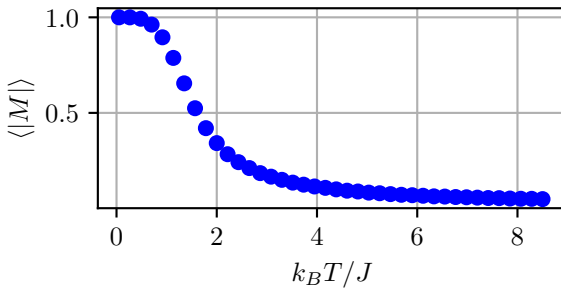


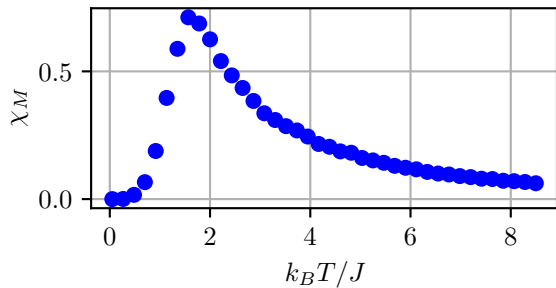
Binary Alloy Thermodynamic Quantities - MCMC Simulation

$$\mu = 0.0, h = 0.5$$

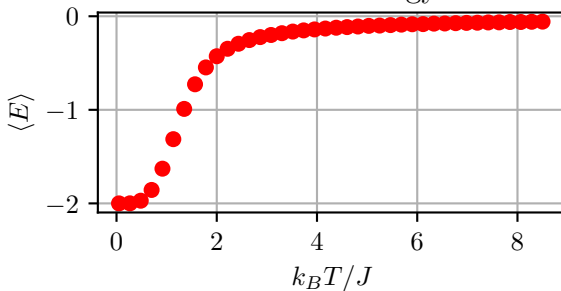
Order Parameter



Derivative of Ord. Parameter



Internal Energy



Specific Heat

