

Generation of maximum spin entanglement induced by a cavity field in quantum-dot systems

Adam Miranowicz,^{1,2,3} Şahin K. Özdemir,^{1,2} Yu-xi Liu,² Masato Koashi,^{1,2} Nobuyuki Imoto,^{1,2,4,5} and Yoshiro Hirayama^{1,5}

¹CREST Research Team for Interacting Carrier Electronics, Japan Science and Technology Corporation, Tokyo, Japan

²School of Advanced Sciences, Graduate University for Advanced Studies (SOKEN), Hayama, Kanagawa 240-0193, Japan

³Nonlinear Optics Division, Institute of Physics, Adam Mickiewicz University, 61-614 Poznań, Poland

⁴University of Tokyo, 7-3-1 Hongo, Bunkyo-ku, Tokyo 113-8654, Japan

⁵NTT Basic Research Laboratories, 3-1 Morinosato Wakamiya, Atsugi, Kanagawa 243-0198, Japan

(Received 5 March 2002; published 19 June 2002)

Equivalent-neighbor interactions of the conduction-band electron spins of quantum dots in the model of Imamoğlu *et al.* [Phys. Rev. Lett. **83**, 4204 (1999)] are analyzed. An analytical solution and its Schmidt decomposition are found and applied to evaluate how much the initially excited dots can be entangled with the remaining dots if all of them are initially disentangled. It is demonstrated that perfect maximally entangled states (MES's) can only be generated in systems of up to six dots with a single dot initially excited. It is also shown that highly entangled states, approximating the MES's with good accuracy, can still be generated in systems of odd numbers of dots with almost half of them excited. A sudden decrease of entanglement is observed on increasing the total number of dots in a system with a fixed number of excitations.

DOI: 10.1103/PhysRevA.65.062321

PACS number(s): 03.65.Ud, 03.67.-a, 68.65.Hb, 75.10.Jm

I. INTRODUCTION

Since the seminal papers of Obermayer, Teich, and Mahler [1], there has been growing interest in the quantum-information properties of quantum dots (QDs) in the quest to implement quantum-dot scalable quantum computers [2–4]. Those high expectations are justified to some extent by recent experimental advances in the coherent observation and manipulation of quantum dots [5,6], including spectacular demonstrations of the quantum entanglement of excitons in a single dot [7] or quantum-dot molecule [8], and observations of Rabi oscillations of excitons in single dots [9]. Among various models of quantum computers based on localized electron spins of quantum dots as qubits [3,4], the scheme of Imamoğlu *et al.* [10] is the first where the interactions between the qubits are mediated by a cavity field. This approach combines the advantages of long-distance optically controlled couplings with long-decoherence times of the spin degrees of freedom. Here, we analyze quantum entanglement in the Imamoğlu *et al.* model.

During the last decade, it has been highlighted that quantum entanglement, being at the heart of quantum mechanics, is also a powerful resource for quantum communication and quantum-information processing. Quantum entanglement in interacting systems is a common phenomenon. It is obvious that any interacting many-body system with defined qubits, if set in a properly chosen state, will evolve through states with entangled qubits. Surprisingly, quantitative descriptions of the entanglement dynamics in multiparticle systems are by no means satisfactory yet [11]. Nevertheless, in a special case of bipartite entanglement, a number of measures have been introduced and studied [12–14]. For example, entanglement of a bipartite system in a pure state, described by the density matrix $\hat{\rho}_{AB} = (|\psi\rangle\langle\psi|)_{AB}$, can be measured by the von Neumann entropy [12,13]

$$E[\hat{\rho}_{AB}] = -\text{Tr}\{\hat{\rho}_A \log_2 \hat{\rho}_A\} = -\text{Tr}\{\hat{\rho}_B \log_2 \hat{\rho}_B\} \quad (1)$$

of the reduced density matrix $\hat{\rho}_A = \text{Tr}_B\{\hat{\rho}_{AB}\}$ or, equivalently, $\hat{\rho}_B = \text{Tr}_A\{\hat{\rho}_{AB}\}$. The entanglement of formation of a mixed state of a bipartite system is often measured by the so-called concurrence proposed by Hill and Wootters [14]. Concurrence has been applied to study entanglement in various models [15] including equivalent-neighbor systems [16,17]. The following two aspects of entanglement are especially important: (i) coherent manipulation of entanglement and (ii) generation of maximum entanglement. The possibility of coherent and selective control of entanglement in a quantum-dot system was analyzed by Imamoğlu *et al.* [10]. Here, we would like to focus on the latter topic, i.e., the generation of the maximally entangled states (MES's) of quantum dots in the model of Imamoğlu *et al.* [10]. MES's are necessary for the majority of quantum information-processing applications. Otherwise, for example, direct application of partly entangled states for teleportation will result in unfaithful transmission, while superdense coding with partly entangled states will cause noise in the resulting classical channel.

