Entangling power and quantum circuit complexity

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Notions of circuit complexity and cost play a key role in quantum computing and simulation where they capture the (weighted) minimal number of gates that is required to implement a unitary. Similar notions also become increasingly prominent in high energy physics in the study of holography. While notions of entanglement have in general little implications for the quantum circuit complexity and the cost of a unitary, in this work, we discuss a simple such relationship when both the entanglement of a state and the cost of a unitary take small values, building on ideas on how values of entangling power of quantum gates add up. This bound implies that if entanglement entropies grow linearly in time, so does the cost. The implications are two-fold: It provides insights into complexity growth for short times. In the context of quantum simulation, it allows to compare digital and analog quantum simulators. The main technical contribution is a continuous-variable small incremental entangling bound.

Introduction

The circuit complexity of a computation captures the number of elementary steps it minimally takes to determine its outcome. A reading of the famous Church-Turing thesis states that all reasonable models of computation give rise to the same class of “easy” problems computable in polynomial time, a statement that can presumably also be applied to processes occurring in nature. Alas, ultimately the world is quantum. Indeed, notions of quantum circuit complexity have long been considered in quantum information science: They provide a quantitative account on the shortest quantum computation that implements a given unitary. Similarly, one can think of the complexity of a quantum state as the circuit complexity of the quantum circuit preparing it, starting from a given fiducial state. Such notions play a similarly central role in quantum as classical circuit complexities do in classical computing. Seminal work \cite{1,4} has introduced a geometric picture of circuit complexities, showing that finding the shortest circuit amounts to identifying the shortest path between two points in a curved geometry. In fact, this program has become so successful that the cost associated with a unitary in such a geometric picture has itself been identified with a notion of circuit complexity.

Yet, it was relatively recently that notions of circuit including those of costs rose to prominence outside the field of quantum computing \cite{5-13}. Again eluding to the physical Church Turing thesis, such an approach is well motivated: One can think of a quantum state – say, one that is being generated by a quantum chaotic Hamiltonian evolution – being highly complex if the quantum circuit that could have prepared it on a quantum computer would have to be long. Since one can argue about how many quantum gates one would have needed to emulate a given Hamiltonian time evolution, such notions also immediately allow to compare the effort in digital and analog quantum simulation \cite{16}. The possibly most compelling application of quantum circuit complexity is in the realm of high energy physics in the context of holography \cite{5-13}.

These thoughts provide fuel for a motivation to actually compute quantum circuit complexities and circuit costs. Yet, to actually quantitatively determine any variant of these quantities is not obvious. After all, there are many ways to decompose a given unitary into a quantum circuit, with the best known algorithms for decomposing given circuits in Clifford and $T$-gates featuring an exponential run-time in the circuit size \cite{20}, and the computation of the complexity requires the optimization over such decompositions. In any decomposition, one may expect cancellations of some sort, with the impact of a unitary gate being partially compensated by the later action of another, rendering naive combinatoric arguments involved. The geometrically motivated notion of a cost of a quantum circuit substantially lessens the technical burden \cite{1,4}, but it is still not obvious how to come up with meaningful lower bounds.

This work provides a compellingly simple lower bound for the cost of a quantum circuit that is tight for small values of the cost. It has indeed rightfully been argued that complexity is not entanglement \cite{12}, and neither is the cost of a circuit. No quantity based on entanglement can accommodate the presumed linear growth of state complexity until a time exponential in the system size \cite{5}, for obvious reasons. That said, for small values of the circuit cost and entanglement there is a simple connection: One can basically add up – if properly put together – potential entangling powers of quantum gates to arrive at tight bounds. The bounds presented are rooted in notions of entanglement capabilities of quantum gates: The argument captures
the insight that quantum gates that are close to the identity in operator norm have little capability to create entanglement from product states (which is very easy to show). They can also add very little entanglement to a given entangled state (which is less obvious to prove), but this can be grasped in terms of the small incremental entangling property \[21\] and which is here freshly proven for Gaussian continuous-variable systems. As such, the simple bound applies both to spin systems as to Gaussian bosonic continuous-variable settings, which are specifically important when approximating non-interacting bosonic quantum fields. Simple as the bound is, it is easily stated and proven (with some of the arguments delegated to the appendix). It can also straightforwardly be applied to important cases of quantum evolutions, for which quenched Hamiltonian many-body dynamics constitute an example.

