

# Measured-induced phase transitions by matrix product states scaling

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## Abstract

We investigate long quantum spin chains under continuous monitoring, employing matrix product states (MPS) and the time-dependent variational principle (TDVP) algorithm. Our findings reveal a notable phase transition in the error rate during monitoring, detectable even with low bond dimensions. This approach efficiently pinpoints critical parameters for measurement-induced phase transitions (MIPTs) in complex many-body quantum systems. Furthermore, we unveil a distinct charge-sharpening (CS) transition in the context of U(1) global spin charge, validating our TDVP approach for identifying phase transitions across diverse system dimensions and sizes.

## Models

We focus on two generic interacting systems in one dimension (whose MIPT has been also studied by different means in [1, 2, 3]) namely a chain of  $L$  spin-1/2 particles, unitarily evolving with  $U(1)$  symmetric (magnetisation conserving) Hamiltonians, defined as

$$\hat{H}_{J\text{-XXX}} = \sum_{i=1}^L \left( \hat{S}_i^x \hat{S}_{i+1}^x + \hat{S}_i^y \hat{S}_{i+1}^y + \hat{S}_i^z \hat{S}_{i+1}^z \right) + \sum_{i=1}^L J \left( \hat{S}_i^x \hat{S}_{i+2}^x + \hat{S}_i^y \hat{S}_{i+2}^y \right), \quad (1)$$

with spin operators  $\hat{S}_i^\alpha$  and with  $J = 0$  (XXX chain) or  $J = 1/2$  (J-XXX chain).

## Basic Concepts

Unitary time evolution tends to entangle the system, whereas measurements serve to disentangle it. In measured-induced phase transitions, measurements disrupt the system's coherent evolution and can drive it across phase boundaries.

### What is charge sharpening?

This transition separates two distinct entangling phases based on the ease or difficulty of determining the charge of the system through measurements.

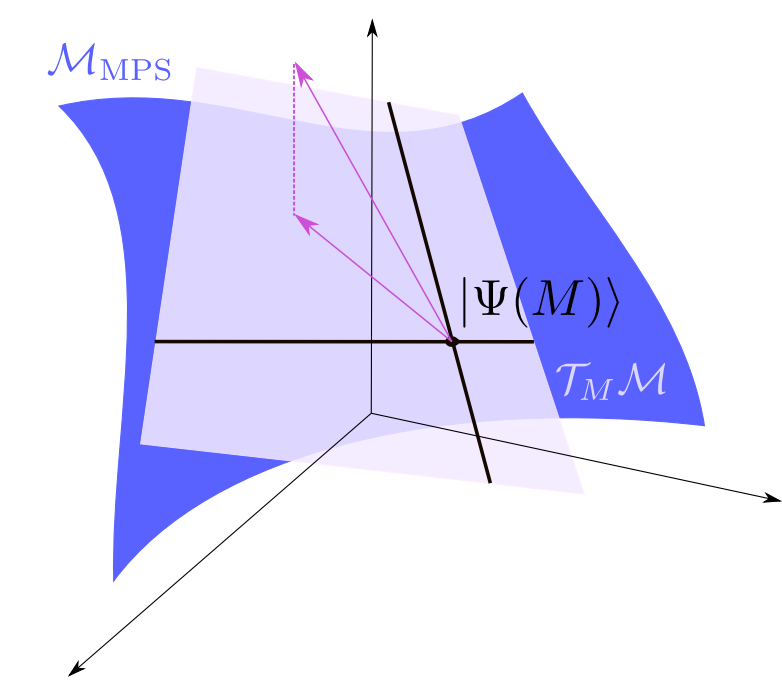
### Time Dependent Variational Principle

The time-evolved quantum state can be brought back to the initial manifold  $\mathcal{M}$  by a projection  $P_{\mathcal{T}_M \mathcal{M}}$ ,

$$i\partial_t |\Psi(M)\rangle = P_{\mathcal{T}_M \mathcal{M}} \hat{H} |\Psi(M)\rangle.$$

The projector  $P_{\mathcal{T}_M \mathcal{M}}$  which projects onto this tangent space is given by

$$P_{\mathcal{T}_M \mathcal{M}} : \mathcal{H} \rightarrow \mathcal{T}_M \mathcal{M}$$

