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Relation between Markovian master equations & CPTP maps

Closed system dynamics described by Schrödinger equation

$$\frac{d}{dt} \rho(t) = \frac{1}{i} [H, \rho(t)] = \mathcal{L}_H(\rho)$$

↑
Liouvillian

Open system dynamics (system coupled to environment) is generally non-local in time:

$$\frac{d}{dt} \rho(t) = \int d\tau \mathcal{L}[\rho(t)]$$

This is due to the memory of the bath. Memory is kept over timescale of bath correlation time τ_{corr} .

If τ_{corr} is much shorter than typical system evolution

time t_{sys} (given by energy distances of occupied states):

$\tau_{\text{corr}} \ll t_{\text{sys}}$, then a time-local (\equiv Markovian)

description of open system dynamics is approximately valid:

$$\frac{d}{dt} \rho(t) = \mathcal{L}_\mathcal{E}[\rho(t)] \leftarrow \text{Liouvillian of CPTP map } \mathcal{E}$$

Evolution over infinitesimal time interval dt is described by CPTP map ($\hat{=}$ quantum channel) \mathcal{E} :

$$\begin{aligned}\rho(t+dt) &= \mathcal{E}_{dt}[\rho(t)] = \sum_a M_a \rho(t) M_a^\dagger \\ &= \rho(t) + dt \mathcal{L}_\mathcal{E}[\rho(t)]\end{aligned}$$

Relation b/w Lindbladian $\mathcal{L}_\mathcal{E}$ and CPTP map \mathcal{E} thus

$$\mathcal{E}_{dt} = \mathbb{I} + \mathcal{L}_\mathcal{E} dt$$



after each infinitesimal time evolution of joint system+environment, the environmental state is discarded (Tr_E) and replaced by a fresh (equilibrium) state of the bath unentangled with the system.

Expanding \mathcal{E}_{dt} in Kraus operators

$$\sum_a M_a \rho(t) M_a^\dagger = (\mathbb{I} + dt \mathcal{L})[\rho(t)]$$

Without loss of generality, we can choose

$$M_0 = \mathbb{I} + O(dt) = \mathbb{I} + (-i \overset{\text{hermitian operators}}{H} + K) dt$$

$$M_a = \sqrt{dt} L_a \quad \text{for } a \neq 0.$$

Operators H, K, L_a are all zeroth order in dt .

We can determine K from the completeness relation

$$\mathbb{I} = \sum_a M_a^\dagger M_a$$

$$= [\mathbb{I} + (iH + K)dt][\mathbb{I} + (-iH + K)dt]$$

$$+ dt \sum_a L_a^\dagger L_a =$$

$$= \mathbb{I} + dt \underbrace{\left[2K + \sum_{a>0} L_a^\dagger L_a \right]}_{=0} + O(dt)^2$$

$$\Rightarrow K = -\frac{1}{2} \sum_{a>0} L_a^\dagger L_a$$

We thus obtain the Lindblad master equation
 (most general Markovian ME describing CPTP evolution):

$$\begin{aligned} \frac{d}{dt} \rho(t) &= \frac{\rho(t+dt) - \rho(t)}{dt} = \frac{1}{dt} \left[\mathcal{E}_{dt}[\rho(t)] - \rho(t) \right] = \\ &= \frac{1}{dt} \left[M_0 \rho M_0^\dagger + \sum_{a>0} M_a \rho M_a^\dagger - \rho(t) \right] = \\ &= \frac{1}{dt} \left\{ \left(\mathbb{I} + (-iH + K)dt \right) \rho \left(\mathbb{I} + (iH + K)dt \right) \right. \\ &\quad \left. + dt \sum_{a>0} L_a \rho L_a^\dagger - \rho(t) \right\} = \end{aligned}$$

$$\Rightarrow \dot{\rho}(t) = -i[H, \rho] + \sum_{a>0} \left(L_a \rho L_a^\dagger - \frac{1}{2} L_a^\dagger L_a \rho - \frac{1}{2} \rho L_a^\dagger L_a \right)$$

Lindblad master equation

Similarly to the Kraus operators, there exist a unitary freedom in choice of Lindblad operators $\tilde{L}_a = \sum_b V_{ba} L_a$ unitary

Different choices of Lindblad "jump" operators correspond to different unravelings of the same Markovian dynamics.

