

Spectroscopy by Tensor Renormalization group method

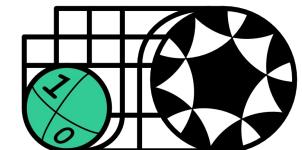
~ how to obtain energy spectrum and determine quantum number
using tensor networks with Lagrangian formalism

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Contents

- Spectroscopy for lattice QCD with Monte Carlo method
- Transfer matrix (TM) formalism
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 - How to determine quantum number
- TM + Tensor network
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- Summary and Future

Spectroscopy for lattice QCD with Monte Carlo method

Energy spectrum

- Schrödinger eq. (e.g., 0^{-+} for pion)
quantum number : J^{PC} , flavors, . . .

$$\hat{H}|n, q\rangle = E_{n,q}|n, q\rangle \quad (n = 0, 1, 2, \dots)$$

QCD Hamiltonian

$$\hat{H}|\Omega\rangle = 0 \quad : \text{vacuum}$$

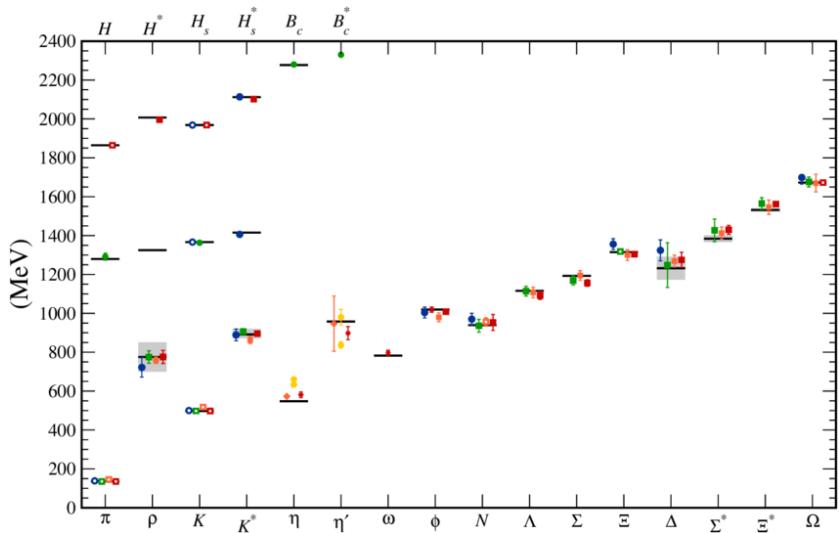
- Two-point function (Euclidean time)

$$\lim_{\beta \rightarrow \infty} \text{Tr} \left[\hat{\mathcal{O}}_q^\dagger(\tau) \hat{\mathcal{O}}_q(0) e^{-\beta \hat{H}} \right] = \sum_{n=0}^{\infty} |\langle \Omega | \hat{\mathcal{O}}_q(0) | n, q \rangle|^2 e^{-\tau E_{n,q}}$$


 $\hat{1} = \sum_{n,q'} |n, q' \rangle \langle n, q'|$

Hadron spectroscopy

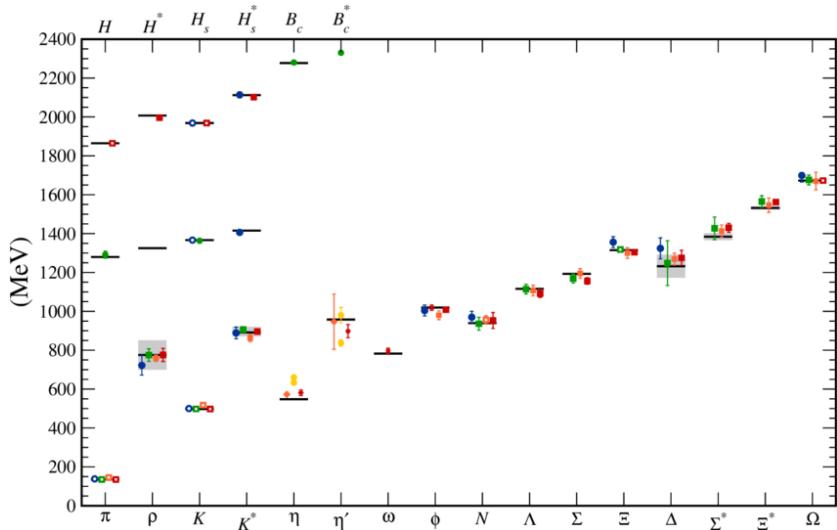
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2013 snowmass report

Hadron spectroscopy

$$\lim_{\beta \rightarrow \infty} \text{Tr} \left[\hat{\mathcal{O}}_q^\dagger(\tau) \hat{\mathcal{O}}_q(0) e^{-\beta \hat{H}} \right] = \sum_{n=0}^{\infty} |\langle \Omega | \hat{\mathcal{O}}_q(0) | n, q \rangle|^2 e^{-\tau E_{n,q}}$$



Problems:

- Need large time extent β and time separation τ
- Need large statistics to extract higher excited states

2013 snowmass report

How to get spectrum by TN?

- Hamiltonian formalism
 - ⇒ Matsumoto's talk arXiv:2307.16655
- Lagrangian formalism
 - Two-point function
 - Large time extent and separation are easily realized
 - nothing new! (just do it)
 - Transfer matrix We use here!
 - No need to extrapolate time extent and separation

Transfer matrix formalism

TM and its spectrum

$$Z = \text{tr} [\mathcal{T} \mathcal{T} \cdots \mathcal{T}]$$

for 2D Ising

$$s_x = \pm 1$$

$$\mathcal{T}_{s' s} = \left(\prod_x e^{\beta s'_x s_x} \right) \left(\prod_x e^{\frac{1}{2}\beta s'_{x+1} s'_x + \frac{1}{2}\beta s_{x+1} s_x} \right)$$

: Hermitian, (semi)positive definite

TM and its spectrum

$$Z = \text{tr} [\mathcal{T} \mathcal{T} \cdots \mathcal{T}]$$

for 2D Ising $s_x = \pm 1$

$$\mathcal{T}_{s' s} = \left(\prod_x e^{\beta s'_x s_x} \right) \left(\prod_x e^{\frac{1}{2}\beta s'_{x+1} s'_x + \frac{1}{2}\beta s_{x+1} s_x} \right)$$

: Hermitian, (semi)positive definite

$$\Rightarrow \mathcal{T} = U e^{-\omega} U^\dagger \quad (\mathcal{T}|a\rangle = e^{-\omega_a} |a\rangle \text{ for } a = 0, 1, 2 \dots)$$
$$\omega_a \geq 0 \quad \omega_0 = 0$$

$$\mathcal{T} \leftrightarrow e^{-\hat{H}} \Rightarrow \omega_a \leftrightarrow E_{n,q}$$

How to identify quantum number of $|a\rangle$?

How to identify quantum number

Answer: compute matrix elements $\langle b | \hat{O}_q | a \rangle$

The diagram consists of two orange arrows. One arrow points from the text "eigenstates of TM" to the state vector $|b\rangle$ in the matrix element. Another arrow points from the text "q is assumed to be well known" to the state vector $|a\rangle$ in the matrix element.

eigenstates of TM
q is assumed to be well known

How to identify quantum number

Answer: compute matrix elements $\langle b | \hat{O}_q | a \rangle$

eigenstates of TM
 q is assumed to be well known

Why:

For conserved charge of **continuous** symmetry \hat{Q} $([\hat{Q}, \hat{H}] = 0)$

$$[\hat{Q}, \hat{X}] = q_X \hat{X}$$

\Downarrow charge of \hat{X} $(\because \hat{Q}\hat{X}|\Omega\rangle = q_X \hat{X}|\Omega\rangle)$ $\hat{Q}|\Omega\rangle = 0$

How to identify quantum number

Answer: compute matrix elements $\langle b | \hat{O}_q | a \rangle$

eigenstates of TM
 q is assumed to be well known

Why:

For conserved charge of **continuous** symmetry \hat{Q} $([\hat{Q}, \hat{H}] = 0)$

$$[\hat{Q}, \hat{X}] = q_X \hat{X}$$

$\langle b | \dots | a \rangle \Downarrow$ charge of \hat{X} $(\because \hat{Q} \hat{X} |\Omega\rangle = q_X \hat{X} |\Omega\rangle)$ $\hat{Q} |\Omega\rangle = 0$

$$(q_b - q_a - q_X) \langle b | \hat{X} | a \rangle = 0$$

$$\hat{Q} |a\rangle = q_a |a\rangle$$

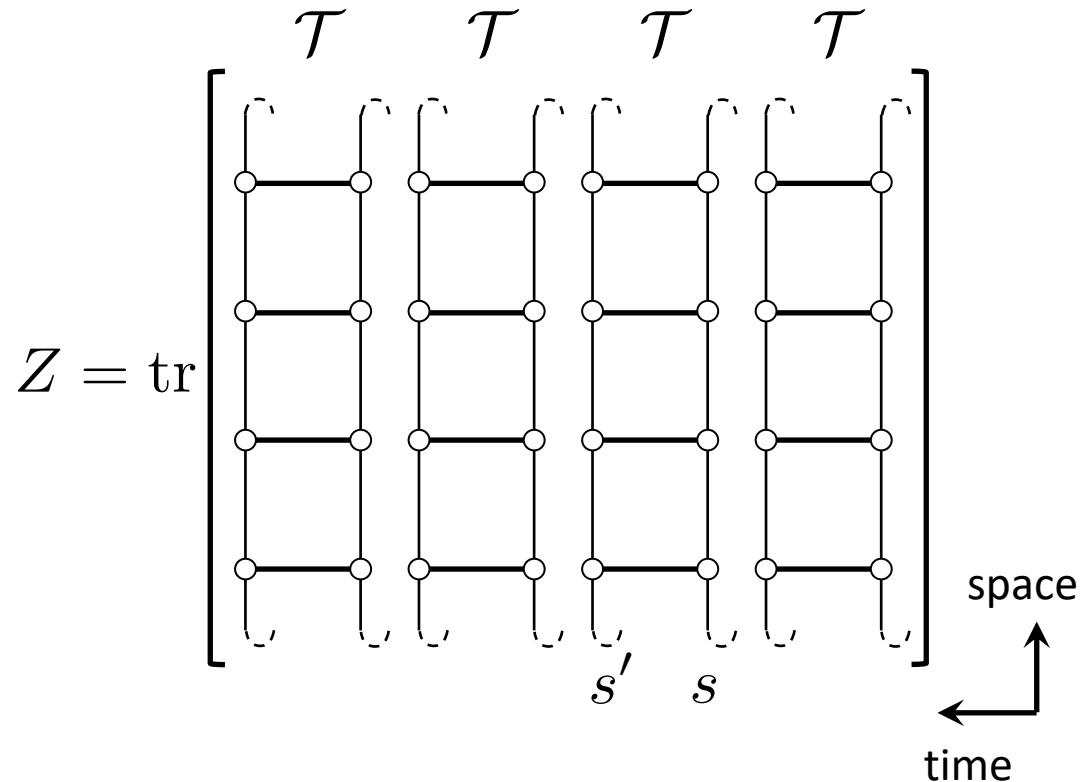
\Rightarrow selection rule: $\langle b | \hat{X} | a \rangle \neq 0 \implies q_b - q_a - q_X = 0$

\Rightarrow for $b = 0 = \Omega$ $\langle \Omega | \hat{X} | a \rangle \neq 0 \implies q_a = q_X$

For discrete symmetry $\hat{D} \hat{X} \hat{D}^{-1} = q_X \hat{X}$ $\langle \Omega | \hat{X} | a \rangle \neq 0 \implies q_a q_X = 1$

$\downarrow \pm$

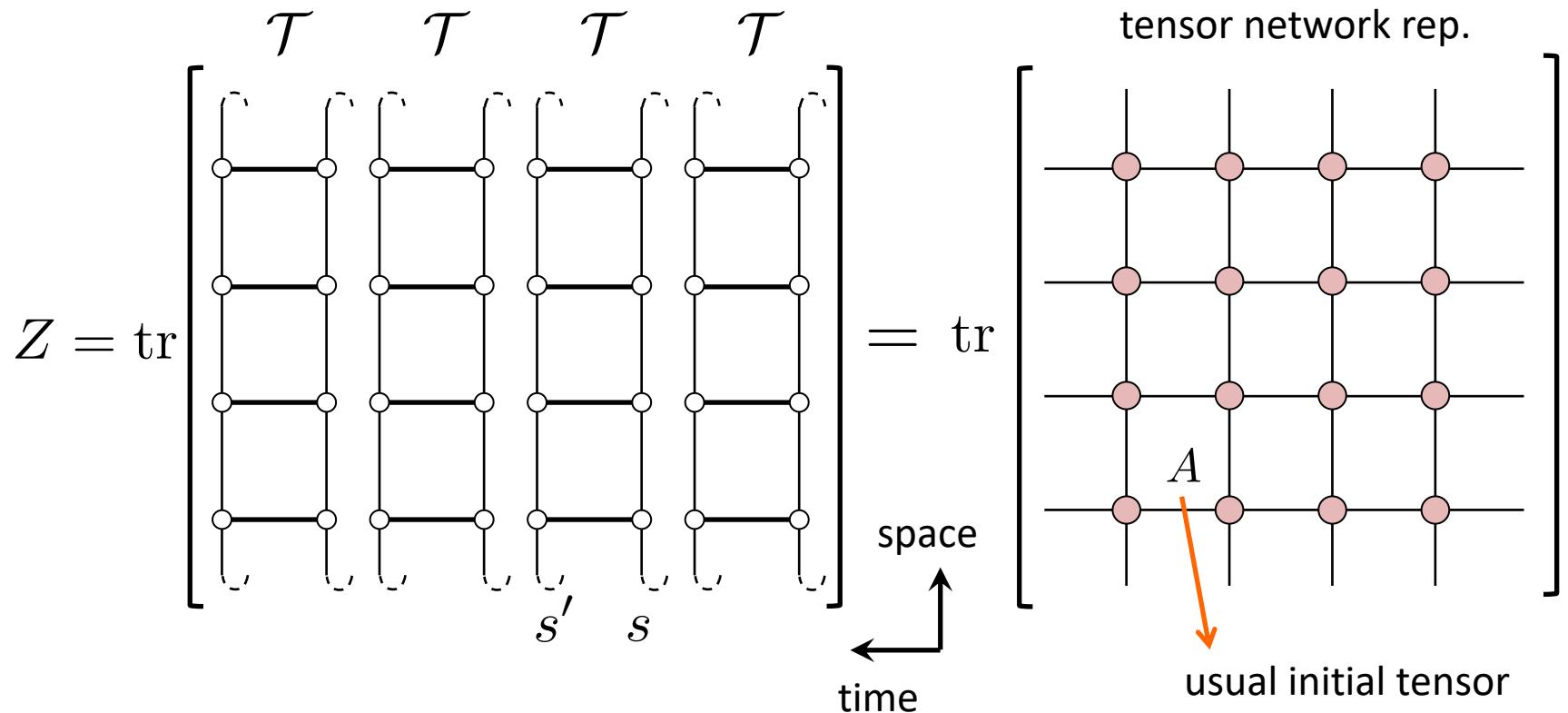
TM + Tensor network



$$\mathcal{T}_{s's} = \left(\prod_x e^{\beta s'_x s_x} \right) \left(\prod_x e^{\frac{1}{2}\beta s'_{x+1}s'_x + \frac{1}{2}\beta s_{x+1}s_x} \right)$$

