

Petz recoverability versus Wilson-loop diagnostics in \mathbb{Z}_2 lattice gauge theory (2+1D)

Exact diagonalization benchmark on small open lattices

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Abstract

We provide a finite-size benchmark testing whether a Petz-type recoverability proxy correlates with Wilson-loop confinement diagnostics in \mathbb{Z}_2 lattice gauge theory in 2+1 dimensions. We compute ground states by sparse exact diagonalization on 2×2 and 2×3 plaquette lattices with open boundary conditions, using qubits on links and enforcing the gauge-invariant sector via a Gauss-law penalty (verified numerically by $\langle G_v \rangle \simeq 1$ for all vertices). For opposite-corner plaquette patches A and C , we define a buffer $B(w)$ by BFS thickening in the link-adjacency graph (with $w = 0$ denoting an empty buffer) and evaluate a trace-renormalized, regularized Petz-type recoverability error $E_{\text{rec}}^{\text{Petz}}(w) = -\log F(\rho_{ABC}, \hat{\rho}_{ABC}(w))$. As a confinement proxy we use Wilson loops and the Creutz ratio $\chi(2, 2)$, treated as an effective string tension σ_{eff} . In 2+1 dimensions the natural confinement-length proxy is σ_{eff}^{-1} . Across couplings g we observe that larger σ_{eff}^{-1} correlates with larger fixed-collar recovery error $E_{\text{rec}}^{\text{Petz}}(w = 1)$ (small-lattice benchmark).

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1 Introduction

Petz-type recoverability provides an operational measure of approximate quantum Markov structure: given a tripartition A – B – C , one asks how well ρ_{ABC} can be reconstructed from ρ_{AB} using a recovery channel acting only on B . In gauge theories, subsystem definitions are subtle due to constraints, and reduced states can depend on the chosen prescription. In companion work we proposed a reproducible protocol for computing Petz-type recoverability on lattices and conjectured links to confinement diagnostics. Here we provide a small-lattice benchmark in a genuine lattice gauge theory.

Scope. We work on very small lattices (2×2 and 2×3 plaquettes) where exact diagonalization is feasible. Finite-size and boundary effects are large. Our goal is a falsifiable, reproducible benchmark rather than a thermodynamic-limit statement.

Dimensional analysis in 2+1 dimensions. In $d+1$ spacetime dimensions, the string tension has dimension $[\sigma] = (\text{length})^{-(d-1)}$. In 2+1 dimensions ($d = 2$), the natural length scale is therefore $\ell_\sigma \sim \sigma^{-1}$, motivating comparisons to σ_{eff}^{-1} .

2 Model and setup

2.1 \mathbb{Z}_2 Hamiltonian lattice gauge theory on links

We work on a spatial $N_x \times N_y$ plaquette lattice with open boundary conditions (OBC), with qubits on links. We use the Hamiltonian

$$H(g) = -g \sum_{\ell} X_{\ell} - \frac{1}{g} \sum_p \prod_{\ell \in \partial p} Z_{\ell} + \Lambda \sum_v (I - G_v),$$

where X_{ℓ} and Z_{ℓ} are Pauli operators on link ℓ , the plaquette term is the product around ∂p , and the gauge generators are

$$G_v := \prod_{\ell \ni v} X_{\ell}.$$

We take $\Lambda = 50$ in the numerics and verify $\langle G_v \rangle \simeq 1$ for all vertices in the computed ground state.

2.2 Tripartition geometry (PARCHES) and collaring rule

Subsystems are defined as subsets of links (an extended-Hilbert-space tensor-product of link degrees of freedom). We choose A and C as *plaquette patches* (four links) in opposite corners: A is the southwest plaquette patch and C is the northeast plaquette patch.

We define the collar $B(w)$ by BFS thickening in the link-adjacency graph (two links adjacent if they share a vertex):

$$B(0) = \emptyset, \quad B(w) = \{\ell : \text{dist}_{\text{link}}(\ell, A) \leq w\} \setminus (A \cup C), \quad w \geq 1.$$

Remark 2.1. On very small lattices, $E_{\text{rec}}^{\text{Petz}}(w)$ need not be monotone in w due to finite-size saturation and the trace normalization used for the regularized Petz-type reconstruction. We therefore emphasize fixed-collar comparisons (notably $w = 1$) and robustness checks, rather than interpreting $E_{\text{rec}}^{\text{Petz}}(w)$ as a clean exponential decay profile at these sizes.