Probability for jump event a to occur during dt :

$$P_{\text{rel}}(a) = dt \langle \psi(t) | L_a^\dagger L_a | \psi(t) \rangle$$

\Rightarrow if jump occurs the updated wavefunction is

$$|\psi(t+dt)\rangle = \frac{L_a |\psi(t)\rangle}{\|L_a |\psi(t)\rangle\|}$$

if no jump occurs the state evolves as

$$|\psi(t+dt)\rangle = \frac{M_0 |\psi(t)\rangle}{\|M_0 |\psi(t)\rangle\|} .$$

Stochastic Schrödinger equation evolution \rightarrow always on many different paths (sampling possible jump event paths).

Quantum state tomography (QST)

Procedure of experimentally determining an unknown quantum state.

Task: Measuring the density operator ρ given many copies of the state.

Reconstruct ρ from measuring its components in a basis of measurement operators E_j , $j = 1, \dots, d^2$ ($d = 2^n$) that span the n -qubit Hilbert space $L(\mathcal{H}_d)$.

Typical choice is the Pauli basis $|E_j\rangle = |j\rangle_{\mathcal{P}}$, where $|j\rangle_{\mathcal{P}} \hat{=}$ vectorized element $P_j \in \{I, X, Y, Z\}^{\otimes n}$.

Generally, one measures the probabilities

Coefficient of ρ in Pauli basis P_j .

$$P_j = \langle\langle E_j | \rho \rangle\rangle = \sum_{\mathcal{R}} \underbrace{\langle\langle E_j | \mathcal{R} \rangle\rangle}_{\text{matrix } A_{j,\mathcal{R}} \text{ is known}} \underbrace{\langle\langle \mathcal{R} | \rho \rangle\rangle}_{\text{measured}}$$

(choice of experiment) and must be invertible.

Can estimate p_j as sample average of measurement outcomes:

$$m_j = \frac{1}{N} \sum_{i=1}^N m_{ij}$$

$m_{ij} \in \{0, 1\}$ is outcome of i -th measurement of E_j .

average over shots (measurements of N copies of S)

Expected value of m_j is

$$\begin{aligned} E(m_j) &= \frac{1}{N} \sum_{i=1}^N E(m_{ij}) = p_j \\ &= p_0 \cdot 0 + p_1 \cdot 1 = p_1 \end{aligned}$$

$$E(X) := \sum_i p_i x_i$$

probabilities random variables

Variance:

$$\text{Var}(m_j) = E[(m_j - E(m_j))^2] =$$

$$= E(m_j^2) - 2E(m_j)^2 + E(m_j)^2 =$$

$$= E(m_j) - 2p_j^2 + p_j^2 = p_j - p_j^2 = p_j(1 - p_j)$$

$$E(m_j^2) = p_0 \cdot 0^2 + p_1 \cdot 1^2 = p_1 = E(m_j)$$

The variance (or error) of the mean:

$$\text{Var}(\bar{X}) = \text{Var}\left(\frac{1}{N} \sum_i X_i\right) = \frac{1}{N^2} \text{Var}\left(\sum_i X_i\right) = \frac{1}{N^2} \sum_i \text{Var}(X_i)$$

general property of the variance

unrelated variables, i.e. uncorrelated (IID) measurements

$$\Rightarrow \text{Var}(\bar{X}) = \frac{\text{Var}(X)}{N} \quad (\text{if all } \text{Var}(X_i) \text{ are the same}).$$

Here:

$$\text{Var}(\bar{m}_j) = \frac{p_j(1-p_j)}{N}$$

Numerator $p_j(1-p_j)$ is also called single-shot variance.

The sample average m_j thus approaches the true probability value p_j as the number of measurements $N \rightarrow \infty$.

Then, \hat{S} is linear inversion estimator of S vector of measurements m_j

$$|\hat{S}\rangle\rangle = A^{-1} |M\rangle\rangle, \quad A_{jg} = \langle\langle E_j | P_g \rangle\rangle$$
$$\Rightarrow \hat{S}_j = \sum_g (A^{-1})_{jg} m_g$$

Note that we can extract expectation value of $P_0 \equiv I$ and $P_j = z$ in same basis measurement $\Rightarrow 3^m - 1$ measurements are necessary to estimate m_j , $P_j \in \{I, x, y, z\}^{\otimes m}$.
One coefficient is fixed by normalization of the state.