The paper is organized as follows. In Sec. II, we describe an equivalent-neighbor quantum-dot model and give its analytical solution. In Sec. III, we analyze the possibilities of generation of the MES's or their good approximations for different initial conditions of the number of excitations and the total number of dots in the system.

II. QUANTUM-DOT MODEL AND ITS SOLUTION

We will apply the model of Imamoğlu *et al.* [10] to describe strong equivalent-neighbor couplings of quantum-dot spins through a single-mode microcavity field. The dots are placed inside a microdisk, put into a microcavity tuned to frequency ω_{cav} , and illuminated selectively by laser fields of frequencies $\omega_n^{(L)}$. Each of N dots with a single electron in the conduction band is modeled by a three-level atom as shown in Fig. 1. The total Hamiltonian for N three-level quantum dots interacting with $N+1$ quantized fields reads

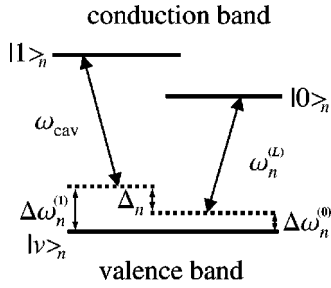


FIG. 1. Three-level atom in V configuration as a model of a semiconductor dot with the conduction-band spin states $|1\rangle_n$ (spin up) of energy $\mathcal{E}_n^{(1)}$ and $|0\rangle_n$ (spin down) of energy $\mathcal{E}_n^{(0)}$, and the effective valence-band state $|v\rangle_n$ of energy $\mathcal{E}_n^{(v)}$ in the n th dot. Key: ω_{cav} , frequency of the common cavity mode; $\omega_n^{(L)}$, frequency of the classical laser field addressed at the n th dot; $\hbar\Delta\omega_n^{(1)} = \mathcal{E}_n^{(1)} - \mathcal{E}_n^{(v)} - \hbar\omega_{\text{cav}}$, $\hbar\Delta\omega_n^{(0)} = \mathcal{E}_n^{(0)} - \mathcal{E}_n^{(v)} - \hbar\omega_n^{(L)}$, and $\Delta_n = \Delta\omega_n^{(1)} - \Delta\omega_n^{(0)}$ are detunings.

$$\hat{H} = \hat{H}_{QD} + \hat{H}_F + \hat{H}_{\text{int}}, \quad (2)$$

$$\hat{H}_{QD} = \sum_n (\mathcal{E}_n^{(0)} \hat{\sigma}_n^{00} + \mathcal{E}_n^{(1)} \hat{\sigma}_n^{11} + \mathcal{E}_n^{(v)} \hat{\sigma}_n^{vv}),$$

$$\hat{H}_F = \hbar\omega_{\text{cav}} \hat{a}_{\text{cav}}^\dagger \hat{a}_{\text{cav}} + \sum_n \hbar\omega_n^{(L)} (\hat{a}_n^{(L)})^\dagger \hat{a}_n^{(L)},$$

$$\hat{H}_{\text{int}} = \sum_n \hbar g_n^{v0} [\hat{a}_n^{(L)} \hat{\sigma}_n^{0v} + (\hat{a}_n^{(L)})^\dagger \hat{\sigma}_n^{v0}] + \sum_n \hbar g_n^{v1} (\hat{a}_{\text{cav}} \hat{\sigma}_n^{1v} + \hat{a}_{\text{cav}}^\dagger \hat{\sigma}_n^{v1}),$$

where \hat{H}_{QD} and \hat{H}_F are the free Hamiltonians of the quantum dots and the fields, respectively; \hat{H}_{int} is the interaction Hamiltonian; \hat{a}_{cav} and $\hat{a}_{\text{cav}}^\dagger$ are the annihilation and creation operators of the cavity mode, respectively; $\hat{a}_n^{(L)}$ and $(\hat{a}_n^{(L)})^\dagger$ are the corresponding operators for the laser modes; $\hat{\sigma}_n^{xy}$ is the n th dot operator given by $\hat{\sigma}_n^{xy} = |x\rangle_n \langle y|$; $\mathcal{E}_n^{(x)}$ is the energy of level $|x\rangle_n$ ($x=0,1,v$); the n th dot levels $|0\rangle_n$ and $|v\rangle_n$ are coupled by dipole interactions with a strength of g_n^{v0} ; analogously, g_n^{v1} is the coupling strength between levels $|1\rangle_n$ and $|v\rangle_n$. There is no direct coupling between levels $|0\rangle_n$ and $|1\rangle_m$ in either the same ($n=m$) or different dots ($n \neq m$). The Hamiltonian (2) simply generalizes, to N dots and $N+1$ fields, models of a three-level atom (dot) interacting with two modes of radiation fields widely discussed in the literature (see, e.g., [19]). By applying an adiabatic elimination method, Imamoglu *et al.* derived the effective interaction Hamiltonian describing the evolution of the conduction-band spins of N quantum dots coupled by a microcavity field in the form [10]

