Energy bands nu=4 as function of U 0K⁺ ↓ 1K⁺ ↓ 40 -2K⁺ ↓ 2K⁺ ↓ 0K_ ↑ Energy bands(meV)

1

C

0

0 20 1K⁻ ↓ -2K⁻↓ 2K-1 0K⁺ ↑ 1K⁺ ↑ -2K⁺ ↑ 2K⁺ ↑ 0K⁻ ↑ 1K⁻ ↑ -40-2K⁻ ↑ alpha H oct int: 1 alpha int H: 1 alpha rand asymmetric calcs: 1 2K⁻ ↑ alpha_reg_asym_calcs: 1 alpha_rho_asymmetric_calcs: 0.05 apha H asym small u: 1 Fermi energy replace_LLm2_LL2_low_u: 1 screening: 0.244 -60uperp meV: -3.2 0.0 meV: 14.0 2.0 1.0 1.5 U(meV)