

Bond-weighting method for the Grassmann tensor renormalization group

Shinichiro Akiyama ^{a), b)}

^{a)} Endowed Chair for Quantum Software (量子ソフトウェア寄付講座), University of Tokyo

^{b)} Institute for Physics of Intelligence, University of Tokyo

Based on SA, JHEP11(2022)030

KEK Theory Workshop 2022

2022.12.8

Advantages of the TRG approach

- ✓ Tensor renormalization group (TRG) is a typical tensor network method and the application to the 4D systems has recently made remarkable progress

Lagrangian (TRG) approach: [Meurice+, Rev. Mod. Phys. 94\(2022\)025005](#)

[Kadoh, PoS\(LATTICE2021\)633](#)

[SA+, PoS\(LATTICE2021\)530](#)

Hamiltonian (TNS) approach: [Bañuls-Cichy, Rep. Prog. Phys. 83\(2020\)024401](#)

- ✓ TRG is a deterministic numerical method based on the idea of the real-space renormalization group
 - **No sign problem**
 - **The computational cost scales logarithmically w. r. t. the system size**
 - **Direct evaluation of the Grassmann integrals**
 - **Direct evaluation of the path integral**

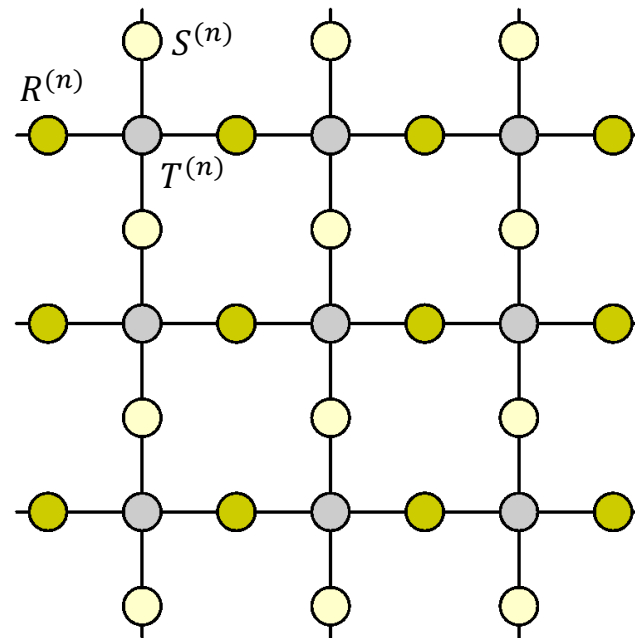
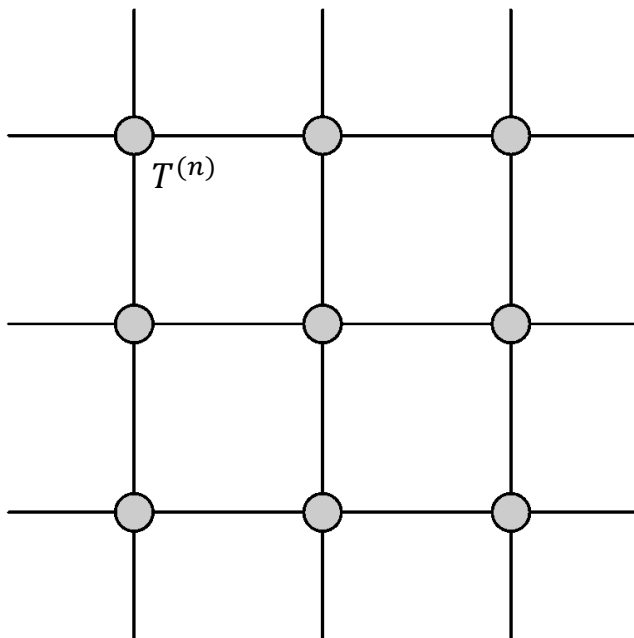
Bond-weighted TRG (BTRG)

Adachi+, PRB105(2022)L060402

✓ The generalized (improved) TRG algorithm

One considers the tensor network w/ **bond weights**.

Introducing the bond weights, one could improve the accuracy of the TRG w/o increasing the computational memory and execution time.



