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Coupling artificial molecular spin states by photon-assisted tunnelling

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Artificial molecules containing just one or two electrons provide a powerful platform for studies of orbital and spin quantum dynamics in nanoscale devices. A well-known example of these dynamics is tunnelling of electrons between two coupled quantum dots triggered by microwave irradiation. So far, these tunnelling processes have been treated as electric-dipole-allowed spin-conserving events. Here we report that microwaves can also excite tunnelling transitions between states with different spin. We show that the dominant mechanism responsible for violation of spin conservation is the spin-orbit interaction. These transitions make it possible to perform detailed microwave spectroscopy of the molecular spin states of an artificial hydrogen molecule and open up the possibility of realizing full quantum control of a two-spin system through microwave excitation.

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n recent years, artificial molecules in mesoscopic systems have drawn much attention owing to a fundamental interest in their quantum properties and their potential for quantum information applications. Arguably, the most flexible and tunable artificial molecule consists of coupled semiconductor quantum dots that are defined in a two-dimensional electron gas using a set of patterned electrostatic depletion gates. Electron spins in such quantum dots exhibit coherence times up to $200\,\mu s$ (ref. 1), about 10^4 – 10^6 times longer than the relevant quantum gate operations^{2–4}, making them attractive quantum bit (qubit) systems.⁵

The molecular orbital structure of these artificial quantum objects can be probed spectroscopically by microwave modulation of the voltage applied to one of the gates that define the dots⁶. In this way, the delocalized nature of the electronic eigenstates of an artificial hydrogen-like molecule was observed^{7,8}. More recently, electrical microwave excitation was used for spectroscopy of single spins⁹⁻¹¹ and coherent single-spin control^{9,11}, through electric-dipole spin resonance.

Here we perform microwave spectroscopy^{7,8,12} on molecular spin states in an artificial hydrogen molecule formed by a double quantum dot (DD) which contains exactly two electrons. In contrast to all previous photon-assisted tunnelling (PAT) experiments, we observe not only the usual spin-conserving tunnel transitions, but also transitions between molecular states with different spin quantum numbers. We discuss several possible mechanisms and conclude from our analysis that these transitions become allowed predominantly through spin-orbit (SO) interaction. The possibility to excite spin-flip tunnelling transitions lifts existing restrictions in our thinking about quantum control and detection of spins in quantum dots, and allows universal control of spin qubits without gate-voltage pulses.

Results

Device and excitation protocol. Figure 1a displays a scanning electron micrograph of a sample similar to that used in the experiments. It shows the metal gate pattern that electrostatically defines a DD and a quantum point contact (QPC) within a GaAs/ (Al,Ga)As two-dimensional electron gas. An on-chip Co micromagnet (u magnet) indicated in blue in Figure 1a generates an inhomogeneous magnetic field across the DD, which adds to the homogeneous external in-plane magnetic field B (Methods), but is not needed for the molecular spin spectroscopy. The sample was mounted in a dilution refrigerator equipped with high-frequency lines. The gate voltages are set so that the DD can be considered as a closed system (the interdot tunnelling rates are 104 times larger than the dot-to-lead tunnelling rates), and the tilt of the DD potential is tuned by the dc-voltages V_L and V_R , applied to the left and right side gates. Working near the turn-on of the first conductance plateau, the current through the QPC, $I_{\rm OPC}$, depends upon the local charge configuration and provides a sensitive meter for the absolute number of electrons (n_L, n_R) in the left and right dot, respectively^{13,14}.