**Quantum circuit complexity and cost**

The exact circuit complexity basically counts the number of quantum gates from a given gate set that is needed to exactly match the given unitary. An approximate reading thereof merely asks for an approximation in operator norm to a given small error. Lower bounds of the circuit complexity are provided by the cost of a given circuit, which is increasingly commonly seen as a notion of circuit complexity in its own right \[1\] [2]. For \( n \) quantum systems of local dimension \( d \) (\( d = 2 \) for spins or qubits), one chooses a collection of 2-local traceless Hamiltonian terms \( O_1, \ldots, O_J \), normalized in operator norm as \( \| O_j \| = 1 \) for \( j = 1, \ldots, J \). We consider both the situation in which \( \{ O_j \} \) are geometrically local and the situation where they are merely locally in their support. For a given \( U \in \text{SU}(d^n) \), one regards the unitary as being generated by a path-ordered integral

\[
U = \mathcal{P} \exp \left( -i \int_0^1 ds H(s) \right),
\]

with

\[
H(s) = \sum_{j=1}^J y_j(s) O_j,
\]

where \( y_j : [0,1] \to \mathbb{R} \) are appropriate continuous cost functions. This path ordered integral can in operator norm arbitrarily well approximated by

\[
V_N = \prod_{k=1}^N \exp \left( -i \frac{1}{N} \sum_{j=1}^J y_j(k/N) O_j \right)
\]

in the limit of \( N \to \infty \), as follows immediately from the definition of the path-ordered integral. The cost of a unitary \( U \in \text{SU}(d^n) \) can then be defined in such terms \[1\] [2].

**Definition 1** (Circuit cost \[1\]). For a given set \( \{ O_1, \ldots, O_J \} \) in the Lie algebra \( \text{su}(d^n) \) of traceless Hermitian matrices normalized as \( \| O_j \| = 1 \) for all \( j = 1, \ldots, J \), the cost of a quantum circuit \( U \in \text{SU}(d^n) \) is the infimum

\[
C(U) := \inf \int_0^1 \sum_{j=1}^J |y_j(s)| ds
\]

over all continuous functions \( y_j : [0,1] \to \mathbb{R} \) so that Eqs. \[2\] are satisfied. We call it the geometrically local circuit cost \( C_{\text{g}}(U) \) if all \( \{ O_j \} \) are geometrically local.

That is to say, the cost of a quantum circuit can be expressed in terms of the limit

\[
\lim_{N \to \infty} \frac{1}{N} \sum_{k=1}^N \sum_{j=1}^J |y_j(k/N)|
\]

of many time steps.

**Potential entangling power**

In what follows, the notion of a potential entangling power of a quantum gate provides some useful intuition. It captures the “coupling strength” and simply takes into account the fact that quantum gates that are close to the identity cannot create much entanglement. A somewhat related, but integer-valued, notion of entangling power has been invoked in Ref. \[23\].

**Definition 2** (Potential entangling power). A unitary \( U \in \text{SU}(d^2) \) has the potential entangling power

\[
e(U) := \log(d) \min \left\{ \| H \| : U = e^{-iH}, H = H^\dagger \right\}.
\]

It is indeed perfectly meaningful to refer to this quantity as the potential entangling power: If \( \rho = \rho_A \otimes \rho_B \), both \( \rho_A \) and \( \rho_B \) being pure and supported on \( \mathbb{C}^d \) each, then the resulting degree of entanglement \( S(t_{\text{BE}}(U \rho U^\dagger)) \) as quantified in terms of the von-Neumann entanglement entropy over the cut \( A : B \) is expected to be small if \( e(U) \) is small, and converging to zero for \( e(U) \to 0 \). Notions of entangling powers of quantum gates have long been connected to coupling strengths of interactions \[24\] [27]. As is well-known, notions of entangling power of unitary gates are altered depending on whether or not auxiliary quantum systems are allowed for: The swap gate obviously has no entangling power, if no auxiliary systems made use of, while it has \( 2 \log_2(d) \) when auxiliary systems are included. It is less obvious to see how much entanglement can be generated, however, if one initially already encounters an intricate entangled state and the unitary acts only on a small subsystem of the total system. The question of how much entanglement can be generated in this fashion has been largely settled in Ref. \[21\], however, which we can make use of here.