A good choice of the hyperparameter

Adachi+, PRB105(2022)L060402

✓ Introducing the hyperparameter $k \in \mathbb{R}$ in the SVD employed in the TRG

$$T_{IJ}^{(n)} \approx \sum_{\alpha=1}^D U_{I\alpha}^{(n)} \left(\sigma_{\alpha}^{(n)}\right)^{\frac{1-k}{2}} \left(\sigma_{\alpha}^{(n)}\right)^k \left(\sigma_{\alpha}^{(n)}\right)^{\frac{1-k}{2}} V_{J\alpha}^{(n)*}$$

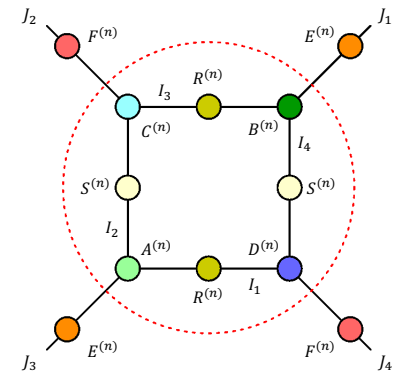
n labels the renormalization steps

With $k = 0$, the BTRG is reduced to be the Levin-Nave TRG.

✓ A good choice of k ?

By the TRG renormalization, $T^{(n+1)} \sim \left[\left(\sigma_{\alpha}^{(n)}\right)^{\frac{1-k}{2}} \left(\sigma_{\alpha}^{(n)}\right)^k \right]^4$

By the SVD of $T^{(n+1)}$, $T^{(n+1)} \sim \sigma_{\alpha}^{(n+1)}$



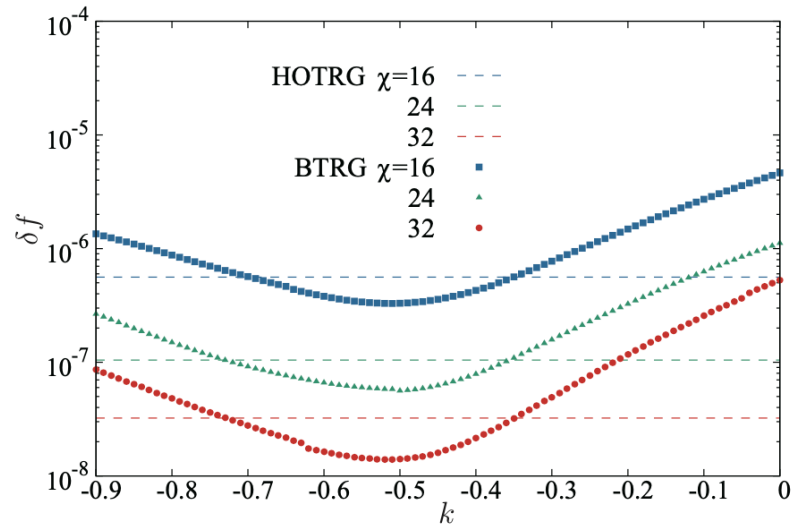
Suppose the singular-value spectrum becomes scale-invariant w/ sufficiently large n , we have

$$\left[\left(\sigma_{\alpha}^{(n)}\right)^{\frac{1-k}{2}} \left(\sigma_{\alpha}^{(n)}\right)^k \right]^4 = \sigma_{\alpha}^{(n)} \Rightarrow \boxed{k = -0.5}$$