First, we excite the DD as indicated in the top panel of Figure 1b, by adding to $V_{\rm R}$ continuous-wave microwave excitation at fixed frequency $v\!=\!11\,\rm GHz$. When the photon energy of the microwaves matches the energy splitting between the ground state and a state with a different charge configuration, a new steady-state charge configuration results, which is visible as a change in the QPC current, $\Delta I_{\rm QPC}$. The excitation is on–off modulated at 880 Hz and lockin detection of $\Delta I_{\rm QPC}$ reveals the microwave-induced change of the charge configuration (see Methods for further experimental details). The lower panel of Figure 1b shows $\Delta I_{\rm QPC}$ as a function of $V_{\rm L}$ and $V_{\rm R}$ near the (1,1) to (0,2) boundary of the charge stability diagram. Sharp red (blue) lines indicate microwave-induced tunnelling of an electron from the right to the left dot (left to right), labelled as $\Delta n = +1$ ($\Delta n = -1$), respectively (Fig. 1c). Sidebands can result from multi-photon absorption. At the boundaries with the (0,1) and (1,2)

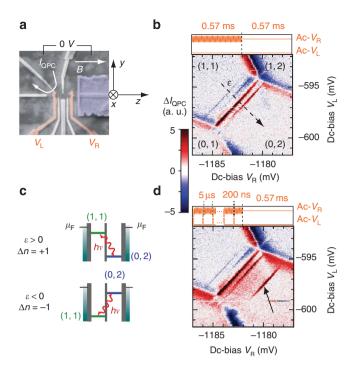


Figure 1 | Photon-assisted tunnelling in a 2-electron double quantum dot. (a) Scanning-electron micrograph top view of the double-dot gate structure with Co micromagnet (blue). The voltages applied to the left V_{\perp} and right V_{\triangleright} side gates (red) control the detuning ε of the double-dot potential. The double-dot charge state is read out by means of the current I_{OPC} running through a nearby quantum point contact (white arrow). (**b**) Charge stability diagram around the 2-electron regime at B = 1.5 T. $(n_{\rm L} n_{\rm R})$ indicate the absolute numbers of electrons in the left and right dot, respectively. During measurements, 11 GHz microwaves with 880 Hz on-off modulation are applied to the right side gate. The top panel displays schematically one cycle of the ac signal applied to the right side gate. Along the detuning axis (dashed arrow) PAT-lines are observed. (c) In the conventional picture of PAT, the first sidebands seen in **b** should appear when the detuning of the (0,2) and (1,1) states matches the photon energy, and interdot tunnelling is induced. Further sidebands are then interpreted as multi-photon transitions. $\mu_{\scriptscriptstyle F}$ is the chemical potential of the left and right electron reservoir. (d) Same as in b, but the microwaves are interrupted every $5 \mu s$ by a 200 ns, $P_s = 2 \,\text{mV}$ detuning pulse applied to the left and right side gates (see the schematic in the top panel). The pulses generate a reference line (black arrow) due to mixing at the ST^+ anti-crossing.

charge states, no energy quantization is observed, because, here, electrons tunnel to and from the electron-state continuum of the leads. At first sight, the observations in Figure 1b thus appear to be well explained by the usual spin-conserving PAT processes.

Surprisingly, the position in gate voltage of the resonant lines exhibits a striking dependence on the in-plane magnetic field, B. This is clearly seen in Figure 2a,b, which display the measured PAT spectrum along the DD detuning ε axis (dashed black arrow in Fig. 1b) as a function of B for 20 and 11 GHz excitation, respectively.

As the gate constitutes an open-ended termination of the transmission line, the excitation produces negligible AC magnetic fields at the DD (estimated in Methods), and is, therefore, expected to give rise to only electric-dipole-allowed spin-conserving transitions, with no B dependence. Furthermore, there is a pronounced asymmetry between the position of the red and blue PAT lines.

