



Motivation

- Characterizing open quantum dynamics is critical for calibrating hardware and designing error correction.
- The Lindblad equation captures both coherent and dissipative dynamics under Markovian noise.
- Prior protocols assume known interaction structure — restrictive when the relevant error mechanisms are unknown.

Can we efficiently learn the full Lindbladian from its time evolution in situ, without structural assumptions?

Lindbladian Model

$$\frac{d}{dt}\rho = \mathcal{L}(\rho) = -i \sum_{P_i \in \mathcal{S}_H} h_i [P_i, \rho] + \sum_{P_i, P_j \in \mathcal{S}_D} a_{ij} \left(P_i \rho P_j - \frac{1}{2} \{P_j P_i, \rho\} \right)$$

- P_i : n -qubit Pauli operators
- $\mathcal{S}_H, \mathcal{S}_D$: unknown Hamiltonian and dissipator supports
- M = total nonzero Lindbladian terms (or an upper bound)
- η = smallest nonzero coefficient (or a lower bound)

Stage 1: Structure Learning

At short times, $e^{\mathcal{L}t}$ is approximated by its Taylor expansion with sparse Pauli structure. The derivatives of Pauli error rates directly reveal the generator structure:

$$\chi_{ii}^{(1)} = a_{ii} \text{ if } a_{ii} > 0 \quad \text{and} \quad \chi_{ii}^{(1)} = 0, \chi_{ii}^{(2)} \geq 2h_i^2 \text{ if } a_{ii} = 0$$

- Estimate all Pauli error rates $\chi_{ii}(t_m)$ at short times via population recovery [Flammia, O'Donnell, 2021].
- Fit Chebyshev interpolants; extract first and second derivatives at $t = 0$.
- Classify dissipator: $\chi_{ii}^{(1)} > \eta/2 \implies \widehat{\mathcal{S}}_D \leftarrow P_i$
Classify Hamiltonian: $\chi_{ii}^{(2)} \geq \eta^2 \implies \widehat{\mathcal{S}}_H \leftarrow P_i$
- Output candidate supports $\widehat{\mathcal{S}}_D$ and $\widehat{\mathcal{S}}_H \leftarrow \widehat{\mathcal{S}}_H \cup \widehat{\mathcal{S}}_D$

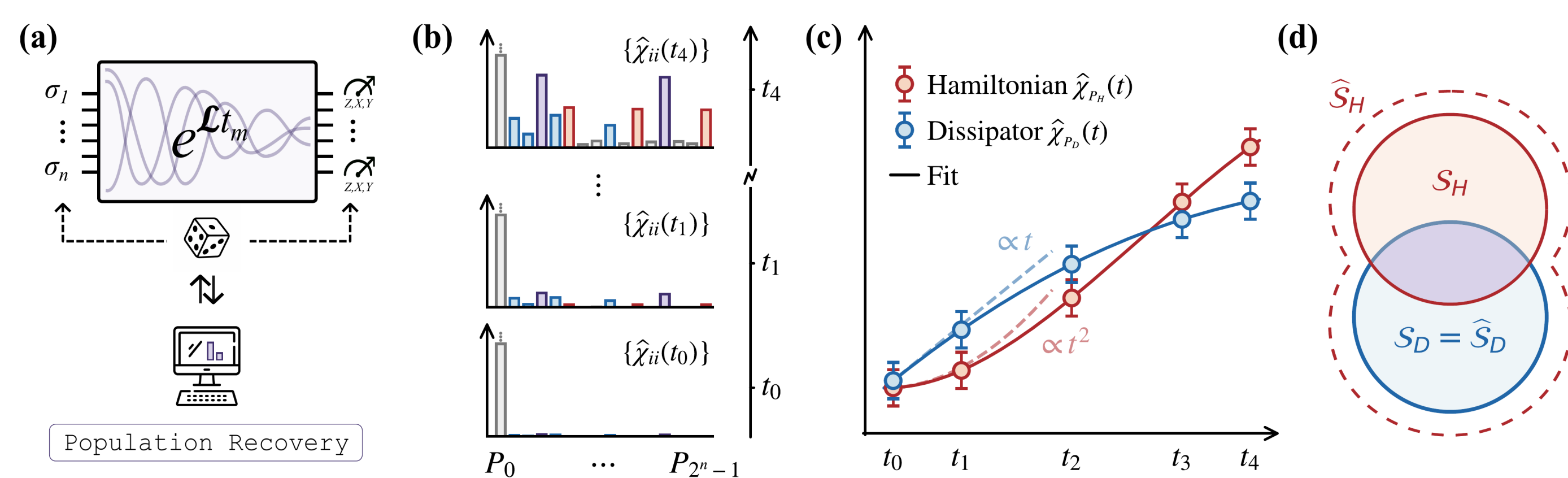


Figure 1. (a) Population recovery, (b) Pauli error rate distributions, (c) Chebyshev derivative fitting separates Hamiltonian ($\propto t^2$) from dissipator ($\propto t$), (d) recovered supports.

Result 1 — Structure Learning

We can recover a polynomially-sized superset of the Lindbladian structure using $\tilde{\mathcal{O}}(M^4/\eta^4)$ queries to $e^{\mathcal{L}t}$ with time resolution $t = \tilde{\Theta}(1/M)$.

