Phase diagram of quantum loop model on triangular lattice

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Collaborators



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X. Ran, Z. Yan, Y-C Wang, R. Samajdar, J. Rong, S. Sachdev, Y. Qi, and Z. Y. Meng. Commun Phys 7, 207 (2024).
 X. Ran, Z. Yan, Y-C Wang, J. Rong, Y. Qi, and Z. Y. Meng. Phys. Rev. B 109, L241109 (2024).
 Z.Yan, Y-C Wang, R. Samajdar, S. Sachdev, Z. Y. Meng, Phys. Rev. Lett. 130, 206501 (2023).

Rydberg atom array

Rydberg atoms on the kagome lattice (links) maps to dimer model on the kagome lattice

2.5





Semeghini G, Levine H, Keesling A, et al. Science, 374, 6572 (2021).

Rydberg atoms on the kagome lattice (sites) maps to dimer\loop model on the triangular lattice



Samajdar R, Ho W W, Pichler H, et al. PNAS, 118, 4 (2021).

Quantum loop model (QLM)



K. Roychowdhury, S. Bhattacharjee, and F. Pollmann, Phys. Rev. B 92, 075141 (2015).





Sweeping cluster quantum Monte Carlo method

Based on stochastic series expansion (SSE) in the path integral $Z = \text{Tr} \{ e^{-\beta H} \}$

Plaquette operators

$$\begin{split} H_{1,p} &= -V\left(|\mathbf{u}\rangle\langle\mathbf{u}| + |\mathbf{z}\rangle\langle\mathbf{z}|\right) + V + C, \\ H_{2,p} &= t\left(|\mathbf{u}\rangle\langle\mathbf{z}| + |\mathbf{z}\rangle\langle\mathbf{u}|\right), \qquad t = 1, \ C = 1 \end{split}$$



Partition function

$$Z = \sum_{\alpha} \sum_{S_M} \frac{\beta^n (M-n)!}{M!} \left\langle \alpha \left| \prod_{i=1}^M H_{a_i, p_i} \right| \alpha \right\rangle$$

Non-zero plaquette matrix elements



|others> plaquettes have 0 or 1 dimer

Sampling in the restricted Hilbert space

Diagonal update

$$P_{\rm ins} = rac{N_p eta \langle lpha | H_{1,p} | lpha
angle}{M-n},$$

 $P_{
m del} = rac{M-n+1}{N_p eta \langle lpha | H_{1,p} | lpha
angle}.$

Z. Yan, Y. Wu, C. Liu, O. F. Syljuasen, J. Lou, and Y. Chen, Phys. Rev. B 99, 165135 (2019).

Off-diagonal (sweeping cluster) update

- (1). Choose a flippable plaquette (FP), create four update lines;
- (2). Three conditions:



(3). Metropolis acceptance probabilities

$$_{accept}(A \to B) = \min(\frac{W(B)P_{select}(B \to A)}{W(A)P_{select}(A \to B)}, 1)$$

= $\min(\frac{N_{FP}}{N_{FP} + \Delta} \left(\frac{2}{1+V}\right)^{\Delta}, 1)$

 N_{FP} : No. of flippable vertex in A $N_{FP} + \Delta : \text{ No. of flippable vertex in B}$

Z. Yan, Y. Wu, C. Liu, O. F. Syljuasen, J. Lou, and Y. Chen, Phys. Rev. B 99, 165135 (2019).

Sketches for two possible cluster updates for QDM:

2 x 2 square lattice







Sketches for a possible cluster updates for QDM:

3 x 3 square lattice







Before update

After update

Phase diagram of quantum loop model

With sweeping cluster QMC up to $L = 20, \beta \propto L$



X. Ran, Z. Yan, Y-C Wang, R. Samajdar, J. Rong, S. Sachdev, Y. Qi, and Z. Y. Meng. Commun Phys 7, 207 (2024).



Real-space vison correlation at V = 0.3

Renormalization-group (RG) analysis

3D Cubic fixed point

$$S = \int dt dx^2 \sum_{i=1}^3 (\partial_\mu \phi_i)^2 + r \sum_{i=1}^3 \phi_i^2 + \mu (\sum_{i=1}^3 \phi_i \phi_i)^2 + \nu_4 \sum_{i=1}^3 (\phi_i)^4 + \dots$$
$$r < 0, \ \nu_4 < 0 \qquad \langle \phi_1 \rangle = \pm v', \quad \langle \phi_2 \rangle = \langle \phi_3 \rangle = 0$$
$$r < 0, \ \nu_4 > 0 \qquad \langle \phi_1 \rangle = \pm v, \quad \langle \phi_2 \rangle = \pm v, \quad \langle \phi_3 \rangle = \pm v.$$

$$\nu_{4} = -\frac{1}{\textit{vol}} \frac{15\sqrt{\pi}}{\langle \phi^{4} \rangle} \left(\langle Y_{4}^{0} \rangle + \frac{\langle \phi^{2} Y_{4}^{0} \rangle \langle \phi^{4} \rangle \langle \phi^{6} \rangle - \langle Y_{4}^{0} \rangle \langle \phi^{6} \rangle^{2}}{\langle \phi^{6} \rangle^{2} - \langle \phi^{4} \rangle \langle \phi^{8} \rangle} \right)$$





rank-2 tensor (or tensorial magnetization) of the (2+1)D O(3)/cubic universality



Cubic* criticality



rank-2 tensor (or tensorial magnetization) of the (2+1)D O(3)/cubic universality

scaler of the (2+1)D O(3)/cubic universality

X. Ran, Z. Yan, Y-C Wang, J. Rong, Y. Qi, and Z. Y. Meng. Phys. Rev. B 109, L241109 (2024).



Rydberg atom array

Rydberg atoms on the kagome lattice (links) maps to dimer model on the kagome lattice



Rydberg atoms on the kagome lattice (sites) maps to dimer\loop model on the triangular lattice



Emergent glassy behavior in a kagome

Rydberg atom array

Z.Yan, Y.-C. Wang, R. Samajdar, S. Sachdev, Z. Y. Meng, Phys. Rev. Lett. 130, 206501 (2023).

Phase diagram

$$\begin{split} H = &\sum_{i=1}^{N} \left[\frac{\Omega}{2} \left(\left| g \right\rangle_{i} \left\langle r \right| + \left| r \right\rangle_{i} \left\langle g \right| \right) - \delta \left| r \right\rangle_{i} \left\langle r \right| \right] \\ &+ \sum_{i,j=1}^{N} \frac{V_{ij}}{2} \left(\left| r \right\rangle_{i} \left\langle r \right| \otimes \left| r \right\rangle_{j} \left\langle r \right| \right), \end{split}$$

Repulsive interaction $V_{ij} = \Omega R_b^6 / R_{ij}^6$ $\Omega = 1$

- Ω : Rabi frequency, transverse field
- δ : detuning, longitudinal field
- R_b : Rydberg blockade radius
- $R_{i,j}$: distance between the sites *i* and *j*



$$S(\mathbf{k}) = \frac{1}{N} \sum_{i,j,\alpha=1,2,3}^{N} e^{i\mathbf{k}\cdot\mathbf{r}_{ij}} \langle \langle n_{i,\alpha}n_{j,\beta} \rangle - \langle n_{i,\alpha} \rangle \langle n_{j,\beta} \rangle)$$



Z Yan, YC Wang, R Samajdar, S Sachdev, ZY Meng, Phys. Rev. Lett. 130, 206501 (2023).

Phase transition

Nonlocal loop operator $< loop > = (-1)^{N_{P_{ij}}} N_{P_{ij}} : #cut dimers$

Edwards-Anderson order parameter

$$q_{\rm EA} = \sum_{i=1}^{N} \langle n_i - \rho \rangle^2 / [N\rho(1-\rho)] \qquad n_i \equiv |r\rangle_i \langle r| \qquad \rho \equiv \sum_{i=1}^{N} \langle n_i \rangle / N_i$$



Glassy dynamics



Triangular lattice quantum dimer model

with variable dimer density

Rydberg atom array





Rydberg atoms on the kagome lattice (sites) maps to dimer\loop model on the triangular lattice



Samajdar R, Ho W W, Pichler H, et al. PNAS, 118, 4 (2021).



