Real Randomized Benchmarking

A. K. Hashagen, S. T. Flammia, D. Gross, and J. J. Wallman

1Department of Mathematics, Technical University of Munich, Germany
2Centre for Engineered Quantum Systems, School of Physics, University of Sydney, Sydney, Australia
3Yale Quantum Institute, Yale University, New Haven, Connecticut 06520, USA
4Institute for Theoretical Physics, University of Cologne, Germany
5Institute for Quantum Computing and Department of Applied Mathematics, University of Waterloo, Canada

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Randomized benchmarking provides a tool for obtaining precise quantitative estimates of the average error rate of a physical quantum channel. Here we define real randomized benchmarking, which enables a separate determination of the average error rate in the real and complex parts of the channel. This provides more fine-grained information about average error rates with approximately the same cost as the standard protocol. The protocol requires only averaging over the real Clifford group, a subgroup of the full complex Clifford group, and makes use of the fact that it forms an orthogonal 2-design. It therefore allows benchmarking of fault-tolerant gates for an encoding which does not contain the full Clifford group transversally. Furthermore, our results are especially useful when considering quantum computations on rebits (or real encodings of complex computations), in which case the real Clifford group now plays the role of the complex Clifford group when studying stabilizer circuits.

1 Introduction

The design of reliable quantum information processing devices requires the quantitative characterization of the average error rate of a physical quantum channel. Full characterization of quantum processes is possible through quantum process tomography [29]. This method is, however, infeasible in practice. Firstly, it relies upon the challenging assumption that the set of measurements and the quantum state preparation admit lower errors than the process itself. Furthermore, the number of experimental configurations required – including quantum state preparation and quantum measurements – grows exponentially with the number of qubits even when employing improvements such as compressed sensing [17, 26].

An alternative approach is randomized benchmarking (RB) and variants thereof [8, 10, 11, 14, 15, 20, 37–39]. An RB protocol gives an estimate of the average fidelity between the realized and ideal implementations of a group of quantum gates by estimating the decay rate of the survival probability over random sequences of varying lengths. The effort of implementing the RB protocol scales efficiently with the number of qubits and it is robust against measurement and state preparation errors. Due to this, RB has become a popular tool to assess the quality of quantum processes [1, 3, 4, 9, 19, 34, 42, 45, 47, 54].

In this work, we study RB protocols in which the quantum gates are taken from the real Clifford group, which we refer to as real randomized benchmarking. We define the
notion of an orthogonal 2-design, and show that the real Clifford group constitutes one. This property allows one to efficiently estimate the average fidelity of an experimental implementation of the real Clifford group.

There are two primary motivations for using alternative groups for randomized benchmarking. First, some gates may be significantly worse than others due to different implementations (such as fault-tolerant implementations of non-transversal gates). Including such gates in the benchmarking group would result in a rapid decay dominated by the worst gate(s), so that little information can be obtained about the majority of gates. Furthermore, some quantum codes do not allow all transversal Clifford gates. The real Clifford group might, however, be accessible. This insight was recently used to do randomized benchmarking inside the code space of the [4,2,2] code using a variant of the protocol discussed in the present manuscript [28]. Second, the average gate fidelity quantifies the error rate over the entire Hilbert space. If an experiment only involves states in a portion of Hilbert space, then the relative figure of merit should only average over the states in that portion of Hilbert space. Real randomized benchmarking allows a direct characterization of the average gate fidelity over real-valued density operators, which is directly relevant to universal quantum computation with rebits [7, 46]. Third, information about which part of the Hilbert space is afflicted by the worst errors provides more information with which to optimize the experimental implementation of a group of quantum gates.

Summary. We analyze real randomized benchmarking, where the quantum gates are taken from the real Clifford group. The real Clifford group acting on \( n \)-qubits is generated by

\[
\mathcal{C}(\mathbb{R}_{2^n}) := \{Z_i, H_i, CZ_{ij}\},
\]

where the subscripts indicate that the gate is acting on the \( i \)th qubit and \( Z \) is the Pauli \( Z \)-gate, \( H \) is the Hadamard gate and \( CZ \) is the controlled \( Z \)-gate, defined respectively as,

\[
Z = \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}, \quad H = \frac{1}{\sqrt{2}} \begin{pmatrix} 1 & 1 \\ 1 & -1 \end{pmatrix}, \quad CZ = \begin{pmatrix} 1 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 \\ 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & -1 \end{pmatrix}.
\]

The protocol that gives an estimate of the average fidelity between the physical and ideal implementations of these gates, denoted as \( \tilde{C} \) and \( C \) respectively, is given in protocol 2 in section 7. We assume that the error quantum channel is gate and time independent throughout. However, we note that the methods of Wallman [50] and Merkel et al. [41] can be used to prove that the gate-dependent assumption can be relaxed with negligible effect on the estimate; we leave a careful and detailed proof of this to future work. The protocol estimates the decay rate of the survival probability over random gate sequences of varying length \( m + 1 \) as illustrated in fig. 1.

The protocol gives an approximation to the average sequence fidelity,

\[
\bar{F}(m, E, \rho) = A + b^m B + c^m C,
\]

where \( A \), \( B \) and \( C \) depend only on the state preparation and measurement, and \( b \) and \( c \) depend on the noise quantum channel. Let \( S_i : \mathcal{M}_d \rightarrow \mathcal{M}_d \), \( i = a, b, c \), be defined as

\[
S_a(\cdot) = \text{Tr}[\cdot] \frac{I}{d},
\]

\[
S_b(\cdot) = \frac{1}{2}(\text{id} + \theta) - \text{Tr}[\cdot] \frac{I}{d}, \quad \text{and}
\]

\[
S_c(\cdot) = \frac{1}{2}(\text{id} - \theta),
\]
Figure 1: The main setup of real randomized benchmarking (see Protocol 2 for more details). For a fixed \( m \in \mathbb{N} \), a sequence of \( m + 1 \) real Clifford gates is applied to an initial quantum state \( \rho \). The sequence is generated, such that in the case of its ideal implementation, it gives the identity operation. A subsequent measurement is performed given by an effect operator of a POVM, \( E \), to measure the survival probability. Averaging over \( M \in \mathbb{N} \) random realizations of sequences of length \( m \) gives the average sequence fidelity.

Then

\[
A = \text{Tr}[ES_a(\rho)],
\]

\[
B = \text{Tr}[ES_b(\rho)] \quad \text{and}
\]

\[
C = \text{Tr}[ES_c(\rho)],
\]

as well as

\[
b = \frac{\text{Tr}[\mathcal{E} \circ S_b(\rho)]}{\text{Tr}[S_b(\rho)]} \quad \text{and}
\]

\[
c = \frac{\text{Tr}[\mathcal{E} \circ S_c(\rho)]}{\text{Tr}[S_c(\rho)]},
\]

where \( \mathcal{E} : \mathcal{M}_d \to \mathcal{M}_d \) denotes the noise quantum channel.

The parameter \( b \) is linearly proportional to the average rebit fidelity, where the average is taken with respect to the orthogonal group, see eq. (35) in section 6. Together with parameter \( c \), they give the average fidelity, see eq. (34). It is thus possible to obtain more fine-grained information about the physical implementation of real Clifford gates.

**Organization of the paper.** This paper starts with an introduction to the group twirl and in particular studies the twirl over the real orthogonal group. In this particular case, the exact form of a twirled quantum channel is derived. We establish the notion of an orthogonal 2-design and, after defining the real Clifford group, in section 3 we show that the real Clifford group is an orthogonal 2-design. The derived insights into the real Clifford group are then, in the following section 4, compared to the complex case. Section 5 gives a protocol on how to obtain a Haar sample from the real Clifford group in an efficient way. In section 6 we discuss the figures of merit of which real RB obtains estimates. In section 7 we give the real RB protocol that shows how to calibrate the average sequence fidelity to experimental data. Section 8 uses the results established in section 3 to derive the estimate for the average fidelity.

2 Group twirl

In this section we introduce the mathematical background necessary for real randomized benchmarking. We will first introduce the notation used throughout this paper followed by a review of the relevant representation theoretic concepts.

