Universal set of quantum gates for double-dot spin qubits with fixed interdot coupling

Ronald Hanson  
Center for Spintronics and Quantum Computation,  
University of California, Santa Barbara, California 93106, USA

Guido Burkard  
Department of Physics and Astronomy, University of Basel,  
Klingelbergstrasse 82, CH-4056 Basel, Switzerland

We propose a set of universal gate operations for the singlet-triplet qubit realized by two electron spins in a double quantum dot, in the presence of a fixed inhomogeneous magnetic field. All gate operations are achieved by switching the potential offset between the two dots with an electrical bias, and do not require time-dependent control of the tunnel coupling between the dots. We analyze the effects of tunnel-coupling dynamics and calculate the effective qubit rotation angle as a function of the applied electric bias. We present explicit gate sequences for single-qubit rotations about two orthogonal axes, and a CNOT gate sequence, completing the universal gate set.

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Electron spins in semiconductor quantum dots (QDs) are promising candidates for encoding and manipulating quantum information in the solid state. Initialization, manipulation and readout of electron spins have already been demonstrated in these systems [1, 2]. Proposals exist for encoding one logical qubit in one [3], two [4, 5, 6], three [7, 8], or even more [9] spins. Although they differ in many respects, a common essential ingredient of all these proposals is electrical control of the two-electron exchange interaction in a double quantum dot, which is characterized by the singlet-triplet energy splitting $J$.

Conventionally, control over $J$ is envisioned through voltage control of the tunnel coupling $t$ between the two dots. However, in many QD systems, such as vertical pillars [10], self-assembled dots [11], nanowires [12], or etched dots in Si [13], $t$ is fixed by growth or fabrication parameters. Even for double QDs (DQDs) in electrically gated systems, such as GaAs dots [14] and carbon nanotubes [15], fast control over the tunnel coupling is challenging and has not been demonstrated thus far.

A possible way around this problem was demonstrated in a recent experiment by Petta et al. [1], where $J$ is controlled by the misalignment $\varepsilon$ between the two QDs. In contrast to the tunnel coupling, the misalignment can easily be changed over a wide range on a subnanosecond timescale by pulsing the source-drain bias [16] or a gate voltage [1]. Building on this result, Taylor et al. [17] proposed a set of universal gates for a logical qubit whose basis states are the two-electron states $|S\rangle = (|\uparrow\uparrow\rangle - |\downarrow\downarrow\rangle)/\sqrt{2}$ and $|T_0\rangle = (|\uparrow\downarrow\rangle + |\downarrow\uparrow\rangle)/\sqrt{2}$. However, their scheme requires $J$ to be tunable to zero, which is not possible by changing $\varepsilon$ alone [18]. Therefore, voltage control of $t$ is still needed in their scheme.

Here, we propose a set of universal quantum gates for the $S-T_0$ qubit in a constant small inhomogeneous field, that eliminates the need for controlling the tunnel coupling $t$. We demonstrate how arbitrary single-qubit rotations can be performed at \textit{finite} $J$, by combining $Z$ rotations with rotations around an axis in the XZ plane. We discuss the experimental requirements for this scheme and compare them to current-day devices. Finally, we outline a two-qubit CNOT operation, which is based on a change in the rotation angle of the target qubit that is conditional on the control qubit through spin-dependent tunneling and the capacitive coupling between qubits.

Tunable spin dynamics in a DQD--Our qubit is realized in the $|0\rangle \equiv |S\rangle$ and $|1\rangle \equiv |T_0\rangle$ states of two electrons in a double quantum dot, where $S$ and $T_0$ are the lowest-energy singlet and triplet states. The dynamics of these states can be described by the Hamiltonian

$$H = \begin{pmatrix} 0 & \delta h/2 & 0 & 0 \\ \delta h/2 & 0 & \sqrt{2t} & \sqrt{2t} \\ 0 & \sqrt{2t} & U - \varepsilon & 0 \\ 0 & \sqrt{2t} & 0 & U + \varepsilon \end{pmatrix},$$

in the basis $|T_0\rangle$, $|(1,1)S\rangle$, $|(0,2)S\rangle$, $|(2,0)S\rangle$ [21]. Here, $(m,n)$ denotes the number $m$ ($n$) of electrons in dot 1 (2), $\delta h$ is the inhomogeneous magnetic field between the dots and $U$ is the difference in Coulomb energy between the $(1,1)S$ and the $(0,2)S$ or $(2,0)S$ state.