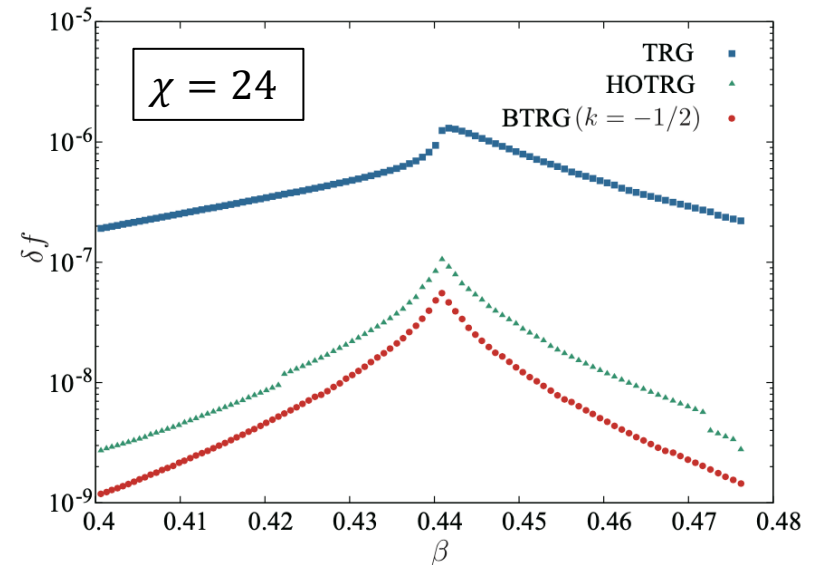
BTRG for the 2D classical Ising model

Adachi+, PRB105(2022)L060402

Relative error vs k



Comparison w/ other TRGs



$k = -0.5$ seems optimal and **the accuracy of the BTRG is higher than the Levin-Nave TRG and the HOTRG with the same bond dimension.**

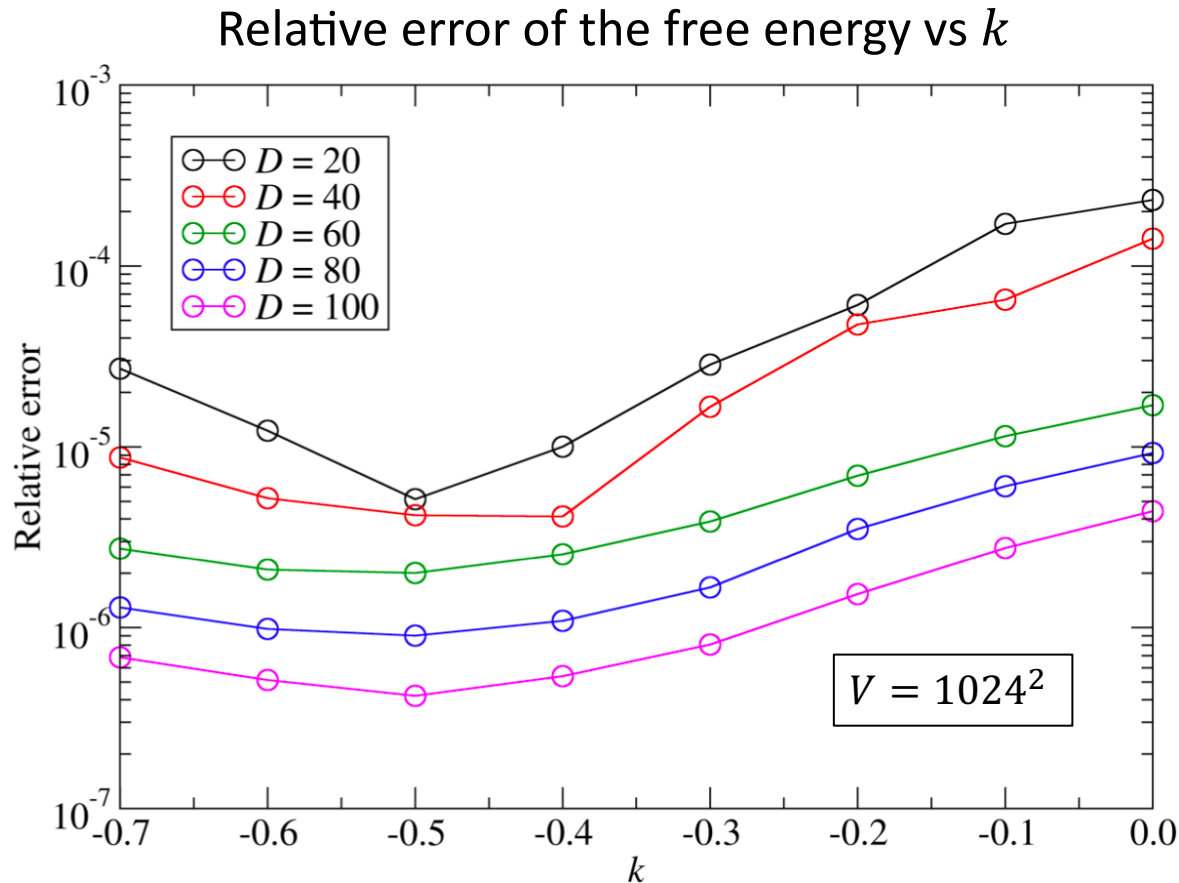
Introduction of k does not increase the computational cost.