We consider an \( n \)-qubit system with an underlying finite-dimensional Hilbert space \( \mathcal{H} \simeq \mathbb{C}^d, d = 2^n \). Denote by \( \mathcal{M}_d(\mathbb{C}) \) the set of complex-valued \( d \times d \)-matrices and as \( \mathcal{M}_d(\mathbb{R}) \) the set of real-valued \( d \times d \)-matrices. Every quantum state is described by a
density matrix $\rho \in \mathcal{M}_d(\mathbb{C})$, with normalization $\text{Tr} [\rho] = 1$ and positivity property $\rho \geq 0$. The set of $d$-dimensional density matrices or quantum states is denoted as $\mathcal{D}_d := \{ \rho \in \mathcal{M}_d(\mathbb{C}) | \rho \geq 0, \text{Tr} [\rho] = 1 \}$. A transformation of a quantum state is described by a quantum channel, which is a completely positive trace preserving linear map $T : \mathcal{M}_d(\mathbb{C}) \rightarrow \mathcal{M}_d(\mathbb{C})$. The Choi-Jamiołkowski representation [33] provides a one-to-one correspondence between linear maps $T : \mathcal{M}_d \rightarrow \mathcal{M}_{d'}$ and operators $\tau_T \in \mathcal{M}_{d'd}$ via

$$\tau_T = (id \otimes T) |\Omega\rangle \langle \Omega|,$$

where $|\Omega\rangle \langle \Omega| = \frac{1}{d} \sum_{i,j=1}^d |ii\rangle \langle jj|$ is the maximally entangled state. This operator $\tau_T$ encodes every property of the linear map $T$ and the representation shows that the set of quantum channels corresponds one-to-one to the set of bipartite quantum states which have one reduced density matrix maximally mixed [29]. This result will be used throughout this work. Denote by $U(d)$ the unitary group acting on $\mathbb{C}^d$ and by $O(d)$ the real orthogonal group acting on $\mathbb{C}^d$. Moreover, $I$ is the identity matrix in $\mathcal{M}_d(\mathbb{C})$.

Throughout this paper, we are interested in group actions on quantum channels. For an extensive review of representations of finite and compact groups, please refer to [48].

To this end consider any finite group $G$ with elements $g \in G$ and a unitary representation $\{ U(g) \}_{g \in G}$ on $\mathbb{C}^d$. Its adjoint representation $U_U : G \rightarrow \text{End}(\mathcal{M}_d(\mathbb{C}))$ is defined through its action on any $X \in \mathcal{M}_d(\mathbb{C})$ as

$$U_U(g)(X) = U(g)XU(g)^* \quad \forall g \in G,$$

and may be represented as a matrix $U \otimes \bar{U} \in \mathcal{M}_{d^2}$.

Indeed, let $H$ be a general matrix group acting on some Hilbert space $\mathcal{K}$ (below, we will be interested in e.g. $H = \{ U(g) \otimes \bar{U}(g) \ | \ g \in G \} = \mathcal{K} = \mathbb{C}^d \otimes \mathbb{C}^d$). The Hilbert space decomposes as

$$\mathcal{K} \simeq \bigoplus_{i=1}^k \mathcal{K}_i \otimes \mathbb{C}^{n_i},$$

where the sum is over irreducible unitary representations of $H$, $\mathcal{K}_i = \mathbb{C}^{d_i}$ carries the $i$th irreducible unitary representation, and $n_i$ is the degeneracy of the irreducible unitary representations in $\mathcal{K}$. Every $U \in H$ is block-diagonal with respect to this decomposition, i.e. of the form

$$U \simeq \bigoplus_{i=1}^k U_i \otimes \mathbb{I}_{n_i \times n_i}.$$  

The commutant $H'$ of $H$ is the algebra $H' = \{ X \ | \ [X,U] = 0 \ \forall U \in H \}$ which commutes with all elements of $H$. By Schur’s Lemma, every $X \in H'$ is of the form

$$X \simeq \bigoplus_{i=1}^k \mathbb{I}_{d_i \times d_i} \otimes X_i,$$

with $d_i = \dim \mathcal{K}_i$, and $X_i$ acting on the $n_i$-dimensional space appearing on the right hand side of eq. (4). We will mainly restrict our attention to the case where all irreducible unitary representations of $H$ on $\mathcal{K}$ are non-degenerate, i.e. $n_i = 1$ for all $i = \{1, \ldots, k\}$. In this case, eq. (5) takes the form

$$X \simeq \sum_{i=1}^k x_i P_i,$$

where the $P_i$ are orthogonal projections onto the $i$th irreducible unitary representation and the $x_i \in \mathbb{C}$. 

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The **group twirl** associated with $H$ is

$$T: A \mapsto \int_H U A U^* dU,$$

where the integration is w.r.t. the Haar measure on $H$. In particular, if $H$ is finite, the integral is the normalized sum over the group. The group twirl is (i) idempotent, (ii) self-adjoint (w.r.t. the Hilbert-Schmidt inner product), and (iii) leaves elements $X \in H'$ of the commutant invariant. It is thus the orthogonal projection onto $H'$. In the non-degenerate case, one can check that this projection is given explicitly by

$$T(A) = \sum_{i=1}^k \frac{1}{d_i} \text{Tr}(P_i A) P_i.$$

Clearly, the group twirl over $H$ only depends on the commutant $H'$. Thus, if a group $S$ is such that $S' = H'$, twirling over $S$ is equivalent to twirling over $H$. In practice, this freedom can be advantageous, if $S$ has e.g. smaller cardinality than $H$, or simpler implementations as a quantum circuit. The notion of a **group design** captures this relation: A **unitary $t$-design** is any group $G$ such that we have the equality of commutants

$$\{U^\otimes t | U \in G\}' = \{O^\otimes t | O \in \mathcal{O}(d)\}'$$

General many-qubit unitaries do not have an efficient gate decomposition, while there are many-qubit unitary 2-designs and 3-designs that do. This was the original motivation for introducing the notion [11].

Phrased this way, it is natural to generalize eq. (9) to arbitrary “reference groups”, beyond the now well-studied case of $\mathcal{U}(d)$. In particular, we will be concerned with the following case:

**Definition 1.** Let $G$ be a matrix group acting on $\mathbb{C}^d$ for some $d$. Then $G$ is an **orthogonal $t$-design** if we have the equality of commutants, i.e.,

$$\{U^\otimes t | U \in G\}' = \{O^\otimes t | O \in \mathcal{O}(d)\}'$$

where $\mathcal{O}(d)$ is the real orthogonal group acting on $\mathbb{C}^d$.

### 2.1 Group twirl over the orthogonal group

Throughout this paper, our main emphasis will be on orthogonal 2-designs. In this case, it is possible to work out the commutant easily [49]. Consider the unitary representation $O^{(2)}: g \mapsto O(g) \otimes O(g)$ of the orthogonal group $\mathcal{O}(d)$ on $\mathcal{H} = \mathbb{C}^d \otimes \mathbb{C}^d$, given as

$$G = \{O \otimes O | O \in \mathcal{O}(d)\}.$$

The commutant $G'$ is spanned by three orthogonal projections [40, 49],

$$P_0 = |\Omega\rangle\langle\Omega|,$$

$$P_1 = \frac{1}{2} (I - F) \text{ and }$$

$$P_2 = \frac{1}{2} (I + F) - |\Omega\rangle\langle\Omega|,$$

where $F = \sum_{i,j=1}^d |i,j\rangle\langle j,i|$ is the flip (or swap) operator and $|\Omega\rangle\langle\Omega| = \frac{1}{d} \sum_{i,j=1}^d |ii\rangle\langle jj|$ is the maximally entangled state. For any symmetric $X \in \mathcal{M}_d(\mathbb{C})$, we have that $FX = \ldots$
X, and for any antisymmetric X ∈ M_d(ℂ), we get FX = −X. The projections thus correspond to multiples of the identity, antisymmetric matrices and traceless symmetric matrices respectively. Every density operator in the commutant must thus be in the convex hull of the corresponding normalized density matrices \( ρ_i = P_i/d_i \) for \( i = 0, 1, 2 \).

