# Lattice quantum electrodynamics in (3+1)-dimensions 

 at finite density with tensor networksG. Magnifico, T. Felser, P. Silvi, and S. Montangero, Nature Commun. 12 (2021) 1

Abstract<br>The first tensor network simulation of<br>(3+1)D LGT with dynamical matter in the Hamiltonian formalism

## Tensor Network (TN) methods

Tensor network = A contraction of many \# of tensors
TN methods are free from the sign problem


Hamiltonian formalism

$$
E_{0} \leq \frac{\langle\psi| \widehat{H}|\psi\rangle}{\langle\psi \mid \psi\rangle}
$$

Prepare $|\psi\rangle$ as a TN

Variational method w/ TN

Ex. DMRG

## TN method based on the Hamiltonian formalism

The goal is to obtain the ground state, $|\Psi\rangle=\sum_{i_{1}, \cdots, i_{N}=1}^{m} \Psi_{i_{1} \cdots i_{N}}\left|i_{1}, \cdots, i_{N}\right\rangle$

We assume a TN form of $|\Psi\rangle$ and variationally tune it

Ex) Matrix Product State (MPS) ansatz

$$
|\Psi\rangle=\sum_{i_{1}, \cdots, i_{N}=1}^{m} \operatorname{Tr}\left[A^{\left(i_{1}\right)} A^{\left(i_{2}\right)} \cdots A^{\left(i_{N}\right)}\right]\left|i_{1}, \cdots, i_{N}\right\rangle \quad A^{\left(i_{k}\right)}: \chi \times \chi \text { matrix }
$$


$m^{N}$


$$
N m \chi^{2} \sim \operatorname{poly}(N)
$$

The bond dimension $\chi$ determines \# of variational parameters

## Examples of TN ansatz



| Type of ansatz | Cost |
| :---: | :---: |
| MPS | $O\left(\chi^{3}\right)$ |
| PEPS | $O\left(\chi^{10}\right)$ |
| MERA | $O\left(\chi^{8}\right)$ |
| Tree Tensor Network (TTN) | $\boldsymbol{O}\left(\chi^{4}\right)$ |

Figure from Bañuls-Cichy, Rep. Prog. Phys. 83(2020)024401
Cf. Nishino san's lecture (11/17, 22, 24)

The following TN simulation is based on three-dimensional TTN
Computationally economic \& Straightforward extension to higher dim.

- Area law in $d$ dimensions $(d \geq 2)$


## The model

$\checkmark$ Lattice QED based on the KS formalism
fermions on even (odd) sites are (anti-)particle w/ positive (negative) charge

$$
\begin{array}{r}
H=-t \sum_{x, \mu}\left(\psi_{x}^{\dagger} U_{x, \mu} \psi_{x+\mu}+\text { h.c. }\right)+m \sum_{x}(-1)^{x} \psi_{x}^{\dagger} \psi_{x} \\
+\frac{g_{\mathrm{e}}^{2}}{2} \sum_{x, \mu} E_{x, \mu}^{2}-\frac{g_{\mathrm{m}}^{2}}{2} \sum_{x}\left(\square_{\mu_{x} \mu_{y}}+\square_{\mu_{y} \mu_{z}}+\square_{\mu_{x} \mu_{z}}+\text { h. c. }\right)
\end{array}
$$

with

$$
\square_{\mu_{\alpha} \mu_{\beta}}=U_{x, \mu_{\alpha}} U_{x+\mu_{\alpha}, \mu_{\beta}} U_{x+\mu_{\beta}, \mu_{\alpha}}^{\dagger} U_{x, \mu_{\beta}}^{\dagger}
$$

$\boldsymbol{V}(1)$ electromagnetic fields are truncated up to a spins representation The original model is restored by the limit $s \rightarrow \infty$
$\checkmark$ TN simulation is implemented selecting $s=1$ (smallest but nontrivial one)

## Transition at zero total charge w/ PBC


$m_{c}$ is shifted by the effect of magnetic coupling

## Quantum capacitor w/ OBC

Charge density along $\mathrm{w} / \mu_{x}$ direction


## Confinement properties



## Surface charge density at finite density

a



## Summary

$\checkmark$ Frist study of (3+1)D LGT with dynamical matter using the TN method based on the Hamiltonian formalism
$\checkmark$ Confinement property has been confirmed
$\boldsymbol{\checkmark}$ TTN works efficiently in this study
$\checkmark$ How about the large-volume calculation w/ TN in the Hamiltonian formalism? $s>1$ ?
"a single simulation for the maximum size that we reached, an $8 \times 8 \times 8$ lattice, can last up to five weeks until final convergence, depending on the different regimes of the model and the control parameters of the algorithms"

## Backup

## Tree tensor network (TTN)



## Convergence



The relative error of the energy is in the range of $\left[10^{-2}, 10^{-4}\right]$

## The model (1/2)

$$
\begin{aligned}
& H=-t \sum_{x, \mu}\left(\psi_{x}^{\dagger} U_{x, \mu} \psi_{x+\mu}+\text { h.c. }\right)+m \sum_{x}(-1)^{x} \psi_{x}^{\dagger} \psi_{x} \\
& +\frac{g_{\mathrm{e}}^{2}}{2} \sum_{x, \mu} E_{x, \mu}^{2}-\frac{g_{\mathrm{m}}^{2}}{2} \sum_{x}\left(\square_{\mu_{x} \mu_{y}}+\square_{\mu_{y} \mu_{z}}+\square_{\mu_{x} \mu_{z}}+\text { h.c. }\right)
\end{aligned}
$$

Relation btw the model parameters and $a, m_{0}, g$ :

$$
t=a^{-1}, m=m_{0}, g_{\mathrm{e}}^{2}=g^{2} a^{-1}, g_{\mathrm{m}}^{2}=8 /\left(g^{2} a\right)
$$

Physical regime of QED: $g_{\mathrm{e}} g_{\mathrm{m}}=2 \sqrt{2} t$

The physical state $|\Phi\rangle$ satisfies $G_{x}|\Phi\rangle=0 \mathrm{w} /$

$$
G_{x}=\psi_{x}^{\dagger} \psi_{x}-\frac{1-(-1)^{x}}{2}-\sum_{\mu} E_{x, \mu}
$$

## The model (2/2)

## Charge operator

Matter occupation operator

$$
Q=\sum_{x}\left(\psi_{x}^{\dagger} \psi_{x}-\frac{1-(-1)^{x}}{2}\right)
$$

$$
n_{x}=\frac{1-(-1)^{x}}{2}-(-1)^{x} \psi_{x}^{\dagger} \psi_{x}
$$

Charge density along w/ $\mu_{x}$ direction

$$
q_{c}(d)=\frac{2}{L^{2}} \sum_{j, k=1}^{L}\langle\mathrm{GS}|(-1)^{x} \psi_{x}^{\dagger} \psi_{x}|\mathrm{GS}\rangle \quad \mathrm{w} / \quad x=(d, j, k)
$$

Average matter density

$$
\rho=\frac{1}{L^{3}} \sum_{x}\langle\mathrm{GS}| n_{x}|\mathrm{GS}\rangle \quad \mathrm{w} / \quad n_{x}=\frac{1-(-1)^{x}}{2}-(-1)^{x} \psi_{x}^{\dagger} \psi_{x}
$$

