

"Phase diagram of the GWW phase transition of the matrix quantum mechanics with a chemical potential"

Takehiro Azuma (Setsunan Univ.) From Strings to LHC IV, Sinclairs Retreat, Chalsa Mar. 10th 2017, 14:25-14:50 with Pallab Basu (ICTS,TIFR) and Prasant Samantray (IIT, Indore)



Thermodynamic aspects of quantum gravity in AdS spacetime:

- Small blackhole (SBH): Unstable. Horizon radius smaller than AdS
  Big blackhole (BBH):
  - Stable. Horizon radius comparable to AdS

Gross-Witten-Wadia (GWW) phase transition of the gauge theory and the blackhole phase transition

[L. Alvarez-Gaume, C. Gomez, H. Liu and S.R. Wadia, hep-th/0502227]

#### 2. The model



Finite-temperature matrix quantum mechanics with a chemical potential  $S = S_b + S_f + S_q$ , where (µ=1,2,...,D, β=1/T) 
$$\begin{split} S_{\rm b} &= N \int_0^\beta \operatorname{tr} \left\{ \frac{1}{2} \sum_{\mu=1}^D (D_t X_\mu(t))^2 - \frac{1}{4} \sum_{\mu,\nu=1}^D [X_\mu(t), X_\nu(t)]^2 \right\} dt \\ D_t X_\mu(t) &= \partial_t X_\mu(t) - i [A(t), X_\mu(t)] \\ S_{\rm f} &= N \int_0^\beta \operatorname{tr} \left\{ \sum_{\alpha=1}^p \bar{\psi}_\alpha(t) D_t \psi_\alpha(t) - \sum_{\mu=1}^D \sum_{\alpha,\eta=1}^p \bar{\psi}_\alpha(t) (\Gamma_\mu)_{\alpha\eta} [X_\mu(t), \psi_\eta(t)] \right\} dt \\ S_{\rm g} &= N \mu (\operatorname{tr} U + \operatorname{tr} U^\dagger) \quad U = \mathcal{P} \exp \left( i \int_0^\beta A(t) dt \right) \end{split}$$

•Bosonic (S=S<sub>b</sub>+S<sub>g</sub>): D is an arbitrary integer D=2,3,... •Fermionic(S=S<sub>b</sub>+S<sub>f</sub>+S<sub>g</sub>): (D,p)=(3,2),(5,4),(9,16) (For D=9, the fermion is Majorana-Weyl ( $\Psi \rightarrow \Psi$ ) In the following, we focus on D=3.)

#### 2. The model



A(t),  $X_{\mu}(t)$ ,  $\Psi(t)$ : N × N Hermitian matrix Boundary conditions:  $\begin{array}{l} A(t+\beta) = A(t), \quad X_{\mu}(t+\beta) = X_{\mu}(t) \\ \psi(t+\beta) = -\psi(t) \end{array}$ 

Static diagonal gauge:

$$A(t) = \frac{1}{\beta} \operatorname{diag}(\alpha_1, \alpha_2, \cdots, \alpha_N) \quad -\pi \leq \alpha_k < \pi$$

 $\Rightarrow$  Add the gauge-fixing term  $S_{g.f.} = -\sum_{k,l=1,k\neq l}^{N} \log \left| \sin \frac{\alpha_k - \alpha_l}{2} \right|$ 

Under this gauge  $u_n = \frac{1}{N} \operatorname{tr} U^n = \frac{1}{N} \sum_{k=1}^{N} e^{in\alpha_k}$ 

Supersymmetry for  $S=S_b+S_f$  (µ=0), broken at µ≠0.

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# 2. The model



Previous works for  $\mu=0$  (without  $S_a$ )



SUSY (S=S<sub>b</sub>+S<sub>f</sub>)



[Quoted for D=9 from N. Kawahara, J. Nishimura and S. Takeuchi, arXiv:0706.3517]

Confinement-deconfinement phase transition at T=T<sub>c0</sub> Absence of phase transition  $<|u_1|>=a_0 \exp(-a_1/T)$ 

[Quoted for D=9 from K.N. Anagnostopoulos, M. Hanada,

J. Nishimura and S. Takeuchi, arXiv:0707.4454]



- Bosonic model without fermion S=S<sub>b</sub>+S<sub>g</sub> [T. Azuma, P. Basu and S.R. Wadia, arXiv:0710.5873]
- Result of D=3 (D=2,6,9 cases are similar)



At  $(\mu_c, T_c)$ ,  $d < |u_{1,2}| > /d\mu$  and  $d < |u_{1,2}| > /dT$  are not smooth  $(d^2 < |u_{1,2}| > /d\mu^2$  and  $d^2 < |u_{1,2}| > /dT^2$  are discontinuous)  $\Rightarrow$  suggests third-order phase transition.