Therefore, the cost of the BTRG is same with the Levin-Nave TRG.

BTRG for the lattice fermion?

- ✓ $k = -0.5$ seems to be just determined by the geometry of the tensor network.
→ **the optimal choice of k does not depend on the details of the model?**
- ✓ Any TRG algorithm can be extended to evaluate the Grassmann path integral.
→ **BTRG is also efficient for the lattice fermions?**
- ✓ We develop the Grassmann BTRG algorithm employing the formulation of [SA-Kadoh, JHEP10\(2021\)188](#), which allows us to extend the given TRG algorithm in a simple way to evaluate the Grassman path integral.

Benchmarking w/ the free massless Wilson fermion in 2D



Again, $k = -0.5$ seems optimal and the bond-weighting method does improve the Grassmann Levin-Nave TRG.

Finite-entanglement scaling

- ✓ In 1+1D, we have the finite-entanglement scaling based on the Matrix Product State (MPS), which is a tensor network method in the Hamiltonian formalism.

w.r.t. the bond dimension, the correlation length scales with $\xi_D \sim D^\kappa$, where

$$\kappa = \frac{6}{c \left(\sqrt{\frac{12}{c} + 1} \right)}. \quad (c: \text{the central charge})$$

Tagliacozzo+, PRB78(2008)024410,
Pollmann+, PRL102(2009)255701

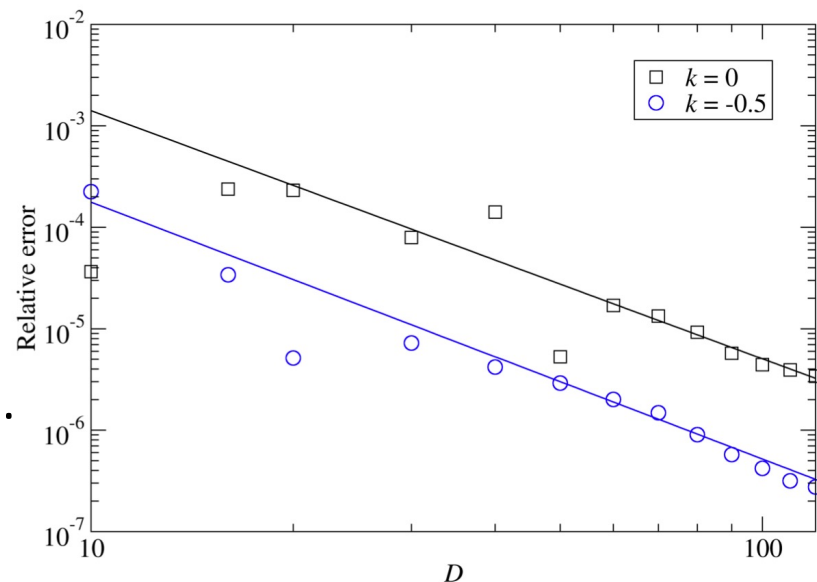
- ✓ Fitting the relative error of the free energy with $aD^{-2\kappa}$,