The theory discussed above therefore applies to quantum channels too. To this end, let \( T \) be a quantum channel on a \( d \)-dimensional quantum system, and let \( G \) be a matrix group on \( ℂ^d \). The twirled channel \( \tilde{T} \) over the full real orthogonal group, \( \tilde{T}(\cdot) = \int_{O(d)} OT (O^* \cdot O) O^* dO, \) (12)
can then be expressed as in eq. (7). Using the state channel duality, we see that \( ρ_0 = P_0/d_0 \) then corresponds to the ideal channel \( T(\cdot) = \text{id} \), \( ρ_1 = P_1/d_1 \) corresponds to the Werner-Holevo channel given by

\[
T(\cdot) = \frac{\text{Tr}[\cdot] - \theta}{d - 1},
\]

where \( \theta \) denotes the usual transposition \( ρ \mapsto \theta(ρ) := ρ^T \), and \( \sum_i P_i/d^2 \) corresponds to the completely depolarizing channel \( T(\cdot) = \text{Tr}[\cdot]/d \). This yields

\[
\tilde{T}(\cdot) = \alpha \text{id} + \beta \frac{\|}{d} \text{Tr}[\cdot] + \gamma \frac{\|}{d - 1} \text{Tr}[\cdot] - \theta,
\]

(13)

with \( \alpha, \beta, \gamma \in ℍ \) satisfying \( \alpha + \beta + \gamma = 1 \) (which makes \( \tilde{T} \) trace-preserving). As before, \( \theta \) denotes the usual transposition \( ρ \mapsto \theta(ρ) := ρ^T \). This immediately follows from the correspondence between a quantum channel and its Jamiołkowski state,

\[
\int_{O(d)} (O \otimes O) \tau_T (O \otimes O)^* dO
= \int_{O(d)} (O \otimes O) (\text{id} \otimes T) |Ω⟩⟨Ω| (O \otimes O)^* dO
= \int_{O(d)} \frac{1}{d} \sum_{i,j=1}^d O |i⟩⟨j| O^* \otimes OT (|i⟩⟨j|) O^* dO
= \int_{O(d)} \frac{1}{d} \sum_{i,j=1}^d |i⟩⟨j| \otimes OT (O^* |i⟩⟨j| O) O^* dO
= \left( \text{id} \otimes \int_{O(d)} OT (O^* \cdot O) O^* dO \right) |Ω⟩⟨Ω|.
\]

A twirl over a quantum channel and the corresponding twirl over its Jamiołkowski quantum state are thus equivalent descriptions. We will, however, focus on the twirling of quantum channels throughout the rest of this paper, where we will use that the theory allows us to give an explicit representation of the image of a channel under an orthogonal twirl.

### 3 Real Clifford group

We now define the real Clifford group and show that it is an orthogonal 2-design. Following [44], we define the real Pauli group \( E(n) \) on \( n \) qubits as the \( n \)-fold tensor power

\[
E(n) := \langle E(1)^{\otimes n} \rangle,
\]

(14)
where $E(1)$ is just the real Pauli group on 1 qubit defined as

$$E(1) := \left\{ X := \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}, Z := \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix} \right\}. \tag{15}$$

$E(n)$ is thus generated by tensor products of the Pauli matrices $X$ and $Z$ with $2 \times 2$ identity matrices $I_2$.

**Definition 2** (Real Clifford group). The real Clifford group $C(n)$ is the normalizer in $O(2^n)$ of the real Pauli group $E(n)$, i.e.

$$C(n) := \{ O \in O(2^n) | OE(n) = E(n)O \}. \tag{16}$$

In the simple case when $n = 1$ the real Clifford group is generated by

$$C(1) = \left\{ Z := \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}, H := \frac{1}{\sqrt{2}} \begin{pmatrix} 1 & 1 \\ 1 & -1 \end{pmatrix} \right\}, \tag{17}$$

and in the case when $n = 2$ the real Clifford group is

$$C(2) = \left\{ C(1) \otimes C(1), CZ := \begin{pmatrix} 1 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 \\ 0 & 0 & 0 & 1 \\ 0 & 0 & 0 & -1 \end{pmatrix} \right\}, \tag{18}$$

where $H$ is the Hadamard gate and $CZ$ is the controlled $Z$-gate. See [44] for a thorough discussion of the real Clifford group.

**Theorem 3.** The representation $O(2^n) : g \mapsto O(g) \otimes O(g)$ of the real Clifford group $C(n)$ decomposes into three non-degenerate irreducible unitary representations.

**Proof.** See [44, theorem 6.8.1.] and proofs therein. \qed

We are now equipped to give the theorem that acts as the main mathematical ingredient for real randomized benchmarking.

**Theorem 4.** The real Clifford group $C(n)$ is an orthogonal 2-design.

**Proof.** The real Clifford group $C(n)$ is a subgroup of the real orthogonal group $O(2^n)$. Its commutant therefore contains the commutant of the real orthogonal group. By theorem 3 these two commutants have the same dimensions, and must thus be equal. This proves the claim. \qed

Theorem 4 will be the main ingredient for real RB. It is, however, possible to also prove the following interesting fact about the real Clifford group.

**Proposition 5.** The real Clifford group $C(n)$ is an orthogonal 3-design, but it is not an orthogonal 4-design.

**Proof.** See [44, notes below theorem 6.8.1] for the fact that the real Clifford group $C(n)$ is an orthogonal 3-design. See [43, corollary 4.13] for the fact that it is not an orthogonal 4-design, where it is shown that in this case the commutant has an additional element to it. \qed
4 Complex Clifford group

The real Clifford group shares the properties observed in the last chapter with its complex counterpart. The complex Clifford group is a unitary 2-design, a unitary 3-design, but fails to be an exact unitary 4-design [27, 31, 36, 53, 55, 56]. This, however, is not a coincidence. To this end, let us first define the complex Clifford group.

The Pauli group on one qubit $\mathcal{P}(1)$ is defined as the group generated by

$$\mathcal{P}(1) := \langle X, Z, iI \rangle,$$

where $I, X, Y, Z$ are the standard Pauli matrices given as

$$I := \begin{pmatrix} 1 & 0 \\ 0 & 1 \end{pmatrix}, \quad X := \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}, \quad Y := \begin{pmatrix} 0 & -i \\ i & 0 \end{pmatrix}, \quad Z := \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}.$$  \hfill (20)

The Pauli group on $n$ qubits is defined to be

$$\mathcal{P}(n) := \mathcal{P}(1)^\otimes n.$$  \hfill (21)

**Definition 6** (Complex Clifford group). The complex Clifford group $\mathcal{X}(n)$ is the group-theoretic normalizer in the unitary group $U(2^n)$ of the Pauli group $\mathcal{P}(n)$, i.e.

$$\mathcal{X}(n) := \{ U \in U(2^n) | UP(n) = P(n)U \}.$$  \hfill (22)

In the simple case where $n = 1$ the complex Clifford group is therefore just given as

$$\mathcal{X}(1) = \langle H, P \rangle,$$

where $H$ is the Hadamard gate defined in eq. (17) and $P$ is the $\pi/4$-phase gate given as

$$P := \begin{pmatrix} 1 & 0 \\ 0 & i \end{pmatrix}.$$  \hfill (24)

In the simple case when $n = 2$ the complex Clifford group is just given as

$$\mathcal{X}(2) = \langle \mathcal{X}(1) \otimes \mathcal{X}(1), CZ \rangle,$$

where $CZ$ is again the controlled $Z$-gate defined in eq. (18).