$$\hat{H}_{\text{eff}} = \frac{\hbar}{2} \sum_{n \neq m} \kappa_{nm}(t) [\hat{\sigma}_n^+ \hat{\sigma}_m^- e^{i(\Delta_n - \Delta_m)t} + \hat{\sigma}_n^- \hat{\sigma}_m^+ e^{-i(\Delta_n - \Delta_m)t}] \quad (3)$$

in terms of the Pauli spin creation $\hat{\sigma}_n^+$ and annihilation $\hat{\sigma}_n^-$ operators acting on the conduction-band spin states of the n th dot. The effective two-dot coupling strength between the spins of the n th and m th dots is given by $\kappa_{nm}(t) = g_n(t)g_m(t)/\Delta_n$, where the effective single-dot coupling of the n th spin to the cavity field is $g_n(t) = g_n^{v0}g_n^{v1}|E_n^{(L)}(t)|/\Delta\omega_n$ with $\Delta\omega_n$ being the harmonic mean of $\Delta\omega_n^{(1)}$ and $\Delta\omega_n^{(0)}$. For simplicity, the laser fields are assumed to be strong and treated classically as described by the complex amplitudes $E_n^{(L)}(t)$. The Hamiltonian (3) was derived by applying adiabatic eliminations of the valence-band states $|v\rangle_n$ and cavity mode \hat{a}_{cav} , which are valid under the assumptions of negligible coupling strength, cavity decay rate, and thermal fluctuations in comparison to $\hbar\Delta_n$ and $\hbar\Delta\omega_n^{(x)}$ ($x=0,1$) and the energy difference $\mathcal{E}_n^{(1)} - \mathcal{E}_n^{(0)}$ (see Fig. 1). Moreover, the valence-band levels $|v\rangle_n$ were assumed to be far off resonance. Although the Hamiltonian (3) describes apparently direct spin-spin interactions, the real physical picture is different: Quantum-dot spins are coupled only indirectly via the cavity and laser fields.

Imamoglu *et al.* [10] applied their model for quantum computing purposes by implementing the conditional phase-flip and controlled-NOT (CNOT) operations between two arbitrary dots addressed selectively by laser fields to satisfy the condition $\Delta_n = \Delta_m$. Here, we are interested in a realization of an equivalent-neighbor model scalable for a large number of dots (even for more than 100 [10]). This goal can readily be achieved by assuming that all dots are identical and illuminated by a single-mode stationary laser field of frequency $\omega_n^{(L)} \equiv \omega^{(L)}$, which implies $\kappa_{nm}(t) = \kappa = \text{const}$. In fact, the condition of equivalent-neighbor interactions can also be assured for nonidentical dots by adjusting the laser-field frequencies $\omega_n^{(L)}$ to get the same detuning $\Delta_n = \text{const}$, and by choosing the proper laser intensities $|E_n^{(L)}|^2$ to obtain the effective coupling constants of $g_n(t) = \text{const}$ or, equivalently, $\kappa_{nm}(t) = \text{const}$ for every pair of dots. Thus, Eq. (3) can be reduced to the effective equivalent-neighbor N -dot Hamiltonian as

$$\hat{H}_{\text{eff}} = \frac{\hbar\kappa}{2} \sum_{n \neq m} (\hat{\sigma}_n^+ \hat{\sigma}_m^- + \hat{\sigma}_n^- \hat{\sigma}_m^+), \quad (4)$$

where κ is the coupling constant. The system described by Eq. (4) is sometimes referred to as the spin-1/2 van der Waals model [20], the infinitely coordinated system [21], the Lipkin or Lipkin-Meshkov-Glick model [22], or just the equivalent-neighbor model [23]. Let us assume that the initial state describing a system of M ($M=0, \dots, N$) dots initially excited (i.e., with conduction-band spins up) and $N-M$ dots in the ground state (conduction-band spins down) is given as

$$\begin{aligned} |\psi(0)\rangle &= \{|1\rangle_A^{\otimes M}\} \{|0\rangle_B^{\otimes (N-M)}\} \\ &\equiv \underbrace{|11 \dots 1\rangle}_M \underbrace{|00 \dots 0\rangle}_{N-M}. \end{aligned} \quad (5)$$

Then, we find the solution of the Schrödinger equation of motion for the model (4) in the form

$$|\psi(t)\rangle = \sum_{m=0}^{M'} C_m^{NM}(t) \{ |1\rangle^{\otimes(M-m)} |0\rangle^{\otimes m} \}_{A} \otimes \{ |1\rangle^{\otimes m} |0\rangle^{\otimes(N-M-m)} \}_{B}, \quad (6)$$