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When  $\mu$ =0, at the critical point T<sub>c0</sub>=1.1, there is a first-order phase transition at small D. [T. Azuma, T. Morita and S. Takeuchi, arXiv:1403.7764]

We fit the susceptibility with (y,p,c) as

$$\chi = N^2 \{ \langle |u_1|^2 \rangle - (\langle |u_1| \rangle)^2 \} = \gamma V^p + c \ (V = N^2)$$

 $p=1 \Rightarrow$  suggests first-order phase transition.

[M. Fukugita, H. Mino, M. Okawa and A. Ukawa, Phys. Rev. Lett.65, 816 (1990)]





The model with fermion (D=3)  $S=S_b+S_f+S_g$  $\Gamma_{\mu}=\sigma_{\mu} (\mu=1,2,3) \qquad \sigma_1 = \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}, \quad \sigma_2 = \begin{pmatrix} 0 & -i \\ i & 0 \end{pmatrix}, \quad \sigma_3 = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}$ 

Non-lattice simulation with Fourier expansion [K.N. Anagnostopoulos, M. Hanada, J. Nishimura and S. Takeuchi, arXiv:0707.4454]

$$\begin{split} X_{\mu}^{kl}(t) &= \sum_{n=-\Lambda}^{\Lambda} \tilde{X}_{\mu,n}^{kl} e^{i\omega nt}, \quad \psi_{\alpha}^{kl}(t) = \sum_{r=-\Lambda+\frac{1}{2}}^{\Lambda-\frac{1}{2}} \tilde{\psi}_{\alpha,r}^{kl} e^{i\omega rt}, \quad \bar{\psi}_{\alpha}^{kl}(t) = \sum_{r=-\Lambda+\frac{1}{2}}^{\Lambda-\frac{1}{2}} \tilde{\psi}_{\alpha,-r}^{kl} e^{i\omega rt}. \quad \left(\omega = \frac{2\pi}{\beta}\right) \\ S_{\text{F,Fourier}} &= N\beta \sum_{r=-\Lambda+\frac{1}{2}}^{\Lambda-\frac{1}{2}} \left\{ i \left\{ r\omega - \frac{\alpha_k - \alpha_l}{\beta} \right\} \tilde{\psi}_{\alpha,r}^{lk} \tilde{\psi}_{\alpha,r}^{kl} - (\sigma_{\mu})_{\alpha\eta} \text{tr} \left\{ [\tilde{\psi}_{\alpha,r} \left( [\tilde{X}_{\mu}, \tilde{\psi}_{\eta}] \right)_r \right\} \right\} \\ \left( f^{(1)} \cdots f^{(p)} \right)_q &= \sum_{k_1+\cdots+k_p=q} f^{(1)}_{k_1} \cdots f^{(p)}_{k_p} \text{ Integrating out } \Psi \Rightarrow N_0 \times N_0 \text{ matrix } \mathcal{M} \\ \left( N_0 = 2 \times (2\Lambda) \times (N^{2}-1) \right) \text{ traceless adjoint } \\ \text{Gamma matrix [4(D=5), 16(D=9)]} \quad r=-\Lambda+1/2, \dots \Lambda-1/2 \end{split}$$

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Result of D=3, N=16, after large- $\Lambda$  extrapolation:



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History of  $R^2 = \frac{1}{N\beta} \int_0^\beta dt \, \text{tr} X_\mu(t)^2$ at  $\Lambda = 3$ 

No instability in the typical  $(\mu,T)$  region.