**Entanglement bounds circuit costs**

Such notions of potential entangling power can be related to tight bounds of circuit costs. In what follows, we denote for
a pure state $\rho$ defined on a spatially one-dimensional system of $n$ constituents for $s \in \{1, \ldots, n-1\}$ with

$$E(\rho : s) := S(\text{tr}_B(\rho))$$

(7)

being the entanglement entropy over the cut $A = \{1, \ldots, s\}$ and $B = \{s+1, \ldots, n\}$.

**Observation 1** (Entanglement lower bounds the cost). The geometrically local circuit cost of a $U \in SU(d^n)$ is lower bounded by

$$C_g(U) \geq \frac{1}{c \log(d)} \sum_{s=1}^{n-1} E(\phi : \phi | U^\dagger : s),$$

(8)

for an absolute constant $c > 0$, where $|\phi\rangle \in (\mathbb{C}^d)^\otimes n$ is a product state vector. For the cost, one finds

$$C(U) \geq \frac{1}{c \log(d)} \max_{\phi} E(\phi : \phi | U^\dagger : s).$$

(9)

The potential cancellation of gates in notions of complexity is faithfully captured in this bound: If most gates in a circuit commute, they will give rise to a lower circuit cost, but at the same time also to a smaller entanglement. So even if the bound is simple indeed, it does capture a key feature of the relationship of the circuit cost to notions of entanglement.

**Proof.** The proof of this observation is straightforward, acknowledging the results of Ref. [21]. We start by decomposing the circuit in a convenient manner. Making use of a Trotter decomposition, we find that $U$ can in operator norm $\|U - W_N\|$ be arbitrarily well approximated as a product

$$W_N := \prod_{k=1}^{N} V_k$$

(10)

with each term being given by

$$V_k := \exp \left( -\frac{i}{N} \sum_{j=1}^{J} y_j(k/N) O_j \right)$$

(11)

$$= \lim_{m \to \infty} \left( V_{k,1}^{1/m} \ldots V_{k,J}^{1/m} \right)^m$$

where

$$V_{k,j} := \exp \left( -\frac{i}{N} y_j(k/N) O_j \right).$$

(12)

Building upon this, let

$$|\psi_l\rangle := \bigotimes_{k=1}^{l} V_k |\psi\rangle$$

(13)

be the state vector after $l \in \{1, \ldots, N\}$ temporal layers, with $|\psi_0\rangle := |\phi\rangle$. Then, for $l = 1, \ldots, N$, using the time integrated instance of Lemma 1, one finds that the entanglement growth over the cut $A = \{1, \ldots, s\}$ and $B = \{s+1, \ldots, n\}$ in each step can at most be

$$E(|\psi_l\rangle |\psi_l\rangle : s) - E(|\psi_{l-1}\rangle |\psi_{l-1}\rangle : s)$$

$$= E(V_l |\psi_{l-1}\rangle |\psi_{l-1}\rangle V_l^\dagger : s) - E(|\psi_{l-1}\rangle |\psi_{l-1}\rangle : s)$$

$$\leq \frac{mc}{N} \sum_{j=1}^{J} y_j(l/N) \|O_j\| \log(d)$$

(14)

which gives

$$E(|\psi_l\rangle |\psi_l\rangle : s) - E(|\psi_{l-1}\rangle |\psi_{l-1}\rangle : s)$$

$$\leq \frac{c \log(d)}{N} \sum_{j=1}^{J} |y_j(l/N)|.$$  

(15)

Iterating this expression, one finds

$$E(U |\phi\rangle \langle \phi | U^\dagger : s) - E(|\phi\rangle \langle \phi | : s)$$

$$\leq \frac{c \log(d)}{N} \sum_{k=1}^{N} \sum_{j=1}^{J} |y_j(l/N)|.$$  

(16)

Acknowledging that the right hand side approximates the circuit cost $C(U)$ arbitrarily well, find finds the statement of Observation 1, by applying the argument to the cut $A = \{1, \ldots, s\}$ and $B = \{s+1, \ldots, n\}$ providing the tightest bound. For the geometrically local circuit cost $C_g(U)$, the argument can be applied to each such cut, leading to the statement of Observation 1. $\square$

In the above statement, the following statement from Ref. [21] has been made use of.