3 Observables

3.1 Regularized Petz-type recoverability error

For each collar width w , we compute reduced states ρ_{ABC} , ρ_{AB} , ρ_B , and ρ_{BC} . We use a regularized Petz-type map

$$\mathcal{R}_{B \rightarrow BC}^{\text{Petz}(\delta)}(X_B) = \rho_{BC}^{1/2} \left(\rho_{B,\delta}^{-1/2} X_B \rho_{B,\delta}^{-1/2} \otimes I_C \right) \rho_{BC}^{1/2}, \quad \rho_{B,\delta} := \rho_B + \delta I,$$

with $\delta = 10^{-12}$, and normalize the reconstructed operator by its trace before computing fidelity. We report

$$E_{\text{rec}}^{\text{Petz}}(w) := -\log F(\rho_{ABC}, \hat{\rho}_{ABC}(w)).$$

3.2 Wilson loops and Creutz proxy

For a closed loop γ we define the \mathbb{Z}_2 Wilson loop operator $W(\gamma) = \prod_{\ell \in \gamma} Z_\ell$. As a practical proxy for string tension on tiny lattices, we compute the Creutz ratio

$$\chi(2, 2) = -\log \frac{\langle W(2, 2) \rangle \langle W(1, 1) \rangle}{\langle W(2, 1) \rangle \langle W(1, 2) \rangle},$$

and set $\sigma_{\text{eff}} := \chi(2, 2)$.

Remark 3.1 (Dimensional analysis in 2+1 dimensions). In 2+1 spacetime dimensions, $[\sigma] = (\text{length})^{-1}$, so the natural confinement-length proxy is σ_{eff}^{-1} (not $\sigma_{\text{eff}}^{-1/2}$).

4 Results

4.1 Summary table

$N_x \times N_y$	g	E_0	σ_{eff}	$1/\sigma_{\text{eff}}$	$E_{\text{rec}}^{\text{Petz}}(w=0)$	$E_{\text{rec}}^{\text{Petz}}(w=1)$
2×2	0.5	-9.5660	0.150780	6.632169	1.566201×10^{-1}	3.685747×10^{-1}
2×2	1.0	-12.4973	1.073026	0.931944	1.202218×10^{-3}	4.432791×10^{-3}
2×2	2.0	-24.0625	2.418666	0.413451	4.281872×10^{-6}	1.848438×10^{-5}
2×3	0.5	-13.9474	0.102507	9.755459	4.304489×10^{-1}	4.261377×10^{-1}
2×3	1.0	-17.7472	1.059645	0.943713	4.315631×10^{-2}	4.311739×10^{-2}
2×3	2.0	-34.0937	2.417752	0.413607	2.869337×10^{-3}	2.867963×10^{-3}

Table 1: Finite-size benchmark for \mathbb{Z}_2 lattice gauge theory (OBC, link qubits). The ground state lies in the gauge-invariant sector (numerically verified by $\langle G_v \rangle \simeq 1$ for all vertices). We report $\sigma_{\text{eff}} := \chi(2, 2)$ as a Creutz proxy and compare fixed-collar recoverability $E_{\text{rec}}^{\text{Petz}}(w=1)$ against the natural 2+1D length proxy $1/\sigma_{\text{eff}}$.

