ON GLOBAL EFFECTS CAUSED BY LOCALLY NONEFFECTIVE UNITARY OPERATIONS

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Given a bipartite quantum state ρ with subsystems A and B of arbitrary dimensions, we study the entanglement detecting capabilities of locally noneffective, or cyclic, unitary operations [L. B. Fu, Europhys. Lett., vol. 75, pp. 1–7, 2006]. Local cyclic unitaries have the special property that they leave their target subsystem invariant. We investigate the distance between ρ and the global state after local application of such unitaries as a possible indicator of entanglement. To this end, we derive and discuss closed formulae for the maximal such distance achievable for three cases of interest: (pseudo)pure quantum states, Werner states, and two-qubit states. What makes this criterion interesting, as we show here, is that it surprisingly displays behavior similar to recent anomalies observed for non-locality measures in higher dimensions, as well as demonstrates an equivalence to the CHSH inequality for certain classes of two-qubit states. Yet, despite these similarities, the criterion is not itself a non-locality measure. We also consider entanglement detection in bound entangled states.

Keywords: entanglement, entanglement detection, cyclic unitary, locally noneffective unitary, non-locality

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1 Introduction

Over the last two decades, quantum entanglement has been the subject of intense research, due to the continuing discoveries of interesting uses for the phenomenon by the quantum computing and information community (see [1] and [2] for surveys). One of the remaining open problems, however, is that of deciding separability of a quantum state - that is, given a (in our case, bipartite) state ρ acting on the Hilbert space $\mathcal{H}^M \otimes \mathcal{H}^N$, where M and N denote the respective dimensions of the subsystems, decide whether ρ can be written in the form

$$\rho = \sum_{k=1}^{n} p_k |a^k\rangle\langle a^k| \otimes |b^k\rangle\langle b^k|, \tag{1}$$

for $p_k \in \mathbb{R}^+$, $\sum_{k=1}^n p_k = 1$, $n \ge 1$, $|a^k\rangle \in \mathcal{H}^M$, $|b^k\rangle \in \mathcal{H}^N$, and $\|a^k\|_2 = \|b^k\|_2 = 1$, for $\|\cdot\|_2$ denoting the Euclidean norm. A state which can be written in this form is called *separable*,

and if additionally we have n > 1, we refer to the state as classically correlated. This problem was proven NP-hard by Gurvits [3] (with subsequent hardness results in Refs. [4, 5]), implying that it is highly unlikely for a general solution to exist for all possible inputs ρ . Another topic of interest which has recently garnered renewed attention is the study of entanglement and non-locality as comprising distinct resources [6, 7, 8, 9]—for generalized Bell inequalities for qutrits, or the ability to close the detection loophole, for example, it has been found that non-maximally entangled quantum states outperform maximally entangled states. Here, we study an approach to detecting entanglement which surprisingly also displays effects similar to those seen for such entanglement versus non-locality arguments.

The approach we consider was originally proposed by Fu [10]. It consists of applying a unitary operation to one of the subsystems, while demanding that the density matrix of the subsystem is invariant under this transformation. However, the global density matrix may be changed, and therefore there may be a non-zero distance between the original (global) state and the one after applying the local unitary operation. We will call this quantity "Fu distance", and will be particularly interested in its maximal value, for a given initial state, where the maximisation is over all locally noneffective unitaries.

Note that the principle underlying this approach has long been implicitly harnessed, for example, in superdense coding [11], where applying a Pauli operator to half of a Bell state can give rise to an orthogonal Bell state.

This paper is organized as follows. In Section 2, we define the Fu distance and discuss its relevant properties. In Sections 3.1 and 3.2 we derive closed formulae for the maximal Fu distance for pseudopure states (i.e. mixtures of a projector onto a pure state with the identity) and Werner states, respectively. In Section 4, we discuss connections between the CHSH inequality and the maximal Fu distance for two-qubit systems. In Section 5 we present the attempt to detect bound entanglement using the Fu distance. Finally, we conclude in Section 6 and pose open questions.

2 Definition and properties of the Fu Distance

Given a quantum state ρ , acting on $\mathcal{H}^M \otimes \mathcal{H}^N$, with density matrices of the subsystems $\rho_A = \text{Tr}_B(\rho)$ and $\rho_B = \text{Tr}_A(\rho)$, define a locally noneffective, or *cyclic* [10], unitary operation U^B , as one satisfying the condition $U^B \rho_B U^{B\dagger} = \rho_B$. This is equivalent to demanding

$$[\rho_B, U^B] = 0. (2)$$

Then, letting $\rho_f = (I \otimes U^B) \rho (I \otimes U^{B\dagger})$, our quantity of interest is

$$d(\rho, U^B) := \frac{1}{\sqrt{2}} \| \rho - \rho_f \|_{F}, \qquad (3)$$

which we dub the Fu distance [10], and where $\|A\|_{\rm F} = \sqrt{{\rm Tr}(A^\dagger A)}$ denotes the Frobenius norm (or Euclidean norm) for matrices. Thus, we are applying a local unitary operation which leaves the target reduced state invariant, and yet may produce a global shift in the joint system, the quantification of which we will study as a possible indicator of entanglement. To this end, we will be most interested in the quantity

$$d_{\max}(\rho) := \max_{\text{cyclic } U^B} d(\rho, U^B), \tag{4}$$

i.e. the maximal possible global distance achieved under any locally noneffective unitary operation.

Let us briefly discuss some relevant properties of $d_{max}(\rho)$. First, note that Eq. (3) can be straightforwardly rewritten in the useful form [10]

$$d(\rho, U^B) = \sqrt{\text{Tr}(\rho^2) - \text{Tr}(\rho\rho_f)},$$
(5)

from which it is easy to see that $0 \leq d(\rho, U^B) \leq 1$, with the latter inequality saturated if and only if ρ is pure and orthogonal to ρ_f . For any product state, i.e. $\rho = \rho_A \otimes \rho_B$, it is clear that $d_{\max}(\rho) = 0$ [10]. It is not known whether $d_{\max}(\rho) > 0$ for all entangled states, although we will later show that this is in indeed the case for all entangled pseudopure and Werner states.