If A is not square since we are using an over-complete measurement basis $|E_j\rangle$ with $j > d^2$, then we obtain \hat{g} as pseudo-inverse (minimizes least-squares)

$$|\hat{g}\rangle\rangle = (A^T A)^{-1} A^T |m\rangle\rangle$$

Quantum process tomography:

Goal: Characterize gate G (instead of state ρ).

Done by measuring process matrix \mathcal{K} (e.g. in Pauli or one-hot $|i\rangle\langle j|$ basis) or PTM $(R_E)_{kl} = \langle\langle l | G | k \rangle\rangle_{\mathcal{E}}$.

This is achieved by measuring the d^4 probabilities

$$P_{ji} = \langle\langle E_j | G | \rho_i \rangle\rangle$$

gate that we want to estimate \mathcal{K} or PTM R_E for.

Basis of measurement operators (fiducial basis, informationally-complete)

the probabilities $P_i = \text{Tr}(\rho E_i)$ completely characterize the state.

One can distinguish measurement outcomes j from POVMs m and label $E_j^{(m)}$, e.g. $m = \{X, Y, Z\}$

basis measurements with outcomes $i = 0, 1$. Then E_i contains all of the $E_j^{(m)}$.

set of vectors $|\rho_i\rangle\rangle$ forms a complete basis for the Hilbert-Schmidt space.

Note that we cannot choose $|\rho_i\rangle\rangle$ as $|i\rangle\rangle_{\mathcal{E}}$ as $\text{Tr}(|\rho_i\rangle\rangle) = 1$, but $\text{Tr}(|i\rangle\rangle_{\mathcal{E}}) = 0$.

Instead, each $|\rho_i\rangle\rangle$ must be chosen as some linear combination of Pauli basis vectors $|\rho_i\rangle\rangle = \sum_j C_{ij} |j\rangle\rangle_{\mathcal{E}}$.

Inserting two complete sets of Pauli basis states

$$P_{ji} = \langle\langle E_j | G | \mathcal{S}_i \rangle\rangle$$

$$\Leftrightarrow P_{ji} = \sum_{\mathcal{R}, \mathcal{L}} \underbrace{\langle\langle E_j | \mathcal{R} \rangle\rangle}_{\substack{\text{known projection} \\ \text{of measurement effects} \\ \langle\langle E_j | \text{ on Pauli basis} \rangle\rangle}} \underbrace{\langle\langle \mathcal{R} | G | \mathcal{L} \rangle\rangle}_{= \text{PTM}} \underbrace{\langle\langle \mathcal{L} | \mathcal{S}_i \rangle\rangle}_{\substack{\text{known coefficients of} \\ |\mathcal{S}_i\rangle \text{ in Pauli basis}}}$$

Now vectorize the PTM to write this as a matrix equation:

$$\mathbf{g}_{\mathcal{R} + (l-1)d^2} = \langle\langle \mathcal{R} | G | \mathcal{L} \rangle\rangle \quad (\text{column stacking})$$

$$\mathbf{S}_{\mathcal{J} + (i-1)d^2, \mathcal{R} + (l-1)d^2} = \langle\langle E_j | \mathcal{R} \rangle\rangle \langle\langle \mathcal{L} | \mathcal{S}_i \rangle\rangle$$

$$P_{\mathcal{J} + (i-1)d^2} = P_{\mathcal{S}_i}$$

$$\Rightarrow \vec{p} = S \vec{g}$$

If S is a square matrix (it has full rank by definition), then

$$\vec{g} = S^{-1} \vec{p}$$

Here, \vec{p} = vector of $\hat{p}_i = \frac{1}{N} \sum_s m_{si}$ (estimates of p_i).

If the sets of measurement operators $\{E_j\}$ or input states $\{\rho_i\}$ are overcomplete, the S matrix has more rows than columns, $i, j > d^2$. Then we obtain \vec{g} by the pseudo-inverse of S (this corresponds to the least-squares solution of $S\vec{g} = \vec{p}$):

$$\vec{g} = (S^T S)^{-1} S^T \vec{p}$$

Note, without vectorization, we can write

$$P_{ji} = \sum_{r,l} \underbrace{A_{jr}}_{\langle\langle E_j | \rho \rangle\rangle} G_{rl} \underbrace{B_{li}}_{\langle\langle \rho | S_i \rangle\rangle}$$