4.2 Recoverability profiles

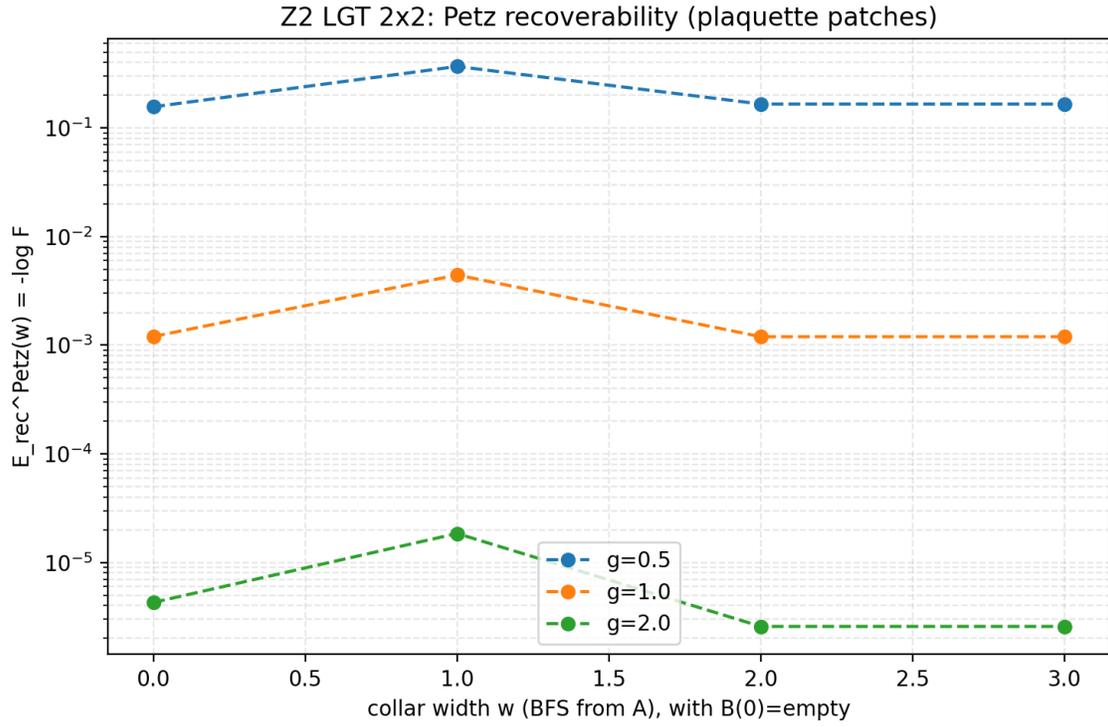


Figure 1: Petz-type recoverability profile $E_{\text{rec}}^{\text{Petz}}(w)$ on the 2×2 plaquette lattice (12 links) for plaquette-patch tripartitions with opposite-corner patches A and C . The collar $B(w)$ is defined by BFS thickening in the link-adjacency graph, with $B(0) = \emptyset$.