Figure 1 shows the energy of the lowest eigenstates as a function of $\varepsilon$ for $\delta h=0$. In this case, the eigenstates are pure spin states for all values of $\varepsilon$ and we define the qubit basis states as the triplet $T_0$ and the lowest-energy singlet $S$, which are separated by an energy $J$. At $\varepsilon = U$, there is an avoided crossing of the $(1,1)S$ and $(0,2)S$ states, and as a consequence $J$ is large at this point. In the presence of an inhomogeneous field $\delta h$ with magnitude much smaller than $t$ (as in the inset of Fig. 1), $S$ and $T_0$ remain eigenstates near $\varepsilon = U$ where $J \approx t \gg \delta h$. Far away from the avoided crossing, however, $J \approx \delta h$ and therefore $S$ and $T_0$ are strongly mixed. As a consequence, the qubit rotates about an axis determined by $\delta h$ and $J$.

The qubit subspace is energetically separated from the
and, in the qubit basis, the Hamiltonian (1) can be reduced to the qubit subspace difference between the qubit states and all other states, always changed adiabatically with respect to the energy not to rely on this fast control of

where we have chosen the zero of energy midway between the qubit states and all other states, the Hamiltonian \( H \) can be reduced to the qubit subspace and, in the qubit basis \( |S\rangle \) and \( |T_0\rangle \), has the general form

\[
H = \frac{1}{2} \begin{pmatrix} J & \delta h \\ \delta h^* & -J \end{pmatrix} = \mathbf{B} \cdot \mathbf{\sigma},
\]

where we have chosen the zero of energy mid-way between the states \(|0\rangle\) and \(|1\rangle\) and introduced a pseudo-spin notation with Pauli matrices \( \mathbf{\sigma} = (\sigma_x, \sigma_y, \sigma_z) \) in the two-dimensional qubit subspace. The pseudo-magnetic field is \( \mathbf{B} = (\text{Re}[\delta h], \text{Im}[\delta h], J)/2 \), where the exchange coupling \( J \) and effective difference field \( \delta h \) are functions of \( t, U, \delta h, \) and \( \varepsilon \). In what follows, \( \delta h \) will be real and thus the pseudo-field always lies in the \( XZ \)-plane (Fig. 1). The angle of the pseudo-field with the \( X \) axis is

\[
\theta = \arctan(J/\delta h),
\]

and therefore \( J \) remains finite. Thus, we cannot reach a point where \( \mathbf{B} \) lies in the equator plane, i.e., we have to work with a finite angle \( \theta > 0 \).

Single-qubit gates—We now show that arbitrary single-qubit rotations can be constructed from the two available elementary operations: (i) rotations about the \( \theta \)-tilted axis by some angle \( \chi = B\tau \), \( \tau \) being the switching time,

\[
U_\theta(\chi) = \begin{pmatrix} \cos \frac{\chi}{2} + i \sin \frac{\chi}{2} \sin \theta & i \sin \frac{\chi}{2} \cos \theta \\ i \sin \frac{\chi}{2} \cos \theta & \cos \frac{\chi}{2} - i \sin \frac{\chi}{2} \sin \theta \end{pmatrix},
\]

and (ii) nearly perfect \( Z \) rotations by \( \phi = J\tau \approx \sqrt{2t}\tau \), given by the diagonal matrix \( U_Z(\phi) \) with diagonal entries \( e^{i\phi/2} \) and \( e^{-i\phi/2} \). Arbitrary single-qubit rotations can be constructed using the Euler angle method, if rotations by arbitrary angles about two orthogonal axes are available. Therefore, it is sufficient to show that arbitrary rotations about the \( X \) axis, \( U_X(\gamma) = \exp(\gamma \sigma_x/2) \), in addition to the \( Z \)-rotations, are feasible. The three-step sequence

\[
U_X(\gamma) = U_\theta(\chi)U_Z(\phi)U_\theta(\chi),
\]

with the rotation angles [23]