In these figures, the detuning axis was calibrated for all magnetic fields by introducing a reference line (Fig. 1d, black arrow) that facilitates interpretation of the spectra despite residual orbital effects of the magnetic field. This line was produced by interspersing

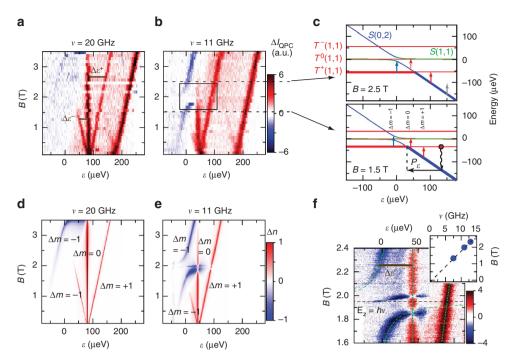


Figure 2 | Photon-assisted-tunnelling spectra and simulations. (a,b) Microwave-induced change of the QPC current $\Delta I_{\rm QPC}$ as a function of the double-dot detuning ε and the external magnetic field B for 20 GHz (a) and 11 GHz (b) frequency, respectively. Singlet-triplet mixing due to 2-mV detuning pulses generates a reference signal that is used to calibrate the detuning axis (see lower panel in Fig. 2f). (c) Eigenenergies versus double-dot detuning ε of the two-electron spin states in the (1,1) and (0,1) charge regime for two external magnetic fields B = 2.5 T (upper panel) and B = 1.5 T (lower panel), respectively. S(0,2), S(1,1) and T(1,1) character of the eigenstates is indicated by blue, green and red colour, respectively. The molecular spin ground state is indicated by thick lines. The vertical arrows indicate PAT transitions for a constant microwave frequency involving spin flips. The transition indicated by a dashed arrow is suppressed, because the initial state lies above the ground state. The red circle in the lower panel indicates the detuning position of the reference signal, that is generated by a detuning pulse with amplitude P_{ε} to the ST^+ anti-crossing. (d,e) Simulated PAT spectra for 20 GHz (a) and 11 GHz (b) frequency, respectively. The colour indicates the change of the population of the steady-state charge state Δn as would be observed in an on-off lock-in detection. A finite temperature of 100 mK and spontaneous relaxation through the phonon bath are taken into account. (f) $I_{\rm QPC}$ scanned with higher resolution in the anti-crossing region (black rectangle in Fig. 2b) at 11 GHz excitation. The green dashed lines indicate the expected detuning positions of the PAT transitions. The horizontal black dashed line indicates the magnetic field, at which the electron spin resonance condition is fulfilled $E_z = g\mu_B(B + b_0) = hv$. The graph is concatenated from two scans that overlap at 1.9 T. The inset displays three magnetic fields (blue dots), at which the ce

the microwaves every $5\,\mu s$ with $200\,ns$ gate-voltage pulses along the detuning axis (see the top panel in Fig. 1d), leading to singlet–triplet mixing as described in ref. 2. The short gate-voltage pulses do not noticeably alter the position of the PAT lines (compare Fig. 1b,d). The reference peaks visible at around $\epsilon{=}\,200\,\mu eV$ in Figure 2a,b were aligned by shifting all data points at a given B by the same amount in detuning (see Supplementary Note 1 'Calibration of the detuning axis' for the full details of this post-processing step and Supplementary Fig. S1a–d for the corresponding spectra).

Interpretation of the photon-assisted tunnelling spectra. The complexity of the PAT spectra shown in Figure 2a,b can be understood in detail, if we allow for non-spin conserving transitions. The two diagrams in Figure 2c show the energies of all relevant DD-states (four (1,1)-states and one singlet S(0,2)-state) as a function of ε for two different (fixed) magnetic fields, that is, the spectrum of the DD along the two horizontal dotted lines in Figure 2b¹⁵. Note that the only difference between the two diagrams is the splitting between the three triplet T(1,1)-states.