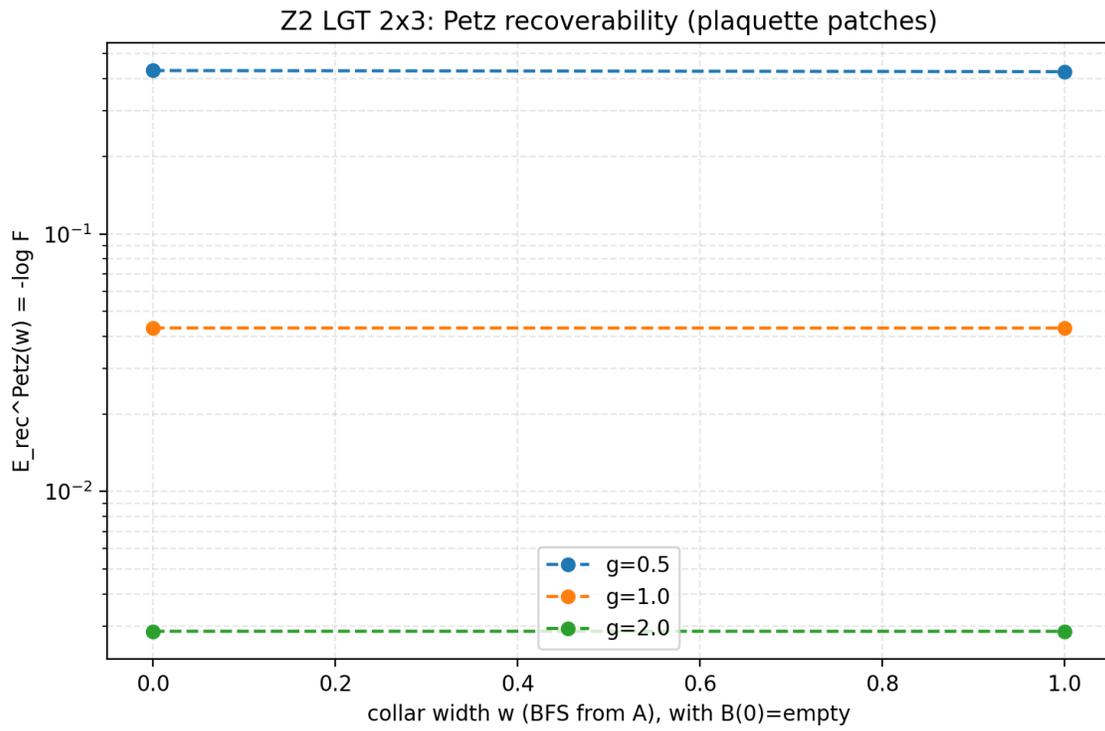


Figure 2: Petz-type recoverability profile $E_{\text{rec}}^{\text{Petz}}(w)$ on the 2×3 plaquette lattice (17 links) for the same plaquette-patch geometry. Dense Petz reconstruction restricts feasible collar widths; here we report $w = 0, 1$.

4.3 Fixed-collar recoverability versus confinement-length proxy

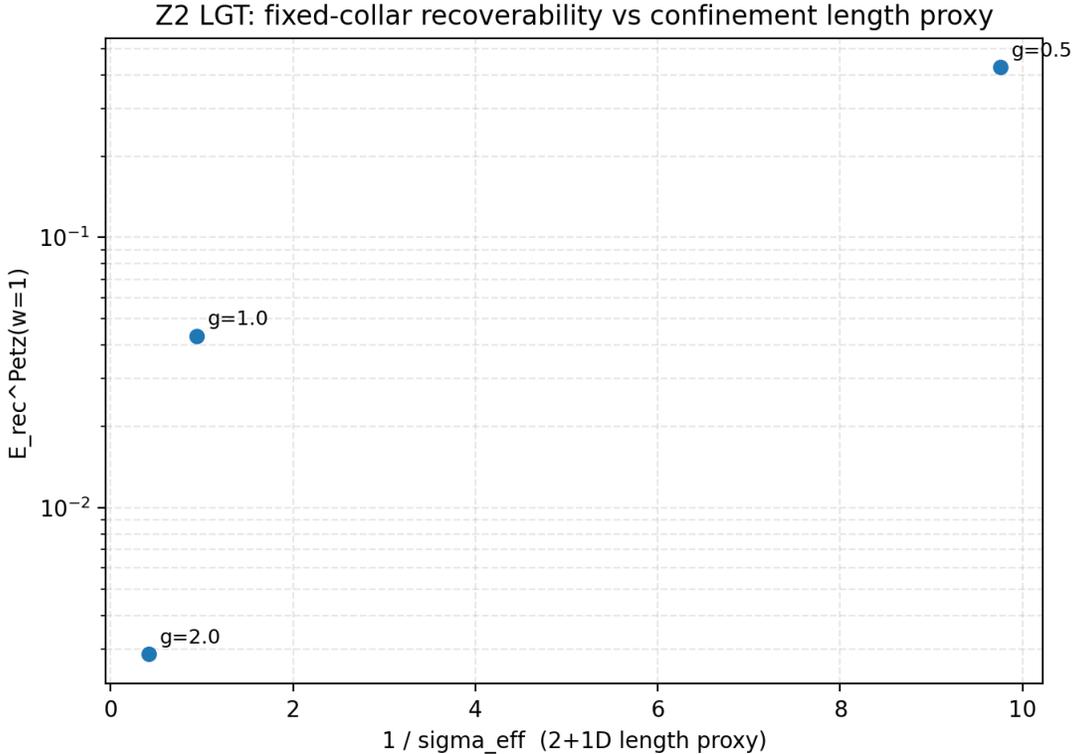


Figure 3: Fixed-collar recoverability versus confinement-length proxy on the 2×3 plaquette lattice. The horizontal axis is σ_{eff}^{-1} with $\sigma_{\text{eff}} = \chi(2, 2)$ (Creutz ratio), and the vertical axis is the fixed-collar recovery error $E_{\text{rec}}^{\text{Petz}}(w = 1)$. Across couplings g , larger σ_{eff}^{-1} correlates with larger $E_{\text{rec}}^{\text{Petz}}(w = 1)$ (finite-size benchmark).