If the inverses of A and B exist, then

$$G = A^{-1} P B^{-1}$$

If not, we can use the pseudo-inverses and write

$$G = (A^T A)^{-1} A^T P B^T (B B^T)^{-1}$$

Example: Quantum process tomography of a single qubit

$$\rho = \frac{1}{2}(\mathbb{I} + \vec{r} \cdot \vec{\sigma})$$

$$\mathcal{E}(\rho) = \sum_a M_a \rho M_a^\dagger, \quad \sum_a M_a^\dagger M_a = \mathbb{I} \quad \begin{array}{l} |i\rangle\langle c \\ \parallel \\ |i\rangle\langle c \end{array}$$

We can write $M_a = \sum_r C_{ar} P_r = \sum_r \tilde{C}_{ar} E_r$

$$\Rightarrow M_a^\dagger = \sum_r C_{ar}^* P_r = \sum_r \tilde{C}_{ar}^* E_r \quad \begin{array}{l} |i\rangle\langle c \\ \parallel \\ |i\rangle\langle c \end{array} \quad \begin{array}{l} E_r = E_{i+(j-1)d^2} \\ = |i\rangle\langle j| \\ = |i,j\rangle \end{array}$$

Thus,

$$\mathcal{E}(\rho) = \sum_a \sum_{r,l} C_{ar} C_{al}^* P_r \rho P_l$$

$$= \sum_{r,l} \chi_{rl} P_r \rho P_l^\dagger \equiv \rho'$$

χ_{rl} process matrix in Pauli basis

with process matrix $\chi_{rl} = \sum_a C_{ar} C_{al}^*$.

Can use any basis, $P_r = |i\rangle\langle j|$, $E_r = |i\rangle\langle c|$, or any other.

Now expand $\rho' = E(\rho)$ in any basis & do state tomography.

Choose $|g_i\rangle$ as follows:

$$g_1 = |0\rangle\langle 0|$$

$$g_2 = |1\rangle\langle 1|$$

$$g_3 = |+\rangle\langle +| \quad \text{with } |+\rangle = \frac{1}{\sqrt{2}}(|0\rangle + |1\rangle)$$

$$g_4 = |-\rangle\langle -| \quad \text{with } |-\rangle = \frac{1}{\sqrt{2}}(|0\rangle + i|1\rangle)$$

Then,

$$g_3 + i g_4 - \frac{1+i}{2} g_1 - \frac{1+i}{2} g_2 =$$

$$= |+\rangle\langle +| + i |-\rangle\langle -| - \frac{1+i}{2} |0\rangle\langle 0| - \frac{1+i}{2} |1\rangle\langle 1| =$$

$$= \frac{1}{2} [\underline{|0\rangle\langle 0|} + \underline{|0\rangle\langle 1|} + \underline{|1\rangle\langle 0|} + \underline{|1\rangle\langle 1|}]$$

$$+ \frac{i}{2} [\underline{|0\rangle\langle 0|} - i \underline{|0\rangle\langle 1|} + i \underline{|1\rangle\langle 0|} + \underline{|1\rangle\langle 1|}]$$

$$- \frac{1+i}{2} [\underline{|0\rangle\langle 0|} + \underline{|1\rangle\langle 1|}] =$$

$$= |0\rangle\langle 1| \quad (|1\rangle\langle 0|).$$

By preparing the states $|0\rangle, |1\rangle, |+\rangle, |-\rangle$, we can thus determine

$$\begin{aligned} \mathcal{E}(|0\rangle\langle 1|) &= \mathcal{E}(|+\rangle\langle +|) + i\mathcal{E}(|-\rangle\langle -|) \\ &\quad - \frac{1+i}{2} [\mathcal{E}(|0\rangle\langle 0|) + \mathcal{E}(|1\rangle\langle 1|)]. \end{aligned}$$

Of course, $\mathcal{E}(|1\rangle\langle 0|) = \mathcal{E}^*(|0\rangle\langle 1|)$ and the diagonal elements are obtained from preparing $|0\rangle$ and $|1\rangle$.