First, we explain the resonances observed along the upper dotted line in Figure 2b (B=2.5 T). In the corresponding (upper) diagram in Figure 2c, we plot the ground state energy for all ε with a thick line. If the microwave excitation is off-resonance with all transitions, the system will be in this ground state. For instance, at ε =150 μ eV, there is no state available 11 GHz above the S(0,2) ground state (grey

arrow) and the system stays in S(0,2). However, when decreasing the detuning, at some point $T^+(1,1)$ becomes energetically accessible (red arrow) and, because we allow for non-spin-conserving transitions, is populated owing to the microwave excitation. For this PAT transition, the spin projection on the quantization axis is changed by $\Delta m = +1$. The resulting change of steady-state charge population (increased population of (1,1), or $\Delta I_{\rm QPC} > 0$) is detected by the QPC and yields the red peak in Figure 2b. Decreasing ε further, there are two more resonances detectable: (i) the S(0,2) - S(1,1) transition (dotted red arrow, $\Delta m = 0$), although the signal will be weakened owing to the fact that S(0,2) is not unambiguously the ground state anymore. Note that a transition $S(0,2) - T^{0}(1,1)$ could appear in nearly the same detuning position, as will be discussed below. (ii) the $T^+(1,1)$ -S transition (blue arrow), where S stands for the hybridized S(0,2) - S(1,1) singlet, results in a negative (blue, $\Delta I_{OPC} < 0$, $\Delta m = -1$) signal from the charge detector as the ground state is now (1,1) and the excited state is partly (0,2). We see that this simple analysis explains both the positions and the signs of the resonances observed in the data.

A similar analysis can be made for other magnetic fields. For instance, for the spectrum plotted in the lower diagram of Figure 2c, we find two resonances with $\Delta I_{\rm QPC} > 0$ (red arrows), and one with $\Delta I_{\rm QPC} < 0$ (blue arrow). Note that the 'blue' transition now connects the ground state to the other branch of the hybridized S compared to the high magnetic field case. Indeed, the singlet anti-crossing is

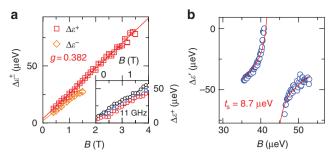


Figure 3 | Analysis of the photon-assisted-tunnelling spectra. (a) The detuning difference $\Delta\varepsilon$ between the spin-conserving line $\Delta m=0$ and the $\Delta m=\pm 1$ line is plotted as a function of the external magnetic field B at 20 GHz (see Fig. 2a), to fit the absolute effective electron g-factor g from the slopes of the linear fits (solid lines). inset, $\Delta\varepsilon$ for the $\Delta m=+1$ line at 11 GHz for different tunnel couplings. The offset of the curves increases with increasing t_s (red to black circles) while g remains constant. (b) $\Delta\varepsilon$ between the $\Delta m=0$ line and the $\Delta m=-1$ (1,1) to (0,2) transition from Figure 2c. The fit of the anti-crossing (red line) allows for a precise determination of t_s .

directly probed, resulting in the two blue curved lines observed in the data around ε =0 (Fig. 2b). The fading out of the blue signal at low fields can be understood from pumping into the metastable state S(1,1): the microwaves excite the system from $T^+(1,1)$ to S(0,2), from where it relaxes quickly to S(1,1). However, relaxation from S(1,1) back to the ground state is slow because of the small energy difference of this transition and the small phonon density of states at low energies ^{14,16}. This pumping weakens the detector signal, as S(1,1) has the same charge configuration as the ground state.

To verify this interpretation, we calculate in Figure 2d,e the position and intensity of the spectral lines at fixed microwave frequency, on the basis of the energy level diagram of Figure 2c. In the simulations, all single-photon transitions between the ground state and the excited states are allowed by including a matrix element $|T^{\pm}(1,1)\rangle\langle S(0,2)|$ (Methods). The input parameters for the calculation of the resonant positions are the interdot S(1,1) to S(0,2)tunnel coupling t_o , the absolute electron g-factor |g|, a magnetic-field contribution b_0 from the μ magnet parallel to B as well as a magnetic field gradient $\Delta \vec{B} = (\Delta B_x^{\perp}, 0, \Delta B^{\parallel})$ between the dots (in Fig. 3, we show how $t_{s_0}[g]$ and b_0 can be extracted from the experimental spectra). The colour scale represents the calculated steady-state Δn that results from microwave excitation, orbital hybridization, and phonon absorption and emission at 100 mK. All the PAT transitions visible in the simulation also appear in the experiment, with excellent agreement in both the position and relative intensity of the spectral lines. Especially, the vanishing signal due to spin pumping is also predicted by the simulations that include phonon relaxation. Surprising at first glance, the intensities of the spin-flip lines $\Delta m = \pm 1$ and the spin-conserved line $\Delta m = 0$ are about the same, which holds true for both simulation and measurement, although $\Delta m = 0$ has a larger matrix element. This is because the same matrix elements enter the spontaneous phonon-mediated relaxation rates, resulting in similar steady-state populations for the two transitions.