One can find an upper bound for the Fu distance of any classically correlated state ρ_{cc} in a bipartite system, by generalizing an argument of Fu [10] to dimensions M and N for the subsystems:

$$d_{\max}(\rho_{cc}) \le \sqrt{\frac{2(M-1)(N-1)}{MN}}.$$
 (6)

The derivation of Eq. (6) is as follows: Any bipartite state ρ can be written in Fano form [12]:

$$\rho = \frac{1}{MN} \left(I^A \otimes I^B + \vec{r}^A \cdot \vec{\sigma}^A \otimes I^B + I^A \otimes \vec{r}^B \cdot \vec{\sigma}^B + \sum_{i=0}^{M^2 - 1} \sum_{j=0}^{N^2 - 1} T_{ij} \sigma_i^A \otimes \sigma_j^B \right), \quad (7)$$

where I denotes the identity matrix, \vec{r}^A denotes the (M^2-1) -dimensional Bloch vector for subsystem A with $r_i^A = \frac{M}{2} \text{Tr}(\sigma_i^A \rho_A)$, $\vec{\sigma}^A$ denotes the $(M^2 - 1)$ -component vector of traceless Hermitian generators for SU(M), and the matrix T is a real matrix known as the correlation matrix, whose entries are given by $T_{ij} = \frac{MN}{4} \text{Tr}(\sigma_i^A \otimes \sigma_j^B \rho)$. The definitions for subsystem B are analogous.

Eq. (5) can now be rewritten via straightforward manipulation as [10]

$$d(\rho, U^B) = \frac{2}{MN} \sqrt{\sum_{ij} T_{ij}^2 - \sum_{ij} T_{ij} T_{ij}^f},$$
 (8)

where T^f is the correlation matrix for ρ_f . Let us derive bounds on each sum in the square root. First, for a separable state $\rho_{cc} = \sum_{k=1}^n p_k |a^k\rangle\langle a^k| \otimes |b^k\rangle\langle b^k|$, one has $T_{ij} = \sum_k p_k r_i^{A_k} r_j^{B_k}$, where \vec{r}^{A_k} and \vec{r}^{B_k} are the Bloch vectors corresponding to states $|a^k\rangle\langle a^k|$ and $|b^k\rangle\langle b^k|$, respectively [10]. Via the Cauchy-Schwarz inequality, we thus have

$$\sum_{ij} T_{ij}^2 = \sum_{ij} \left(\sum_k p_k r_i^{A_k} r^{B_k} \right)^2 \le \sum_k p_k \left(\sum_i r_i^{A_k} \right) \left(\sum_j r_j^{B_k} \right) \le \frac{MN(M-1)(N-1)}{4}, \tag{9}$$

where the last inequality follows from the fact that $\|r^{A_k}\|_2 \leq \sqrt{M(M-1)/2}$ (resp. for $||r^{B_k}||_2$) [13]. A lower bound of $\sum_{ij} T_{ij} T_{ij}^f \ge -\sum_{ij} T_{ij}^2$ is also easily found using the Cauchy-Schwarz inequality. Substituting these into Eq. (8) gives Eq. (6), as desired.

Note that the bound in Eq. (6) is only non-trivial for small dimensions M and N, as $d(\rho, U^B) \leq 1$ has to hold. For M = N = 2, Eq. (6) gives a tight bound of $d_{\text{max}}(\rho_{\text{cc}}) \leq 1/\sqrt{2}$,

 $[\]overline{{}^{a}\text{Consider}}$, for example, $\rho_{cc} = \frac{1}{2}|00\rangle\langle 00| + \frac{1}{2}|11\rangle\langle 11|$ and $U^{B} = \sigma_{x}$, the Pauli X operator.

and for M = N = 3, a possibly loose bound of $d_{\text{max}}(\rho_{\text{cc}}) \leq \sqrt{8/9}$. Determining a tight upper bound on $d_{\text{max}}(\rho_{\text{cc}})$ for arbitrary dimensions remains at present an intriguing open problem.

3 Maximizing the Fu Distance

3.1 Pseudopure States

We now derive a closed formula for $d_{max}(\rho)$ for pseudopure states, and follow with a discussion of its implications. Specifically, consider a pseudopure quantum state ρ acting on $\mathcal{H}^M \otimes \mathcal{H}^N$, such that

$$\rho = \epsilon \sigma + \frac{1 - \epsilon}{MN} I,\tag{10}$$

for $\sigma = |\psi\rangle\langle\psi|$ a pure state of dimension MN, and $0 < \epsilon \le 1$. (The case $\epsilon = 0$ leads trivially to $\mathrm{d_{max}}(\rho) = 0$.) Without loss of generality, we assume $M \ge N$ in the following. Using the Schmidt decomposition [14], we can write $|\psi\rangle = \sum_{k=0}^{N-1} a_k |k\rangle_A \otimes |k\rangle_B$, where a_k are non-negative real numbers (Schmidt coefficients) with $\sum_{k=0}^{N-1} a_k^2 = 1$, and $\{|k\rangle_A\}_{k=0}^{N-1}$ are the elements of the Schmidt basis for subsystem A (analogously for B). We first prove the following useful lemma.

Lemma 1 Let ρ , acting on $\mathcal{H}^M \otimes \mathcal{H}^N$, be a pseudopure quantum state as defined in Eq. (10). Then, for any k such that $a_k \neq a_j$ for all $j \neq k$, and for any unitary U^B , if $[\rho_B, U^B] = 0$, then $|U_{k,k}^B| = 1$.

Proof. We first write $U^B = \sum_{ln} \langle l|U^B|n\rangle |l\rangle\langle n|$, for $\{|n\rangle\}_{n=0}^{N-1}$ the Schmidt basis for subsystem B. Then:

$$[\rho_B, U^B] = \epsilon \sum_{ln} (a_l^2 - a_n^2) \langle l | U^B | n \rangle | l \rangle \langle n |.$$
(11)

If two Schmidt coefficients of $|\psi\rangle$ differ in value, it therefore follows that the corresponding entry in U^B must be 0 in order for ρ_B and U^B to commute. Thus, for unique a_k , row k and column k of U^B must be all zeroes, except for position $U^B_{k,k}$, for which $|U^B_{k,k}| = 1$, since U^B is unitary \Box .

We now show the main result of this section. For the remainder of our discussion, let us denote the maximal Schmidt coefficient of $|\psi\rangle$ as $a_m := \max_k a_k$.