where $M' = \min(M, N-M)$. The states in curly brackets $\{ |1\rangle^{\otimes(n-m)} |0\rangle^{\otimes m} \}$ denote the sum of all n -dot states with $(n-m)$ excitations. For example, $\{ |1\rangle^{\otimes 2} |0\rangle^{\otimes 2} \}$ stands for $|0011\rangle + |0101\rangle + |0110\rangle + |1001\rangle + |1010\rangle + |1100\rangle$. The number of states in the superposition $\{ |1\rangle^{\otimes(n-m)} |0\rangle^{\otimes m} \}$ (or equivalently $\{ |1\rangle^{\otimes m} |0\rangle^{\otimes(N-m)} \}$) is equal to the binomial coefficient $\binom{n}{m}$. Thus, for given N and M , the solution (6) contains $\binom{N}{M}$ terms. The energy of the QD system described by Eq. (4) is conserved; thus all the superposition states in Eq. (6) have the same number M of excitations. We find the time-dependent superposition coefficients in Eq. (6) as

$$C_m^{NM}(t) = \sum_{n=0}^{M'} b_{nm}^{NM} \exp\{i[n(N+1-n) - M(N-M)]\kappa t\} \quad (7)$$

in terms of

$$b_{nm}^{NM} = \sum_{k=0}^m (-1)^k \binom{m}{k} \binom{N-2k}{M-k}^{-1} \left[\binom{N+1-2k}{n-k} - 2 \binom{N-2k}{n-k-1} \right], \quad (8)$$

where $\binom{x}{y}$ are binomial coefficients. Our solution can be represented in a biorthogonal form via the Schmidt decomposition

$$|\psi(t)\rangle = \sum_{m=0}^{M'} \sqrt{P_m^{NM}(t)} |\phi_m(t)\rangle_A \otimes |\varphi_m(t)\rangle_B, \quad (9)$$

where $|\phi_m(t)\rangle_A$ and $|\varphi_m(t)\rangle_B$ are the orthonormal basis states of subsystems A and B , respectively. We find that the real and positive Schmidt coefficients can be related to the squared module of superposition coefficients (7) as follows:

$$P_m^{NM}(t) = \binom{M}{m} \binom{N-M}{m} |C_m^{NM}(t)|^2, \quad (10)$$

while the phases of $C_m^{NM}(t)$ are absorbed into the definitions of the basis states $|\phi_m(t)\rangle_A$ and $|\varphi_m(t)\rangle_B$. The Schmidt coefficients are normalized to unity. The evolutions of all P_m^{NM} for systems with single and two excitations are given in Figs. 2 and 3, respectively. We observe that the evolution of Schmidt coefficients is periodic with the period of $\kappa T = 2\pi/N$ for systems with a single ($M=1$ or, equivalently, $M=N-1$) excitation (Fig. 2), and π -periodic (2π periodic) for systems of even (odd) numbers of dots with higher numbers of excitations (see Fig. 3). For brevity, only half of the period is depicted in the right-hand panels of Fig. 3.

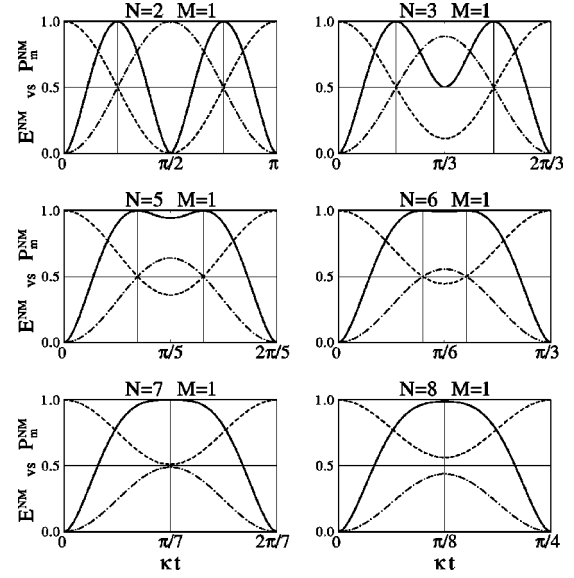


FIG. 2. Evolution of the quantum entanglement of $E^{N1}(t)$ (solid) and the Schmidt coefficients of $P_0^{N1}(t)$ (dashed) and $P_1^{N1}(t)$ (dot-dashed curves) for systems of $N=2, \dots, 8$ quantum dots with only one ($M=1$) of them initially excited. Figure illustrates that the exact maximally entangled states can be generated in systems of N up to 6 dots only.