5 Discussion

On 2×3 , the near-equality of $E_{\text{rec}}^{\text{Petz}}(w = 0)$ and $E_{\text{rec}}^{\text{Petz}}(w = 1)$ indicates rapid finite-size saturation of the buffer effect for this geometry. On 2×2 , mild non-monotonicity in w can occur; we therefore emphasize fixed-collar comparisons and treat collar dependence as a diagnostic at these sizes.

The primary observation is a clear correlation between fixed-collar recoverability $E_{\text{rec}}^{\text{Petz}}(w = 1)$ and the confinement-length proxy σ_{eff}^{-1} . Extending to larger collars and lattices will require tensor-network contractions or alternative recovery implementations.

A Reproducibility: code to regenerate the figures

The figures in this paper can be regenerated by running the Python script below. It computes ground states by sparse exact diagonalization, checks the gauge constraint $\langle G_v \rangle \simeq 1$ in the resulting state, evaluates the Petz-type recoverability profile for the plaquette-patch geometry, and produces the PNG plots used in the main text.

Dependencies. Python 3 with numpy, scipy, and matplotlib.

Run. Save the listing to `paper_d_reproduce.py` and run:

```
python paper_d_reproduce.py
```

This generates: `z2_2x2_recoverability_v2.png`, `z2_2x3_recoverability_v2.png`, `z2_2x3_fixed_w` and `paper_d_results.csv`.

Listing 1: `paper_d_reproduce.py` (reproducibility script)

```
1 # NOTE: This is the same code used to generate the CSV/PNGs for Paper D v1.
2 import csv
3 import numpy as np
4 import scipy.sparse as sp
5 import scipy.sparse.linalg as spla
6 import scipy.linalg as la
7 import matplotlib
8 matplotlib.use("Agg")
9 import matplotlib.pyplot as plt
10
11 def idx_h(x, y, Nx, Ny): return y * Nx + x
12 def idx_v(x, y, Nx, Ny):
13     Nh = Nx * (Ny + 1)
14     return Nh + y * (Nx + 1) + x
15
16 def build_lattice(Nx, Ny):
17     def vid(x, y): return y * (Nx + 1) + x
18     link_endpoints = []
19     for y in range(Ny + 1):
20         for x in range(Nx):
21             link_endpoints.append((vid(x, y), vid(x + 1, y)))
22     for y in range(Ny):
23         for x in range(Nx + 1):
24             link_endpoints.append((vid(x, y), vid(x, y + 1)))
25     Nlinks = len(link_endpoints)
26
27     plaquettes = []
28     for y in range(Ny):
29         for x in range(Nx):
30             pl = [idx_h(x, y, Nx, Ny), idx_v(x + 1, y, Nx, Ny), idx_h(x, y + 1, Nx, Ny), idx_v(x, y,
31                 Nx, Ny)]
32             plaquettes.append(pl)
33
34     Nv = (Nx + 1) * (Ny + 1)
35     stars = [[] for _ in range(Nv)]
36     for ell, (a, b) in enumerate(link_endpoints):
37         stars[a].append(ell); stars[b].append(ell)
38
39     v2links = [[] for _ in range(Nv)]
40     for ell, (a, b) in enumerate(link_endpoints):
41         v2links[a].append(ell); v2links[b].append(ell)
42     adjacency = [set() for _ in range(Nlinks)]
43     for v in range(Nv):
44         Ls = v2links[v]
45         for i in Ls:
46             for j in Ls:
47                 if i != j:
48                     adjacency[i].add(j)
```

```

48
49     return Nlinks, plaquettes, stars, adjacency
50
51 def mask_from_links(link_list):
52     m = 0
53     for ell in link_list:
54         m |= (1 << int(ell))
55     return m
56
57 def parity_u64(x): return x.bit_count() & 1
58
59 def build_hamiltonian_z2(Nlinks, plaquette_masks, star_masks, g, Lambda=0.0):
60     dim = 1 << Nlinks
61     states = np.arange(dim, dtype=np.uint32 if Nlinks <= 32 else np.uint64)
62     invg = 1.0 / float(g)
63     diag = np.zeros(dim, dtype=np.float64)
64
65     for mp in plaquette_masks:
66         par = np.fromiter((parity_u64(int(s) & mp) for s in states), count=dim, dtype
67                             =np.int8)
68         eig = 1.0 - 2.0 * par
69         diag += -(invg) * eig
70
71     if Lambda != 0.0:
72         diag += float(Lambda) * len(star_masks)
73
74     rows = []; cols = []; data = []
75
76     rows.extend(range(dim)); cols.extend(range(dim)); data.extend(diag.tolist())
77
78     for ell in range(Nlinks):
79         mask = 1 << ell
80         tgt = states ^ mask
81         rows.extend(states.tolist()); cols.extend(tgt.tolist())
82         data.extend((-float(g)) * np.ones(dim, dtype=np.float64).tolist())
83
84     if Lambda != 0.0:
85         for mv in star_masks:
86             tgt = states ^ mv
87             rows.extend(states.tolist()); cols.extend(tgt.tolist())
88             data.extend((-float(Lambda)) * np.ones(dim, dtype=np.float64).tolist())
89
90     H = sp.coo_matrix((np.array(data, dtype=np.float64),
91                       (np.array(rows), np.array(cols))), shape=(dim, dim)).tocsr()
92     H.sum_duplicates()
93     return H
94
95 def ground_state(H, tol=1e-10, maxiter=30000):
96     vals, vecs = spla.eigsh(H, k=1, which="SA", tol=tol, maxiter=maxiter)
97     v = vecs[:, 0].astype(np.complex128)
98     v /= np.linalg.norm(v)
99     return float(vals[0]), v
100
101 def plaquette_patch_A_C(Nx, Ny):
102     A = [idx_h(0,0,Nx,Ny), idx_v(1,0,Nx,Ny), idx_h(0,1,Nx,Ny), idx_v(0,0,Nx,Ny)]
103     C = [idx_h(Nx-1,Ny-1,Nx,Ny), idx_v(Nx,Ny-1,Nx,Ny), idx_h(Nx-1,Ny,Nx,Ny), idx_v(
104         Nx-1,Ny-1,Nx,Ny)]
105     return sorted(set(A)), sorted(set(C))

