

# CP(2) Model at Non-Zero Chemical Potential

via dimensional reduction of a quantum spin system

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# Motivation

- i) The  $(1+1)$ -d  $\mathbb{C}P(2)$  model is a popular toy model for QCD, since it exhibits:
  - Asymptotic freedom
  - Dynamically generated mass gap
  - Non-trivial topology
  
- ii) Obtaining the  $\mathbb{C}P(2)$  model from a  $(2+1)$ -d system of quantum spins provides experimentalists with a clear analogy to cold atoms in an optical lattice: C. Laflamme et al., *Annals of Physics* (2016), pp. 117-127
  - The ultimate goal is to quantum simulate QCD at finite chemical potential

# Formulation Overview

i) Microscopic (2+1)-d antiferromagnetic  $SU(3)$  spin system:

$$H = -J \sum_{\substack{\langle ij \rangle \\ i \in A}} T_i^a T_j^{a*} - \mu^a \left( \sum_{i \in A} T_i^a - \sum_{j \in B} T_j^{a*} \right), \quad J > 0.$$

ii) Spontaneous symmetry breaking:  $SU(3)/U(2) = \mathbb{C}P(2)$ .  
Emergent Nambu-Goldstone bosons are governed by:

$$S[P] = \int_0^\beta dt \int_0^L dx \int_0^{L'} dy \operatorname{Tr} \left[ \rho_s \partial_i P \partial_i P + \frac{\rho_s}{c^2} D_t P D_t P \right] - i\theta Q[P].$$

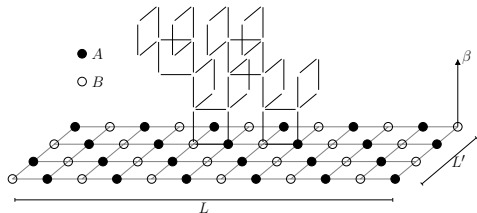
iii) Dimensional reduction:  $\xi \propto \exp[4\pi L' \rho_s / 3c] \gg L'$

$$S[P] = \int_0^\beta dt \int_0^L dx \operatorname{Tr} \left[ \rho_s L' \partial_x P \partial_x P + \frac{\rho_s L'}{c^2} D_t P D_t P \right] - i\theta Q[P].$$

# Microscopic (2+1)-d Quantum Spin System

$$H = -J \sum_{\substack{\langle ij \rangle \\ i \in A}} T_i^a T_j^{a*} - \mu^a \left( \sum_{i \in A} T_i^a - \sum_{j \in B} T_j^{a*} \right)$$

- $[T_x^a, T_y^a] = i\delta_{xy} f_{abc} T_x^c$ ,  $\bar{T}_x^a = -T_x^{a*}$ .
- Global  $SU(3)$  symmetry at  $\mu = 0$ .
- Total spin conservation,  $\left[ H, \left( \sum_{x \in A} T_x^a - \sum_{y \in B} T_y^{a*} \right) \right] = 0$ .
- $J > 0$  gives anti-ferromagnet.
- Spins live in the fundamental and antifundamental representations of  $SU(3)$  on sublattice A and B respectively.



# Spontaneous Symmetry Breaking

- In the thermodynamic limit:  $\beta, L, L' \rightarrow \infty$ ,

$$\text{Global } SU(3) \rightarrow U(2).$$

K. Harada, N. Kawashima, M. Troyer, PRL 90, 117203.

- Consequently 4 Nambu-Goldstone bosons emerge with degrees of freedom in the coset space:

$$\frac{SU(3)}{U(2)} = \mathbb{C}P(2).$$

- These massless fields can be described by complex  $3 \times 3$  Hermitean projection matrices  $P$  using:

$$P(x) = P(x)^2, \quad \text{Tr } P(x) = 1, \quad P(x) = P(x)^\dagger.$$

- Effective action for  $\mathbb{C}P(2)$  model,

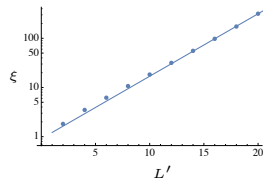
$$S[P] = \int_0^\beta dt \int_0^L dx \int_0^{L'} dy \text{Tr} \left[ \rho_s \partial_i P \partial_i P + \frac{\rho_s}{c^2} D_t P D_t P \right] - i\theta Q[P],$$

$$i\theta Q[P] = \frac{1}{\pi} \int_0^\beta dt \int_0^L dx \int_0^{L'} dy \text{Tr} [P \partial_x P \partial_t P], \quad D_t P = \partial_t P + [\mu_a N_a, P].$$