The real Clifford group is therefore a subgroup of the complex Clifford group, $\mathcal{C}(n) \subset \mathcal{X}(n)$, and we necessarily have that for any $t \in \mathbb{N}$,

$$\{ O^\otimes t | O \in \mathcal{C}(n) \}^\prime \supset \{ U^\otimes t | U \in \mathcal{X}(n) \}^\prime.$$

Similarly, the orthogonal group is a subgroup of the unitary group, $\mathcal{O}(2^n) \subset U(2^n)$, and we thus have that

$$\{ O^\otimes t | O \in \mathcal{O}(2^n) \}^\prime \supset \{ U^\otimes t | U \in U(2^n) \}^\prime.$$

In the special case $t = 2$ (and in fact $t = 3$), we get the following correspondence:

$$\{ O^\otimes 2 | O \in \mathcal{C}(n) \}^\prime \supset \{ U^\otimes 2 | U \in \mathcal{X}(n) \}^\prime$$

\hspace{1cm} \parallel \hspace{1cm} \parallel

$$\{ O^\otimes 2 | O \in \mathcal{O}(2^n) \}^\prime \supset \{ U^\otimes 2 | U \in U(2^n) \}^\prime$$
In [44], using the language of invariant harmonic polynomials, it was shown that any harmonic polynomial \( p \) that is invariant w.r.t. the complex Clifford group, must be invariant w.r.t. any of its subgroups, including the real Clifford group. If we decompose the harmonic polynomial into its real and imaginary parts, then the restrictions \( p_{\text{Re}(A)} \) and \( p_{\text{Im}(A)} \) must be invariant harmonic polynomials of the real Clifford group. Unfortunately, the resulting real polynomials may turn out to be zero. It is therefore not possible to infer the absence of harmonic invariants of \( \mathcal{X}(n) \) from the absence of real harmonic invariants of \( \mathcal{C}(n) \). However, it explains the observation regarding the real and complex Clifford group and their \( t \)-design properties for \( t = 2, 3 \).

If some family of matrix groups \( G \) acting on \( \mathbb{C}^d \) fails to be a unitary \( t \)-design, it might still turn out to be useful for RB. If its commutant, \( \{U^\otimes t | U \in G\}' \), has \( l \) additional elements to it than the commutant of the unitary group, \( \{U^\otimes t | U \in \mathcal{U}(d)\}' \), and there are only \( l = O(1) \) additional elements to the commutant, then we may term the family of matrix groups \( G \) an algebraic almost \( t \)-design.

An example of an algebraic almost unitary \( t \)-design is the family of Clifford groups, which form an algebraic almost 4-design with one additional generator in its commutant compared to the unitary group [30, 31, 56]. As another example, the real Clifford group is an algebraic almost unitary 2-design (and an exact orthogonal 2-design, as discussed above). The dihedral-CNOT family of groups provides yet another example [10].

The \( l \) additional factors in the commutant are useful for RB because they yield \( l \) additional decay terms in the average fidelity. A successful fit to the multi-exponential decay in a benchmarking experiment would then yield finer-grained information about the average error rate by finding the average fidelity associated to the projections onto each of the commutant algebras. In these scenarios, however, stability is an issue in the case of large \( l \) as fitting a multi-exponential decay is in general poorly conditioned [12]. For the case discussed in the most detail in this paper, the real Clifford group, there is only one extra decay term, and successful data processing methods can be employed to fit the model that we derive below with an efficient number of measurements [22, 24]. Another avenue for dealing with multi-exponential decays is to consider state preparations and measurements that optimize the contrast between various competing terms, perhaps even canceling all but one (or a constant fraction of) the exponential decay terms [8, 18]. We explore this idea in more detail in section 8.

The next section answers the question of how to obtain a Haar sample from the real Clifford group in an efficient way. This is an important ingredient for the real RB protocol.

5 Haar sample from the real Clifford group

5.1 Structure of the real Clifford group, and orthogonal transformations

Here, we summarize results and notions from [6] and [2, Chapter 7].

Consider the phase space \( V = \mathbb{F}_2^n \). Elements of phase space will often be written as \( (p, q) \in \mathbb{F}_2^n \), with \( p, q \in \mathbb{F}_2^n \). An important piece of structure for the real Clifford group [6] is the quadratic form

\[
Q((p, q)) = p \cdot q = \sum_{i=1}^{n} p_i q_i.  \tag{26}
\]

For \( x = (p, q), x' = (p', q') \), one checks that

\[
Q(x + x') - Q(x) - Q(x') = p \cdot q' - p' \cdot q = [x, x'], \tag{27}
\]
where the square brackets denote the standard symplectic form on phase space. (While over $\mathbb{F}_2$, $-1 = +1$, we occasionally use negative signs when these would appear for analogous calculations in odd characteristic). The form $Q$ turns $V$ into an orthogonal space.

A vector $x \in V$ is singular if $Q(x) = 0$. A hyperbolic pair is a set of two singular vectors $e, f \in V$ such that $[e, f] = 1$. A two-dimensional subspace is a hyperbolic plane if it is spanned by a hyperbolic pair.

A space $U \subset V$ is totally singular if $Q$ and $\cdot, \cdot$ vanish on $U$. Clearly, $U = \{(p, 0) \mid p \in \mathbb{F}_2^n\}$ is totally singular and has dimension half of $V$. This, by definition, means that $Q$ has Witt index $n$, or, equivalently sign $+1$. A $2n$-dimensional orthogonal space in characteristic two has sign $+1$ if and only if it is isometric to the orthogonal sum of $n$ hyperbolic planes:

$$V = \bigoplus_{i=1}^n \langle \{e_i, f_i\} \rangle, \quad Q(e_i) = Q(f_i) = [e_i, e_j] = [f_i, f_j] = 0, \quad [e_i, f_j] = \delta_{i,j}. \quad (28)$$

The set of linear transformations $GL(V)$ preserving such a quadratic form of positive sign (and hence, by eq. (27), the form $\cdot, \cdot$) is the group $O^+(2n, 2)$. By eq. (28), a matrix $S$ represents an element of $O^+(2n, 2)$ with respect to a hyperbolic basis $\{e_1, f_1, \ldots, e_n, f_n\}$ if and only if it is symplectic and its columns are singular – i.e. if and only if its columns form again a hyperbolic basis.

Recall that the complex Clifford group up to Pauli operators is isomorphic to the symplectic group $\mathcal{X}(n)/P(n) \simeq Sp(2n, 2)$ [35]. That is true in the sense that for each $U \in \mathcal{X}(n)$, there exists a $S \in Sp(2n, 2)$ such that

$$UP(x)U^* \propto P(Sx),$$

where for $x = (p, q) \in \mathbb{F}_2^{2n}$, we have defined the Pauli operator

$$P(x) = i^{Q(x)} \bigotimes_{i=1}^n X^{p_i} Z^{q_i}. \quad (29)$$

Any two Cliffords $U$ that differ by the left- or right-action of an element of the Pauli group induce the same $S$.

If $U/P(n)$ contains a real-valued matrix, then $S \in O^+(2n, 2)$, i.e., in addition to being symplectic, its columns are singular. Conversely, any element of the real Clifford group is associated with such an $S$, which again does change under left- or right-multiplication with elements from $E(n)$.

Thus, to sample from the real Clifford group, one can proceed by 1) drawing a random element from $O^+(2n, 2)$ (see below), 2) use one of the known constructions (e.g. [13, 32]) for generating a gate sequence that implements a given symplectic matrix as a Clifford operation, and 3) multiply with a randomly chosen element of $E(n)$.

5.2 Efficient sampling from $O^+(2n, 2)$

It remains to describe an efficient protocol for drawing an element $S$ from $O^+(2n, 2)$ uniformly at random. This will be achieved as follows:

| **Protocol 1:** Sampling($O^+(2n, 2)$) |
|---------------------------------------------------------|
| **Initialization** Choose a basis $\mathcal{B}_1$ of $\mathbb{F}_2^{2n}$. |
| **Iterate for** $i = 1$ to $n$ **Choose random linear combinations** $x$ of the vectors in $\mathcal{B}_i$ |

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until $x$ is non-zero and singular. Set $e_i = x$. Choose random linear combinations $y$ of the vectors in $B_i$ until $[e_i, y] = 1$. If $y$ is singular, set $f_i = y$. Else, set $f_i = e_i + y$. Choose a basis $B_{i+1}$ for $(\{e_j, f_j\}_{j=1}^i)^\perp$.