```

```

105 def bfs_distances(adjacency, sources):
106     n = len(adjacency)
107     dist = np.full(n, fill_value=-1, dtype=np.int32)
108     q = []
109     for s in sources:
110         dist[s] = 0; q.append(s)
111     head = 0
112     while head < len(q):
113         u = q[head]; head += 1
114         for v in adjacency[u]:
115             if dist[v] < 0:
116                 dist[v] = dist[u] + 1
117                 q.append(v)
118     return dist
119
120 def buffer_Bw(adjacency, A, C, w):
121     if w <= 0:
122         return []
123     distA = bfs_distances(adjacency, A)
124     B = [i for i, d in enumerate(distA) if (d >= 0 and d <= w and i not in A and i
125         not in C)]
126     return sorted(B)
127
128 def reduce_pure_to_density_qubits(psi, N, keep_qubits):
129     keep = list(keep_qubits)
130     trace = [i for i in range(N) if i not in keep]
131     psi_t = psi.reshape([2]*N)
132     perm = keep + trace
133     psi_perm = np.transpose(psi_t, axes=perm)
134     d_keep = 1 << len(keep)
135     d_tr = 1 << len(trace)
136     M = psi_perm.reshape(d_keep, d_tr)
137     rho = M @ M.conj().T
138     return 0.5*(rho + rho.conj().T)
139
140 def partial_trace_rho(rho, dims, keep):
141     dims = list(dims)
142     N = len(dims)
143     keep = list(keep)
144     trace = [i for i in range(N) if i not in keep]
145     d_keep = int(np.prod([dims[i] for i in keep]))
146     d_tr = int(np.prod([dims[i] for i in trace]))
147     rho = np.asarray(rho, dtype=np.complex128)
148     rho_t = rho.reshape(dims + dims)
149     perm = keep + trace + [i+N for i in keep] + [i+N for i in trace]
150     rho_p = np.transpose(rho_t, axes=perm).reshape(d_keep, d_tr, d_keep, d_tr)
151     out = np.trace(rho_p, axis1=1, axis2=3)
152     return 0.5*(out + out.conj().T)
153
154 def sqrtm_psd(M, eps=0.0):
155     w, V = la.eigh(np.asarray(M))
156     w = np.maximum(w, eps)
157     return (V * np.sqrt(w)) @ V.conj().T
158
159 def invsqrtm_pd(M):
160     w, V = la.eigh(np.asarray(M))
161     if np.min(w) <= 0:
162         raise ValueError("Matrix_not_strictly_positive;_increase_delta.")
163     return (V * (1.0/np.sqrt(w))) @ V.conj().T