# Dimensional Reduction

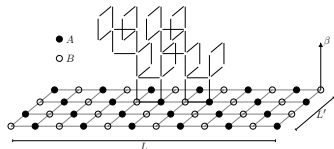
- (1+1)-d  $\mathbb{CP}(2)$  model is asymptotically free:

$$\xi \propto \exp \left[ \frac{4\pi L' \rho_s}{3c} \right].$$



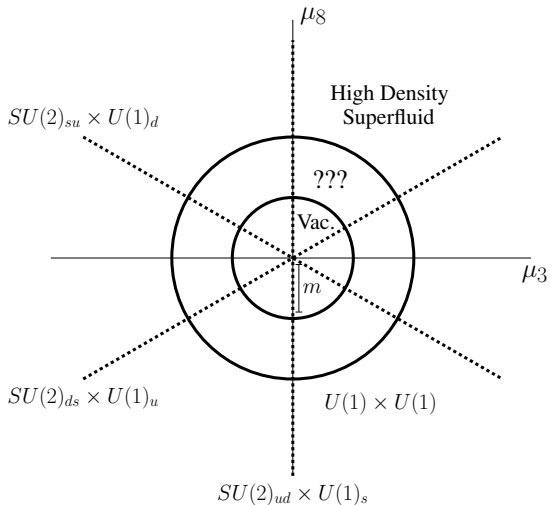
- For finite  $L'$ , Mermin-Wagner theorem states Nambu-Goldstone bosons must pick up a mass  $m = 1/\xi$ .
- If  $\xi \gg L'$  then system undergoes dimensional reduction in the  $L'$  extent.
- Effective action after dimensional reduction,

$$S[P] = \frac{c}{g^2} \int_0^\beta dt \int_0^L dx \text{Tr} \left[ \partial_x P \partial_x P + \frac{1}{c^2} D_t P D_t P \right] - i\theta Q[P],$$



$$g^2 = \frac{c}{\rho_s L'}.$$

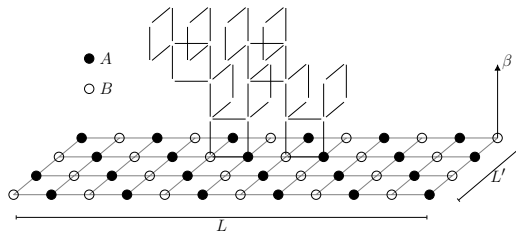
# (1+1)-d $\mathbb{CP}(2)$ Phase Diagram



$\lambda_8 \pm \frac{3}{\sqrt{3}} \lambda_3$  has  $SU(2) \times U(1)$  symmetry.

# Simulation

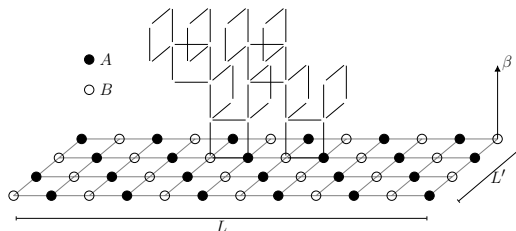
- The (2+1)-d spin system is quantized in the  $T_3$ - $T_8$  basis.
- We have  $u$ ,  $d$ , or  $s$  on sublattice A and  $\bar{u}$ ,  $\bar{d}$ , or  $\bar{s}$  on sublattice B.
- The time extent is discretized with the Trotter decomposition to give four time-slices for each Euclidean time step  $\epsilon$ .
- The system is updated using a worm algorithm. Typically  $10^5$  thermalization updates and  $10^6$  measurement updates.





# Worm Algorithm

- Randomly pick lattice site, we note the spin color at this site.
- Randomly choose one of the two spin colors not at starting site.
- Flip start site spin to the color chosen, creates the worm head or defect.
- Worm head propagates according to the rules derived from the Hamiltonian, flipping spins as it goes.
- Worm closes once the worm head returns to the starting site where the worm tail remained.
- Defect is eliminated and system has been updated.