Theorem 1 Let ρ , acting on $\mathcal{H}^M \otimes \mathcal{H}^N$, be a pseudopure quantum state as defined in Eq. (10). Then,

$$d_{\max}(\rho) = \begin{cases} \epsilon & \text{if } a_m^2 \le \frac{1}{2}, \\ 2\epsilon a_m \sqrt{1 - a_m^2} & \text{otherwise.} \end{cases}$$
 (12)

Proof. Inserting ρ of Eq. (10) into Eq. (5) leads for arbitrary U^B (not necessarily cyclic), to

$$d(\rho, U^B) = \epsilon \sqrt{1 - \left| \sum_{k=0}^{N-1} a_k^2 \langle k | U^B | k \rangle \right|^2}, \tag{13}$$

from which it follows that $d(\rho, U^B)$ depends only on the diagonal entries of U^B . Let us hence first assume that U^B is a diagonal unitary matrix with eigenvalue $e^{i\theta_k}$ on row k (for

 θ_k to be chosen), and subsequently show that choosing U^B in this way is always optimal. By Eq. (13), we then find that maximizing $d(\rho, U^B)$ reduces to minimizing $\left|\sum_{k=0}^{N-1} a_k^2 e^{i\theta_k}\right|$. Since ρ_B is diagonal, any choice of θ_k 's constitutes a commuting unitary operation U^B , and so this minimization problem has a simple geometric solution as follows.

If $a_m^2 \leq \sum_{k \neq m} a_k^2$ (or equivalently, $a_m^2 \leq 1/2$), then one can always construct a closed polygon using vectors of the lengths a_k^2 each exactly once. Hence, $\min_{\{\theta_k\}_k} \left| \sum_{k=0}^{N-1} a_k^2 e^{i\theta_k} \right| = 0$, and so $d_{\max}(\rho) = \epsilon$. If, however, $a_m^2 > \sum_{k \neq m} a_k^2$, then no such polygon can be constructed, and the best minimization strategy is simply to set $\theta_m = 0$ and $\theta_k = \pi$, for all $k \neq m$. Substitution into our simplified expression for Eq. (13) and using the normalization $\sum_{k=0}^{N-1} a_k^2 = 1$ then gives $d_{\max}(\rho) = 2\epsilon a_m \sqrt{1 - a_m^2}$, as desired.

Finally, to see that choosing U^B diagonal is always optimal, note that if $a_m^2 \leq \sum_{k \neq m} a_k^2$, then our strategy for diagonal U^B achieves the maximum possible value for Eq. (13). If $a_m^2 > \sum_{k \neq m} a_k^2$, on the other hand, clearly $a_m \neq a_k$ for all $k \neq m$, and so it follows from Lemma 1 that $|\langle m|U^B|m\rangle|=1$ in Eq. (13). Thus, the best minimization strategy is again the same as that outlined for the diagonal case. This concludes the proof \Box .

Let us point out some consequences of Theorem 1. First, from Eq. (6), we know that a necessary condition for using $d_{\max}(\rho)$ to detect entanglement in ρ , acting on $\mathcal{H}^M \otimes \mathcal{H}^N$, is $d_{\text{max}}(\rho) > 1/\sqrt{2}$. Thus, from Theorem 1 and straightforward calculation, one finds that $d_{max}(\rho)$ may be used to detect entanglement in pseudopure states ρ only if

$$\frac{1}{2} \left(1 - \sqrt{1 - \frac{1}{2\epsilon^2}} \right) \le a_m^2 \le \frac{1}{2} \left(1 + \sqrt{1 - \frac{1}{2\epsilon^2}} \right). \tag{14}$$

For two-qubit pure states, this becomes an if and only if condition, and the corresponding upper bound on a_m simplifies to approximately 0.924 (the lower bound in this case is implicitly given by $1/\sqrt{2}$, by definition of a_m). Second, for general dimensions, if one knows that $|\psi\rangle$ is a maximally entangled state of the form $|\psi\rangle = \frac{1}{\sqrt{D}} \sum_{k=0}^{D-1} |kk\rangle$ for D = M = N, then it follows from Theorem 1 that one can always reliably detect the entanglement of the pseudopure state ρ , since in this case ρ is entangled if and only if $\epsilon > 1/(D+1)$ [15]. Third, it is clear from Theorem 1 that $d_{max}(\rho) > 0$ for any entangled pseudopure ρ . Finally, for pure ρ , Theorem 1 implies that one achieves $d_{\max}(\rho) = 1$ as long as $a_m \leq 1/\sqrt{2}$. Hence, ρ need not be maximally entangled in order to achieve a maximal shift. This surprising behavior is plotted in Figure 1. One possible interpretation for this is as follows: Observe that any pure state ρ with $a_m \leq 1/\sqrt{2}$ is majorized by the singlet (i.e. a pure state with Schmidt rank 2 and both amplitudes $1/\sqrt{2}$). It hence follows by Nielsen's majorization theorem [16] that a pure state achieves $d_{max}(\rho) = 1$ if and only if it can be deterministically converted to the singlet via local operations and classical communication (LOCC). It similarly follows that a pure state ρ with $d_{max}(\rho) < 1$ cannot be converted via LOCC to a state with Fu distance 1 — the converse, however, is possible provided the required majorization condition [16] holds.

Is there any connection between the Fu distance and the concurrence? This is the case for a two-qubit pure state, i.e. $|\psi\rangle = a_0|00\rangle + a_1|11\rangle$ and $\epsilon = 1$. Then Eq. (12) reduces to $d_{\max}(\rho) = 2a_0a_1 = C(\rho)$, where $C(\rho)$ denotes the concurrence of ρ [17]. Used by Wootters [18] to derive an analytic formula for the entanglement of formation of two-qubit states, and an

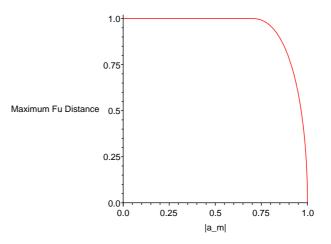


Fig. 1. A plot of $d_{\max}(\rho)$ for pure states as a function of a_m , as given by Eq. (12). Note that only the region $a_m \geq 1/\sqrt{N}$ is accessible, where N is the smaller dimension of the two subsystems.