We thus have used the input states \mathcal{S}_i to determine

$$\text{and can express } |i\rangle_c = \sum_j C_{ij} \mathcal{S}_j$$

column stacked basis $\begin{pmatrix} 1 \\ 0 \\ 0 \\ 0 \end{pmatrix}, \begin{pmatrix} 0 \\ 1 \\ 0 \\ 0 \end{pmatrix}, \begin{pmatrix} 0 \\ 0 \\ 1 \\ 0 \end{pmatrix}, \begin{pmatrix} 0 \\ 0 \\ 0 \\ 1 \end{pmatrix}$.

with coefficients $\langle\langle j | \mathcal{S}_i \rangle\rangle = C_{ij}$

$$C_{1j} = (1, 0, 0, 0)$$

$$C_{2j} = \left(-\frac{1-i}{2}, -\frac{1-i}{2}, 1, -i\right)$$

$$C_{3j} = \left(-\frac{1+i}{2}, -\frac{1+i}{2}, 1, i\right)$$

$$C_{4j} = (0, 1, 0, 0).$$

We can also easily express the Pauli basis in terms of the S_i i.e., $\langle\langle j | S_i \rangle\rangle$ or $|j\rangle\rangle_{S_i} = \sum_i C_{ji} S_i$.

$$P_0 = I = \begin{pmatrix} 1 & 0 \\ 0 & 1 \end{pmatrix} = S_1 + S_2$$

$$P_1 = X = \left(S_3 + i S_4 - \frac{1+i}{2} S_1 - \frac{1+i}{2} S_2 \right) + c.c. =$$

$$= 2 S_3 - S_1 - S_2 =$$

$$P_2 = Y = -i \left(S_3 + i S_4 - \frac{1+i}{2} S_1 - \frac{1+i}{2} S_2 \right) + c.c.$$

$$= 2 S_4 - S_1 - S_2$$

$$P_3 = Z = S_1 - S_2$$

Then measure these states $E(S_i)$ in the Pauli basis P_j to experimentally obtain P_{ji} :

$$P_{ji} = \text{Tr} [P_j E(S_i)] = A G B$$

Alternatives.

$$|s_i\rangle = \sum_{\mu} B_{\mu i} |\mu\rangle$$

$$P_{ji} = \sum_{\mu, \nu} A_{j\mu} G_{\mu\nu} B_{\nu i}$$

\parallel
 $\langle E_j | \mu \rangle$

$\langle \mu | s_i \rangle \rightarrow$ from above, we can read off:

$$S_1 = \frac{1}{2}(I+Z); S_2 = \frac{1}{2}(I-Z)$$

$$S_3 = \frac{1}{2}(I+X); S_4 = \frac{1}{2}(I+Y)$$

$$\Rightarrow A_{j\mu} = \delta_{j\mu}$$

$$B_{\ell i} = \begin{pmatrix} \frac{1}{\sqrt{2}} & \frac{1}{\sqrt{2}} & \frac{1}{\sqrt{2}} & \frac{1}{\sqrt{2}} \\ 0 & 0 & \frac{1}{\sqrt{2}} & 0 \\ \frac{1}{\sqrt{2}} & -\frac{1}{\sqrt{2}} & 0 & 0 \\ 0 & 0 & 0 & \frac{1}{\sqrt{2}} \end{pmatrix}$$

Since

$$P_{\mu} \in \left\{ \frac{I}{\sqrt{2}}, \frac{X}{\sqrt{2}}, \frac{Y}{\sqrt{2}}, \frac{Z}{\sqrt{2}} \right\}^{\otimes n}$$

$$\Rightarrow S_1 = \frac{1}{\sqrt{2}}(P_0 + P_3)$$

$$S_2 = \frac{1}{\sqrt{2}}(P_0 - P_3)$$

$$S_3 = \frac{1}{\sqrt{2}}(P_0 + P_1)$$

$$S_4 = \frac{1}{\sqrt{2}}(P_0 + P_2)$$

$$\Rightarrow B^{-1} = \begin{pmatrix} \frac{1}{\sqrt{2}} & -\frac{1}{\sqrt{2}} & -\frac{1}{\sqrt{2}} & \frac{1}{\sqrt{2}} \\ \frac{1}{\sqrt{2}} & \frac{1}{\sqrt{2}} & -\frac{1}{\sqrt{2}} & -\frac{1}{\sqrt{2}} \\ 0 & \frac{1}{\sqrt{2}} & 0 & 0 \\ 0 & 0 & \frac{1}{\sqrt{2}} & 0 \end{pmatrix}$$

$$\Rightarrow G = A^{-1} P B^{-1} = P B^{-1}$$

$$\uparrow$$
$$A^{-1} = I$$

$$\text{where } P = P_{ji} = \text{Tr} [P_j E(s_i)].$$

Example: Depolarizing noise channel with $\rho \equiv \lambda = 0.1$.