\[
\chi = \arccos \frac{\cos \frac{\phi}{2} \sqrt{1 - \tan^2 \theta \sin^2 \frac{\gamma}{2}}} {\cos^2 \frac{\phi}{2} + \cos^2 \theta \sin^2 \frac{\gamma}{2}},
\]

\[
\phi = -2 \arctan \frac{\sin \chi \sin \theta} {\cos^2 \frac{\phi}{2} + \cos 2\theta \sin^2 \frac{\gamma}{2}},
\]

generates a rotation about the \( X \)-axis by an arbitrary angle \( \gamma \), as long as \( 0 \leq \theta \leq \pi/4 \). One can intuitively understand this sequence by following the state on the Bloch sphere; Fig. 2a depicts the three steps for a rotation from \( |S\rangle \) to \( |T_0\rangle \) (\( \gamma = \pi \)). We note that switching between the working points has to be performed non-adiabatically with respect to \( J \). The rotation angles \( \chi \) and \( \phi \) are plotted as a function of \( \theta \) in Fig. 2b for three different \( X \)-rotation angles \( \gamma \). For a \( \pi \)-flip about the \( X \) axis, \( \chi = \pi \), we find the simpler expressions,

\[
\chi = \arccos(-\tan^2 \theta), \quad \phi = -2 \arctan \frac{\sin \theta}{\sqrt{\cos 2\theta}}.
\]

We note that the sequence Eq. 5 is not simply one of the known NMR sequences. Actually, in NMR it is usually not a problem to perform rotations about an axis in the equator plane of the Bloch sphere [24].

Doing nothing—A convenient “idle” position would be close to the avoided crossings \( \varepsilon = \pm U \), i.e., close to the \( Z \)-gate operation point, as here only \( Z \)-rotations need to be accounted for. A disadvantage of this position is that the qubit is more susceptible to decoherence from charge fluctuations, due to the different orbital characters of the basis states close to the avoided crossing [25]. The best waiting position in terms of coherence is probably the symmetric point \( \varepsilon = 0 \). However, since \( \delta h \) and \( J \) are of the same order at \( \varepsilon = 0 \), the spin constantly rotates.
for effect, one could always wait for an integer number of full periods, \( \tau \). General methods exist for creating an inhomogeneous field \( \epsilon \), \( \phi \), \( \theta \) parameters to anything between 1 and time, and orbitals yield a lower \( \epsilon \), finally, can be controlled in all quantum dot systems listed in the introduction, by pulsing the source, drain or gate voltage \( |\bar{\epsilon}/\epsilon| \lesssim U/\hbar \), away from the avoided crossing.

The three-step sequence Eq. (5) for \( X \)-rotations can be obtained for \( |U \pm \epsilon| \gg t \) by way of a Schrieffer-Wolff transformation \([26]\). Since we assume that \( \delta h \) is fixed, the angle \( \theta \) will never be exactly \( \pi/2 \), which is required for perfect \( Z \) rotations. The desired values of \( t \) and \( \delta h \) therefore depend on the error that can be tolerated (see Fig. 3(b)), with \( t \) typically exceeding \( \delta h \) by more than an order of magnitude. For the \( X \) rotations we need \( \theta \leq \pi/4 \), which gives \( \delta h \sim J \), which can be satisfied by moving away from the avoided crossing. The minimum value of \( \delta h \) needed for the \( X \) rotation is given in Fig. 3(b) for two typical values of \( U \). We note that more detailed calculations including higher orbitals yield a lower (\(<20\%) \) value of \( J \) \([22]\).

In most systems, \( t \) can be set by gates or fabrication parameters to anywhere between 1 \( \mu \)eV and 1 meV. Several methods exist for creating an inhomogeneous field \( \delta h \): (i) application of an inhomogeneous magnetic field, (ii) different \( g \)-factors in the two dots—either by composition or confinement \([27]\) in combination with a homogeneous magnetic field, and (iii) inhomogeneous nuclear polarizations \([28]\). Note that the effect of a fluctuating nuclear field can be diminished by bringing it into an eigenstate \([29]\). The electrical bias \( \epsilon \), finally, can be controlled in all quantum dot systems listed in the introduction, by pulsing the source, drain or gate voltage \([30]\).