entanglement measure in its own right, the concurrence has a number of generalizations to higher dimensions [19, 20, 21, 22, 23], two of which have simple closed forms for the case of pure states, which we shall compare to Eq. (12) of Theorem 1 here. To do so, let ρ , acting on $\mathcal{H}^M \otimes \mathcal{H}^N$, be pure, and set $D = \min\{M, N\}$. Then, Rungta et al. [20] define the concurrence $C_R(\rho)$ for a pure state ρ as

$$C_R(\rho) = \sqrt{1 - \text{Tr}(\rho_A^2)},\tag{15}$$

where $0 \le C_R(\rho) \le \sqrt{2(D-1)/D}$. This expression can be rewritten as

$$C_R(\rho) = \sqrt{2\left(1 - \sum_{k=0}^{D-1} a_k^4\right)},$$
 (16)

from which it is clear that Eq. (12) does not reduce to (a normalized version of) $C_R(\rho)$. As a supporting example, consider $|\psi\rangle=\frac{1}{\sqrt{2}}|00\rangle+\frac{1}{2}|11\rangle+\frac{1}{2}|22\rangle$, for which $d_{\max}(\rho)=1$, but $C_R(\rho)\approx 0.9682$, where we have normalized the latter by the maximum value possible for qutrits, $2/\sqrt{3}$. Next, consider the generalization of Audenaert et al. [21], which states that for pure ρ , we have $C_A(\rho)=2a_ma_{m2}$, where a_m and a_{m2} are the first and second largest Schmidt coefficients in the Schmidt decomposition of $|\psi\rangle$. Again, it is clear that Eq. (12) does not reduce to this definition $C_A(\rho)$ either. As an example, consider the maximally entangled two qutrit state $|\psi\rangle=\frac{1}{\sqrt{3}}(|00\rangle+|11\rangle+|22\rangle)$, for which $d_{\max}(\rho)=1$, but $C_A(\rho)=2/3$.

Note that in Figure 1 it is already evident that for $D \ge 3$ several non-maximally entangled states lead to the same maximal Fu distance, and therefore the Fu distance can in general not be used to define an entanglement measure.

3.2 Werner States

We now turn our attention to bipartite Werner states ρ_W acting on $\mathcal{H}^D \otimes \mathcal{H}^D$ with $D \geq 2$, for which we derive a closed formula for $d_{\max}(\rho_W)$. Denoting as $\{|i\rangle\}_{i=0}^{D-1}$ an arbitrary

orthonormal basis for \mathcal{H}^D , the Werner state ρ_W can be defined as follows [24]:

$$P = \sum_{ij} |i\rangle\langle j| \otimes |j\rangle\langle i| \tag{17}$$

$$P_{sym} = \frac{1}{2}(I_{D^2} + P) \tag{18}$$

$$P_{as} = \frac{1}{2}(I_{D^2} - P) \tag{19}$$

$$\rho_W = p \frac{2}{D^2 + D} P_{sym} + (1 - p) \frac{2}{D^2 - D} P_{as}, \tag{20}$$

where I_{D^2} is the D^2 -dimensional identity matrix and $0 \le p \le 1$. The state ρ is invariant under operation $U \otimes U$, for any choice of unitary U, and is entangled for p < 1/2, and separable otherwise. Investigating in terms of Fu distance, we find the following result.

Theorem 2 Let ρ_W , acting on $\mathcal{H}^D \otimes \mathcal{H}^D$, be a Werner state, as defined in Eq. (20). Then

$$d_{\max}(\rho_W) = \frac{|2pD - D - 1|}{D^2 - 1},\tag{21}$$

obtained using any traceless $D \times D$ choice of unitary U^B .

Proof. Consider substitution of ρ_W and arbitrary U^B into Eq. (5). Observing that Tr(P) = D, $\text{Tr}(P^2) = D^2$, and defining $\beta := \text{Tr}(P(I \otimes U^B)P(I \otimes U^{B^{\dagger}})) = \text{Tr}(U^B)\text{Tr}(U^{B^{\dagger}})$, straightforward manipulation leads us to

$$d(\rho_W, U^B) = \frac{\sqrt{(2pD - D - 1)^2(D^2 - \beta)}}{D(D^2 - 1)}.$$
(22)

Examining the boundary and critical points of the first derivative of Eq. (22) with respect to β , we find that the two cases of interest are $\beta = -D$ and $\beta = 0$. Note, however, that $\beta = -D$ implies $\text{Tr}(U^B)\text{Tr}(U^{B\dagger}) = -D$, which is impossible, since $aa^* \geq 0$ for all $a \in \mathbb{C}$. Hence, the maximum Fu distance is achieved when $\beta = 0$, implying that U^B is traceless, giving the desired result \square .

We remark that for a two-qubit Werner state ρ_W , Eq. (21) reduces to Eq. (12), as required, since in this case ρ_W can be written as the pseudopure state

$$\rho_W = \left| 1 - \frac{4}{3}p \right| \cdot |\psi^-\rangle\langle\psi^-| + \left(1 - \left| 1 - \frac{4p}{3} \right| \right) \cdot \frac{I}{4}$$
 (23)

where $|\psi^{-}\rangle = 1/\sqrt{2}(|01\rangle - |10\rangle)$ and $0 \le p \le 1$.

We now direct the reader's attention to Figure 2, which graphically depicts $d_{max}(\rho_W)$ for various dimensions D. For D=2, we find from Eq. (6) that one can detect entanglement in Werner states if $p \leq (3\sqrt{2}-3)/4\sqrt{2} \approx 0.220$. For arbitrary D, if we are promised that ρ_W is a Werner state, but not given the value of p, then by Figure 2, attaining $d_{max}(\rho_W) > 1/(D+1)$ is sufficient to conclude that ρ_W is entangled, seen by considering the case p=1 which leads to the maximal Fu distance for a separable Werner state. From Eq. (21), it is also clear that $d_{max}(\rho_W) > 0$ for all entangled Werner states.

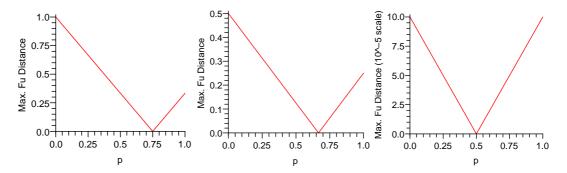


Fig. 2. $d_{\text{max}}(\rho_W)$ plotted for D=2,3,10000, respectively, as given by Eq. (21). Note the scale of 10^{-5} for the vertical axis for the case of D=10000.

We remark that examining the critical points of the first derivative of Eq. (21) as $D \to \infty$ suggests that there is a "kink" in the graph at p=1/2, which is precisely the boundary between entangled and separable Werner states. We also see that as $D \to \infty$, $d_{\max}(\rho_W) \to 0$, such that the possibility of distinguishing between classical and quantum correlations vanishes. To explain these phenomena (at least in a mathematical sense), we invoke the following theorem.

Theorem 3 ([25]) For any state
$$\rho$$
 acting on $\mathcal{H}^M \otimes \mathcal{H}^N$, $d_{\max}(\rho) \leq \sqrt{2 \left[\operatorname{Tr}(\rho^2) - \frac{1}{MN} \right]}$.