for 2D Ising

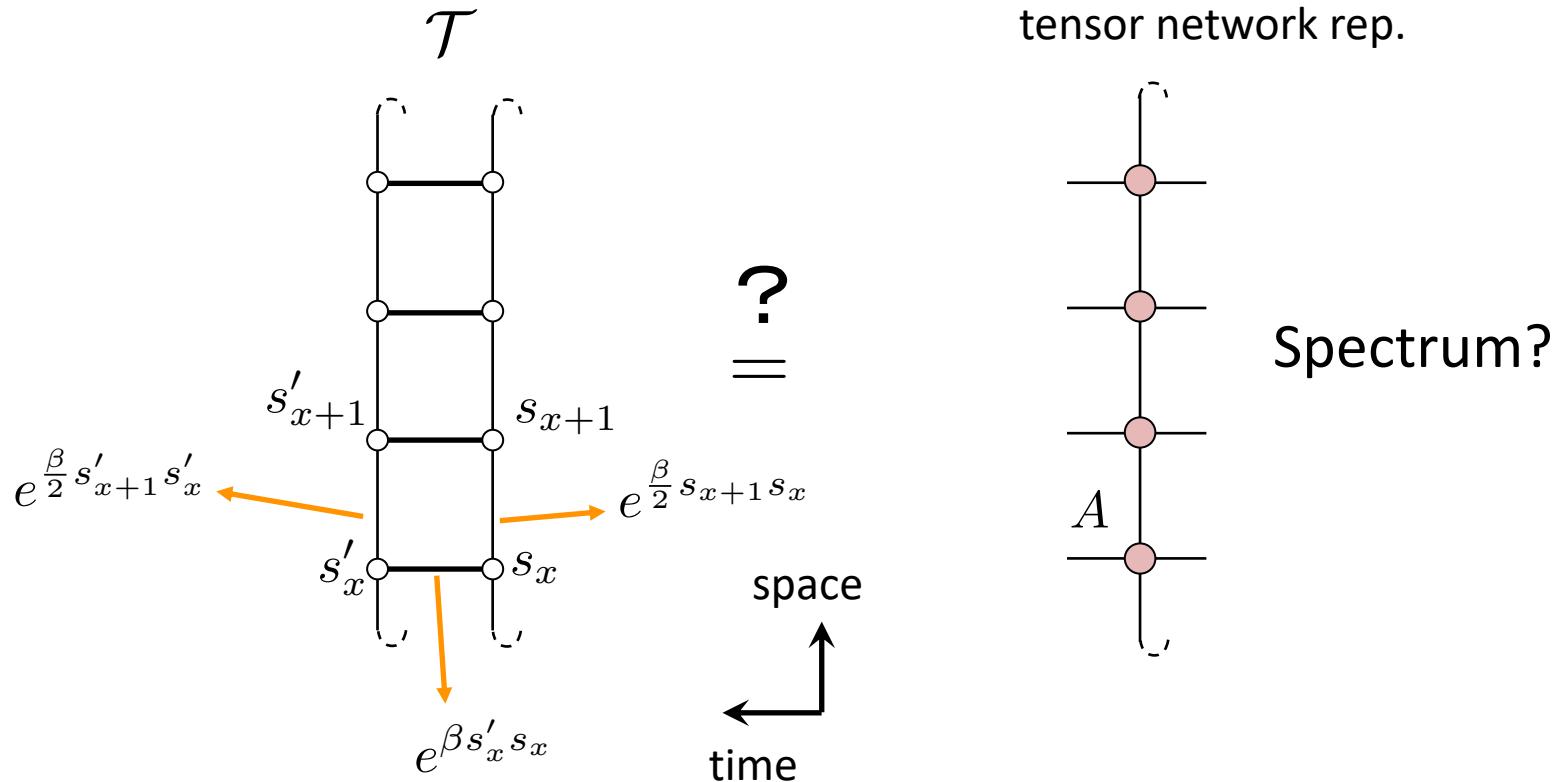
TM + Tensor network



$$\mathcal{T}_{s's} = \left(\prod_x e^{\beta s'_x s_x} \right) \left(\prod_x e^{\frac{1}{2}\beta s'_{x+1}s'_x + \frac{1}{2}\beta s_{x+1}s_x} \right)$$

for 2D Ising

TM + Tensor network

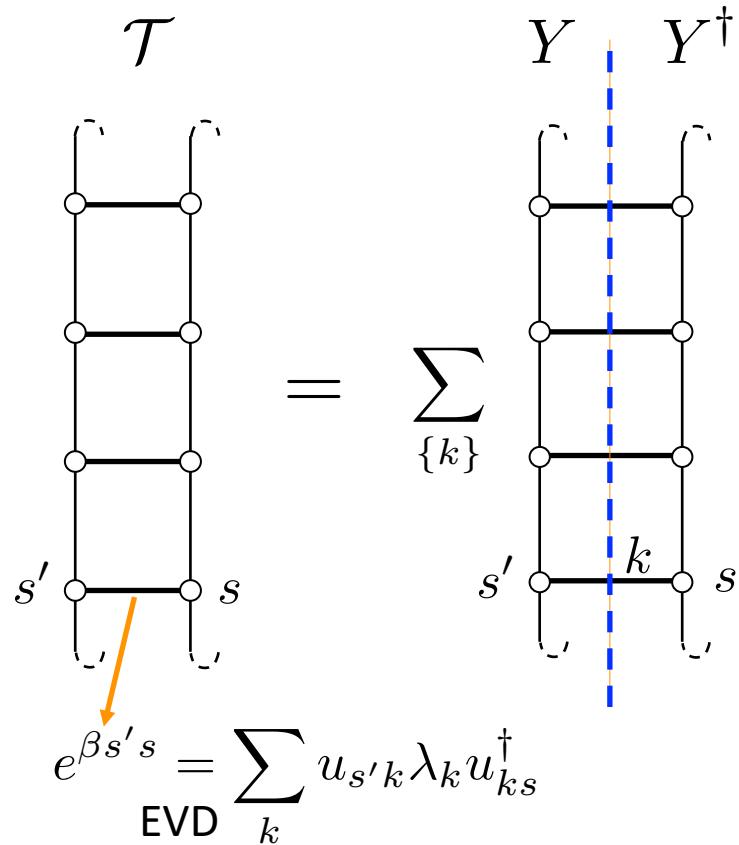


$$\mathcal{T}_{s's} = \left(\prod_x e^{\beta s'_x s_x} \right) \left(\prod_x e^{\frac{1}{2}\beta s'_{x+1}s'_x + \frac{1}{2}\beta s_{x+1}s_x} \right)$$

for 2D Ising

TM + Tensor network

$$Z = \text{tr}[\mathcal{T}^n] = \text{tr}[(YY^\dagger)^n]$$



TM + Tensor network

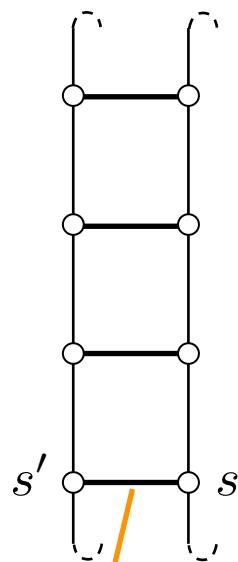
$$Z = \text{tr}[\mathcal{T}^n]$$

$$= \text{tr}[(YY^\dagger)^n]$$

$$= \text{tr}[(Y^\dagger Y)^n]$$

$$= \text{tr}[\mathcal{A}^n]$$

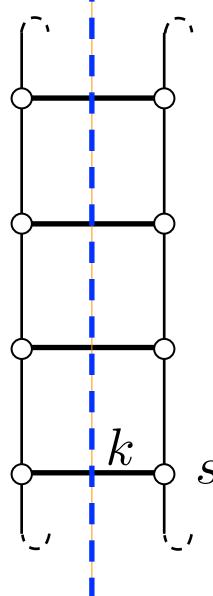
\mathcal{T}



$$e^{\beta s' s} = \sum_k u_{s'k} \lambda_k u_{ks}^\dagger$$

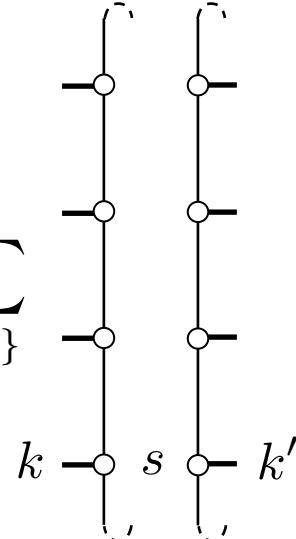
EVD

$Y \quad Y^\dagger$

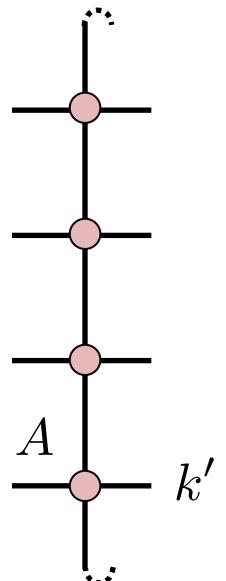


\Rightarrow

$Y^\dagger \quad Y$



\mathcal{A} TN rep.