```

```

163
164 def fidelity_squared(rho, sigma):
165     sr = sqrtm_psd(rho)
166     inner = sr @ sigma @ sr
167     return (np.trace(sqrtm_psd(inner)).real)**2
168
169 def petz_recoverability_from_rhoABC(rho_ABC, dA, dB, dC, delta=1e-12):
170     dims = [dA, dB, dC]
171     rho_AB = partial_trace_rho(rho_ABC, dims, keep=[0,1])
172     rho_B = partial_trace_rho(rho_ABC, dims, keep=[1])
173     rho_BC = partial_trace_rho(rho_ABC, dims, keep=[1,2])
174     rho_B_delta = rho_B + delta*np.eye(dB, dtype=np.complex128)
175     O_B = invsqrtm_pd(rho_B_delta)
176     sqrt_rho_BC = sqrtm_psd(rho_BC)
177     IA = np.eye(dA, dtype=np.complex128)
178     IC = np.eye(dC, dtype=np.complex128)
179     sandwich = np.kron(IA, O_B)
180     middle = sandwich @ rho_AB @ sandwich
181     middle = 0.5*(middle + middle.conj().T)
182     left = np.kron(IA, sqrt_rho_BC)
183     sigma = left @ np.kron(middle, IC) @ left
184     sigma = 0.5*(sigma + sigma.conj().T)
185     tr_sigma = float(np.trace(sigma).real)
186     sigmaN = sigma / tr_sigma
187     F = float(np.clip(fidelity_squared(rho_ABC, sigmaN), 0.0, 1.0))
188     E = -np.log(F + 1e-16)
189     return E, F, tr_sigma
190
191 def expval_Z_string(psi, mask, Nlinks):
192     dim = 1 << Nlinks
193     states = np.arange(dim, dtype=np.uint32 if Nlinks <= 32 else np.uint64)
194     par = np.fromiter((parity_u64(int(s) & mask) for s in states), count=dim, dtype=
195                       np.int8)
196     eig = 1.0 - 2.0 * par
197     prob = (psi.conj() * psi).real
198     return float(np.sum(prob * eig))
199
200 def expval_X_string(psi, mask, Nlinks):
201     dim = 1 << Nlinks
202     s = np.arange(dim, dtype=np.uint32 if Nlinks <= 32 else np.uint64)
203     t = s ^ mask
204     return float(np.vdot(psi, psi[t]).real)
205
206 def gauge_diagnostics(psi, star_masks, Nlinks):
207     vals = [expval_X_string(psi, mv, Nlinks) for mv in star_masks]
208     return float(np.min(vals)), float(np.mean(vals))
209
210 def rectangle_loop_mask(Nx, Ny, x0, y0, R, T):
211     links = set()
212     for x in range(x0, x0+R):
213         links.add(idx_h(x, y0, Nx, Ny))
214         links.add(idx_h(x, y0+T, Nx, Ny))
215     for y in range(y0, y0+T):
216         links.add(idx_v(x0, y, Nx, Ny))
217         links.add(idx_v(x0+R, y, Nx, Ny))
218     return mask_from_links(sorted(links))
219
220 def creutz_ratio(W_RT, W_Rm1Tm1, W_Rm1T, W_RTm1):
221     eps = 1e-16