III. ENTANGLEMENT IN QUANTUM-DOT SYSTEMS

We address the following questions: How much can the initially excited dots (say, subsystem A) be entangled with the remaining dots (subsystem B) in the equivalent-neighbor system of initially all disentangled dots if the evolution is governed by Hamiltonian (4)? And whether the maximally entangled states can be generated exactly or, at least, ap-

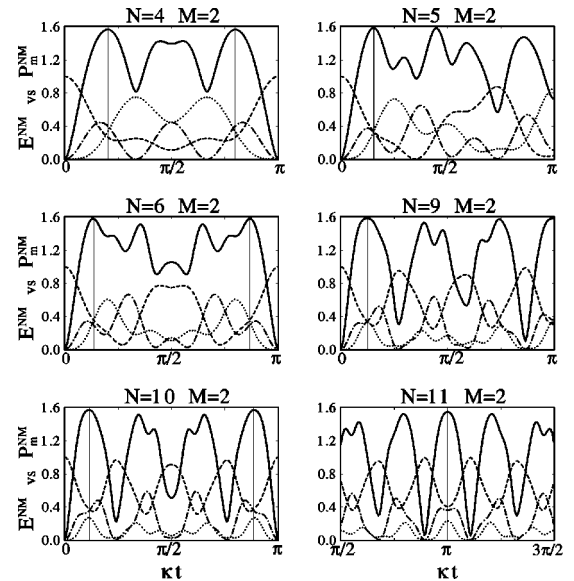


FIG. 3. Evolution of the entanglement of $E^{N2}(t)$ (solid) and all Schmidt coefficients: $P_0^{N2}(t)$ (dashed), $P_1^{N2}(t)$ (dot-dashed), and $P_2^{N2}(t)$ (dotted curves), in systems with two ($M=2$) dots initially excited.

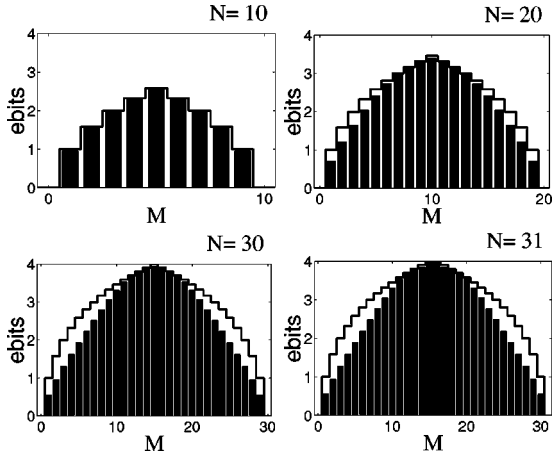


FIG. 4. Maximum entanglement $E_{\max}^{NM} = \max_t E^{NM}(t)$ (solid bars), measured in ebits, as a function of the excitation number M generated in systems of $N = 10, 20, 30,$ and 31 dots. The empty staircase corresponds to entanglement of E_{MES}^{NM} for the MES's. The figure illustrates that the highest entanglement, closest to E_{MES}^{NM} , can be generated in systems with $M = \lfloor N/2 \rfloor$ excitations. On decreasing M or $(N-M)$, the entanglement decreases. The discrepancy between E_{\max}^{NM} and E_{MES}^{NM} becomes more pronounced with increasing N especially for $0 < M \ll \lfloor N/2 \rfloor$.

proximately in systems of an arbitrary number N of dots while M of them are excited.

With the help of an explicit form of the Schmidt decomposition, it is convenient to calculate the entanglement (1) via the Shannon entropy

$$E^{NM}(t) \equiv E[|\psi(t)\rangle\langle\psi(t)|] = - \sum_{m=0}^{M'} P_m^{NM}(t) \log_2 P_m^{NM}(t) \quad (11)$$

of the Schmidt coefficients given for our system by Eq. (10). By applying Eq. (11), we can determine the maximum entanglement given by $E_{\max}^{NM}(t) \equiv \max_t E^{NM}(t)$, which can periodically be generated during the evolution of N -dot systems with M excitations. The coefficients (10), as well as (7), possess the symmetry of $P_m^{NM}(t) = P_m^{N, N-M}(t)$, which implies equal evolutions of entanglement

$$E^{NM}(t) = E^{N, N-M}(t) \quad (12)$$

in the N -dot system with M and $N-M$ excitations. Figure 4 shows this symmetry in a special case for maximum entanglement of $\max_t E^{NM}(t) = \max_t E^{N, N-M}(t)$.

To solve the second problem proposed at the beginning of this section, we have to determine the quantum correlations of the maximally entangled state of two subsystems having d equally weighted terms in its Schmidt decomposition. According to the theorem of Bennett *et al.* [12], the MES has $\log_2 d$ ebits of entanglement, where d is the Hilbert space dimension of the smaller subsystem. Thus, in our case, the MES of the subsystem A consisting of M dots and the subsystem B of $N-M$ dots has

$$E_{\text{MES}}^{NM} = \log_2[\min(M, N-M) + 1] \quad (13)$$

ebits of entanglement. In particular, the MES in the N -dot system with a single initial excitation has only 1 ebit independent of N . The empty staircase in Fig. 4 and solid lines in Fig. 5 correspond to E_{MES}^{NM} . To show a deviation of a given state from the MES, it is convenient to use the relative (or scaled) entanglement defined to be

$$e_{\max}^{NM} \equiv \frac{E_{\max}^{NM}}{E_{\text{MES}}^{NM}} = \max_t \frac{E^{NM}(t)}{E_{\text{MES}}^{NM}}. \quad (14)$$