**Result** Return matrix $S$, with columns given by $e_1, f_1, e_2, f_2, \ldots, e_n, f_n$. 

The following statements are true for this construction:

1. By definition, the span of $B_1$ is an orthogonal space of sign $+1$ and dimension $2n$.

2. Assume $V_i = \langle B_i \rangle$ spans an orthogonal space of sign $+1$ and dimension $2(n - i + 1)$. Then, in expectation, one will find a non-zero singular $x$ after no more than 4 attempts. To see this, let $\{e'_j, f'_j\}_{j}$ be a hyperbolic basis for $V_i$. Then $x$ will be of the form

   \[ x = \sum_{j=1}^{n-i+1} (p_j e'_j + q_j f'_j), \]

   with the $p_j, q_j$ drawn uniformly at random. Hence

   \[ Q(x) = \sum_{j=1}^{n-i+1} (p_j Q(e'_j) + q_j Q(f'_j) + p_j q_j [e'_j, f'_j]) = \sum_{j=1}^{n-i+1} p_j q_j, \]

   which is zero with probability at least $\frac{1}{2}$, while at least $1 - 2^{-n} \geq \frac{1}{2}$ of these cases correspond to non-zero $x$.

3. Under the same assumption as before, a vector $y$ with $[e_i, y] = 1$ will be found after an expected number of 2 attempts. This is because $|\ker(y \mapsto [e_i, y])| = 2^\dim(V_i) - 1$ and thus exactly half of all vectors in $V$ do not lie in the kernel.

4. The vectors $\{e_i, f_i\}$ form a hyperbolic pair. Indeed, $e_i$ is singular by construction. If $y$ is singular, so is $f_i$. If $Q(y) = 1$, then

   \[ Q(f_i) = Q(e_i + y) = Q(e_i) + Q(y) + [e_i, y] = 0 + 1 + 1 = 0. \]

   Also, $[e_i, y] = [e_i, y + e_i] = 1$.

5. The basis $B_{i+1}$ of $(\{e_j, f_j\}_{j=1}^i)^\perp$ describes the solution space of a set of linear equations over $\mathbb{F}_2$, and can thus be found efficiently. By eq. (28), it spans an orthogonal space of sign $+1$ and dimension $2(n - (i + 1) + 1)$.

   Hence, by induction, the columns of $S$ form a hyperbolic basis of $\mathbb{F}_2^{2n}$, and every such basis is equally likely to arise this way. The above procedure thus samples uniformly from $O^+(2n, 2)$.

   The next chapters use the result that the real Clifford group is an orthogonal 2-design to analyze real RB using gates from the real Clifford group.

### 6 Figures of merit

The main goal of RB is to quantify how close a physical quantum channel $\tilde{C}$ is to the ideal quantum gate $C$. In order to do so, we seek a figure of merit assessing this quality in an
efficient way. We can always write the physical quantum channel as a composition of the ideal quantum channel with an error quantum channel, \( \tilde{C} = C \circ E \) [29]. We assess the quality of the physical quantum channel using the average fidelity

\[
\bar{F}(C, \tilde{C}) = \int_{U(d)} \text{Tr} \left[ C(U |0\rangle \langle 0|U^*) \tilde{C}(U |0\rangle \langle 0|U^*) \right] dU,
\]

and the average rebit fidelity

\[
\bar{F}_R(C, \tilde{C}) = \int_{O(d)} \text{Tr} \left[ C(O |0\rangle \langle 0|O^*) \tilde{C}(O |0\rangle \langle 0|O^*) \right] dO.
\]

Please notice that in the case of the average rebit fidelity, the average is taken with respect to the orthogonal group. The fidelity is thus averaged over rebits. For quantum gates \( C(\cdot) = C \cdot C^* \) with \( C \) unitary, this simplifies to

\[
\bar{F}(C, \tilde{C}) = \int_{U(d)} \text{Tr} \left[ C(U |0\rangle \langle 0|U^*) C^* C E(U |0\rangle \langle 0|U^*) C^* \right] dU
\]

\[
= \int_{U(d)} \langle 0|U^* E(U |0\rangle \langle 0|U^*) U |0\rangle dU = \bar{F}(E, \text{id}),
\]

and similarly for the average rebit fidelity to

\[
\bar{F}_R(C, \tilde{C}) = \int_{O(d)} \langle 0|O^*_E(O |0\rangle \langle 0|O^*) O |0\rangle dO = \bar{F}_R(E, \text{id}).
\]

These are the quantities that are related to the two parameters, which RB can estimate,

\[
\bar{F}(E, \text{id}) = \frac{b(d^2 + d - 2) + cd(d - 1) + 2(d + 1)}{2d(d + 1)}
\]

as well as

\[
\bar{F}_R(E, \text{id}) = \frac{b(d - 1) + 1}{d}.
\]

For real density matrices the action of the twirled channel thus reduces to the action of the depolarizing channel with parameter \( b \), and the above average gate fidelities can be interpreted as fidelities restricted to the respective commutant spaces. This makes our analysis especially interesting when considering quantum computations on rebits. It has been shown that universality holds in this case [46] and the real Clifford group now plays the role of the complex Clifford group when studying stabilizer circuits [7].

The next chapter gives the real RB protocol, which has to be executed to quantitatively describe the quality of a sequence of physical quantum gates taken from the real Clifford group.

7 Real randomized benchmarking protocol

The real randomized benchmarking protocol is given in the following. The real RB protocol considers quantum gates taken from the real Clifford group. We assume that the error quantum channel is both gate and time independent. We follow the notation of [38].

We will first describe the protocol for a given state preparation \( \rho \), sequence length \( m \), and final measurement \( E \). This defines a protocol that can be repeated many times to obtain data with those labels, \((m, E, \rho)\). Averaging these data gives a fidelity decay curve, which in expectation is a function only of these three data labels and the noise channel,
which is assumed to be gate and time independent. This forms the core of real RB. In
the subsequent section, we will show one way to process these data to obtain accurate
estimates of the parameters of the decay curve without having to fit multi-exponential
decays, which is possible but in practice quite challenging. In the following, we give the
real RB protocol:

**Protocol 2: RealRB**$(m, E, \rho)$

**Step 1** Fix a positive integer $m \in \mathbb{N}$ that varies with every loop.

**Step 2** Generate a sequence of $m + 1$ quantum gates taken from the real Clifford
group, i.e., $C_1, \ldots, C_{m+1}$, where $C_j(\cdot) = C_j \cdot C_j^* \in \mathcal{C}(n)$, for $j = 1, \ldots, m+1$. The first $m$ quantum gates, $C_1, \ldots, C_m$, are chosen independent and uniformly at random from the real Clifford group. The final quantum gate, $C_{m+1}$, is chosen from the real Clifford group, such that the net sequence (if realized without errors) is the identity operation,

$$C_{m+1} \circ C_m \circ \ldots \circ C_2 \circ C_1 = \text{id},$$

where $\circ$ represents composition. The entire sequence is therefore given by

$$S_m = \bigcirc_{j=1}^{m+1} C_j \circ E,$$

where $E$ is the associated error, a completely positive trace preserving linear map.

**Step 3** For each sequence, measure the survival probability given by the fidelity

$$F(m, E, \rho) = \text{Tr} [E S_m(\rho)],$$

where $\rho$ is the initial quantum state, taking into account preparation errors, and $E$ is an effect operator of a POVM taking into account measurement errors.

**Step 4** Repeat steps 2-3 and average over $M$ random realizations of the sequence of length $m$ to find the averaged sequence fidelity

$$\bar{F}(m, E, \rho) = \text{Tr} \left[ E \bar{S}_m(\rho) \right],$$

where

$$\bar{S}_m = \frac{1}{M} \sum_m S_m$$

is the average sequence operation.

Repeating these steps many times and averaging the results gives a good approximation
to the average sequence fidelity,

$$\bar{F}(m, E, \rho) = A + b^m B + c^m C,$$

where $A$, $B$ and $C$, given in eq. (58) below, depend only on the state preparation and measurement, and $b$ and $c$ depend on the average noise channel.

Please note that the final quantum gate $C_{m+1}$ can be found efficiently by the Gottesman-Knill theorem [21].
The question of how to choose the sequence lengths $m$ and the number of sequences at each length $M$ is still unanswered, but is addressed in [16, 23, 30, 51]. The sequence length $m$ should be exponentially spaced from 4, in order to avoid gate-dependent effects [41, 52], to around $1/(1 - F^R)$, for optimal information gain [23].

In the next section, we analyze RB focusing on the real Clifford group. The fact that it is an orthogonal 2-design will prove that it obeys the model given in eq. (41) when the assumptions of the noise model hold. Finally, we will discuss parameter estimation.