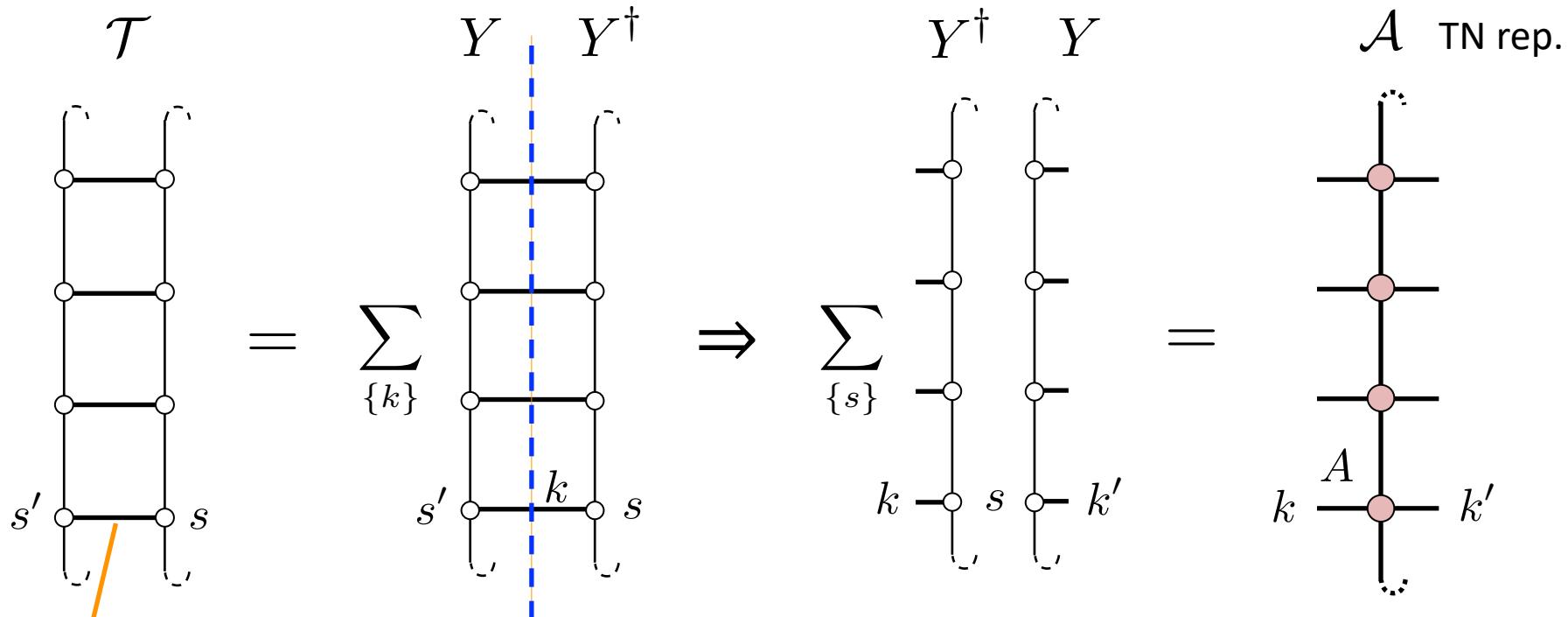
TM + Tensor network

$$Z = \text{tr}[\mathcal{T}^n]$$

$$= \text{tr}[(YY^\dagger)^n]$$

$$= \text{tr}[(Y^\dagger Y)^n]$$

$$= \text{tr}[\mathcal{A}^n]$$



$$e^{\beta s' s} = \sum_k u_{s' k} \lambda_k u_{ks}^\dagger$$

EVD

$$Y \xrightarrow{\text{SVD}} U \sigma W^\dagger$$

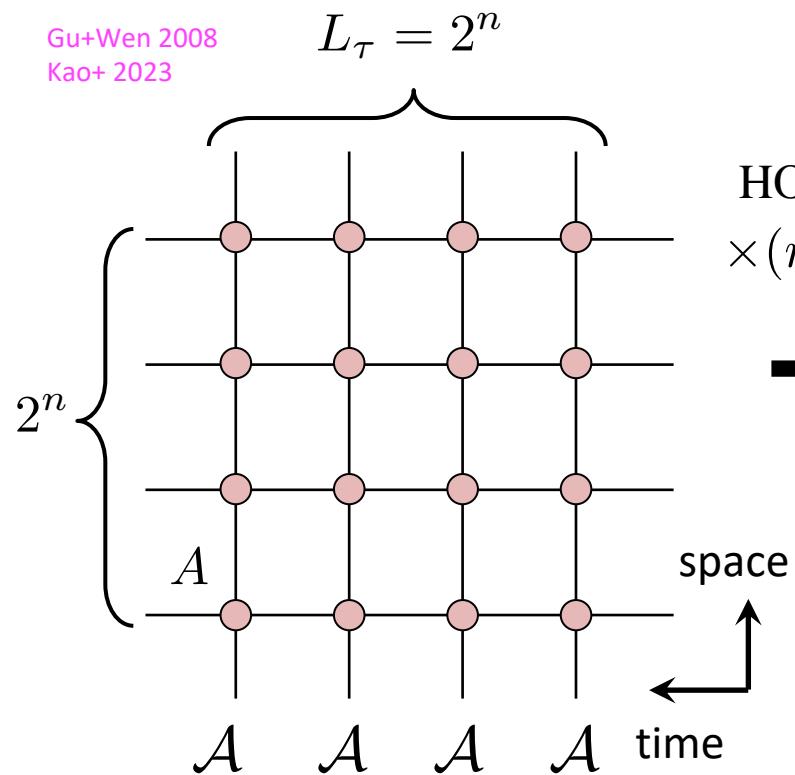
$$\mathcal{T} = YY^\dagger = U \underline{\sigma^2} U^\dagger$$

$$Y^\dagger Y = \mathcal{A} = W \underline{\sigma^2} W^\dagger$$

Identical spectrum!

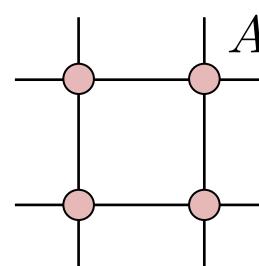
How to obtain spectrum using HOTRG

Gu+Wen 2008
Kao+ 2023

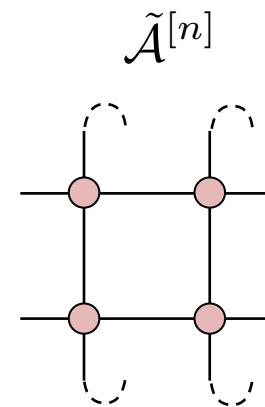


HOTRG
 $\times(n - 1)$

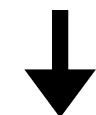
→



PBC



diagonalization

$$\tilde{A}^{[n]} = W^{[n]} \lambda^{[n]} W^{[n]\dagger}$$


$$\omega_a^{[n]} \equiv \frac{1}{L_\tau} \log \left(\frac{\lambda_0^{[n]}}{\lambda_a^{[n]}} \right) \approx \omega_a$$

$(\mathcal{T} = U e^{-\omega} U^\dagger)$

$\tilde{A}^{[n]} \approx \mathcal{A}^{L_\tau}$

How to compute matrix elements

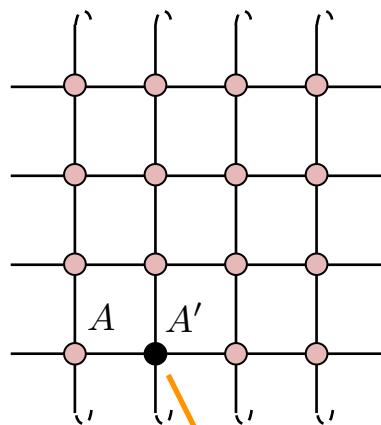
$$\langle b | \hat{\mathcal{O}}_q | a \rangle = (U^\dagger \mathcal{O}_q U)_{ba} \quad (\mathcal{T} = U e^{-\omega} U^\dagger)$$