Phase diagram for D=2,3,6,9 (boson) and D=3(fermion) . Some phase transitions at ( $\mu_c$ ,T<sub>c</sub>) where <|u\_1|>=0.5



Fitting of the	
critical point	by
$T_c = a(0.5 - \mu_c)^{b}$	

	D	2(boson)	3(boson)	6(boson)	9(boson)	3(fermion)
/	а	1.36(12)	1.01(15)	0.91(9)	0.90(8)	1.39(72)
	b	0.55(6)	0.34(7)	0.25(4)	0.23(4)	2.30(59)

# 5. Summary



- We have studied the matrix quantum mechanics with a chemical potential  $S_g = N\mu(trU + trU^{\dagger})$
- •bosonic model  $\Rightarrow$  GWW-type third-order phase transition (except for very small  $\mu$ )
- •phase diagram of the bosonic/fermionic model

#### backup: GWW phase transition





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[D.J. Gross and E. Witten, Phys. Rev. D21 (1980) 446, S.R. Wadia, Phys. Lett. B93 (1980) 403]

# Gross-Witten-Wadia (GWW) type third-order phase transition



#### backup: GWW phase transition



#### Continuous at $\mu = 1/2$ .

$$\langle |u_1| \rangle = \begin{cases} \mu & \left( 0 \leq \mu \leq \frac{1}{2} \right) \\ 1 - \frac{1}{4\mu} & \left( \mu \geq \frac{1}{2} \right) \end{cases}, \quad \langle |u_2| \rangle = \begin{cases} 0 & \left( 0 \leq \mu \leq \frac{1}{2} \right) \\ \left( 1 - \frac{1}{2\mu} \right)^2 & \left( \mu \geq \frac{1}{2} \right) \end{cases}$$
$$\frac{d\langle |u_1| \rangle}{d\mu} = \begin{cases} 1 & \left( 0 \leq \mu \leq \frac{1}{2} \right) \\ \frac{1}{4\mu^2} & \left( \mu \geq \frac{1}{2} \right) \end{cases}, \quad \frac{d^2 \langle |u_1| \rangle}{d\mu^2} = \begin{cases} 0 & \left( 0 \leq \mu \leq \frac{1}{2} \right) \\ -\frac{1}{2\mu^3} & \left( \mu \geq \frac{1}{2} \right) \end{cases}$$

Discontinuous at  $\mu = 1/2$ .  $\Rightarrow$  For free energy, d<sup>3</sup>F/d $\mu$ <sup>3</sup> is discontinuous.



#### backup: RHMC



Simulation via Rational Hybrid Monte Carlo (RHMC) algorithm. [Chap 6,7 of B.Ydri, arXiv:1506.02567, for a review]



## backup: RHMC



Hot spot (most time-consuming part) of RHMC:  $\Rightarrow$ Solving  $(\mathscr{D} + b_k)\chi_k = F$   $(k = 1, 2, \dots, Q)$ by conjugate gradient (CG) method.

Multiplication  $\mathcal{M}\chi_k \Rightarrow$ 

 $\mathscr{M}$  is a very sparse matrix. No need to build  $\mathscr{M}$  explicitly.  $\Rightarrow$  CPU cost is O(N<sup>3</sup>) per CG iteration The required CG iteration time depends on T. (while direct calculation of  $\mathscr{M}^{-1}$  costs O(N<sup>6</sup>).)

Multimass CG solver: [B. Jegerlehner, hep-lat/9612014] Solve  $(\mathscr{D} + b_k)\chi_k = F$  only for the smallest  $b_k$  $\Rightarrow$ The rest can be obtained as a byproduct, which saves O(Q) CPU cost.

#### backup: RHMC

Conjugate Gradient (CG) method:

Iterative algorithm to solve the linear equation Ax=b (A: symmetric, positive-definite n × n matrix)

Initial config.  $\mathbf{x}_0 = 0$   $\mathbf{r}_0 = \mathbf{b} - A\mathbf{x}_0$   $\mathbf{p}_0 = \mathbf{r}_0$ (for brevity, no preconditioning on  $\mathbf{x}_0$  here)

$$\mathbf{x}_{k+1} = \mathbf{x}_k + \boldsymbol{\alpha}_k \mathbf{p}_k \quad \mathbf{r}_{k+1} = \mathbf{r}_k - \boldsymbol{\alpha}_k A \mathbf{p}_k \quad \boldsymbol{\alpha}_k = \frac{(r_k, r_k)}{(p_k, A p_k)}$$
$$\mathbf{p}_{k+1} = \mathbf{r}_{k+1} + \frac{(\mathbf{r}_{k+1}, \mathbf{r}_{k+1})}{(\mathbf{r}_k, \mathbf{r}_k)} \mathbf{p}_k$$
Iterate this until  $\sqrt{\frac{(\mathbf{r}_{k+1}, \mathbf{r}_{k+1})}{(\mathbf{r}_0, \mathbf{r}_0)}} < (\text{tolerance}) \simeq 10^{-4}$ 

The approximate answer of Ax=b is  $x=x_{k+1}$ .

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