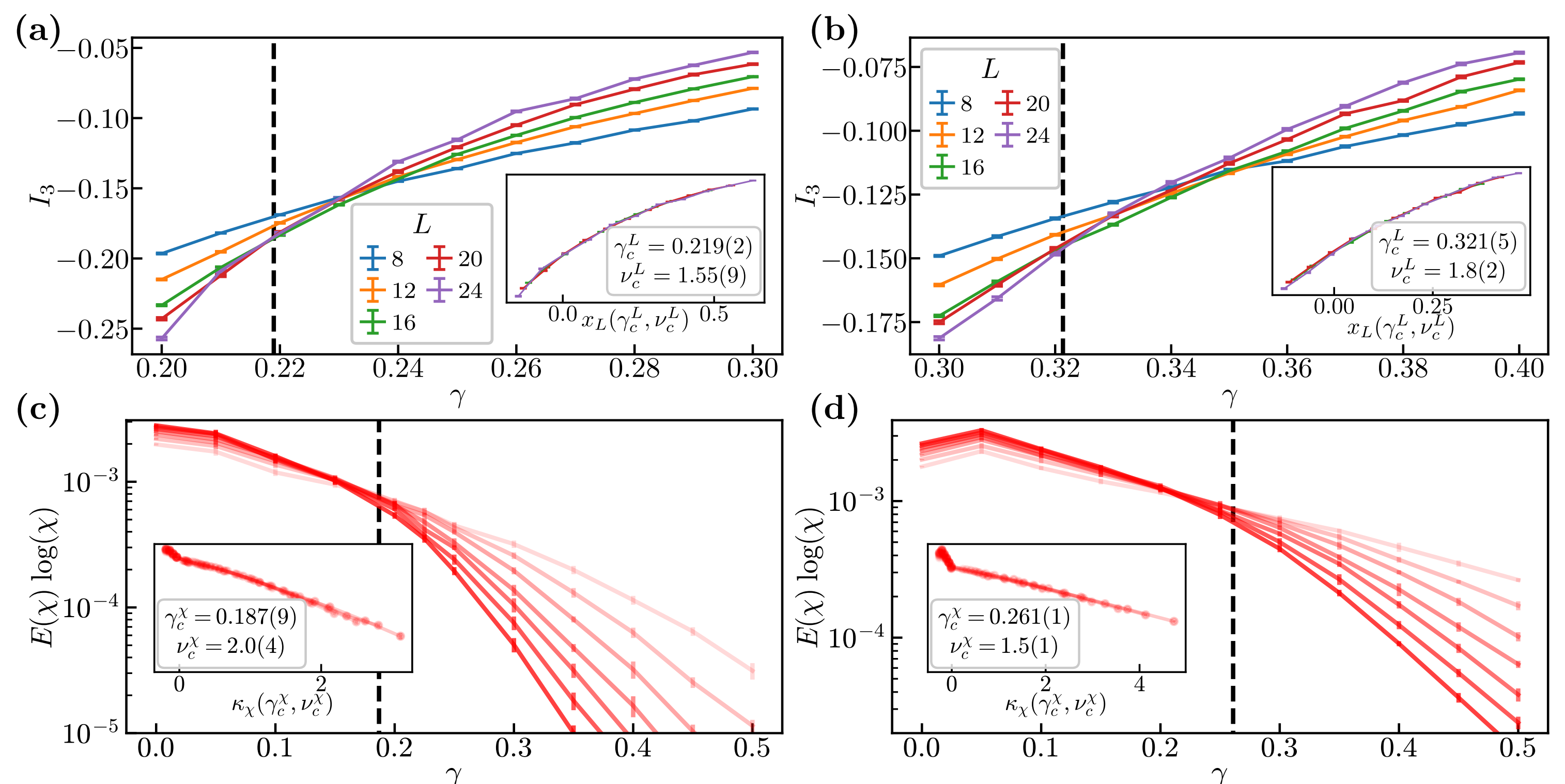


## References

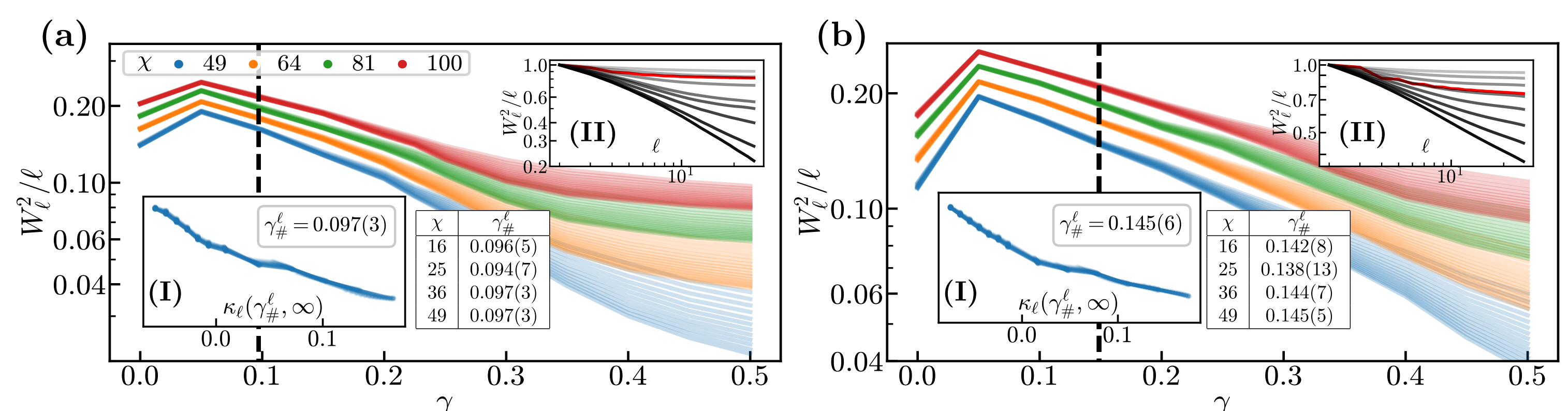
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## Numerical Results

Entanglement/Error phase transitions in (a),(c) XXX model, and (b),(d) in J-XXX model.



Charge-sharpening phase transitions for (a) the XXX model and (b) the J-XXX model.



## Methods

The associated monitored dynamics of the quantum state are described by the following stochastic Schrödinger equation (SSE) [4, 5, 6] for the many-body state  $|\psi\rangle$ ,

$$d|\psi_t\rangle = -iHdt|\psi_t\rangle + \sum_{i=1}^L \left[ \sqrt{\gamma}(\hat{S}_i^z - \langle \hat{S}_i^z \rangle_t) dW_t^i - \frac{\gamma}{2}(\hat{S}_i^z - \langle \hat{S}_i^z \rangle_t)^2 dt \right] |\psi_t\rangle, \quad (2)$$

with the expectation value of the local magnetisation given by the state at a given time  $\langle \hat{S}_i^z \rangle_t = \langle \psi_t | \hat{S}_i^z | \psi_t \rangle$ . Eq. (2) can be easily simulated by alternating its unitary and measurement terms via a Trotter splitting,

$$|\psi_{t+\delta t}\rangle \approx C e^{\sum_{j=1}^L [\delta W_t^j + 2\langle \hat{S}_j^z \rangle_t \gamma \delta t] \hat{S}_j^z} e^{-i\hat{H}\delta t} |\psi_t\rangle, \quad (3)$$

where the set of  $\delta W^i$  are generated each time step from a normal distribution with variance  $\sqrt{\gamma\delta t}$  and zero mean and  $C$  is a normalizing constant. In order to extend the CS protocol to large system sizes, we here consider the variance of the fluctuations of the local magnetization on a sub-system of size  $\ell$ , e.g. by defining  $Q_\ell = \sum_{j \in \ell} S_j^z$ , we introduce its variance as