How to compute matrix elements

$$\langle b | \hat{\mathcal{O}}_q | a \rangle = (U^\dagger \mathcal{O}_q U)_{ba} \quad (\mathcal{T} = U e^{-\omega} U^\dagger) \quad m = L_\tau / 2$$
$$= (U^\dagger \underline{\mathcal{T}^{-m}} \underline{\mathcal{T}^m \mathcal{O}_q \mathcal{T}^m} \underline{\mathcal{T}^{m+1}} \underline{\mathcal{T}^{-(m+1)}} U)_{ba} \quad (\because \mathcal{T} \mathcal{T}^{-1} = 1)$$

$$\mathcal{T}^{-1} = U e^{+\omega} U^\dagger$$

roughly speaking



one-point function

impurity tensor

How to compute matrix elements

$$\begin{aligned} \langle b | \hat{\mathcal{O}}_q | a \rangle &= (U^\dagger \mathcal{O}_q U)_{ba} & (\mathcal{T} = U e^{-\omega} U^\dagger) \\ &= (U^\dagger \underbrace{\mathcal{T}^{-m}}_{\mathcal{T} = YY^\dagger} \underbrace{\mathcal{T}^m}_{\mathcal{T}^{-1} = U e^{+\omega} U^\dagger} \mathcal{O}_q \underbrace{\mathcal{T}^{m+1}}_{\mathcal{T}^{-1}} \underbrace{\mathcal{T}^{-(m+1)}}_{\mathcal{T}^{-1}} U)_{ba} & m = L_\tau / 2 \\ && (\because \mathcal{T}\mathcal{T}^{-1} = 1) \end{aligned}$$

$$\begin{aligned} &= e^{\omega(m-1/2)} W^\dagger \tilde{\mathcal{A}}' W e^{\omega(m+1/2)} \\ &= \mathcal{A} = W e^{-\omega} W^\dagger \end{aligned}$$

one-point function

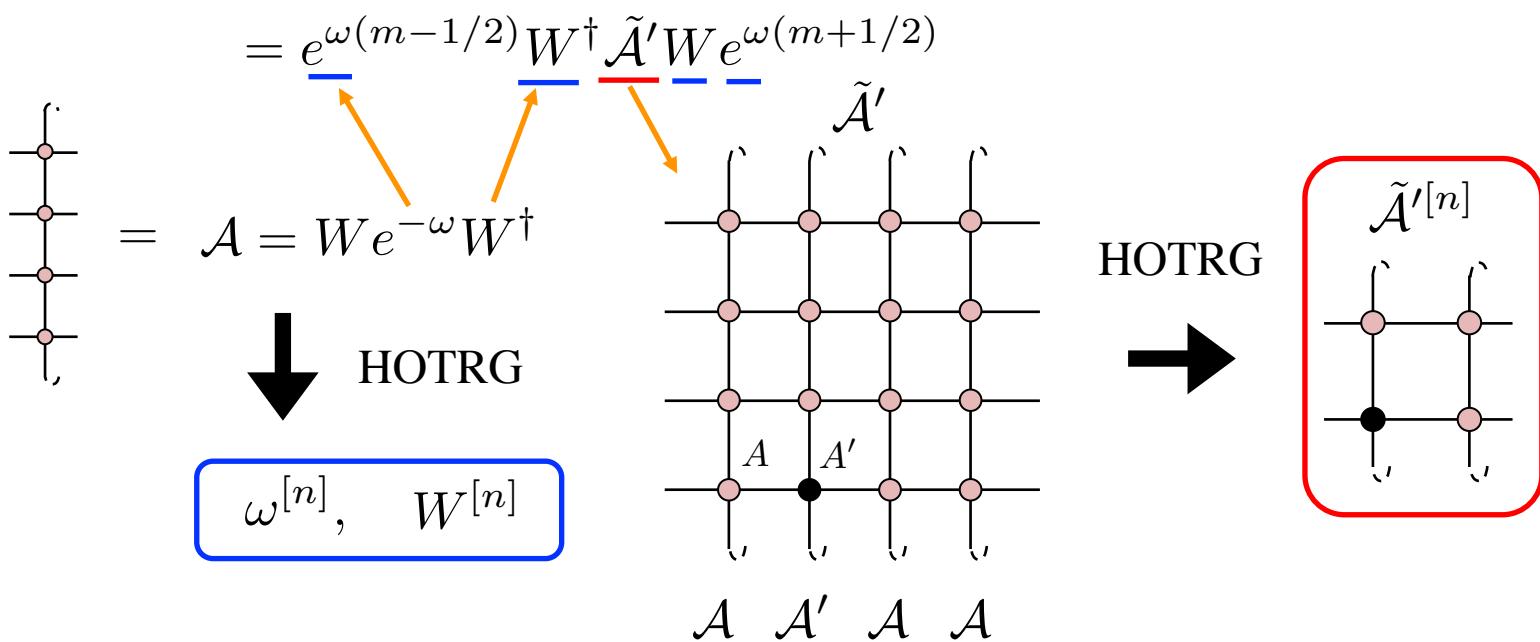
How to compute matrix elements

$$\langle b | \hat{\mathcal{O}}_q | a \rangle = (U^\dagger \mathcal{O}_q U)_{ba} \quad (\mathcal{T} = U e^{-\omega} U^\dagger)$$

$$= (U^\dagger \underbrace{\mathcal{T}^{-m}}_{\mathcal{T} = YY^\dagger} \underbrace{\mathcal{T}^m}_{\mathcal{T}^{-1} = U e^{+\omega} U^\dagger} \mathcal{O}_q \underbrace{\mathcal{T}^{m+1}}_{\mathcal{T}^{-1}} \underbrace{\mathcal{T}^{-(m+1)}}_{\mathcal{T}^{-1}} U)_{ba}$$

$\mathcal{T} = YY^\dagger \quad \mathcal{T}^{-1} = U e^{+\omega} U^\dagger$

$$m = L_\tau / 2 \quad (\because \mathcal{T} \mathcal{T}^{-1} = 1)$$



How to compute matrix elements

$$\begin{aligned}\langle b | \hat{\mathcal{O}}_q | a \rangle &= (U^\dagger \mathcal{O}_q U)_{ba} & (\mathcal{T} = U e^{-\omega} U^\dagger) \\ &= (U^\dagger \underbrace{\mathcal{T}^{-m}}_{\mathcal{T} = YY^\dagger} \underbrace{\mathcal{T}^m}_{\mathcal{A}' = \tilde{\mathcal{A}}' W^{[n]}} \mathcal{O}_q \underbrace{\mathcal{T}^{m+1}}_{\mathcal{T}^{-1} = U e^{+\omega} U^\dagger} \underbrace{\mathcal{T}^{-(m+1)}}_{\mathcal{T}^{-1}} U)_{ba} & m = L_\tau / 2 \\ &= e^{\omega(m-1/2)} \underbrace{W^\dagger}_{\mathcal{A}' = \tilde{\mathcal{A}}' W^{[n]}} \underbrace{\tilde{\mathcal{A}}'}_{\mathcal{A}' = \tilde{\mathcal{A}}' W^{[n]}} \underbrace{W e^{\omega(m+1/2)}}_{\mathcal{A}' = \tilde{\mathcal{A}}' W^{[n]}} \\ &\approx e^{\omega^{[n]}(m-1/2)} \underbrace{W^{[n]\dagger}}_{\mathcal{A}' = \tilde{\mathcal{A}}' W^{[n]}} \underbrace{\tilde{\mathcal{A}}'^{[n]}}_{\mathcal{A}' = \tilde{\mathcal{A}}' W^{[n]}} \underbrace{W^{[n]}}_{\mathcal{A}' = \tilde{\mathcal{A}}' W^{[n]}} e^{\omega^{[n]}(m+1/2)}\end{aligned}$$

