

# Quantum flexible quasi-molecular chains in optical lattice

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We consider dipolar molecules in layered optical lattice with  $N$  layers (which can be 2d planes or 1d "cigars") and the polarization perpendicular to the layers<sup>1,2</sup>. First-principle quantum Monte Carlo canonical simulations of  $N$  molecules arranged one per layer in the absence of inter-layer tunneling find formation of a single  $N$ -molecular chain for strong enough dipole interaction  $V$  (in units of the intra-layer band width  $t$ ). Such flexible quantum chains are analogous to classical chains found in magnetic and electric dipolar colloids<sup>3</sup>. Single chain formation is seen by columnar imaging of the density distribution map: For small  $V$  such map shows mostly single molecules distributed independently in each layer. Above some critical value  $V_c \approx 0.75$  configurations where all molecules are atop of each other dominate the density map. Gyration radius exhibits log-growth with number of layers in a narrow range of  $V$  (up to  $\sim 1.25$  in 2d), which indicates quantum rough state of single chain. At higher  $V$  this chain exhibits KT transition into autolocalized stiff (smooth) chain. The question of the nature of multi-chain state is addressed by J-current model simulations<sup>2</sup>. Superfluid phase of chains for fixed  $N$  is found, with the range of its existence shrinking as  $N$  increases.

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<sup>1</sup> D.-Wei Wang, M. D. Lukin, E. Demler Phys. Rev. Lett. **97**, 180413 (2006).

<sup>2</sup> A.B. Kuklov, talk at KITPC program " Condensed matter physics of cold atoms", Beijing, 10/08/2009: <http://www.kitpc.ac.cn/program.jsp?id=PC20090921&i=sched>

<sup>3</sup> Yu.E.Lofovik, V.A.Mandelshtam, Phys. Lett. **A138**, 204(1989).