Intuitively, this means that the maximal Fu distance for a state is upper bounded by a dependence on the state's mixedness. In our case, straightforward calculation yields for the Werner state ρ :

$$Tr(\rho^2) = p^2 \frac{2}{D^2 + D} + (1 - p)^2 \frac{2}{D^2 - D}.$$
 (24)

The first derivative of this reveals that $\text{Tr}(\rho^2)$ has a minimum at $\frac{1+1/D}{2}$. Thus, as $D \to \infty$, ρ is most mixed at p approaching 1/2, explaining the first phenomenon above. Similarly, one finds from Eq. (24) that $\text{Tr}(\rho^2) \to 0$ as $D \to \infty$, explaining the second phenomenon.

4 Connections to the CHSH Inequality

We now investigate connections between the CHSH inequality and the Fu distance for twoqubit systems. Our motivation stems from the following observation. From Eqs. (12) and (6), it is immediate that in order to use $d_{\max}(\rho)$ to detect entanglement in the two-qubit Werner state $\rho_W = p|\psi^-\rangle\langle\psi^-| + \frac{1-p}{4}I$, where $|\psi^-\rangle = \frac{1}{\sqrt{2}}(|01\rangle - |10\rangle)$, we require $p > 1/\sqrt{2}$. On the other hand, define the following quantity for any quantum state ρ , acting on $\mathcal{H}^2 \otimes \mathcal{H}^2$:

$$M(\rho) := \tau_1(T^T T) + \tau_2(T^T T),$$
 (25)

where T is the correlation matrix of ρ from Eq. (7), T^T its transpose, and $\tau_1(T^TT)$ and $\tau_2(T^TT)$ the first and second largest eigenvalues of T^TT , respectively. Then, ρ can violate the CHSH inequality if and only if [26]

$$M(\rho) > 1. \tag{26}$$

For the Werner state ρ_W , one has $M(\rho_W) = 2p^2$, and so in order to detect entanglement in ρ_W using the CHSH inequality, one requires $p > 1/\sqrt{2}$, which is the same bound obtained above for the Fu distance. Thus, we will pose and answer the question of whether there is a connection between the ability to detect entanglement via the CHSH inequality versus the Fu distance.

Our approach is as follows. We first show that, without loss of generality, one can take the correlation matrix T of ρ to be diagonal. We then derive a closed formula for $d_{max}(\rho)$ for any two-qubit state ρ with diagonal T. Using this formula, we compare $d_{\max}(\rho)$ and $M(\rho)$.

To begin, we follow [27, 28] and note that applying a unitary operation $U_1 \otimes U_2$ to ρ is the equivalent of applying orthogonal rotation matrices O_1 and O_2 to \vec{r}^A , \vec{r}^B , and T, such that:

$$\vec{r}_f^A = O_1 \vec{r}^A \qquad \qquad \vec{r}_f^B = O_2 \vec{r}^B \qquad \qquad T_f = O_1 T O_2^{\dagger}$$
 (27)

Thus, given any ρ , we can find some $U_1 \otimes U_2$ such that $\rho' = U_1 \otimes U_2 \rho U_1^{\dagger} \otimes U_2^{\dagger}$ has diagonal T [28]. Further, by the following lemma, application of $U_1 \otimes U_2$ to ρ leaves $d_{\max}(\rho)$ invariant, implying we can assume without loss of generality that T is diagonal, as desired.

Lemma 2 For any state ρ , acting on $\mathcal{H}^M \otimes \mathcal{H}^N$, $d_{\max}(\rho)$ is invariant under local unitary operations applied to ρ .

Proof. Let U_1 and U_2 be arbitrary unitary operations acting on subsystems A and B, respectively. Then, straightforward manipulation of Eq. (5) yields $d(U_1 \otimes U_2 \rho U_1^{\dagger} \otimes U_2^{\dagger}, U^B) =$ $\mathrm{d}(\rho, U_2^{\dagger} U^B U_2)$. It is easy to see that $[U_2 \rho_B U_2^{\dagger}, U^B] = 0$ if and only if $[\rho_B, U_2^{\dagger} U^B U_2] = 0$. Thus, since the set of unitary matrices of fixed dimension forms a group, cycling through all possible choices of U^B in $d(\rho, U_2^{\dagger}U^BU_2)$ gives the desired result \square .

We can now derive a closed formula for $d_{max}(\rho)$. Henceforth, assuming that T is diagonal, denote $\lambda_i := T_{ii}$.

Lemma 3 Given a quantum state ρ , acting on $\mathcal{H}^2 \otimes \mathcal{H}^2$, with diagonal correlation matrix T, we have

$$d_{\max}(\rho) = \frac{1}{\sqrt{2}} \sqrt{\sum_{i=0}^{2} \lambda_i^2 (1 - n_i^2)}.$$
 (28)

Here, if $\rho_B \neq I/2$, then $\vec{n} := \vec{r}_B / \|\vec{r}_B\|_2$, and otherwise $n_i := 1$ for $\lambda_i = \min_k \lambda_k$ (with $\vec{n} := (n_0, n_1, n_2)^T$ and $\|\vec{n}\|_2 = 1$).

Proof. Assume first that $\rho_B \neq I/2$. We shall manipulate Eq. (8) to achieve the claimed form. Specifically, let U^B be a unitary operation corresponding to a rotation of angle $\theta \in [0, 2\pi)$ about axis $\vec{n} = (n_0, n_1, n_2)^T$, with $||\vec{n}||_2 = 1$ (θ and \vec{n} to be chosen as needed). We can characterize T^f in terms of T and U^B by applying Eq. (27) for $U_1 = I$ and $U_2 = U^B$, and

^bThis holds in the two-qubit case due to the existence of a surjective homomorphism from SU(2) to SO(3).

utilizing the following formula for O_2 in terms of U^B [29]:

$$O_{2} = I + \sin \theta A + (1 - \cos \theta) A^{2}, \text{ where}$$

$$A = \begin{pmatrix} 0 & -n_{2} & n_{1} \\ n_{2} & 0 & -n_{0} \\ -n_{1} & n_{0} & 0 \end{pmatrix}.$$
(29)

This simplifies Eq. (8) to:

$$d(\rho, U^B) = \frac{1}{2} \sqrt{\sum_{i=0}^{2} \lambda_i^2 (1 - \cos \theta) (1 - n_i^2)}.$$
 (30)

To choose \vec{n} , observe that demanding $[U^B, \rho_B] = 0$ requires U^B to induce a rotation about the Bloch vector of ρ_B (unless $U^B = I$ or $\vec{r}_B = \vec{0}$). Thus, set $\vec{n} = \vec{r}_B / \|\vec{r}_B\|_2$. Since $\|\vec{n}\|_2 = 1$, we have $(1 - n_i^2) \ge 0$ for all i, and so the expression above is maximized for $\cos \theta = -1$, or $\theta = \pi$, giving the desired result.