all building blocks are computed from
tensor network representations

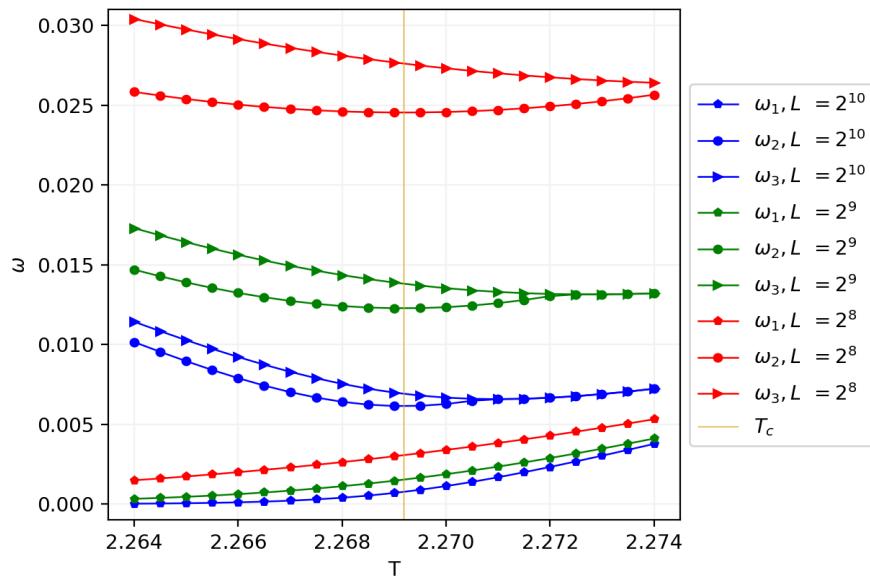
Numerical results for 2D Ising model

Energy spectrum

after n HOTRG steps

$$\begin{array}{c}
 \text{Diagram of a } 2 \times 2 \text{ grid of nodes with dashed arcs connecting them.} \\
 = W^{[n]} \lambda^{[n]} W^{[n]\dagger} \quad \rightarrow \quad \omega_a^{[n]} \equiv \frac{1}{L_\tau} \log \left(\frac{\lambda_0^{[n]}}{\lambda_a^{[n]}} \right) \approx \omega_a
 \end{array}$$

$\tilde{\mathcal{A}}^{[n]}$



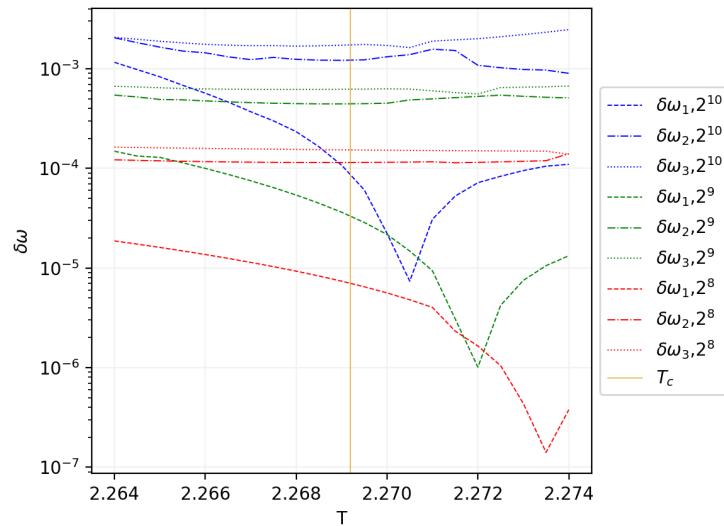
$\chi = 100$

$L = L_\tau$

Energy spectrum

Kaufmann 1949

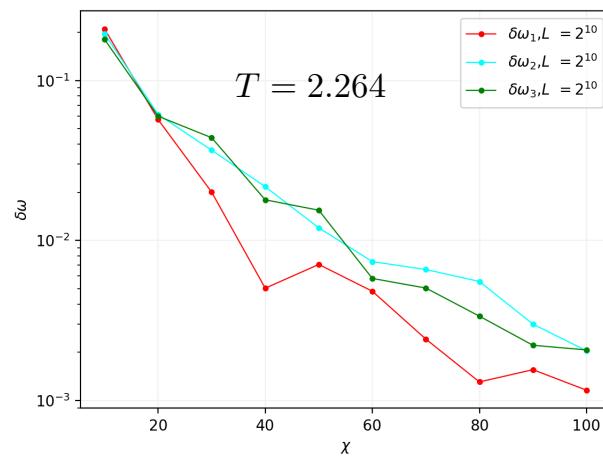
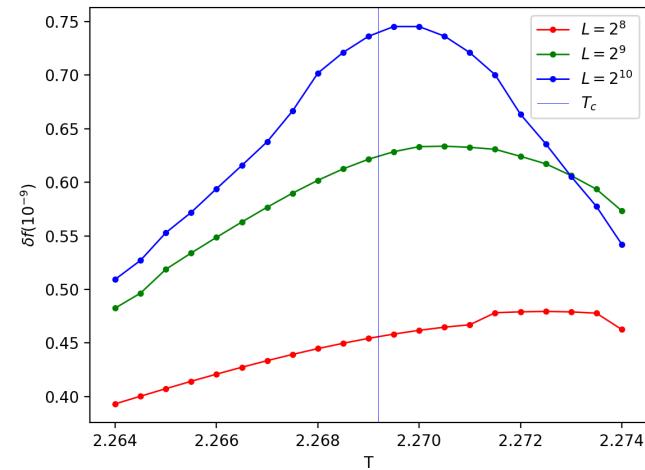
$$\delta\omega = \left| \frac{\omega - \omega_{\text{exact},L}}{\omega_{\text{exact},L}} \right|$$



$\chi = 100$

Kaufmann 1949

$$\delta f = \left| \frac{f - f_{\text{exact},L}}{f_{\text{exact},L}} \right|$$



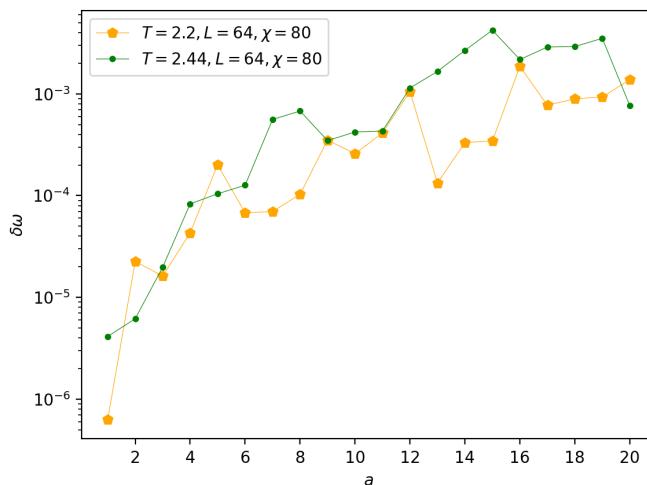
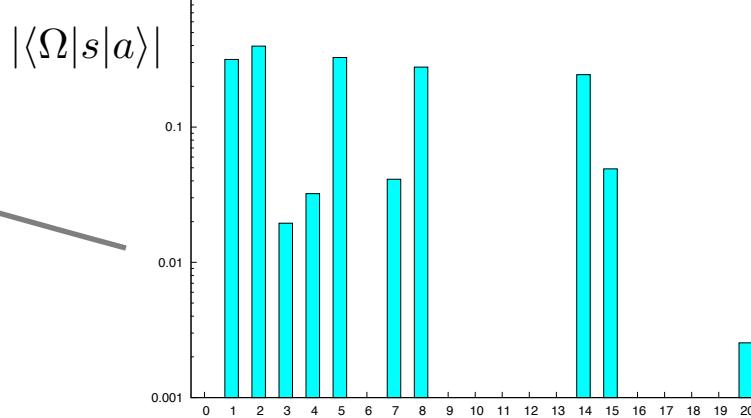
Determination of quantum number

$$Z_2 : + \quad - \quad q_a$$

$$\langle \Omega | s_{x=0} | a \rangle \neq 0 \implies q_a = -$$

| a | $\omega^{[\text{exact}]}$ | q_a | $\omega^{[\text{hotrg}]}$ | q_a | $\delta\omega$ |
|-----|---------------------------|-------|---------------------------|-------|----------------|
| 1 | 0.1262302 | - | 0.1262307 | - | 0.000004 |
| 2 | 0.1597880 | - | 0.1597889 | - | 0.000006 |
| 3 | 0.1597880 | - | 0.1597911 | - | 0.000020 |
| 4 | 0.2326853 | - | 0.2327046 | - | 0.000083 |
| 5 | 0.2326853 | - | 0.2327095 | - | 0.000104 |
| 6 | 0.2708016 | + | 0.2708359 | + | 0.000127 |
| 7 | 0.3181546 | - | 0.3183329 | - | 0.000560 |
| 8 | 0.3181546 | - | 0.3183705 | - | 0.000679 |
| 9 | 0.3290037 | + | 0.3291180 | + | 0.000347 |
| 10 | 0.3290037 | + | 0.3291425 | + | 0.000422 |
| 11 | 0.3290037 | + | 0.3291456 | + | 0.000431 |
| 12 | 0.3290037 | + | 0.3293794 | + | 0.001142 |
| 13 | 0.3872058 | + | 0.3878486 | + | 0.001660 |
| 14 | 0.4073042 | - | 0.4083937 | - | 0.002675 |
| 15 | 0.4073042 | - | 0.4090231 | - | 0.004220 |
| 16 | 0.4100181 | + | 0.4109090 | + | 0.002173 |
| 17 | 0.4100181 | + | 0.4112006 | + | 0.002884 |
| 18 | 0.4100181 | + | 0.4112120 | + | 0.002912 |
| 19 | 0.4100181 | + | 0.4114574 | + | 0.003510 |
| 20 | 0.4457831 | - | 0.4461242 | - | 0.000765 |