**Lemma 1** (Small incremental entanglement [21]). For a pure state $\rho$ and a Hamiltonian $h$ supported on a $d \times d$-dimensional subspace acting over the cut $\{1, \ldots, s\}$ and $\{s+1, \ldots, n\}$, the entangling rate defined as

$$\Gamma(h, \rho) := \frac{d}{dt} E(e^{-ith} \rho e^{ith} : s) \bigg|_{t=0}$$

(17)

is upper bounded by

$$\Gamma(h, \rho) \leq c \log(d) \|h\|.$$  

(18)

The constant presented in the proof is $c = 22$, but numerical evidence is shown that rather $c = 2$ actually provides a tight bound. Interpreted in terms of the above notion of an potential entangling power of a unitary $X \in U(d^2)$ acting on two constituents connecting the subsystems over the cut, one can argue that

$$|E(X \rho X^\dagger : s) - E(\rho : s)| \leq ce(X),$$

(19)

so that up to an absolute constant, the maximum increase of entanglement is indeed nothing but the potential entangling power: In each application, a quantum gate with a certain potential entangling power can increase the value of entanglement only to some extent, no matter how entangled the initial state has been. From the above Trotter decomposition it also follows that the circuit cost is nothing but the weighted quantum circuit complexity, weighted by the potential entangling power of each quantum gate.
Corollary 1 (Weighted quantum circuit complexity). For a given $U \in SU(2^n)$, the infimum of the sum of weights $e(U_j)$ of a circuit consisting of quantum gates $\{U_j\}$ generated by $\{O_j\}$ is given by $C(U)$.

Gaussian circuit cost

In fact, there is a small incremental entanglement bound as well as a harmonic equivalent of the above relationship between entanglement and quantum circuit cost for Gaussian bosonic settings $\{11, 13\}$, including ones motivated by evolutions of non-interacting bosonic quantum fields. For such bosonic systems, characterized by canonical coordinates $R = \{x_1, p_1, x_2, p_2, \ldots, x_n, p_n\}$, the appendix presents the proof of the following small incremental entanglement statement for such continuous-variable systems.

Theorem 1 (Gaussian small incremental entanglement). For a pure Gaussian state $\rho$ and a Hamiltonian $H = RH^\dagger$ supported on one of the modes each of $A = \{1, \ldots, s\}$ and $B = \{s + 1, \ldots, n\}$, the entangling rate defined as

$$\Gamma(h, \rho) := \frac{d}{dt} \mathbb{E} \left( e^{-itH} \rho e^{itH} : s \right) \bigg|_{t=0}$$

is upper bounded by

$$\Gamma(h, \rho) \leq \|h\| f(\|\gamma(0)\|), \quad (20)$$

where $f : [1, \infty) \rightarrow \mathbb{R}$ is a monotone increasing function.

Interestingly, it is not the operator norm of the Hamiltonian as such (which would make little sense anyway and would not be finite) but that of the kernel matrix when expressed as a polynomial in canonical coordinates that features in this small incremental entanglement statement. In the same way as above, and elaborated upon in the appendix, we can conclude the following.

Observation 2 (Gaussian entanglement lower bounds Gaussian circuit cost). The geometrically local Gaussian quantum circuit cost of a bosonic Gaussian unitary $U$ that prepares $U|\phi\rangle$ from the product state vector $|\phi\rangle$ associated with the covariance matrix $\gamma(0)$ is lower bounded by

$$G_\text{g} \geq \frac{1}{f(\|\gamma(0)\|)} \sum_{s=1}^{n-1} \mathbb{E}(U|\phi\rangle \langle \phi|U^\dagger : s). \quad (22)$$

For the Gaussian quantum circuit cost one finds

$$G \geq \frac{1}{f(\|\gamma(0)\|)} \max_s \mathbb{E}(U|\phi\rangle \langle \phi|U^\dagger : s). \quad (23)$$

Making use of these statements, one can infer about non-interacting bosonic theories in largely the same way as for spin systems, despite unbounded operators featuring in the problem.