$$W_\ell^2 = \langle Q_\ell^2 \rangle - \langle Q_\ell \rangle^2 = \sum_{i,j \in \ell} \langle S_i^z S_j^z \rangle^c, \quad (4)$$

In order to fit the MPS transition, given the much quicker convergence in  $\chi$  on the left of the critical point, we need to introduce a *piece-wise ansatz*, namely

$$x = \kappa_A(\gamma_0, \nu_0) = \begin{cases} (\gamma - \gamma_0) & \gamma < \gamma_0 \\ (\gamma - \gamma_0) A^{1/\nu_0} & \gamma > \gamma_0 \end{cases}, \quad (5)$$

and we cross-checked this scaling with the one where only the right side of the transition is fitted, giving analogous results.

## Discussion

Starting with the initial Néel state  $|\Psi(0)\rangle = |\downarrow\uparrow\downarrow\uparrow \dots \downarrow\uparrow\rangle$ , monitored evolution at different  $\chi$  and  $L$  values is performed. Analysis of error rate evolution reveals saturation after a timescale of order  $L$ , indicating a transition between volume-law and area-law phases. Precisely determined critical parameters suggest a smaller critical measurement rate  $\gamma_c^\chi$  than that of ED simulations. Unitary terms dominate for small  $L$ , shifting critical measurement rates leftward for larger  $L$ . Although critical exponents differ slightly between ED and MPS, well-converged results are obtained for small bond dimensions. Transition to the CS phase is observed through magnetization variance values, showing a shift from extensive to sub-linear scaling with increasing  $\gamma$ . Cross-checking with ED simulations confirms the observed transition, highlighting the method's importance in studying large systems with continuous time.