In the simplest nontrivial case, for $M=1$, the Schmidt coefficients reduce to

$$P_1^{N1}(t) = 4 \frac{(N-1)}{N^2} \sin^2\left(\frac{N}{2} \kappa t\right) \quad (15)$$

and $P_0^{N1}(t) = 1 - P_1^{N1}(t)$, which enable a direct calculation of the entanglement $E^{N1}(t)$ with the help of Eq. (11). The evolutions of entanglement and the Schmidt coefficients of $P_m^{N1}(t)$ for $m=0,1$ are depicted in Fig. 2. The quantum-dot systems evolve into the MES's at evolution times, that are the roots of the equation

$$0 = \dot{E}^{N1}(t) = 2\kappa \frac{N-1}{N} \sin(N\kappa t) \times \log_2 \left[\frac{N^2}{4(N-1)} \csc^2\left(\frac{N}{2} \kappa t\right) - 1 \right]. \quad (16)$$

Thus, we get

$$\kappa t' = \frac{2}{N} \operatorname{arccsc} \left(\frac{2}{N} \sqrt{2(N-1)} \right) \quad (17)$$

and $\kappa t'' = \pi/N$. We find that the maximum entanglement, equal to $E^{N1}(t') = 1$ ebit, can be achieved at evolution times t' for $N \leq 6$ only. For $N > 6$, a real solution for t' does not exist. Another explanation of this result, as illustrated in Fig. 2, can be given as follows: The maximum entanglement corresponds to the Schmidt coefficients mutually equal or, in general, the least different. But the MES corresponds solely to the former case. As is seen in Fig. 2, the condition $P_0^{N1}(t') = P_1^{N1}(t')$ is strictly satisfied for $N \leq 6$. The entanglement for $N > 6$ reaches its maximum at the evolution times t'' . This maximum value is given by

$$E^{N1}(t'') = \frac{2}{N^2} \{ N^2 \log_2 N - (N-2)^2 \log_2(N-2) - 2(N-1) \log_2[4(N-1)] \}, \quad (18)$$

which is less than unity and monotonically decreases with increasing N as clearly illustrated in Figs. 5 and 6 for $M=1$. Thus, the perfect MES's cannot be generated in systems of $N > 6$ dots. Nevertheless, a good approximation of the MES can also be obtained for $N=7$. On the scale of Fig. 2, $\max_t E^{7,1}(t) = E^{7,1}(\pi/7) = 0.9997$ is close to unity since $P_0^{N1}(\pi/7)$ and $P_1^{N1}(\pi/7)$ are almost the same. It is worth noting that a critical value of $N=6$ was also found, although

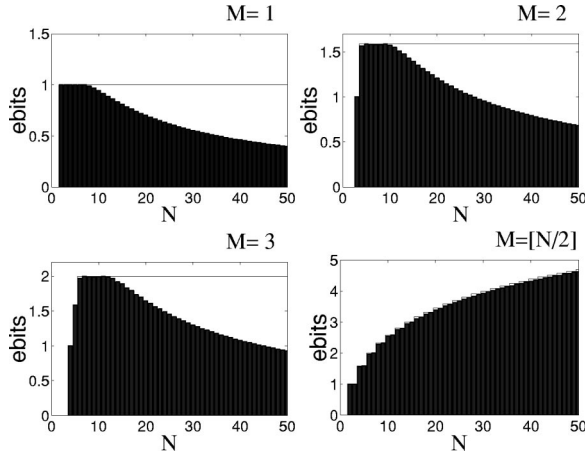


FIG. 5. Maximum entanglement E_{\max}^{NM} as a function of the total number N of dots generated in systems with $M=1, 2, 3$, and $[N/2]$ excitations. The solid lines and empty staircase correspond to E_{MES}^{NM} . On the scale of the figure, an apparent plateau occurs for N smaller than some critical value N_M . For N higher than N_M and fixed M , a monotonic decrease of the maximum entanglement is clearly visible. One concludes that arbitrary high entanglement can be achieved by increasing N and keeping half $M=[N/2]$ of the system excited.

in the different context of the pairwise entanglement measured by the concurrence [14], for an equivalent-neighbor model of entangled webs in Ref. [16]. In comparison, a critical value of $N=6$ for the concurrence in the equivalent-neighbor isotropic or anisotropic Heisenberg models was not observed (see, e.g., [17]). Similarly, generation of the MES in an equivalent-neighbor quantum-dot model of Reina *et al.* was discussed only in two special cases of the Bell ($N=2$) and Greenberger-Horne-Zeilinger (GHZ) ($N=3$) entangled states [18]. Thus, no critical behavior of entanglement as a function of N was reported there.