Quenched quantum many-body systems

Simple as the above bounds are, they provide tight and relevant bounds to circuit costs and complexities for small times in a number of settings. An interesting insight along these lines of thought is the point that whenever a quantum many-body system undergoing non-equilibrium dynamics leads to a linear increase in the entanglement entropy over suitable cuts, so does the quantum state complexity. This is in particular true for quenched quantum many-body systems, for which the linear growth of entanglement entropies is generic $[28–30]$. In fact, both upper $[18, 19]$ and lower bounds $[31]$ for the entanglement entropy as a function of time have readily been established. That is to say, whenever the right hand side of Eq. (9) grows linearly in time, so does the left hand side, as an immediate corollary (see Fig. 1). We state this explicitly for the Ising Hamiltonian, but it should be clear that the same behaviour is expected for any local Hamiltonian (not featuring disorder).

Observation 3 (Growth of circuit cost in dynamics). For any time $T > 0$ there exists a system size $n$ for a translationally invariant Ising Hamiltonian such that the unitary dynamics $e^{-iHT}$ applied to a product state vector $|\phi\rangle$ leads to $C(e^{-iHT}) > \delta t$ for an absolute constant $\delta > 0$, for all times $t \in [0, T]$.

The upper bound in time $T$ is merely accommodating the possibility of having a finite system of finitely many degrees of freedom $n$, for which at some point, the respective entanglement entropies will no longer grow in time (rendering the bound then uninteresting). The result stated here is a corollary of Observation 1, together with the results of Ref. [22]. Since the model is translationally invariant, any cut serves to show the linear growth of the quantum state complexity in time. For the geometrically local circuit cost, one also finds a growth linear in time, but now the largest value of $C_\text{g}(e^{-iHT})$ attained at intermediate times scales as $\Theta(n^2)$ in the system size $n$, instead of the essentially linear scaling $\Theta(n)$ in case of the quantity $C(e^{-iHT})$.

Summary and outlook

In this work, we have carefully and quantitatively revisited the connection between entanglement and notions of circuit cost and complexity. While there is in general no tight connection between these quantities, for small values, there actually is, as this work shows: Indeed, one arrives at compellingly simple bounds. The usefulness of such bounds is manifest. One can argue, for example, how deep a weighted quantum circuit has to be to give rise to a given entanglement pattern in a desired final state; this is true at least for pure states, but it seems perfectly conceivable that similar techniques can be established for mixed quantum states. Also, it helps assessing the power and capabilities of analog quantum simulators $[16]$. Using such tools, one can argue that a digital quantum simulator would have required a precisely defined computational effort to produce the same results as a given analog quantum...
simulator. In this sense, it makes the computational effort of
digital and analog quantum simulators comparable. It is the
hope that this simple bound provides a useful and versatile tool
in various studies of this kind.

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APPENDIX

Preliminaries

Consider a quantum system of \(n\) bosonic modes,
equipped with the canonical coordinates \(R = (x_1, p_1, x_2, p_2, \ldots, x_n, p_n)\) reflecting positions and
momenta. For a quantum system comprising of \(n\) modes, the
canonical commutation relations give rise to a symplectic form

\[
\sigma = \bigoplus_{j=1}^{n} \begin{bmatrix} 0 & 1 \\ -1 & 0 \end{bmatrix}.
\]  

(24)

We consider Gaussian states \(\rho\) \([33, 34]\) with vanishing first
moments (which can be assumed to be the case without loss of
generality in the context considered) and second moments that
can be captured in the covariance matrix \(\gamma \in \mathbb{R}^{2n \times 2n}\) with entries

\[
\gamma_{j,k} = \text{tr}(\rho(R_j R_k + R_k R_j)).
\]  

(25)

Covariance matrices always satisfy the Heisenberg uncertainty
principle \(\gamma + i\sigma \geq 0\): It takes a moment of thought that this
is nothing but the standard Heisenberg uncertainty principle
written in a way that is manifestly invariant under symplectic
transformations that preserve the symplectic form \(\sigma\).