$$T = 2.44 \quad L = 64 \quad \chi = 80$$



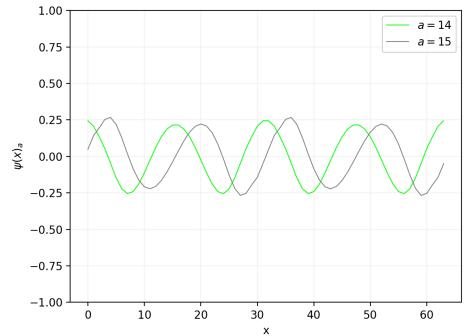
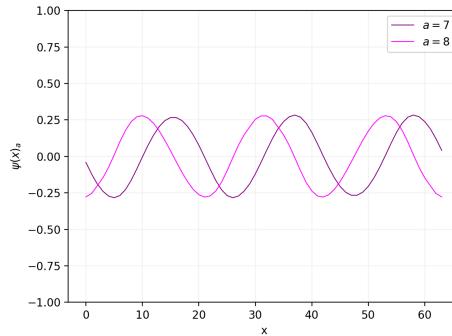
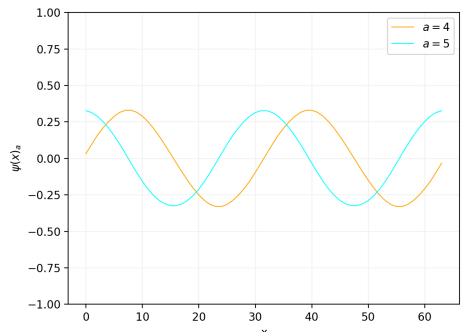
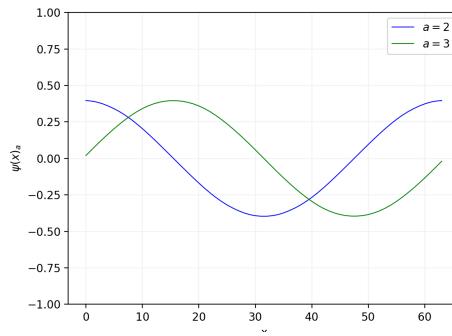
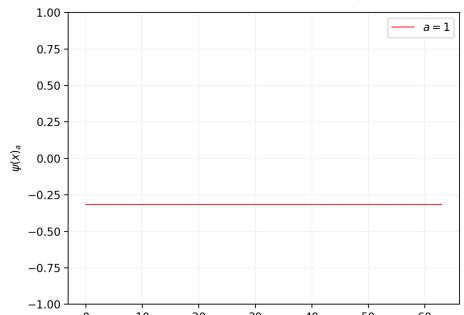
Bethe-Salpeter wave function

$$\psi_a(x) = \langle \Omega | s_x | a \rangle$$

$$(0 \leq x < L)$$

| a | ω^{exact} | q_a | ω^{hotrg} | q_a | $\delta\omega$ |
|-----|-------------------------|-------|-------------------------|-------|----------------|
| 1 | 0.1262302 | — | 0.1262307 | — | 0.000004 |
| 2 | 0.1597880 | — | 0.1597889 | — | 0.000006 |
| 3 | 0.1597880 | — | 0.1597911 | — | 0.000020 |
| 4 | 0.2326853 | — | 0.2327046 | — | 0.000083 |
| 5 | 0.2326853 | — | 0.2327095 | — | 0.000104 |
| 6 | 0.2708016 | + | 0.2708359 | + | 0.000127 |
| 7 | 0.3181546 | — | 0.3183329 | — | 0.000560 |
| 8 | 0.3181546 | — | 0.3183705 | — | 0.000679 |
| 9 | 0.3290037 | + | 0.3291180 | + | 0.000347 |
| 10 | 0.3290037 | + | 0.3291425 | + | 0.000422 |
| 11 | 0.3290037 | + | 0.3291456 | + | 0.000431 |
| 12 | 0.3290037 | + | 0.3293794 | + | 0.001142 |
| 13 | 0.3872058 | + | 0.3878486 | + | 0.001660 |
| 14 | 0.4073042 | — | 0.4083937 | — | 0.002675 |
| 15 | 0.4073042 | — | 0.4090231 | — | 0.004220 |
| 16 | 0.4100181 | + | 0.4109090 | + | 0.002173 |
| 17 | 0.4100181 | + | 0.4112006 | + | 0.002884 |
| 18 | 0.4100181 | + | 0.4112120 | + | 0.002912 |
| 19 | 0.4100181 | + | 0.4114574 | + | 0.003510 |
| 20 | 0.4457831 | — | 0.4461242 | — | 0.000765 |