Finally, if $\rho_B = I/2$, one can choose any axis of rotation \vec{n} . By Eq. (30), choosing $n_i = 1$ for $\lambda_i = \min_k \lambda_k$ is the optimal choice, as claimed \square .

Lemma 3 confirms that for any pure state $|\psi\rangle = a_0|00\rangle + a_1|11\rangle$ with $|a_0|^2 + |a_1|^2 = 1$, $d_{\max}(\psi) > 0$ if and only if $|\psi\rangle$ violates the CHSH inequality [10]. To see this, note that for $\rho = |\psi\rangle\langle\psi|$, T is diagonal with $T_{00} = 2a_0a_1$, $T_{11} = -2a_0a_1$, $T_{22} = 1$, and $\vec{r}_B = (0, 0, a_0^2 - a_1^2)^T$. Eq. (28) hence reduces to Eq. (12), yielding $d_{\max}(\psi) = 2|a_0a_1|$. The maximum violation of the CHSH inequality for a pure state is given by [30]

$$B_{\text{max}}(\psi) = 2\sqrt{1 + 4|a_0 a_1|^2} = 2\sqrt{1 + d_{\text{max}}^2(\psi)},$$
 (31)

where the inequality is violated if and only if $B_{\text{max}}(\psi) > 2$. The claim immediately follows. We now show the main results of this section.

Theorem 4 Given a quantum state ρ , acting on $\mathcal{H}^2 \otimes \mathcal{H}^2$, with diagonal correlation matrix T, the implication $d_{\max}(\rho) > 1/\sqrt{2} \Rightarrow M(\rho) > 1$ holds. The converse is not true, i.e. $M(\rho) > 1 \neq d_{\max}(\rho) > 1/\sqrt{2}$.

Proof. Assume first that $d_{\max}(\rho) > 1/\sqrt{2}$. Without loss of generality, let $|\lambda_0| \ge |\lambda_1| \ge |\lambda_2|$, and let $\vec{n} = (\sqrt{\epsilon_0}, \sqrt{\epsilon_1}, \sqrt{\epsilon_2})^T$, where $\epsilon_0 + \epsilon_1 + \epsilon_2 = 1$. Then, substitution into Eq. (28) gives

$$(1 - \epsilon_0)\lambda_0^2 + (1 - \epsilon_1)\lambda_1^2 + (1 - \epsilon_2)\lambda_2^2 > 1,$$
(32)

from which it follows that $M(\rho) = \lambda_0^2 + \lambda_1^2 > 1$, since setting $\epsilon_2 = 1$ can only increase the left hand side of Eq. (32).

To show that the converse does not hold, consider the following counterexample: Given a pure state $|\psi\rangle=a_0|00\rangle+a_1|11\rangle$ (with real coefficients a_0 and a_1 , normalized via $a_0^2+a_1^2=1$), for all $a_0\leq 0.3827$ or $a_0\geq 0.9239$, we have $M(\rho)>1$, but $\mathrm{d_{max}}(\rho)\leq 1/\sqrt{2}$ \square .

Theorem 4 implies that the Fu distance is generally a weaker entanglement criterion (at least in the two-qubit case) than the CHSH inequality. We next ask if there are specific classes

of two-qubit states for which the Fu distance is "equivalent" to the CHSH inequality, in the sense that $M(\rho) > 1 \Leftrightarrow d_{\max}(\rho) > \frac{1}{\sqrt{2}}$ holds? It turns out that this is indeed the case, as we will show now.

Theorem 5 Given a quantum state ρ , acting on $\mathcal{H}^2 \otimes \mathcal{H}^2$, with diagonal correlation matrix T and its entries λ_i , with i = 0, 1, 2, consider the following conditions:

- 1. $\lambda_i = \min_k |\lambda_k|$, and $|n_i| = 1$, where $\vec{n} := \vec{r}_B / ||\vec{r}_B||_2$, and $\vec{r}_B \neq (0,0,0)^T$.
- 2. $|\lambda_0| = |\lambda_1| = |\lambda_2|$.
- 3. $\rho_B = I/2$.

Then, $M(\rho) > 1 \Leftrightarrow d_{\max}(\rho) > \frac{1}{\sqrt{2}}$ if and only if one of the above conditions holds.

Proof. We proceed case by case.

1. Suppose without loss of generality that $\lambda_2 = \min_k |\lambda_k|$, and $\vec{n} = (0,0,1)^T$. Then, Eq. (28) simplifies to

$$d_{\text{max}}(\rho) = \frac{1}{\sqrt{2}} \sqrt{\lambda_0^2 + \lambda_1^2} = \frac{1}{\sqrt{2}} \sqrt{M(\rho)},$$
 (33)

from which we have $M(\rho) > 1 \Leftrightarrow d_{\max}(\rho) > 1/\sqrt{2}$.

2. Suppose $|\lambda_0| = |\lambda_1| = |\lambda_2|$. Then, since $||\vec{n}||_2 = 1$, Eq. (28) simplifies to:

$$d_{\max}(\rho) = \frac{1}{\sqrt{2}} \sqrt{\lambda_0^2 (3 - n_1^2 - n_2^2 - n_3^2)} = \frac{1}{\sqrt{2}} \sqrt{\lambda_0^2 + \lambda_1^2} = \frac{1}{\sqrt{2}} \sqrt{M(\rho)}, \quad (34)$$

and we arrive at the same conclusion as in Case 1.

3. Suppose $\rho_B = I/2$. Then by Lemma 3, it straightforwardly follows that we are reduced to Case 1.

Finally, in order to show that the demonstrated equivalence holds if and only if one of these conditions hold, assume without loss of generality that $|\lambda_0| \ge |\lambda_1| \ge |\lambda_2|$. Then, unless $|\lambda_0| = |\lambda_1| = |\lambda_2|$ (Case 2), the only way to guarantee the equivalence is to have in Eq. (28) the equalities $(1-r_0^2)=1$ and $(1-r_1^2)=1$, which implies that $\vec{r}_B=(0,0,\pm 1)^T$. But such a choice of \vec{r}_B can only correspond to a cyclic unitary operation if we have Case 1 or 3 above, as required \square .

Thus, there exist certain classes of two-qubit states for which the CHSH inequality and $d_{max}(\rho)$ are equally capable of detecting entanglement. Specifically, note that the Werner state ρ_W that we considered at the start of this section falls into such a class, since $\rho_B = \frac{I}{2}$ for ρ_W , explaining the observed coincidence.