In the encore of bounding quantum circuit complexities,
the above quantification in terms of operator norms no longer
makes sense. However, when assessing notions of circuit
cost and complexity, similar bounds can still be derived when
appropriately evaluated for Hamiltonian terms. To this goal,
let \(\{O_j : j = 1, \ldots, J\}\) be a collection of operators – with
no assumption on the cardinality \(J\) of the set being made – each
of which being of the form

\[
O_j = (X_{k,j}, P_{k,j}, X_{l,j}, P_{l,j}) h_j (X_{k,j}, P_{k,j}, X_{l,j}, P_{l,j})^T,
\]  

(26)

where \(k \in \{1, \ldots, s\}\) and \(l \in \{s + 1, \ldots, n\}\) are labels refer-
ing to modes in \(A\) and \(B\), respectively. The fact that the
indices \(k, l\) are from the set of modes reflects the feature that
the operators are 2-local, but as before, geometric locality may
or may not be assumed. The matrices \(h_j \in \mathbb{R}^{2 \times 2}\) satisfy

\[
h_j = h_j^T, \quad ||h_j|| = 1.
\]  

(27)

That is to say, it is no longer the operators \(O_j\) as such that
have unit operator norm, but rather the kernels of quadratic
operators in the canonical coordinates. Equipped with this
preparation, we can again think of Hamiltonians

\[
H(s) = \sum_{j=1}^{J} y_j(s) O_j,
\]  

(28)

with as before \(y_j : [0, 1] \to \mathbb{R}\) being arbitrary continuous
functions. Just in the same way considered above, the Gaussian
circuit cost of a Gaussian unitary \(U\) becomes

\[
G(U) = \lim_{N \to \infty} \frac{1}{N} \sum_{k=1}^{N} \sum_{j=1}^{J} |y_j(k/N)|,
\]  

(29)

or analogously \(G_g(U)\), depending on whether the generators
have been chosen merely local or additionally geometrically
local. We are now in the position to prove Theorem 1.

Proof of Theorem 1

Proof. In what follows, we prove the Gaussian bosonic small
incremental entanglement statement summarized in Theorem
1 of the main text. As before, we consider a bi-partite system
consisting of parts \(A = \{1, \ldots, s\}\) and \(B = \{s + 1, \ldots, n\}\) with
two-local Hamiltonian supported on modes labeled
\(k \in A\) and \(l \in B\), with the rest of the systems

\[
a = \{1, \ldots, s\} \setminus \{k\}
\]  

(30)

and \(b = \{s + 1, \ldots, n\} \setminus \{l\}\) acting as auxiliary systems. The
entanglement rate for a given state \(\rho\) is then

\[
\Gamma(H, \rho) := \left. \frac{d}{dt} E(\rho(t) : s) \right|_{t=0}
\]  

(31)

where

\[
\rho(t) = e^{-itH} \rho e^{itH}.
\]  

(32)

On an abstract level, this entangling rate has been shown in Ref. \([21]\) to be given by

\[
\Gamma(H, \rho) = -it \text{tr}(\mathbb{I}_a \otimes H) \rho A, \{\{i\}\}, \log(\rho A) \otimes \mathbb{I}_{\{i\}}).
\]  

(33)

The Hamiltonian

\[
H = RhR^T,
\]  

(34)

which is a quadratic polynomial in the bosonic operators \(R\),
acts non-trivially on two modes in \(A\) and \(B\) each, so that \(h \in \mathbb{R}^{2n \times 2n}\) features non-vanishing elements only within a 4 \(\times\) 4 block reflecting pairs of canonical coordinates of two modes.
and has rank($h$) = 4 (more common is a coupling in position only, so that the respective block in the momentum sector is proportional to $I$). The generalization to $k$-local Hamiltonians with Hamiltonian kernels of rank $2k$ is immediate and only omitted for notational convenience.

The subsequent steps can be proven entirely on the level of level of second moments. The covariance matrix $\gamma$ of a pure Gaussian state $\rho$ takes the form

$$\gamma = \begin{bmatrix} \gamma_A & \Xi \\ \Xi^T & \gamma_B \end{bmatrix},$$

(35)

where the principal sub-matrices $\gamma_A$ and $\gamma_B$ reflect the reduced quantum states of the sub-systems labeled $A$ and $B$, respectively. The entanglement entropy of $\rho$ with respect to the split of $A$ versus $B$ [33] can be computed as

$$E(\rho : s) = \sum_{k=1}^s \left( \frac{\sigma_k + 1}{2} \right) \log \left( \frac{\sigma_k + 1}{2} \right) - \left( \frac{\sigma_k - 1}{2} \right) \log \left( \frac{\sigma_k - 1}{2} \right),$$