8 Fidelity decay and parameter estimation

The main setup of real RB is illustrated in fig. 1. A sequence of $m + 1$ quantum gates $\tilde{C}_j$ acts on an initial quantum state $\rho \in D_d$, followed by a measurement represented by a POVM with effect operators $E$. Consider the physical quantum channel

$$\tilde{C}_j = C_j \circ \mathcal{E},$$

where $C_j(\cdot) = C_j \cdot C_j^*$, $C_j \in \mathcal{C}(n)$, is a real Clifford gate and $\mathcal{E} : \mathcal{M}_d(\mathbb{C}) \to \mathcal{M}_d(\mathbb{C})$ is the associated error, a completely positive trace preserving linear map. We assume that the error quantum channel is both gate and time independent.

Let us first derive the decay curve of the expected data, eq. (41).

**Theorem 7.** The protocol RealRB$(m, E, \rho)$ has an expected fidelity of the form

$$\tilde{F}(m, E, \rho) = A + b^m B + c^m C,$$

where $A$, $B$ and $C$ are functions only of $(E, \rho)$, and $b$ and $c$ depend on the average noise channel.

**Proof.** The expected fidelity of the above described sequence is given by

$$F(m, E, \rho) = \text{Tr} \left[ E \left[ \tilde{C}_{m+1} \circ \tilde{C}_m \circ \ldots \circ \tilde{C}_2 \circ \tilde{C}_1 \right] (\rho) \right]$$

$$= \text{Tr} \left[ E \left[ \mathcal{C}_{m+1} \circ \mathcal{E} \circ \mathcal{C}_m \circ \mathcal{E} \circ \ldots \circ \mathcal{C}_2 \circ \mathcal{E} \circ \mathcal{C}_1 \circ \mathcal{E} \right] (\rho) \right].$$

Absorbing the first error quantum channel into the state as a preparation error gives

$$F(m, E, \rho) = \text{Tr} \left[ E \left[ \mathcal{C}_{m+1} \circ \mathcal{E} \circ \mathcal{C}_m \circ \mathcal{E} \circ \ldots \circ \mathcal{C}_2 \circ \mathcal{E} \circ \mathcal{C}_1 \right] (\rho) \right]$$

$$= \text{Tr} \left[ E \left[ S_m \right] (\rho) \right],$$

with

$$S_m = \mathcal{C}_{m+1} \circ \mathcal{E} \circ \mathcal{C}_m \circ \mathcal{E} \circ \ldots \circ \mathcal{C}_2 \circ \mathcal{E} \circ \mathcal{C}_1$$

$$= \mathcal{C}_{m+1} \circ (\mathcal{C}_m \circ \ldots \circ \mathcal{C}_1 \circ \mathcal{C}_1^* \circ \ldots \circ \mathcal{C}_m^*) \circ \mathcal{E} \circ \mathcal{C}_m \circ \mathcal{E} \circ \ldots \circ \mathcal{C}_2 \circ \mathcal{E} \circ \mathcal{C}_1$$

$$= \mathcal{D}_m \circ \mathcal{E} \circ \mathcal{D}_m \circ \ldots \circ \mathcal{D}_m \circ \mathcal{E} \circ \mathcal{D}_m \circ \mathcal{E} \circ \mathcal{D}_m$$

$$= \bigcap_{j=1}^{2^m} \left( \mathcal{D}_j \circ \mathcal{E} \circ \mathcal{D}_j \right),$$

where we have used the fact that $\mathcal{C}(n)$ is a group and defined a new quantum gate

$$\mathcal{D}_j := \bigcap_{i=1}^{2^j} \mathcal{C}_l,$$
with $\mathcal{D}(\cdot) = D_j \cdot D^*_j$, $D_j \in \mathcal{C}(n)$. Please note that all $D_j$ are independent uniformly distributed real Clifford gates.

The sequence fidelity can then be written as

$$F(m, E, \rho) = \text{Tr} \left[ E \bigotimes_{j=1}^m \left[ D^*_j \circ \mathcal{E} \circ D_j \right] (\rho) \right]. \quad (48)$$

Taking the average over the real Clifford group yields an average sequence fidelity given by

$$\bar{F}(m, E, \rho) = \text{Tr} \left[ E \left[ \frac{1}{|\mathcal{C}(n)|} \sum_{D_j \in \mathcal{C}(n)} S_m(\rho) \right] \right]$$

$$= \text{Tr} \left[ E \left[ \frac{1}{|\mathcal{C}(n)|} \sum_{D_j \in \mathcal{C}(n)} \bigotimes_{j=1}^m \left[ D^*_j \circ \mathcal{E} \circ D_j \right] (\rho) \right] \right]$$

$$= \text{Tr} \left[ E \left[ \frac{1}{|\mathcal{C}(n)|} \sum_{D_j \in \mathcal{C}(n)} D^*_j \circ \mathcal{E} \circ D_j \right]^\otimes_m (\rho) \right], \quad (49)$$

where we have used the fact that all $D_j$ are independent uniformly distributed Clifford gates. Because the real Clifford group is an orthogonal 2-design, see theorem 4, we have that

$$\frac{1}{|\mathcal{C}(n)|} \sum_{D_j \in \mathcal{C}(m)} D^*_j \circ \mathcal{E} \circ D_j (\cdot) = \int_{\mathcal{O}(d)} \mathcal{O}^* \circ \mathcal{E} \circ \mathcal{O} (\cdot) d\mathcal{O}, \quad (50)$$

where $\mathcal{O}(\cdot) = O \cdot O^*$, $O \in \mathcal{O}(d)$, is the orthogonal quantum channel. By eq. (13), the averaged sequence fidelity is

$$\bar{F}(m, E, \rho) = \text{Tr} \left[ E \left( \alpha \text{id} + \beta \frac{\Pi}{d} \text{Tr} [\cdot] + \gamma \frac{\Pi \text{Tr} [\cdot] - \theta}{d-1} \right)^\otimes_m (\rho) \right], \quad (51)$$

with $\alpha, \beta, \gamma \in \mathbb{R}$. Consider the following change of variables,

$$\alpha = \frac{1}{2} (b + c), \quad (52a)$$

$$\beta = a - b + \frac{1}{2} (b - c)d \quad \text{and} \quad (52b)$$

$$\gamma = \frac{1}{2} (c - b)(d - 1), \quad (52c)$$

such that

$$a = \alpha + \beta + \gamma, \quad (53a)$$

$$b = \alpha - \frac{\gamma}{d-1} \quad \text{and} \quad (53b)$$

$$c = \alpha + \frac{\gamma}{d-1}. \quad (53c)$$

Note that $a = \alpha + \beta + \gamma = 1$. The resulting quantum channel is

$$\hat{T}(\cdot) = a \text{Tr} [\cdot] \frac{\Pi}{d} + b \left( \frac{1}{2} (\text{id} + \theta) - \text{Tr} (\cdot) \frac{\Pi}{d} \right) + c \frac{1}{2} (\text{id} - \theta), \quad (54)$$

with $\alpha, \beta, \gamma \in \mathbb{R}$. Its $m$-fold concatenation is given by

$$\hat{T}^m (\cdot) = a^m \text{Tr} [\cdot] \frac{\Pi}{d} + b^m \left( \frac{1}{2} (\text{id} + \theta) - \text{Tr} (\cdot) \frac{\Pi}{d} \right) + c^m \frac{1}{2} (\text{id} - \theta). \quad (55)$$
Therefore, the averaged sequence fidelity is 

\[ \bar{F}(m, E, \rho) = \text{Tr} \left[ E \left( \alpha \text{id} + \frac{1}{d} \text{Tr} [\cdot] + \frac{1}{d-1} \right)^m \rho \right] \]

\[ = \text{Tr} \left[ E \left( a \text{Tr} [\cdot] \frac{1}{d} + b \left( \frac{1}{2} (\text{id} + \theta) - \text{Tr} (\cdot) \frac{1}{d} \right) + c \frac{1}{2} (\text{id} - \theta) \right)^m \rho \right] \]

\[ = \text{Tr} \left[ E \left( a^m \text{Tr} [\cdot] \frac{1}{d} + b^m \left( \frac{1}{2} (\text{id} + \theta) - \text{Tr} (\cdot) \frac{1}{d} \right) + c^m \frac{1}{2} (\text{id} - \theta) \right) \rho \right] \]

\[ = a^m \text{Tr} \left[ E \frac{1}{d} \right] + b^m \text{Tr} \left[ E \left( \frac{1}{2} \left( \rho + \rho^T \right) - \frac{1}{d} \right) \right] + c^m \text{Tr} \left[ E \frac{1}{2} \left( \rho - \rho^T \right) \right]. \quad (56) \]