Controlled-\( NOT \) gate—To complete our universal set of quantum gates, we require a suitable two-qubit operation, e.g., the controlled-\( NOT \) (CNOT, or quantum XOR) gate that flips the target qubit \((|0\rangle \leftrightarrow |1\rangle)\) if the control qubit is in state \(|1\rangle\), and otherwise leaves the target unchanged. This can be achieved by applying a bias voltage \( \epsilon_{\text{control}} \) on the control qubit, such that its charge state partly shifts to \((0, 2)\) if the qubit state is \(|S\rangle\), but remains mostly in \((1, 1)\) if the state is \(|T_0\rangle\) because the \((0, 2)\) triplet state is far away in energy \([1,31]\). Due to the Coulomb interaction between the control and the target qubit, the target qubit will experience a conditional bias shift (see Fig. 3), that can be of the same order as the interdot Coulomb energy within a single logical qubit.

The CNOT is a conditional \( X \)-rotation by \( \gamma = \pi \), thus it is natural to use a sequence analogous to Eq. (9):

\[
\hat{W}_X(\gamma) = W_\theta(\chi)U_Z(\phi)W_\theta(\chi),
\]

where \( \chi \) and \( \phi \) are given in Eq. (8) in terms of \( \theta(\epsilon) \) and \( \gamma \) at the conditional bias point \( \epsilon \), induced by the charge movement in the control qubit, and \( U_Z \) is the single-qubit \( Z \)-rotation. The conditional rotations \( W_\theta(\chi) \) about the \( \theta \)-axis are analogous to \( U_\theta(\chi) \), but instead of being induced by a direct manipulation of the bias \( \epsilon \), they are...
FIG. 4: A CNOT is performed by electrically biasing the control qubit, shifting its charge distribution toward the target qubit (upper panel) if it is in a singlet state (a), but leaving the charge distribution unchanged if it is in the triplet state (b). This leads to a conditional shift of the working point of the target qubit (lower panels) and to a conditional operation.

controlled by applying $\varepsilon_{\text{control}}$ to the control qubit, which results in a conditional bias $\varepsilon$ at the target qubit.

The sequence Eq. (1) for $\gamma = \pi$ is not a true CNOT yet, because (i) the Z-rotation $U_Z(\phi)$ is not conditional on the control qubit being in state $|1\rangle$, but is in fact always carried out, and (ii) the conditional $W_\delta(\chi)$ rotations also perform a Z-rotation in case the control qubit is in state $|0\rangle$. In summary, $\tilde{W}_X(\pi)$ does a NOT operation (X-rotation by $\pi$) on the target if the control is in $|1\rangle$ and a Z-rotation by $2\chi + \phi$ if the control is in state $|0\rangle$. The true CNOT operation does nothing on the target qubit if the control is in $|0\rangle$; it can be obtained with the sequence

$$U_{\text{CNOT}} = \tilde{W}_X(\pi/2)U_X(\pi)\tilde{W}_X(\pi/2)U_X(\pi),$$

(12)
canceling the undesired phases if the control qubit is $|0\rangle$.

In conclusion, we have proposed a universal set of quantum gates for the $S - T_0$ qubit, consisting of single-qubit rotations about two orthogonal axes X and Z about arbitrary angles combined with the CNOT gate. The electrical bias $\varepsilon$ is the only parameter that needs to be tuned fast and with high precision, which considerably relieves the experimental requirements compared to previous spin-based qubit control proposals and makes our scheme applicable to virtually any quantum dot system.

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[18] Note that $J$ is predicted to go to zero in a finite perpendicular magnetic field [19, 20], but such a relatively large field also closes the essential single-dot singlet-triplet gap.
[21] We only consider the states with spin projection $S_z = 0$. The triplet states $|T_\pm\rangle$ with $S_z = \pm 1$ can be efficiently decoupled from the $S_z = 0$ subspace by choosing $\delta h$ purely along $z$ and by applying a magnetic field $B_z$ along $z$. Note that choosing $z$ in the plane of the dots is preferable to minimize effects of $B_z$ on the orbitals. See [22] for details.
[23] The angles $\chi$ and $\phi$ are obtained by solving the transcendental equations Eq. [3].
[30] For some systems, $t$ might have a weak dependence on $\varepsilon$ which could easily be accounted for in Eq. (1).
[31] The controlled-phase gate proposed in Ref. [17] can also be used here, as it does not require control over $t$. 