$$\mathcal{E}(\mathcal{B}_1) = \mathcal{E}(|0\rangle\langle 0|) = \frac{1}{2}(\mathbb{I} + \vec{\tau}' \cdot \vec{\sigma})$$

$$\text{with } \vec{\tau}' = \begin{pmatrix} 0 \\ 0 \\ 1 - \frac{4}{3}\lambda \end{pmatrix} = \frac{1}{2}(\mathbb{I} + (1 - \frac{4}{3}\lambda)\mathbb{Z})$$

$$\Rightarrow P_{j1} = \left(1, 0, 0, 1 - \frac{4}{3}\lambda \right)^T \quad (j=0, \dots, 3)$$

$$\mathcal{E}(\mathcal{B}_2) = \mathcal{E}(|1\rangle\langle 1|) = \frac{1}{2}(\mathbb{I} + \vec{\tau}' \cdot \vec{\sigma})$$

$$\text{with } \vec{\tau}' = \begin{pmatrix} 0 \\ 0 \\ -(1 - \frac{4}{3}\lambda) \end{pmatrix}$$

$$\Rightarrow P_{j2} = \left(1, 0, 0, -(1 - \frac{4}{3}\lambda) \right)^T$$

$$\mathcal{E}(\mathcal{B}_3) = \mathcal{E}(|+\rangle\langle +|) \stackrel{\vec{\tau} = (1, 0, 0)}{=} \frac{1}{2}(\mathbb{I} + \vec{\tau}' \cdot \vec{\sigma})$$

$$\text{with } \vec{\tau}' = \begin{pmatrix} 1 - \frac{4}{3}\lambda \\ 0 \\ 0 \end{pmatrix}$$

$$\Rightarrow P_{j3} = \left(1, 1 - \frac{4}{3}\lambda, 0, 0 \right)^T$$

$$\Sigma(\beta_4) = \Sigma(1-\lambda < -1) \stackrel{\text{GL}}{=} \frac{1}{2}(\mathbb{I} + \tilde{F}' \cdot \tilde{F})$$

$$\text{with } \tilde{F}' = \begin{pmatrix} 0 \\ 1 - \frac{4}{3}\lambda \\ 0 \end{pmatrix}$$

$$\Rightarrow P_{j4} = (1, 0, 1 - \frac{4}{3}\lambda, 0)^T$$

$$\Rightarrow P = (P_{j4}) = \begin{pmatrix} 1 & 1 & 1 & 1 \\ 0 & 0 & 1 - \frac{4}{3}\lambda & 0 \\ 0 & 0 & 0 & 1 - \frac{4}{3}\lambda \\ 1 - \frac{4}{3}\lambda & -1 + \frac{4}{3}\lambda & 0 & 0 \end{pmatrix} \frac{1}{\sqrt{2}}$$



due to normalization

$$P_{j4} \in \left\{ \frac{\mathbb{I}}{\sqrt{2}}, \frac{x}{\sqrt{2}}, \frac{y}{\sqrt{2}}, \frac{z}{\sqrt{2}} \right\}$$

$$\Rightarrow G = P B^{-1} =$$

$$= \begin{pmatrix} 1 & 0 & 0 & 0 \\ 0 & 1 - \frac{4}{3}\lambda & 0 & 0 \\ 0 & 0 & 1 - \frac{4}{3}\lambda & 0 \\ 0 & 0 & 0 & 1 - \frac{4}{3}\lambda \end{pmatrix}$$

as expected.

Randomized benchmarking

Model-free benchmarking technique that yields the average fidelity F_{av} over the full set of Clifford gates Cliff_m . The Clifford group is the normalizer of the Pauli group \mathcal{P} , i.e., it maps Pauli operators $P_m \in \mathcal{P}$ to Pauli operators: $g P_m g^{-1} = P_m \quad \forall g \in \text{Cliff}_m$.