Fu Distances for some Bound Entangled States

Let us now investigate $d_{max}(\rho_{BE})$ for three distinct constructions of bound entangled (BE) states of two qutrits in order to determine whether $d_{max}(\rho_{BE})$ can be used to detect bound entanglement. Throughout this section, we denote the computational basis for qutrits as $\{|0\rangle, |1\rangle, |2\rangle\}.$

5.1 P. Horodecki Construction

Denote by $P_{\psi} = |\psi\rangle\langle\psi|$ the projector onto a state $|\psi\rangle$, and define [31]:

$$Q = I \otimes I - \left(\sum_{k=0}^{2} P_k \otimes P_k\right) - P_2 \otimes P_0, \tag{35}$$

$$|\Psi\rangle = \frac{1}{\sqrt{3}} (|0\rangle \otimes |0\rangle + |1\rangle \otimes |1\rangle + |2\rangle \otimes |2\rangle),$$
 (36)

$$\rho_{\text{ent}} = \frac{3}{8} P_{\Psi} + \frac{1}{8} Q,$$
(37)

$$|\Phi_a\rangle = |2\rangle \otimes \left(\sqrt{\frac{1+a}{2}}|0\rangle + \sqrt{\frac{1-a}{2}}|2\rangle\right),$$
 (38)

$$\rho_a = \frac{8a}{8a+1}\rho_{\text{ent}} + \frac{1}{8a+1}P_{\Phi_a}, \tag{39}$$

for $0 \le a \le 1$. Note that ρ_{ent} is entangled, as its partial transpose has a negative eigenvalue, and P_{Φ_a} is separable, since $|\Phi_a\rangle$ is a product state. The state of interest, ρ_a , is bound entangled for 0 < a < 1 [31]. Let us now determine $d(\rho_a, U^B)$ in terms of a.

Theorem 6 Given ρ_a , such that 0 < a < 1, the maximal Fu distance is $d_{\max}(\rho_a) = 2\sqrt{2}a/(8a+1)$. It is obtained, for example, by using any diagonal unitary matrix $U^B \in \mathbb{C}^{3\times 3}$ with $U^B_{0,0} = -U^B_{1,1} = U^B_{2,2}$.

Proof. Let U^B be an arbitrary complex 3×3 matrix, i.e.

$$U^{B} = \begin{pmatrix} u_{1} & u_{2} & u_{3} \\ u_{4} & u_{5} & u_{6} \\ u_{7} & u_{8} & u_{9} \end{pmatrix}. \tag{40}$$

Defining $\gamma := \sqrt{1 - a^2}$, we have

$$[\rho_B, U^B] = \frac{1}{2(8a+1)} \begin{pmatrix} \gamma(u_7 - u_3) & (1-a)u_2 + \gamma u_8 & \gamma(u_9 - u_1) \\ u_4(a-1) - \gamma u_6 & 0 & u_6(a-1) - \gamma u_4 \\ \gamma(u_1 - u_9) & (1-a)u_8 + \gamma u_2 & \gamma(u_3 - u_7) \end{pmatrix}.$$
(41)

Since 0 < a < 1, it is easy to see that for U^B to be cyclic it must therefore be of the form:

$$U^{B} = \begin{pmatrix} u_{1} & 0 & u_{3} \\ 0 & u_{5} & 0 \\ u_{3} & 0 & u_{1} \end{pmatrix}. \tag{42}$$

Inserting this into Eq. (5) and enforcing unitary constraints on U^B leads to (where Re(x) denotes the real part of $x \in \mathbb{C}$)

$$d(\rho_a, U^B) = \frac{a}{8a+1} \sqrt{6 - 2(|u_1|^2 + 2\operatorname{Re}(u_1^* u_5))}.$$
 (43)

To maximize $d(\rho_a, U^B)$, we need to minimize $|u_1|^2 + 2\operatorname{Re}(u_1^*u_5)$. To do so, set $u_5 = -1$ and let $u_1 = re^{i\theta}$ for $0 \le r \le 1$, $\theta \in [0, 2\pi)$. Then,

$$|u_1|^2 + 2\operatorname{Re}(u_1^*u_5) = r^2 - 2r\cos(\theta),\tag{44}$$

which achieves a minimum at $\theta = 0$ and r = 1, or equivalently for $u_1 = 1$ and $u_5 = -1$, yielding $d_{\text{max}}(\rho_a) = (2\sqrt{2}a)/(8a+1)$, as claimed. It is easy to see that any diagonal U^B with entries $U_{0,0}^B = -U_{1,1}^B = U_{2,2}^B$ gives the same optimum value \square .

The limiting value for $d_{\max}(\rho_a)$ as a approaches 1 is $\frac{2\sqrt{2}}{9}$. By embedding a two-qubit state in a higher-dimensional space, one finds that the value $d_{max}(\rho_{cc}) = 1/\sqrt{2}$ can be reached for a classically correlated state in any dimension. Thus, we conclude that by using $d_{max}(\rho_a)$. one cannot detect entanglement in bound entangled states of the above construction.

$Horodecki^{\otimes 3}$ Construction

Consider now a second bipartite one-parameter qutrit bound entangled class of states due to Paweł, Michał, and Ryszard Horodecki [32]. Define, for $2 \le \alpha \le 5$:

$$\sigma_{+} = \frac{1}{3}(|01\rangle\langle 01| + |12\rangle\langle 12| + |20\rangle\langle 20|),$$
 (45)

$$\sigma_{-} = \frac{1}{3}(|10\rangle\langle10| + |21\rangle\langle21| + |02\rangle\langle02|),$$
 (46)

$$\sigma_{+} = \frac{1}{3}(|01\rangle\langle01| + |12\rangle\langle12| + |20\rangle\langle20|),$$

$$\sigma_{-} = \frac{1}{3}(|10\rangle\langle10| + |21\rangle\langle21| + |02\rangle\langle02|),$$

$$\rho_{\alpha} = \frac{2}{7}|\phi_{+}\rangle\langle\phi_{+}| + \frac{\alpha}{7}\sigma_{+} + \frac{5-\alpha}{7}\sigma_{-},$$
(45)

where $|\phi_{+}\rangle = \frac{1}{\sqrt{3}}(|00\rangle + |11\rangle + |22\rangle)$. The state of interest, ρ_{α} , is separable for $\alpha \in [2,3]$, bound entangled for $\alpha \in (3,4]$, and free entangled for $\alpha \in (4,5]$. Determining an analytical form for $d_{max}(\rho_{\alpha})$ proves difficult, but if one is promised that the input state ρ is of the form ρ_{α} , but does not know α , then choosing any U^{B} with an all-zero diagonal (observing that $\operatorname{Tr}_A(\rho_\alpha) = I/3$) gives

$$d(\rho_{\alpha}, U^B) = \frac{1}{7}\sqrt{\alpha^2 - 5\alpha + 9}.$$
(48)