The case for $M=1$ is the only one where the general formula (10) for the Schmidt coefficients simplifies to a com-

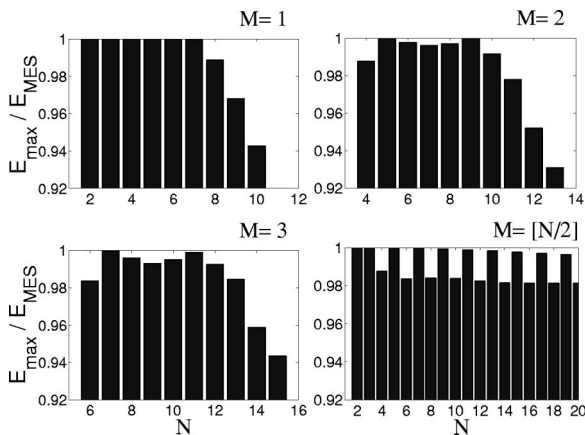


FIG. 6. The same as in Fig. 5 but for the relative maximum entanglement $e_{\max}^{NM} = E_{\max}^{NM} / E_{\text{MES}}^{NM}$. The figure shows that the apparent plateau for finite M actually occurs for $M=1$ only. The first and second highest maxima of entanglement correspond to N equal to $2M+1$ and $2M+5$, respectively.

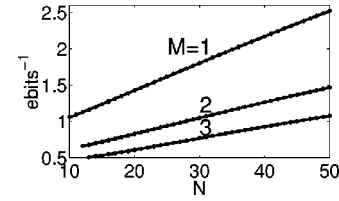


FIG. 7. The inverse of the maximum entanglement, $(E_{\max}^{NM})^{-1}$ (dots) measured in ebits^{-1} , and its approximation (solid lines) as a function of $N > 2M+5$ generated in systems with $M=1, 2, 3$ excitations.

pact form for arbitrary evolution times. Thus, for clarity, we present mainly numerical results for $M \geq 2$. For example, Fig. 3 illustrates that the exact MES cannot be generated in systems with $M=2$ excitations at any evolution time. This conclusion can be drawn from the observation that $P_m^{N2}(t)$ for $m=0, 1, 2$ do not cross simultaneously at any times in the period. Nevertheless, the MES can be approximated with good precision. The highest possible entanglement, corresponding to the least mutually different P_m^{N2} , is observed for $N=5$ and 9 , where the relative entanglement deviates from unity at the order of 10^{-5} and 10^{-4} , respectively (see Fig. 6 for $M=2$). The states generated in N -dot systems with three excitations can be entangled up to $e_{\max}^{7,3} = 0.9996$ (first) and $e_{\max}^{11,3} = 0.9990$ (second maximum) for the relative entanglement (see Fig. 6 for $M=3$). It is interesting to compare the relative entanglement of e_{\max}^{NM} , depicted in Fig. 6, with the “absolute” entanglement of E_{\max}^{NM} presented in Fig. 5. By analyzing the numerical data given, in part, in Fig. 6, we find the following rule: The maximally or almost maximally entangled states can be generated in systems of $N=2M+1$ dots with M excitations. Slightly worse entanglement can be achieved in systems of $N=2M+5$ dots with M excitations. Thus, systems composed of odd rather than even numbers of dots enable generation of the entangled states better approximating the MES for $M > 1$. This is clearly illustrated in Fig. 6 for $M=[N/2]$, i.e., the integer part of $N/2$. We observe that the system of odd and large numbers ($N > 2M+5$ for $M > 1$) of dots is the most entangled at the evolution times $\kappa t = (1+2k)\pi$ for $k=0, 1, \dots$ (see, e.g., Fig. 3 for $N=11$). In this special case, the Schmidt coefficients can be written compactly via

$$\left| C_m^{NM} \left(\frac{\pi}{\kappa} \right) \right| = 2^m m! (N-2M) \frac{(N-2m-2)!!}{N!!}. \quad (19)$$

For $\kappa t = k\pi$ and even N , in contrast to odd N , the entanglement vanishes. The maximum entanglement of E_{\max}^{NM} for $N > N_M \equiv 2M+5 - \delta_{1M}$ can be well fitted by the inverse of linear functions as shown in Fig. 7.

IV. CONCLUSION

We studied the evolution of the conduction-band spins of quantum dots in the model of Imamoglu *et al.* [10]. We found the analytical solution and its Schmidt decomposition for the equivalent-neighbor model and applied them in our study of bipartite entanglement in quantum-dot systems with

arbitrary numbers of dots and their excitations. We have raised and solved the problem to what extent the initially excited dots can be entangled with the remaining dots if all of them are initially disentangled in the equivalent-neighbor energy-conserving model. We have shown that the perfect maximally entangled states can only be generated in systems of $N=2, \dots, 6$ dots with a single dot initially excited. Nevertheless, highly entangled states, being excellent approximations of the MES's, can periodically be generated in systems of odd numbers N of dots with the number M of excitations equal to $M=(N-1)/2$ (leading to the best approximation) and $M=(N-5)/2$ (giving a slightly worse approximation). If we increase N beyond $N_M=2M+5$

– δ_{1M} , the entanglement decreases monotonically as described by the inverse of linear functions.