Stage 2: Coefficient Learning

Given candidate supports $\widehat{\mathcal{S}}_D$ and $\widehat{\mathcal{S}}_H$, estimate all Lindbladian coefficients $\{h_i, a_{ij}\}$ to additive accuracy ε .

$$\text{Linear response relation: } \left. \frac{d}{dt} \langle O(t) \rangle \right|_{t=0} = \text{tr}(\rho \mathcal{L}^\dagger(O))$$

- Select Pauli probes (ρ, O) via patchwise tomography on candidate supports.
- For each probe, sample $f(t) = \text{tr}\{\rho e^{t\mathcal{L}^\dagger}(O)\}$ at Gauss–Chebyshev nodes; fit and extract df/dt at $t = 0$.
- Assemble the linear system $\mathbf{d} = C\mathbf{x}$ from the derivative estimates.
- Solve classically; error amplification controlled by $\nu = \|C^{-1}\|_{\infty \rightarrow \infty}$

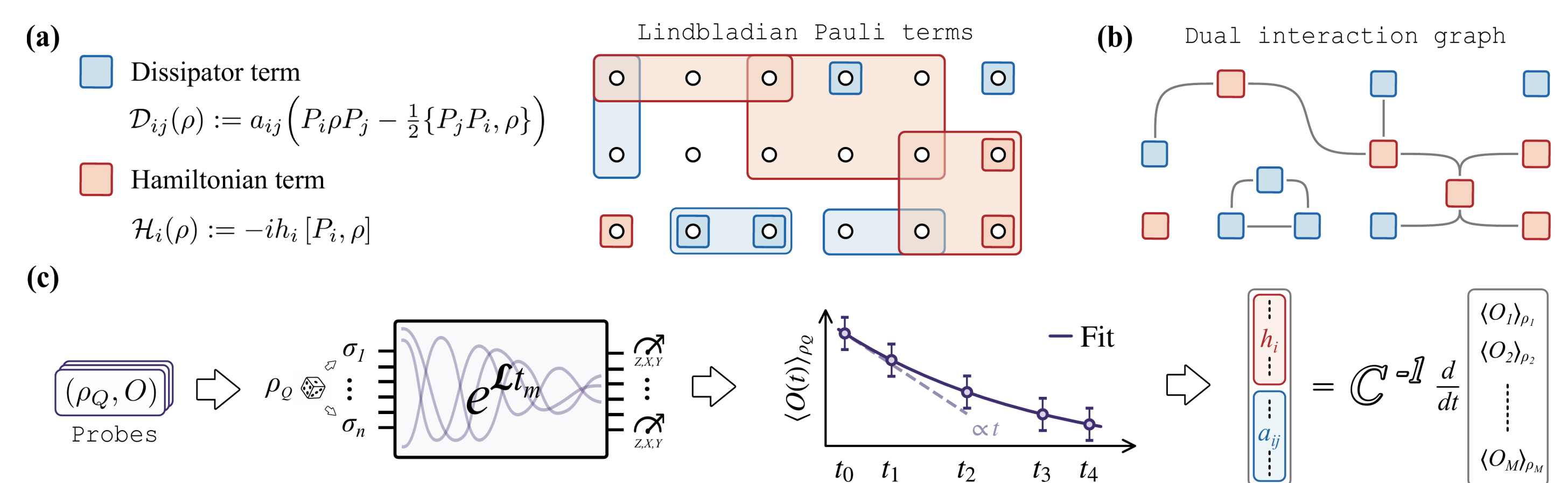


Figure 2. (a) Lindbladian Pauli terms, (b) dual interaction graph; degree $\mathfrak{d}$ controls the derivatives of Pauli observables, (c) linear system from short-time derivatives via Chebyshev interpolation.

Result 2 — Coefficient Learning

Given a superset of the Lindbladian structure, we can recover all coefficients $\{h_i, a_{ij}\}$ to ε -accuracy using

$$\tilde{\mathcal{O}}\left(\min(9^k, \widehat{M}) \mathfrak{d}^2 \nu^2 / \varepsilon^2\right)$$

queries to $e^{\mathcal{L}t}$ with time resolution $t = \tilde{\Theta}(1/\mathfrak{d})$.

Time Resolution

- Our protocol uses minimum evolution time $t_{\min} = \tilde{\Theta}(1/M)$.
- This is provably near-optimal. Coarser resolution forces any protocol to use exponentially many samples. Key idea: rapid mixing drives all input states exponentially close to the maximally mixed state, erasing structural information.

Result 3 — Optimality of Time Resolution

At resolution $t_0 = \Theta(M^{2/\kappa-1})$, $\kappa \geq 2$ there exist M -sparse Lindbladians requiring $\Omega(e^n)$ queries.

Takeaways

- First ansatz-free Lindbladian learning protocol in situ.
- No prior structure, no ancillas — product-state preparations and Pauli measurements only.
- Total sample complexity $m_{\text{tot}} = \tilde{\mathcal{O}}\left(\frac{M^4 \nu^2}{\varepsilon^4}\right)$