$$k = -0.5: a = 0.06(4), \kappa = 1.26(7)$$

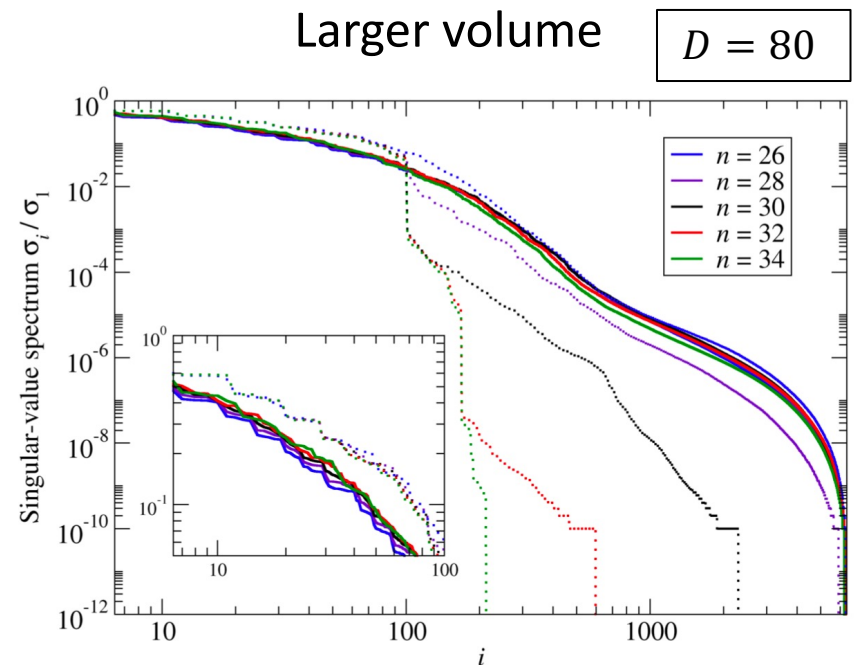
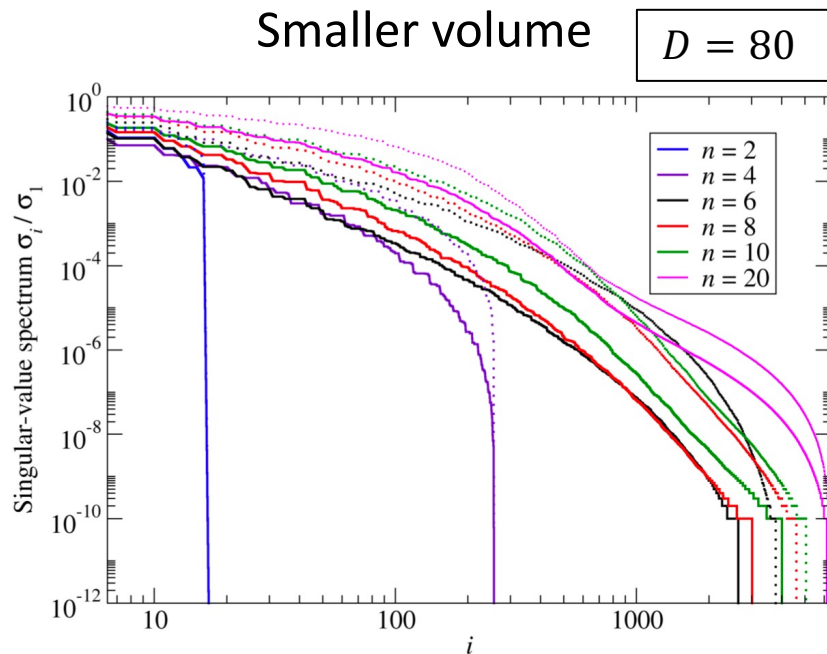
$$k = 0: a = 0.4(3), \kappa = 1.22(8)$$

The Grassmann BTRG reproduces $\kappa = 1.344 \dots$, corresponding to $c = 1$.

Relative error of the free energy vs D



Hierarchy of the singular values



n labels the renormalization steps or the lattice volume via $V = 2^n$

Grassmann BTRG (solid lines)

Grassmann Levin-Nave TRG (dotted lines)

In smaller volumes, the spectrum obtained by the Grassmann BTRG rapidly decays compared with the Grassmann Levin-Nave TRG.

The Levin-Nave algorithm does not reproduce the scale-invariant structure in the local Grassmann tensor, but the Grassmann BTRG does.

Summary

- Bond-weighting method is a new way to improve the TRG algorithm.
- The method was originally proposed for the spin system.
We numerically confirmed that the bond-weighting method is also efficient for the lattice fermions.
- Benchmarking with the 2D free Wilson fermions, we have found that the accuracy of the TRG is highly improved. The scale-invariant structure of the local Grassmann tensor is also reproduced.
- **A sample code of the Grassmann BTRG is available on GitHub.**
2D single-flavor Gross-Neveu-Wilson model at finite density is provided as an example.
see <https://github.com/akiyama-es/Grassmann-BTRG>