ACKNOWLEDGMENTS

We thank J. Bajer, T. Cheon, T. Kobayashi, H. Matsueda, and I. Tsutsui for their stimulating discussions. Y.L. acknowledges support from the Japan Society for the Promotion of Science (JSPS). This work was supported by a Grant-in-Aid for Scientific Research (B) (Grant No. 12440111) and a Grant-in-Aid for Encouragement of Young Scientists (Grant No. 12740243) by the Japan Society for the Promotion of Science.

-
- [1] K. Obermayer, W.G. Teich, and G. Mahler, *Phys. Rev. B* **37**, 8096 (1988); **37**, 8111 (1988).
- [2] A. Barenco *et al.*, *Phys. Rev. Lett.* **74**, 4083 (1995); H. Matsueda and S. Takeno, *IEICE Trans. Fundamentals* **E79-A**, 1707 (1996); H. Matsueda, *Superlattices Microstruct.* **24**, 423 (1998); G.D. Sanders *et al.*, *Phys. Rev. A* **60**, 4146 (1999); E. Biolatti *et al.*, *Phys. Rev. Lett.* **85**, 5647 (2000).
- [3] D. Loss and D.P. DiVincenzo, *Phys. Rev. A* **57**, 120 (1998); G. Burkard *et al.*, *Phys. Rev. B* **59**, 2070 (1999).
- [4] S.N. Molotkov, *JETP Lett.* **64**, 237 (1996); S. Bandyopadhyay, *Phys. Rev. B* **61**, 13 813 (2000); T. Ohshima, *Phys. Rev. A* **62**, 062316 (2000); J. Levy, *ibid.* **64**, 052306 (2001).
- [5] N.H. Bonadeo *et al.*, *Science* **282**, 1473 (1998).
- [6] T.H. Oosterkamp *et al.*, *Nature (London)* **395**, 873 (1998).
- [7] G. Chen *et al.*, *Science* **289**, 1966 (2000).
- [8] M. Bayer *et al.*, *Science* **291**, 451 (2001).
- [9] T.H. Stievater *et al.*, *Phys. Rev. Lett.* **87**, 133603 (2001); H. Kamada *et al.*, *ibid.* **87**, 246401 (2001).
- [10] A. Imamoglu *et al.*, *Phys. Rev. Lett.* **83**, 4204 (1999); *Fortschr. Phys.* **48**, 987 (2000).
- [11] M. Muraio *et al.*, *Phys. Rev. A* **57**, R4075 (1998); A. Thapliyal, *ibid.* **59**, 3336 (1999); C.H. Bennett *et al.*, *ibid.* **63**, 012307 (2000); W. Dür *et al.*, *Phys. Rev. Lett.* **83**, 3562 (1999); N. Linden *et al.*, *ibid.* **83**, 243 (1999); H.A. Carteret *et al.*, *Found. Phys.* **29**, 527 (1999).
- [12] C.H. Bennett *et al.*, *Phys. Rev. A* **53**, 2046 (1996).
- [13] V. Vedral *et al.*, *Phys. Rev. Lett.* **78**, 2275 (1997); V. Vedral and M.B. Plenio, *Phys. Rev. A* **57**, 1619 (1998).
- [14] S. Hill and W.K. Wootters, *Phys. Rev. Lett.* **78**, 5022 (1997); W.K. Wootters, *ibid.* **80**, 2245 (1998).
- [15] W.K. Wootters, e-print quant-ph/0001114; K.M. O'Connor and W.K. Wootters, *Phys. Rev. A* **63**, 052302 (2001); D. Gunlycke *et al.*, *ibid.* **64**, 042302 (2001); M.C. Arnesen, S. Bose, and V. Vedral, *Phys. Rev. Lett.* **87**, 017901 (2001).
- [16] M. Koashi, V. Bužek, and N. Imoto, *Phys. Rev. A* **62**, 050302 (2000).
- [17] X. Wang and K. Mølmer, *Eur. Phys. J. D* **18**, 385 (2002).
- [18] J.H. Reina *et al.*, *Phys. Rev. A* **62**, 012305 (2000).
- [19] C.C. Gerry and J.H. Eberly, *Phys. Rev. A* **42**, 6805 (1990); M. Alexanian and M. Bose, *ibid.* **52**, 2218 (1995); Y. Wu, *ibid.* **54**, 1586 (1996).
- [20] R. Dekeyser and M.H. Lee, *Phys. Rev. B* **19**, 265 (1979); **43**, 8123 (1991); **43**, 8131 (1991).
- [21] R. Botet *et al.*, *Phys. Rev. Lett.* **49**, 478 (1982).
- [22] H.J. Lipkin, N. Meshkov, and A.J. Glick, *Nucl. Phys.* **62**, 188 (1965).
- [23] J.-M. Liu and G. Müller, *Phys. Rev. A* **42**, 5854 (1990); *Phys. Rev. B* **44**, 12 020 (1991).