Straightforward calculation shows that the range of $d(\rho_{\alpha}, U^B)$ is disjoint for domains $\alpha \in [2,3], \alpha \in (3,4],$ and $\alpha \in (4,5],$ and so one can distinguish between all three cases using the Fu distance. This is illustrated in Figure 3. If one does not know that ρ is of the form ρ_{α} , on the other hand, we find via Theorem 3 that

$$d_{\max}(\rho_{\alpha}) \le \frac{2\sqrt{3\alpha^2 - 15\alpha + 31}}{21},\tag{49}$$

which is also plotted in Figure 3. For $\alpha \in (3,4]$, this gives $d_{\max}(\rho_{\alpha}) \leq 2\sqrt{19}/21 \approx 0.415$, and for $\alpha \in (4,5]$, we have $d_{\max}(\rho_{\alpha}) \leq 2\sqrt{31}/21 \approx 0.530$, and so in both cases we cannot detect bound entanglement using $d_{max}(\rho)$.

Unextendible Product Bases Construction

We next consider the construction of Bennett et al. [33], which requires the following definition.

Definition 1 (Unextendible Product Basis (UPB) [33]) Consider a bipartite quantum system in $\mathcal{H}=\mathcal{H}_1\otimes\mathcal{H}_2$ with subsystems of arbitrary dimension. Define an incomplete orthogonal product basis (PB) as a set S of pure orthogonal product states spanning a proper subspace \mathcal{H}_S of \mathcal{H} . Then an unextendible product basis (UPB) is a PB whose complementary subspace $\mathcal{H} - \mathcal{H}_{\mathcal{S}}$ contains no product state.

^cThis definition extends straightforwardly to the multipartite case [33].

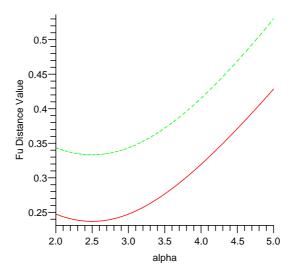


Fig. 3. Plots of the Fu distance for the Horodecki et al. bound entangled state [32]. The bottom solid line depicts $d(\rho_{\alpha}, U^B)$ for U^B with an all-zero diagonal (see Eq. (48)). The top dashed line plots the bound on $d_{\max}(\rho_{\alpha})$ given by Theorem 3 (see Eq. (49)).

Using a UPB, one can systematically construct BE states using the following theorem.

Theorem 7 (Bennett et al. [33]) Given UPB $\{|\psi_i\rangle\}_{i=0}^{n-1}$ in a Hilbert space of total dimension D, the following state is bound entangled:

$$\rho = \frac{1}{D-n} \left(I - \sum_{k=0}^{n-1} |\psi_k\rangle\!\langle\psi_k| \right). \tag{50}$$

The UPB we shall use with Theorem 7 is the following, given by [33]:

$$|\psi_0\rangle = \frac{1}{\sqrt{2}}|0\rangle(|0\rangle - |1\rangle),$$
 (51)

$$|\psi_1\rangle = \frac{1}{\sqrt{2}}(|0\rangle - |1\rangle)|2\rangle,$$
 (52)

$$|\psi_2\rangle = \frac{1}{\sqrt{2}}|2\rangle(|1\rangle - |2\rangle),$$
 (53)

$$|\psi_3\rangle = \frac{1}{\sqrt{2}}(|1\rangle - |2\rangle)|0\rangle,$$
 (54)

$$|\psi_4\rangle = \frac{1}{3}(|0\rangle + |1\rangle + |2\rangle)(|0\rangle + |1\rangle + |2\rangle),$$
 (55)

$$\rho = \frac{1}{4} \left(I - \sum_{k=0}^{4} |\psi_k\rangle\!\langle\psi_k| \right). \tag{56}$$

Using Theorem 3, we find that for ρ ,

$$d_{\max}(\rho) \le \sqrt{\frac{2n}{D(D-n)}}.$$
(57)

It thus follows that $d_{\max}(\rho) \leq 1/\sqrt{2}$ for any UPB for which $n \leq D^2/(D+4)$. Specifically, this rules out the possibility of detecting bound entanglement with the UPB chosen above, for which D=9 and n=5, yielding $d_{\max}(\rho) \leq \sqrt{10}/6 \approx 0.527$.

6 Conclusions and Open Problems

We have investigated locally noneffective unitary operations in connection with the detection of quantum entanglement. Specifically, we have derived and discussed closed formulae for $d_{max}(\rho)$ (the maximal distance between the original state ρ and the state after a locally noneffective unitary operation) for the bipartite cases of (pseudo)pure quantum states, Werner states, and two-qubit states. The first of these reveals the existence of non-maximally entangled states capable of achieving a maximal global shift. Thus, no entanglement measure based on $d_{max}(\rho)$ can be defined, and similarities to anomalies seen in non-locality measures are revealed. Since $d_{max}(\rho)$ is neither an entanglement measure, nor a non-locality measure, yet as demonstrated here possesses clear connections to the CHSH inequality in the two-qubit case, it would be interesting to have a better intuitive understanding of the correlations (both classical and quantum) that $d_{max}(\rho)$ is quantifying, and if and how the anomalies mentioned above are related to those seen in non-locality measures.

There are a number of questions which remain open. First, despite the fact that our formula for $d_{max}(\rho)$ for Werner states demonstrates diminishing distinguishability between classical and quantum correlations in Werner states as the dimension grows, it remains for a tight upper bound on $d_{max}(\rho)$ to be found for classically correlated states of total dimension D>4 in order to conclusively state the efficacy of $d_{max}(\rho)$ as an entanglement detection criterion. Second, although we have demonstrated that any entangled pseudopure or Werner state achieves $d_{max}(\rho)>0$, it is still not known whether this holds for all entangled bipartite states. Third, it would be of interest to determine whether a closed formula for $d_{max}(\rho)$ can be derived for mixed states in general, the existence of which would not contradict known hardness results for the quantum separability problem [3, 4, 5]. Finally, as mentioned briefly in Section 1, the principle behind the Fu distance is implicitly applied in superdense coding, and we would be curious to know whether there exist any other applications in quantum computing and information.

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