The Clifford group is generated by $\{H, S, \text{CNOT}\}$ and circuits composed only out of Clifford gates can be efficiently simulated classically (Gottesman-Knill theorem). Clifford group plays important role in QEC using stabilizer codes and forms a unitary 2-design, which implies

$$\frac{1}{|\text{Cliff}_m|} \sum_{j=1}^{|\text{Cliff}_m|} (C_j \wedge (C_j^\dagger \rho C_j) C_j^\dagger) =$$

w/o proof; we show a similar statement explicitly later for Pauli twirls

uniform Haar measure of $U(d)$

$$= \int_{U(d)} dU (U \wedge (U^\dagger \rho U) U^\dagger) = \Lambda_{\text{dep}}(\rho)$$

Λ_{dep} is depolarizing channel with same average fidelity

F_{ave} as Λ . Now, $\Lambda_{\text{dep}} = (1-\tilde{p})\rho + \tilde{p}\frac{I}{d}$ and thus

$$F_{\text{ave}} = 1 - \tilde{p} + \frac{\tilde{p}}{d} \quad (\text{shown below}).$$

RB is scalable in the number of qubits n ($O(n^2)$ gates and $O(n^4)$ classical preprocessing costs). However, RB only provides partial information about the noise (= error channel), specifically the average fidelity F_{ave} . This is related to the X_{00} element of the χ process matrix as

$$F_{\text{ave}} = \frac{X_{00} d + 1}{d + 1}.$$

RB is unaffected by SPAM errors (unlike quantum process tomography), where SPAM refers to state preparation & measurement errors. RB in fact gives estimates for SPAM errors.

Fidelity of depolarizing channel:

$$\begin{aligned} E(\rho) &= (1-p)\rho + \frac{p}{3}(X\rho X + Y\rho Y + Z\rho Z) = \\ &= \left(1 - \frac{3}{4}\tilde{p}\right)\rho + \frac{\tilde{p}}{4}(X\rho X + Y\rho Y + Z\rho Z) = \\ &= (1-\tilde{p})\rho + \tilde{p}\frac{\mathbb{I}}{d} \quad (d=2 \text{ here,} \\ &\quad \text{generally } 2^n). \end{aligned}$$

$$\text{Here, } \frac{3}{4}\tilde{p} = p \Rightarrow \tilde{p} = \frac{4}{3}p.$$

The average fidelity of the depolarizing channel is defined as

$$\overline{F}_{\text{av}}(E) = \int d\psi \underset{\substack{\uparrow \\ \text{Haar measure}}}{F(|\psi\rangle, E(|\psi\rangle\langle\psi|))}$$

with state-dependent channel

$$F(|\psi\rangle, E(|\psi\rangle\langle\psi|)) = \langle\psi| E(|\psi\rangle\langle\psi|) |\psi\rangle.$$

Inserting the depolarizing channel yields (use def. w/o square root here)

$$\langle\psi| \left[(1-\tilde{p})\underbrace{\rho}_{=|\psi\rangle\langle\psi|} + \tilde{p}\frac{\mathbb{I}}{d} \right] |\psi\rangle = 1 - \tilde{p} + \underbrace{\frac{\tilde{p}}{d}}_{=\tilde{p}/2} = F(|\psi\rangle, E(|\psi\rangle\langle\psi|))$$

RB protocol:

Generate random sequences of Clifford gates $C_i \in \text{Cliff}_m$ of length $m \leq M-1$. Fix initial state $|\psi\rangle$ and maximal sequence length M .

Step 1: fix $m \leq M-1$ and generate K_m sequences consisting of $(m+1)$ quantum operations.

The first m are chosen uniformly random from Cliff_m .

The $(m+1)$ th operation is chosen such that (possible as Cliff_m is a group)

$$C_{i_{m+1}} \circ [C_{i_m} \circ \dots \circ C_{i_1}] = \mathbb{I}$$

Each Clifford gate is associated with an error channel

$\Lambda_{i_j, j}$ such that the sequence K_m corresponds to the quantum operation

$$S_{\vec{i}_m} = \bigcirc_{j=1}^{m+1} (\Lambda_{i_j, j} \circ C_{i_j}), \quad \vec{i}_m = (i_1, \dots, i_m)$$

Note that i_{m+1} is uniquely determined by \vec{i}_m .