# Observables

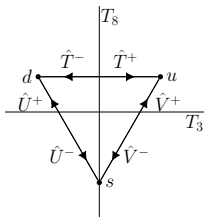
- Particle Number Density

$$\langle n \rangle = \frac{\langle M_3 \rangle}{L},$$

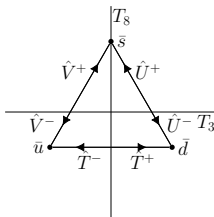
$$M_3 = \sum_{x \in A} T_x^a - \sum_{y \in B} T_y^{a*},$$

$M_3$  is calculated in a particular time-slice.

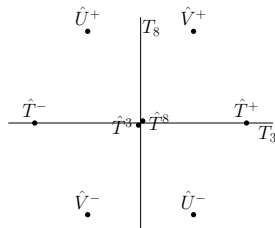
- Shift Operator Correlators



$$\langle U_0^+ U_x^- + U_0^- U_x^+ \rangle$$



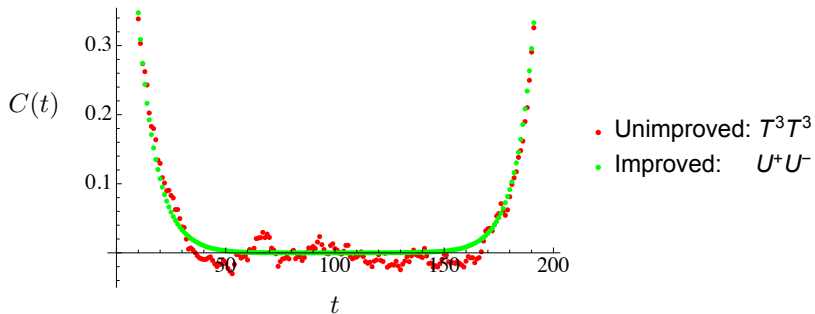
$$\langle T_0^+ T_x^- + T_0^- T_x^+ \rangle$$



$$\langle V_0^+ V_x^- + V_0^- V_x^+ \rangle$$

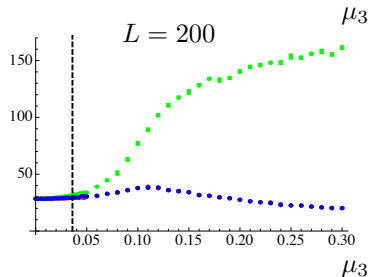
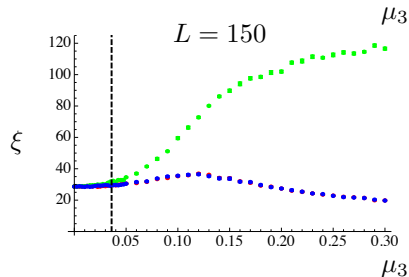
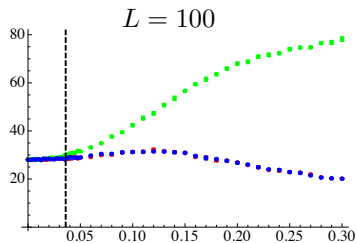
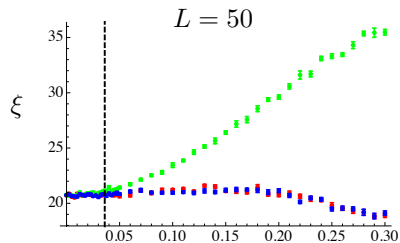
# Improved vs. Unimproved Correlator Estimator

$$\beta = 10, \quad L = 10, \quad L' = 2.$$



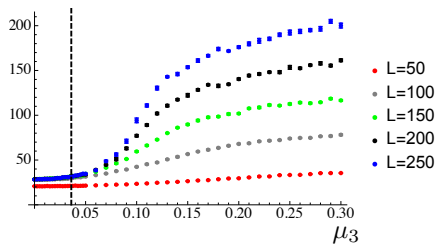
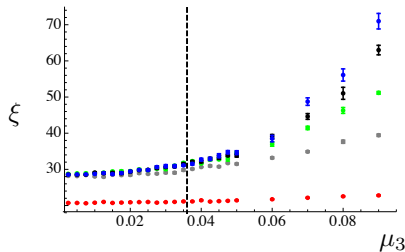
# Correlation Length vs. Chemical Potential

$\beta c \approx L$ ,  $L' = 10$ ,  $\xi = 28.5(2)$  at  $\mu = 0$ .