```

```

221     return -np.log((W_RT*W_Rm1Tm1 + eps) / (W_Rm1T*W_RTm1 + eps))
222
223 def run_lattice(Nx, Ny, g_list, Lambda, delta, w_list, max_abc_qubits, out_prefix):
224     Nlinks, plaquettes, stars, adjacency = build_lattice(Nx, Ny)
225     plaquette_masks = [mask_from_links(pl) for pl in plaquettes]
226     star_masks = [mask_from_links(st) for st in stars if len(st) > 0]
227
228     A, C = plaquette_patch_A_C(Nx, Ny)
229     Bw_dict = {w: buffer_Bw(adjacency, A, C, w) for w in w_list}
230
231     rows = []
232     per_g = []
233
234     for g in g_list:
235         H = build_hamiltonian_z2(Nlinks, plaquette_masks, star_masks, g=g, Lambda=
                Lambda)
236         E0, psi = ground_state(H)
237         mnG, avG = gauge_diagnostics(psi, star_masks, Nlinks)
238
239         # W(1,1) avg
240         W11_vals = []
241         for y0 in range(Ny):
242             for x0 in range(Nx):
243                 W11_vals.append(expval_Z_string(psi, rectangle_loop_mask(Nx, Ny, x0, y0,
                1, 1), Nlinks))
244         W11_avg = float(np.mean(W11_vals))
245
246         # sigma_eff = chi(2,2) at (0,0) (simple benchmark)
247         sigma_eff = np.nan
248         inv_sigma = np.nan
249         if Nx >= 2 and Ny >= 2:
250             W22 = expval_Z_string(psi, rectangle_loop_mask(Nx, Ny, 0, 0, 2, 2), Nlinks
                )
251             W21 = expval_Z_string(psi, rectangle_loop_mask(Nx, Ny, 0, 0, 2, 1), Nlinks
                )
252             W12 = expval_Z_string(psi, rectangle_loop_mask(Nx, Ny, 0, 0, 1, 2), Nlinks
                )
253             W11 = expval_Z_string(psi, rectangle_loop_mask(Nx, Ny, 0, 0, 1, 1), Nlinks
                )
254             sigma_eff = creutz_ratio(W22, W11, W21, W12)
255             inv_sigma = 1.0 / sigma_eff if sigma_eff != 0 else np.nan
256
257         Erec_dict = {}
258         for w in w_list:
259             B = Bw_dict[w]
260             keep = A + B + C
261             k = len(keep)
262
263             if k > max_abc_qubits:
264                 rows.append([Nx, Ny, g, E0, mnG, avG, w, k, np.nan, np.nan, np.nan, W11_avg,
                sigma_eff, inv_sigma])
265                 continue
266
267             rho_ABC = reduce_pure_to_density_qubits(psi, Nlinks, keep)
268             dA = 1 << len(A)
269             dB = 1 << len(B)
270             dC = 1 << len(C)
271
272             Erec, F, tr_sigma = petz_recoverability_from_rhoABC(rho_ABC, dA, dB, dC,

```

```

273         delta=delta)
274         Erec_dict[w] = Erec
275         rows.append([Nx,Ny,g,E0,mnG,avG,w,k,Erec,F,tr_sigma,W11_avg,sigma_eff,
276                     inv_sigma])
277
278     per_g.append((g, Erec_dict, sigma_eff, inv_sigma))
279
280 # plot recoverability
281 plt.figure(figsize=(7.5,5))
282 for (g, Erec_dict, _, _) in per_g:
283     ws = sorted(Erec_dict.keys())
284     Es = [max(Erec_dict[w], 1e-16) for w in ws]
285     plt.plot(ws, Es, "o--", label=f"g={g}")
286 plt.yscale("log")
287 plt.xlabel("collar_width_w_(BFS_from_A),_with_B(0)=empty")
288 plt.ylabel("E_rec^Petz(w)_=-log_F")
289 plt.title(f"Z2_LGT_{Nx}x{Ny}:_Petz_recoverability_(plaque_patches)")
290 plt.grid(True, which="both", ls="--", alpha=0.3)
291 plt.legend()
292 plt.tight_layout()
293 plt.savefig(f"{out_prefix}_recoverability_v2.png", dpi=200)
294
295 return rows
296
297 def main():
298     g_list = [0.5, 1.0, 2.0]
299     Lambda = 50.0
300     delta = 1e-12
301     w_list = [0,1,2,3]
302     max_abc_qubits = 13
303
304     rows = []
305     rows += run_lattice(2,2,g_list,Lambda,delta,w_list,max_abc_qubits,"z2_2x2")
306     rows += run_lattice(2,3,g_list,Lambda,delta,w_list,max_abc_qubits,"z2_2x3")
307
308     header = ["Nx", "Ny", "g", "E0", "min_Gv", "mean_Gv", "w", "ABC_qubits", "E_recP", "F", "
309               tr_sigma", "W11_avg", "sigma_eff", "inv_sigma_eff"]
310     with open("paper_d_results.csv", "w", newline="", encoding="utf-8") as f:
311         wr = csv.writer(f); wr.writerow(header); wr.writerows(rows)
312
313 if __name__ == "__main__":
314     main()

```

References

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