$$T = 2.44 \quad L = 64 \quad \chi = 80$$



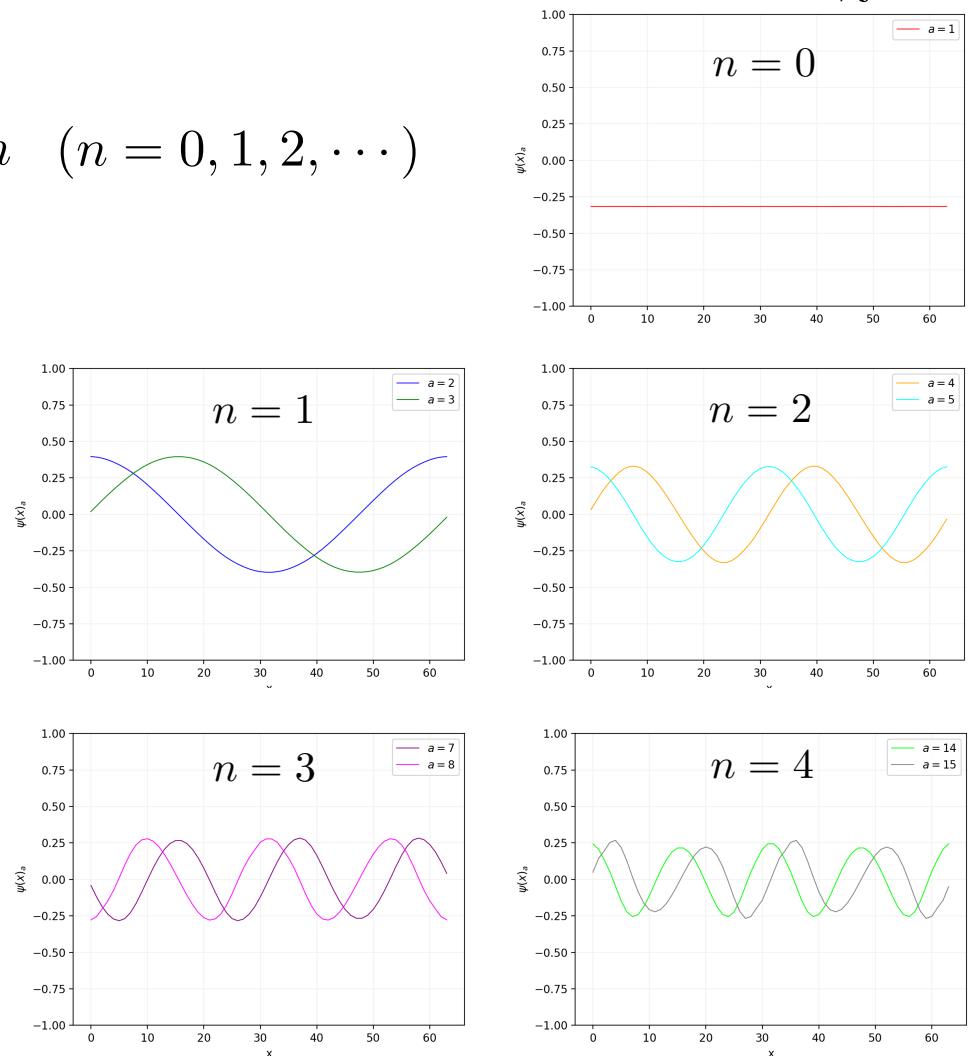
Bethe-Salpeter wave function

$$\psi_a(x) = \langle \Omega | s_x | a \rangle \propto e^{ipx}$$

$$(0 \leq x < L) \quad p = \frac{2\pi}{L}n \quad (n = 0, 1, 2, \dots)$$

| a | ω^{exact} | q_a | ω^{hotrg} | q_a | $\delta\omega$ |
|-----|-------------------------|-------|-------------------------|-------|----------------|
| 1 | 0.1262302 | — | 0.1262307 | — | 0.000004 |
| 2 | 0.1597880 | — | 0.1597889 | — | 0.000006 |
| 3 | 0.1597880 | — | 0.1597911 | — | 0.000020 |
| 4 | 0.2326853 | — | 0.2327046 | — | 0.000083 |
| 5 | 0.2326853 | — | 0.2327095 | — | 0.000104 |
| 6 | 0.2708016 | + | 0.2708359 | + | 0.000127 |
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| 12 | 0.3290037 | + | 0.3293794 | + | 0.001142 |
| 13 | 0.3872058 | + | 0.3878486 | + | 0.001660 |
| 14 | 0.4073042 | — | 0.4083937 | — | 0.002675 |
| 15 | 0.4073042 | — | 0.4090231 | — | 0.004220 |
| 16 | 0.4100181 | + | 0.4109090 | + | 0.002173 |
| 17 | 0.4100181 | + | 0.4112006 | + | 0.002884 |
| 18 | 0.4100181 | + | 0.4112120 | + | 0.002912 |
| 19 | 0.4100181 | + | 0.4114574 | + | 0.003510 |
| 20 | 0.4457831 | — | 0.4461242 | — | 0.000765 |

$$T = 2.44 \quad L = 64 \quad \chi = 80$$

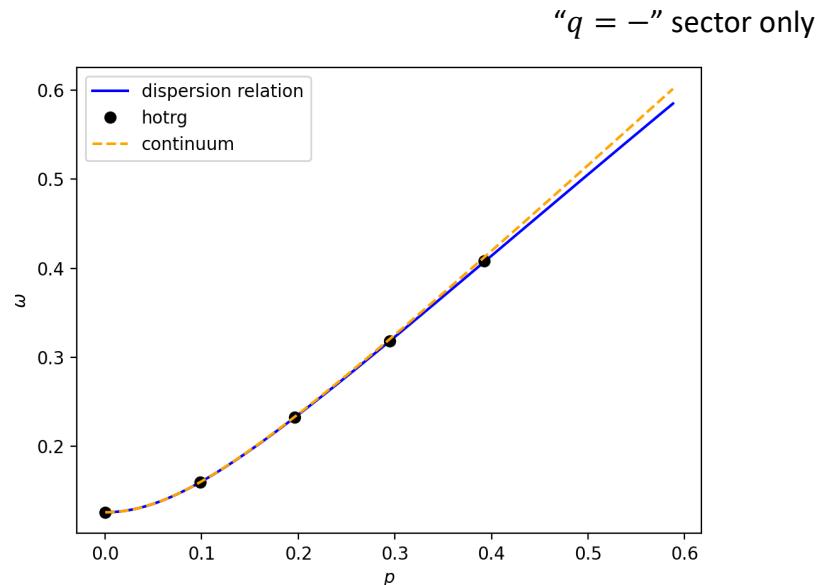


Momentum identification

$$T = 2.44 \quad L = 64 \quad \chi = 80$$

$$\sum_x e^{-ipx} \langle \Omega | s_x | a \rangle \neq 0 \implies p : \text{momentum of } |a\rangle$$

| a | ω^{exact} | q_a | ω^{hotrg} | q_a | $ p $ |
|-----|-------------------------|-------|-------------------------|-------|----------|
| 1 | 0.1262302 | - | 0.1262307 | - | 0 |
| 2 | 0.1597880 | - | 0.1597889 | - | $2\pi/L$ |
| 3 | 0.1597880 | - | 0.1597911 | - | $2\pi/L$ |
| 4 | 0.2326853 | - | 0.2327046 | - | $4\pi/L$ |
| 5 | 0.2326853 | - | 0.2327095 | - | $4\pi/L$ |
| 6 | 0.2708016 | + | 0.2708359 | + | * |
| 7 | 0.3181546 | - | 0.3183329 | - | $6\pi/L$ |
| 8 | 0.3181546 | - | 0.3183705 | - | $6\pi/L$ |
| 9 | 0.3290037 | + | 0.3291180 | + | * |
| 10 | 0.3290037 | + | 0.3291425 | + | * |
| 11 | 0.3290037 | + | 0.3291456 | + | * |
| 12 | 0.3290037 | + | 0.3293794 | + | * |
| 13 | 0.3872058 | + | 0.3878486 | + | * |
| 14 | 0.4073042 | - | 0.4083937 | - | $8\pi/L$ |
| 15 | 0.4073042 | - | 0.4090231 | - | $8\pi/L$ |
| 16 | 0.4100181 | + | 0.4109090 | + | * |
| 17 | 0.4100181 | + | 0.4112006 | + | * |
| 18 | 0.4100181 | + | 0.4112120 | + | * |
| 19 | 0.4100181 | + | 0.4114574 | + | * |
| 20 | 0.4457831 | - | 0.4461242 | - | 0 |



lattice: $\omega(p) = \cosh^{-1}(1 - \cos p + \cosh m)$

continuum: $\omega(p) = \sqrt{m^2 + p^2}$

Summary

- We develop a spectroscopy scheme of Lagrangian tensor network approach
- Quantum number of eigen state is judged by looking at associated matrix elements (one-point function)
- BS wave function for low-lying states can be computed
- Relatively higher momentum states and dispersion relation are clearly observed

Future

- Application to other quantum field theories
- Two-particle state
- Scattering phase shift

Backup slides

How to compute matrix elements

$$\begin{aligned}
 \langle b | \hat{\mathcal{O}}_q | a \rangle &= (U^\dagger \mathcal{O}_q U)_{ba} & (\mathcal{T} = U e^{-\omega} U^\dagger) \\
 &= (U^\dagger \underbrace{\mathcal{T}^{-m}}_{\mathcal{T} = YY^\dagger} \underbrace{\mathcal{T}^m}_{\mathcal{T}^{-1} = U e^{+\omega} U^\dagger} \mathcal{O}_q \underbrace{\mathcal{T}^{m+1}}_{\mathcal{T}^{-1}} \underbrace{\mathcal{T}^{-(m+1)}}_{\mathcal{T}^{-1}} U)_{ba} & m = L_\tau / 2 \\
 &= (U^\dagger (U e^\omega U^\dagger)^m (Y Y^\dagger)^m \mathcal{O}_q (Y Y^\dagger)^{m+1} (U e^\omega U^\dagger)^{m+1} U)_{ba} & (\because \mathcal{T} \mathcal{T}^{-1} = 1) \\
 &= (e^{m\omega} \underbrace{U^\dagger Y}_{e^{-\omega/2} W^\dagger} \underbrace{(Y^\dagger Y)^{m-1}}_{\mathcal{A}'} \underbrace{Y^\dagger \mathcal{O}_q Y}_{\mathcal{A}} \underbrace{(Y^\dagger Y)^m}_{\mathcal{A}} \underbrace{Y^\dagger U e^{(m+1)\omega}}_{W e^{-\omega/2}})_{ba} & (\because Y = U e^{-\omega/2} W^\dagger) \\
 &= (e^{(m-1/2)\omega} W^\dagger \underbrace{\mathcal{A}^{m-1} \mathcal{A}' \mathcal{A}^m}_{\tilde{\mathcal{A}'}} W e^{(m+1/2)\omega})_{ba}
 \end{aligned}$$