Given that \( a = 1 \), the averaged sequence fidelity simplifies to 

\[ \bar{F}(m, E, \rho) = A + b^m B + c^m C, \quad (57) \]

where

\[ A = \text{Tr} \left[ E \frac{1}{d} \right], \quad (58a) \]

\[ B = \text{Tr} \left[ E \left( \frac{1}{2} \left( \rho + \rho^T \right) - \frac{1}{d} \right) \right] \quad \text{and} \quad (58b) \]

\[ C = \text{Tr} \left[ E \frac{1}{2} \left( \rho - \rho^T \right) \right], \quad (58c) \]

for any fixed prepared quantum state \( \rho \) and any fixed quantum measurement represented by a POVM with effect operators \( E \). As claimed, \( A \), \( B \) and \( C \) are independent of the noise channel.

The experimental data may thus be calibrated to this model using the real RB protocol discussed in section 7 by varying the parameter \( m \) and fitting the parameters \( b \) and \( c \). It is important to notice that any quantum state preparation errors and quantum measurement errors are absorbed by \( A \), \( B \) and \( C \) and the calibration of the model to experimental data is thus not affected by these errors. However, direct fitting is sometimes poorly conditioned in multi-exponential decays, and obtaining high accuracy can be challenging. We next describe a way to improve the contrast by choosing several different states and measurements and fitting to simpler decay curves by clever averaging. We make the simplifying assumption that the noise on the state preparations and measurements is independent of the particular state preparation, or the particular measurement. While this will not hold exactly in practice, in the regimes of interest there should still be a marked increase in statistical contrast when applying this idea [8, 18].

The standard approach in RB is to prepare eigenstates of \( Z \) on each qubit and then measure the POVM element that corresponds to \(+1\) eigenstates of \( Z \) on each qubit. In the ideal case, this choice eliminates \( C \) (since \( \rho \) is symmetric). This reduces the number of parameters that need to be fit for the same data, which typically leads to a higher quality fit. We can generalize this idea so that we can accurately estimate \( b \) and \( c \) separately without coupling to all of the nuisance parameters \( A \), \( B \), and \( C \) at once.

The idea is to randomly compile in an extra gate that effectively prepares different Pauli eigenstates, rather than just a \(+1\) eigenstate of \( Z \) on each qubit. First notice that every Pauli eigenstate is either complex symmetric \( (\rho = \rho^T) \) or antisymmetric \( (\rho = -\rho^T) \). The latter only holds if the Pauli eigenstate has an odd number of \( Y \) gates. So we should
choose from these initial states if we want to isolate the parameters $b$ and $c$. Similarly, the POVM element $E$ can be randomly chosen to be the $+1$ eigenstate or the $-1$ eigenstate of each given Pauli input eigenstate. For a given sequence length $m$, this defines four possible data sets. Letting $\beta_m(\rho, E)$ denote the data collected at sequence length $m$ with state preparations $\rho_\pm$ that are symmetric or antisymmetric, and with $\pm 1$ eigenstate projectors $E_\pm$, we have the four data sets

$$
\beta_m(\rho_+, E_+), \quad \beta_m(\rho_+, E_-), \quad \beta_m(\rho_-, E_+), \quad \beta_m(\rho_-, E_-). \tag{59}
$$

Each of these obeys (in expectation) the respective form of the expected fidelity decay eq. (57). Now we can look at the symmetries of the nuisance parameters $A$, $B$, and $C$ from eq. (58) and we see that linear combinations of the data sets can be chosen to eliminate one or more of the above parameters in the ideal case. In fact, we have that the following differences in data approach the following fidelity difference curves:

$$
\beta_m(\rho_+, E_+) - \beta_m(\rho_-, E_-) \to \bar{F}(m, \rho_+, E_+) - \bar{F}(m, \rho_-, E_-) = \Delta B b^m \tag{60a}
$$

$$
\beta_m(\rho_-, E_+) - \beta_m(\rho_+, E_-) \to \bar{F}(m, \rho_-, E_+) - \bar{F}(m, \rho_+, E_-) = \Delta C c^m, \tag{60b}
$$

where $\Delta B = \text{Tr}[E \rho_+]$ and $\Delta C = \text{Tr}[E \rho_-]$. In practice, there will not be exact cancellation, but this transformation will still greatly enhance the contrast. Therefore, collecting data at various values of $m$ and fitting to the model $\Delta B b^m$ or $\Delta C c^m$ respectively yields a much simpler exponential fit model, and standard tools from regression can be applied. Each of these fits then yields a separate estimate of the parameters $b$ and $c$ without a strong covariance between the estimates. Estimates of the average rebit fidelity as well as the average fidelity are thus obtained by eq. (35) and eq. (34) respectively. Real RB thus provides finer-grained information about the channel.

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After this paper was completed, we learned of closely related independent work by Brown and Eastin [5] that derives randomized benchmarking protocols for certain subgroups of the Clifford group.

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Appendix

Following [25], it is also possible to use the frame potential and establish that the real Clifford group is an orthogonal 2-design.

**Theorem 8 (Orthogonal frame potential).** Let \( D = \{ O_k \in \mathcal{O}(d) \}_{k=1,...,K} \) be a set of orthogonal real matrices. Define the frame potential of \( D \) to be

\[
\mathcal{P}(D) = \frac{1}{K^2} \sum_{O_k, O_k' \in D} |\text{Tr}[O_k^* O_k']|^4.
\]

The set \( D \) is an orthogonal 2-design if and only if \( \mathcal{P}(D) = 3 \).

**Proof.** Following the idea and notation of [25], the group \( D \) is an orthogonal 2-design if and only if \( \Delta := \tau_D - \tau_{\text{twirl}} = 0 \), where

\[
\tau_D = (\text{id}_{d^2} \otimes T_D) |\Omega\rangle\langle \Omega|,
\]

\[
\tau_{\text{twirl}} = (\text{id}_{d^2} \otimes T_{\text{twirl}}) |\Omega\rangle\langle \Omega|,
\]

\[
= \left( \text{id}_{d^2} \otimes \int_{\mathcal{O}(d)} (O \otimes O) \cdot (O \otimes O)^* \, dO \right) |\Omega\rangle\langle \Omega|.
\]

In order to see what this means in terms of the frame potential, let us introduce a basis with regards to the minimal projections given in eq. (11). Within the subspace of \( P_0 \), we introduce an orthonormal basis \( \{ |i_0\rangle \}_{j=1}^{d_0} \) with dimension \( d_0 = \text{dim } P_0 \). Similarly, we introduce an orthonormal basis \( \{ |i_1\rangle \}_{j=1}^{d_1} \) within the subspace of \( P_1 \) with dimension \( d_1 = \text{dim } P_1 \) as well as an orthonormal basis \( \{ |i_2\rangle \}_{j=1}^{d_2} \) within the subspace of \( P_2 \) with dimension \( d_2 = \text{dim } P_2 \). These then form an orthonormal basis \( \{ |i\rangle \}_{j=1}^{d^2} \) in \( \mathbb{C}^d \otimes \mathbb{C}^d \). The maximally entangled state is then given by \( |\Omega\rangle\langle \Omega| \) with

\[
|\Omega\rangle = \frac{1}{d} \sum_{m=0}^{2} \sum_{i_m=1}^{d_m} |i_m\rangle \otimes |i_m\rangle.
\]