(36)

where $\sigma_1, \ldots, \sigma_s \geq 1$ are the symplectic eigenvalues of the $s$ modes of $\gamma_A$. These symplectic eigenvalues are the positive square roots of the eigenvalues of $-\gamma_A\sigma\gamma_A^T$, where

$$\sigma_A := \bigoplus_{j=1}^s \begin{bmatrix} 0 & 1 \\ -1 & 0 \end{bmatrix}$$

(37)

is the symplectic form of the $s$ modes constituting $A$, so that the entanglement entropy is found to be

$$S_A(\rho) = \text{tr}\left[ \left( \frac{M_A^{1/2} + \mathbb{I}}{2} \right) \log \left( \frac{M_A^{1/2} + \mathbb{I}}{2} \right) - \left( \frac{M_A^{1/2} - \mathbb{I}}{2} \right) \log \left( \frac{M_A^{1/2} - \mathbb{I}}{2} \right) \right],$$

(38)

where $M_A$ is the Hermitian $s \times s$-matrix defined as

$$M_A := \gamma_A^{1/2} (i\sigma_A) \gamma_A (i\sigma_A) \gamma_A^{1/2}$$

(39)

in terms of matrix square roots of covariance matrices. In the same way, $M_A(t)$ can be defined for all times $t \geq 0$, derived from the second moments of $\rho_A(t)$ given by

$$\gamma_A(t) = e^{-\sigma_A t} e^{\sigma_A t},$$

(40)

$\sigma$ denoting the symplectic form of the entire system involving all $n$ modes as defined in Eq. (24). Prepared in this fashion, we can turn to actually upper bounding the incremental entanglement rate. An explicit calculation shows that

$$\Gamma(H, \rho) = \frac{1}{2} \text{tr}\left[ \frac{d}{dt} M_A^{1/2}(t) \right]_{t=0} \times \left( \log \left( \frac{M_A^{1/2} + \mathbb{I}}{2} \right) - \log \left( \frac{M_A^{1/2} - \mathbb{I}}{2} \right) \right).$$

(41)

The rank

$$\text{rank} \left[ \frac{d}{dt} M_A^{1/2}(t) \right]_{t=0} = \text{rank} \left[ \frac{d}{dt} M_A(t) \right]_{t=0} \leq 3 \text{rank}(h)$$

(42)

is upper bounded by an absolute constant, as is seen by explicitly computing the derivative in time. What is more, we find the upper bound

$$\left\| \frac{d}{dt} M_A^{1/2}(t) \right\|_{t=0} \leq \| M_A(0) \|^{-1/2} \left\| \frac{d}{dt} M_A(t) \right\|_{t=0} \leq \| M_A(0) \|^{-1/2} \frac{d}{dt} \| \gamma_A(t) \|_{t=0}.$$ (43)

Using the sub-multiplicativity of the operator norm several times. Using the property that $\| M_A(0) \| \geq 1$, this gives

$$\left\| \frac{d}{dt} M_A^{1/2}(t) \right\|_{t=0} \leq \frac{d}{dt} \| \gamma_A(t) \|_{t=0}.$$ (44)

Introducing the projection $\pi := \mathbb{1}_A \oplus 0_B$, from this, we can upper bound the operator norm of the derivative and conclude that

$$\left\| \frac{d}{dt} M_A^{1/2}(t) \right\|_{t=0} \leq \| (\sigma_A \gamma_A - \gamma_A \sigma_A) \mathbb{1} \|$$ (45)

$$\leq \| (\sigma_A \gamma_A(0) - \gamma_A(0) \sigma_A) \mathbb{1} \|$$

$$\leq 2 \| \gamma_A(0) \| \| h \|,$$

again using the sub-multiplicativity of the operator norm. Then, the second term above can be bounded from above as

$$\left\| \log \left( \frac{M_A^{1/2} \pm \mathbb{I}}{2} \right) \right\| = \log \left( \frac{\| M_A^{1/2} \|}{2} \pm 1 \right) \leq \log \left( \frac{\| \gamma_A(0) \| + 1}{2} \right).$$ (46)

Putting these results together lets us arrive at the claim of the theorem, giving rise to the monotone increasing function $f : [1, \infty) \to \mathbb{R}$ that lets the small incremental entangling bound depends only on the coupling strength $\| h \|$. \hfill \Box

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