Step 2: For each of K_m sequences, measure the survival probability

$$\text{Tr}[E_\psi S_{\vec{i}_m}(\rho_\psi)]$$

$\rho_\psi = \text{noisy version of } |\psi\rangle\langle\psi| \text{ (state prep. error)}$ } SPAM
 $E_\psi = \text{noisy measurement of } |\psi\rangle\langle\psi| \text{ (measurement error)}$ } error

Step 3: Average over K_m random realizations to find the averaged sequence fidelity

$$F_{\text{seq}}(m, \psi) = \text{Tr}[E_\psi S_{K_m}(\rho_\psi)]$$

where

$$S_{K_m} = \frac{1}{K_m} \sum_{\vec{i}_m} S_{\vec{i}_m}$$

Step 4: Repeat steps 1 through 3 for different values of m and fit results to the model

$$F_g(m, \psi) = A_0 (1-\rho)^m + B_0$$

A_0 : state preparation error
 B_0 : measurement error
 } SPAM error

The average error rate $\tau = 1 - F_{\text{ave}} =$

$$= 1 - \left[1 - \tilde{\rho} + \frac{\tilde{\rho}}{d} \right] =$$

$$= \tilde{\rho} - \frac{\tilde{\rho}}{d}. \quad (d = 2^n \text{ for } n \text{ qubits})$$

Note that it becomes obvious that we measure the twisted channel on the Clifford group, i.e., a depolarizing channel with the same average fidelity as the original channel.

For this define: (1) $D_{i_1} = C_{i_1}$
 (2) $D_{i_2} = C_{i_2} \circ C_{i_1} \Rightarrow C_{i_2} = D_{i_2} \circ D_{i_1}^\dagger$

(3) given if C_{i_1}, \dots, C_{i_j} have been chosen and D_{i_1}, \dots, D_{i_j} have been defined accordingly, define $D_{i_{j+1}}$ such that $C_{i_{j+1}} = D_{i_{j+1}} \circ D_{i_j}$ i.e.,

$$D_{i_{j+1}} = C_{i_j} \circ \dots \circ C_{i_1}.$$

Then,

$$\begin{aligned} S_{\vec{i}_m} &= \Lambda_{i_{m+1}} \circ C_{i_{m+1}} \circ \Lambda_{i_m} \circ C_{i_m} \circ \dots \circ \Lambda_{i_1} \circ C_{i_1} \\ &= \Lambda_{i_{m+1}} \circ D_{i_{m+1}} \circ D_{i_m}^+ \circ \Lambda_{i_m} \circ D_{i_m} \circ D_{i_{m-1}}^+ \circ \Lambda_{i_{m-1}} \circ D_{i_{m-1}} \\ &\quad \circ \dots \circ \underbrace{D_{i_1}^+ \circ \Lambda_{i_1} \circ D_{i_1}}. \end{aligned}$$

Clifford trine channel

\Rightarrow depolarizing noise channel with same average fidelity p as original channel

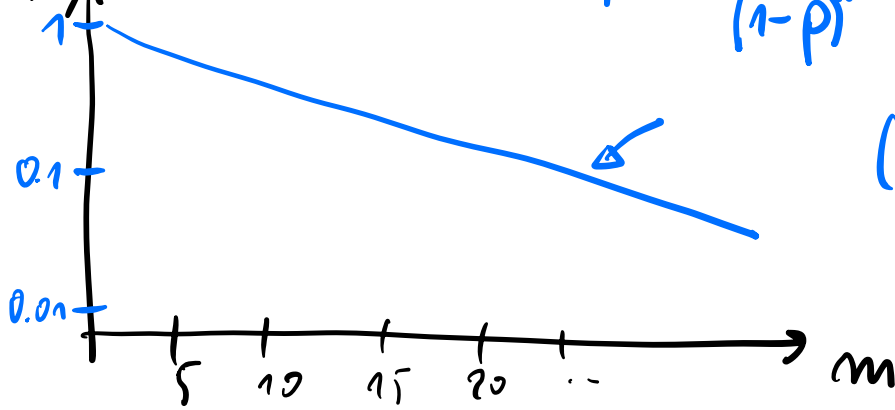
Λ_{i_j} .

$F_{\text{req}}(m, \psi) - B_0$

Semi-log plot

$$(1-p)^m = \exp[m \log(1-p)]$$

$$(1-p) \approx 1 - \epsilon.$$



Fast benchmarking rate

Ideal of interleaved RB: determine av. error of particular Clifford gate G

(1) run original RB: R_m sequences $C_{i_{m+1}} \circ C_{i_m} \circ \dots \circ C_{i_1}$

(2) for every sequence \vec{i}_m , run the circuit

$$\tilde{C}_{i_{m+1}} \circ G \circ C_{i_m} \circ \dots \circ G \circ C_{i_2} \circ G \circ C_{i_1}$$

↑

needs to be updated compared to standard RB (1)

to insert the sequence containing G .

\Rightarrow compute σ for (1) & (2) \Rightarrow difference is σ_G .