Using this decomposition and using eq. (8), we see that

\[
\tau_{\text{twirl}} = \left( \text{id}_{d^2} \otimes \frac{2}{\text{Tr}[P_m]} \frac{\text{Tr}[[i_m]_m j_m] P_m}{\text{Tr}[P_m]} \right) |\Omega\rangle\langle \Omega|,
\]

\[
= \frac{1}{d^2} \sum_{m=0}^{2} \sum_{i_m, j_m=1}^{d_m} |i_m\rangle \langle j_m| \otimes \frac{\text{Tr}[[i_m]_m j_m] P_m}{\text{Tr}[P_m]} P_m
\]

\[
= \frac{1}{d^2} \sum_{m=0}^{2} \frac{1}{\text{Tr}[P_m]} P_m \otimes P_m.
\]

Furthermore,

\[
\tau_D = \frac{1}{d^2} \sum_{i,j=1}^{d^2} |i\rangle \langle j| \otimes \frac{1}{K} \sum_{O_k \in D} (O_k \otimes O_k) |i\rangle \langle j| (O_k \otimes O_k)^*.
\]
We will now show that $||\Delta||^2_2 := \text{Tr} ||\Delta||^2 = 0$, from which the claim follows. We thus want to compute

$$\text{Tr} [\Delta^* \Delta] = \text{Tr} [\tau^{*_{\text{twirl}}}_D \tau^{*_{\text{twirl}}} - \tau^{*_{\text{twirl}}} D - \tau^{*_{\text{twirl}}}_D \tau^{*_{\text{twirl}}} + \tau^{*_{\text{twirl}}} D \tau^{*_{\text{twirl}}}].$$

Its first term is easily calculated to give

$$\text{Tr} [\tau^{*_{\text{twirl}}}_D \tau^{*_{\text{twirl}}}] = \frac{1}{d^4} \text{Tr} \left( \sum_{m=0}^{2} \frac{1}{\text{Tr}[P_m]} P_m \otimes P_m \right) = \frac{3}{d^4}.$$

The second and third term yield

$$\text{Tr} [\tau^{*_{\text{twirl}}} D] = \frac{1}{d^4} \text{Tr} \left( \sum_{m=0}^{2} \frac{1}{\text{Tr}[P_m]} P_m \otimes P_m \right) \left( \sum_{i,j=1}^{d^2} |i\rangle \langle j| \otimes \frac{1}{K} \sum_{O_k \in D} (O_k \otimes O_k) |i\rangle \langle j| (O_k \otimes O_k)^* \right)$$

$$= \frac{1}{d^4 K} \sum_{m=0}^{2} \frac{1}{\text{Tr}[P_m]} \text{Tr} [P_m] |i\rangle \langle j| \text{Tr} \left( \sum_{O_k \in D} (O_k \otimes O_k) |i\rangle \langle j| (O_k \otimes O_k)^* \right),$$

where we have used the fact that $P_m$ commutes with $(O_k \otimes O_k)$. We then have that

$$\text{Tr} [\tau^{*_{\text{twirl}}}_D D] = \frac{1}{d^4 K} \sum_{m=0}^{2} \frac{1}{\text{Tr}[P_m]} \text{Tr} \left( \sum_{O_k \in D} (O_k \otimes O_k) P_m |i_m\rangle \langle i_m| (O_k \otimes O_k)^* \right)$$

$$= \frac{1}{d^4} \sum_{m=0}^{2} \frac{1}{\text{Tr}[P_m]} \text{Tr} [P_m] |i_m\rangle \langle i_m|$$

$$= \frac{3}{d^4}.$$

The last term is

$$\text{Tr} [\tau^{*}_{D} D] = \frac{1}{d^4 K^2} \sum_{i,j=1}^{d^2} \sum_{v,w=1}^{d^2} \text{Tr} [\langle j| \langle i| v \rangle \langle w|]$$

$$= \sum_{O_k, O_{k'} \in D} \sum_{O_{k'}, O_{k'} \in D} (O_k \otimes O_{k'}) |j\rangle \langle i| (O_k \otimes O_{k'})^* (O_{k'} \otimes O_{k'}) |v\rangle \langle w| (O_{k'} \otimes O_{k'})^*$$

$$= \frac{1}{d^4 K^2} \sum_{i,j=1}^{d^2} \sum_{O_k, O_{k'} \in D} \text{Tr} [(O_k \otimes O_{k'}) |j\rangle \langle i| (O_k \otimes O_{k'})^* (O_{k'} \otimes O_{k'}) |i\rangle \langle j| (O_{k'} \otimes O_{k'})^*]$$

$$= \frac{1}{d^4 K^2} \sum_{O_k, O_{k'} \in D} |\text{Tr} O_k^* O_{k'}|^4$$

$$= \frac{\mathcal{P}(D)}{d^4}.$$

The group $D$ is an orthogonal 2-design if and only if $\Delta = 0$, which gives $\mathcal{P}(D) = 3$ and thus the claim follows.

**Theorem 9.** Let $D = \{O_k \in \mathcal{O}(d)\}_{k=1,\ldots,K}$ be a set of real orthogonal matrices with the symmetry $G = \{O_k \otimes O_{k'}|O_k \in D\}$ affording the character $\zeta_G$. Then the following are equivalent:

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1. The set \(D\) is an orthogonal 2-design.

2. The symmetry has no more than three irreducible representations.

3. It holds that \(\langle \zeta_G, \zeta_G \rangle = 3\).

Before we can prove this theorem, let us recall the notion of a character, which plays a major role in the analysis of unitary representations. We mainly follow \[48\].

**Definition 10 (Character).** Given any unitary representation \(U : g \mapsto U(g)\) of a group \(G\), we define its character by
\[
\xi(g) := \text{Tr}[U(g)].
\] (62)

A character is called irreducible, if the unitary representation under consideration is irreducible.

Denote by \(\{V^{(i)}\}_i\) the irreducible unitary representations of \(G\) and the corresponding irreducible characters by \(\{\chi_i\}_i\). The irreducible characters are orthonormal in the group algebra \[48\] with the inner product given by
\[
\langle \chi_i, \chi_j \rangle := \frac{1}{|G|} \sum_{g \in G} \chi_i(g) \overline{\chi_j(g)} = \delta_{ij}.
\] (63)

If a representation is the direct sum of subrepresentations, then the corresponding character is the sum of the characters of those subrepresentations. This holds true especially for the decomposition into irreducible representations.

**Theorem 11** ([48, theorem III.2.4.]). *Every character \(\xi\) is of the form*
\[
\xi = \sum_i n_i \chi_i
\] (64)

*for nonnegative integers \(n_i\) and every such sum is the character of some representation.*

**Proof.** The proof is given in [48], but is reproduced here for the convenience of the reader. It is a standard result in representation theory that any finite-dimensional unitary representation can be written as a direct sum of finite irreducible unitary representations [48, theorem II.2.3.],
\[
U = \oplus_i n_i V^{(i)},
\] (65)
for some \(n_i\). But then we necessarily have that
\[
\xi = \sum_i n_i \chi_i,
\] (66)
which yields the claim. \(\square\)

**Corollary 12.** *If the character \(\xi\) has a decomposition as in theorem 11 given by eq. (64), then*
\[
\langle \xi, \xi \rangle = \sum_i n_i^2.
\] (67)
Proof. Substituting the decomposition into the inner product gives

\[
\langle \xi, \xi \rangle = \left\langle \sum_i n_i \chi_i, \sum_j n_j \chi_j \right\rangle = \sum_{i,j} n_i n_j \langle \chi_i, \chi_j \rangle = \sum_{i,j} n_i n_j \delta_{i,j} = \sum_i n_i^2,
\]

where we have used the orthogonality relation of irreducible characters given in eq. (63).

With these definitions and results at hand, we may now prove theorem 9, following the idea of [25].

Proof of theorem 9. Consider the unitary representation associated to the symmetry $G$ as given in the theorem 9 and its afforded character denoted by $\zeta_G$. The frame potential can be related to the character $\zeta_G$ by

\[
\mathcal{P}(D) = \langle \zeta_G, \zeta_G \rangle.
\]

Due to corollary 12 this must equal $\langle \zeta_G, \zeta_G \rangle = 3$ if and only if $G$ has exactly three irreducible components. This in turn is equivalent to $D$ being an orthogonal 2-design by theorem 8. The claim thus follows.

Theorem 13. The real Clifford group $C(n)$ is an orthogonal 2-design.

Proof. Given that the real Clifford group has three irreducible representations by theorem 3, it must be an orthogonal 2-design by theorem 9.