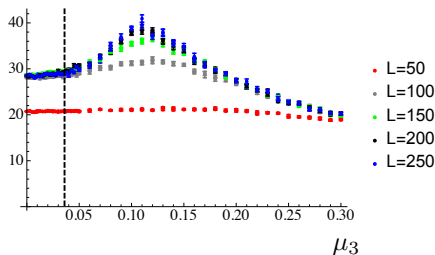
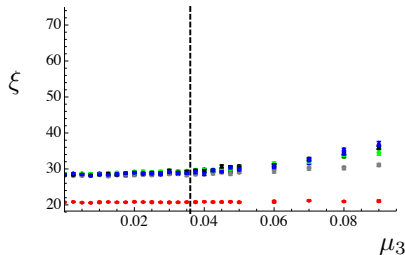


# Correlation Length vs. Chemical Potential

$$\langle T_0^+ T_x^- + T_0^- T_x^+ \rangle, \quad \beta c \approx L, \quad L' = 10.$$

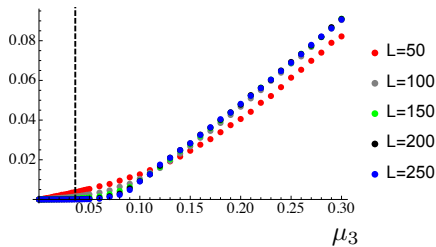
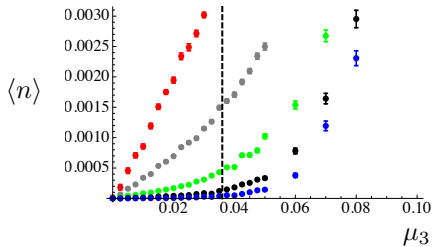
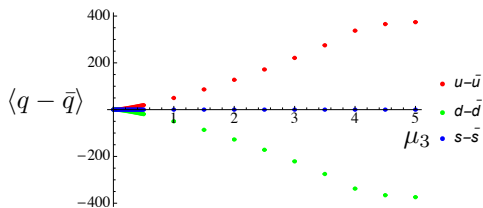


$$\langle U_0^+ U_x^- + U_0^- U_x^+ \rangle, \text{ and } \langle V_0^+ V_x^- + V_0^- V_x^+ \rangle.$$



# Magnetization vs. Chemical Potential

$$\beta c \approx L, \quad L' = 10.$$



## Conclusions and Outlook

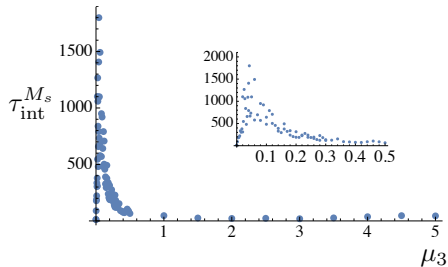
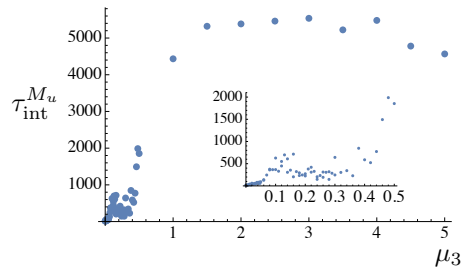
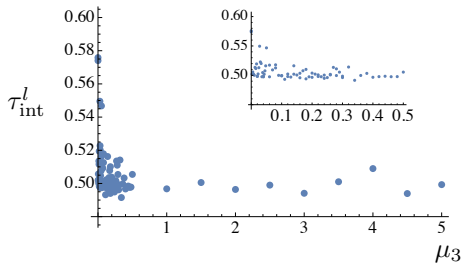
- Behaviour of the magnetization at  $\mu_3 \neq 0$  implies dimensional reduction occurs.
- At  $\mu_8 = 0$ ,  $\mu_3 \neq 0$  the behaviour of the correlation functions suggests  $U(1) \times U(1) \rightarrow U(1)$ .
- Study  $\mu_8 \neq 0$ , and  $\mu_3, \mu_8 \neq 0$ .
- Cold atom experiments.

# Autocorrelation vs. Chemical Potential

$$L = 150$$

$$\beta c \approx L$$

$$L' = 10$$





# Correlation Length vs. Chemical Potential

$$\langle T_0^+ T_x^- + T_0^- T_x^+ \rangle, \quad \beta c \approx L, \quad L' = 10.$$