Soft constraint: one or two dimer(s) per site

$$H = -t \sum_{r} \left(\left| \mathbf{I} \right\rangle \left\langle \mathbf{I} \right| + \text{h.c.} \right) \\ +V \sum_{r} \left(\left| \mathbf{I} \right\rangle \left\langle \mathbf{I} \right| + \left| \mathbf{I} \right\rangle \left\langle \mathbf{I} \right| \right) \\ -h \sum_{l} \left(\left| \mathbf{I} \right\rangle \left\langle \mathbf{I} \right| + \text{h.c.} \right) \\ -\mu \sum_{l} \left(\left| \mathbf{I} \right\rangle \left\langle \mathbf{I} \right\rangle \right\rangle \left\langle \mathbf{I} \right| \right),$$

t = 1

h: transverse- field term of strength

 μ : chemical potential



Z Yan, R Samajdar, YC Wang, S Sachdev, ZY Meng, Nat Commun 13, 5799 (2022).

Phase diagram

$$H = -t \sum_{r} \left(\left| \mathbf{I} \right\rangle \left\langle \mathbf{I} \right| + \text{h.c.} \right) \qquad t = 1$$
$$+V \sum_{r} \left(\left| \mathbf{I} \right\rangle \left\langle \mathbf{I} \right| + \left| \mathbf{I} \right\rangle \left\langle \mathbf{I} \right| \right)$$
$$-h \sum_{l} \left(\left| \mathbf{I} \right\rangle \left\langle \mathbf{I} \right| + \text{h.c.} \right)$$
$$-\mu \sum_{l} \left(\left| \mathbf{I} \right\rangle \left\langle \mathbf{I} \right| \right),$$

- h: transverse- field term of strength
- μ : chemical potential



Only (a) allowed to create/annihilate a dimer on the centre link

Equal-time dimer-structure factor $\tau = 0$ $\mu = -3$ $\mu = 3$ $\mu = -3$ $\mu = 3$ $\mu = 0$ V = 0.9V = 0.9V = -0.5V = -0.5V = 0.9(e) 1/25 $D(\mathbf{k},\tau) = \frac{1}{N} \sum_{i,j}^{L^3} e^{i\mathbf{k}\cdot\mathbf{r}_{ij}} \left(\langle n_{i,\alpha}(\tau)n_{j,\alpha}(0) \rangle - \langle n_{i,\alpha} \rangle \langle n_{j,\alpha} \rangle \right)$ π π π 0 0 0 0 0 a=1,2,3 $-\pi$ $-\pi$ $-\pi$ $-\pi$ 0 π 0 0 π 0 π 0 π -π π $-\pi$ $-\pi$ -π $-\pi$ Odd QSL PM Even QSL Columnar Nematic

Phase transitions between QSLs and the PM phase

V = 0.9 h = 0.4

Polarization

String operator



Dynamical dimer spectra



Phase diagram of quantum loop model on

the square lattice

Phase diagram

$$egin{aligned} H &= - t \sum_{ ext{plaq}} \left(\left| egin{aligned} igstarrow
ight
angle \left\langle igstarrow igstarrow
ight
angle + ext{H.c.}
ight
angle \ &+ V \sum_{ ext{plaq}} \left(\left| igstarrow
ight
angle \left\langle igstarrow igstarrow igstarrow
ight
angle \left\langle igstarrow igstarrow igstarrow
ight
angle \left\langle igstarrow igstarrow$$

Order parameter for rotational symmetry breaking





Translational symmetry breaking



X. Ran, Z. Yan, Y.-C. Wang, J. Rong, Y. Qi, and Z. Y. Meng, Phys. Rev. B 107, 125134 (2023).

Effective height field action

$$\mathcal{L} = \frac{1}{2} (\partial_{\tau} h)^2 + \frac{1}{2} \rho_2 (\nabla h)^2 + \frac{\kappa^2}{2} (\nabla^2 h)^2 + \lambda \cos(4\pi h)$$



NG

e RP

 $\tilde{\rho}_2$

homogenous h





L = 16