When we zoom in on the boxed region of Figure 2b, we see an extra horizontal blue feature at $B\approx2\,\mathrm{T}$ (Fig. 2f) that also appears in the calculated spectra of Figure 2e. This feature is due to a triplet resonance from $T^+(1,1)$ to $T^0(1,1)$ that becomes detectable by relaxation into the meta-stable S(0,2) state. In the detuning range where this line appears, the S(0,2) state lies energetically only slightly above the $T^+(1,1)$ state, so relaxation back to the $T^+(1,1)$ ground state is suppressed, again by the small phonon density of states at low energies (see Supplementary Fig. S2a for a schematic energy diagram). The triplet resonance is expected to appear at $E_z = g\mu_B(B+b_0) = hv$,

where h is Planck's constant and $\mu_{\rm B}$ the Bohr magneton. The inset of Figure 2f shows the magnetic fields corresponding to the centre of the measured triplet resonance line for three excitation frequencies (see Supplementary Note 2 'Triplet spin resonance' for details and Supplementary Fig. S2b,c for extra spectra taken at different microwave frequencies), which are in good agreement with the expected positions (black dashed lines in Fig. 2f) based on the values |g| and b_0 determined in the next section from other features of the PAT spectra. Surprisingly, the measured triplet resonance exhibits a finite slope in the $B(\varepsilon)$ spectra. A longitudinal magnetic field gradient ΔB^{\parallel} gives rise to such a detuning dependence, but the ΔB^{\parallel} required in our simulations to reproduce the observed slope is $\Delta B^{\parallel} {\gtrsim} 80\,{\rm mT}$ per 50 nm, an order of magnitude larger than the gradient we calculate for the μ magnet. The magnitude of the slope remains a puzzle.

Extracting artificial molecule parameters. We now show how |g|, and b_0 , the parameters used for all simulations, can be extracted independently from the experimental spin-flip PAT spectra. For this analysis, we only use the relative distance $\Delta\varepsilon$ between PAT lines at fixed magnetic field, to be independent from the calibration of the detuning axis by means of the reference line. Figure 3a shows $\Delta\varepsilon^\pm$ as a function of B using the 20 GHz data. $\Delta\varepsilon^\pm$ is defined as the difference in detuning between the red $\Delta m = \pm 1$ and $\Delta m = 0$ PAT lines as plotted in Figure 2a. For a fixed v, both $\Delta\varepsilon^\pm$ increase linearly with the Zeeman energy and therefore allow fitting of |g|. (Note that for $\Delta\varepsilon^+$ the linearity is only exact for sufficiently large B, at which the singlet anti-crossing does not affect the detuning position of the S(0,2) to T^+ (1,1) transition. A least-squares fit to the $\Delta\varepsilon^+$ data gives $|g| = 0.382 \pm 0.004$ (Fig. 3a). From the linear behaviour of $\Delta\varepsilon^+$, we also deduce that there is most likely negligible dynamic nuclear polarization in the experiment.

Knowing |g| precisely, we make use of the blue anti-crossing in Figure 2b,f to determine t_s (and b_0). Figure 3b shows the difference in detuning $\Delta \varepsilon$ between the blue $\Delta m = -1$ and the red $\Delta m = 0$ lines in Figure 2f. Assuming the $\Delta m = 0$ line corresponds to the S(0,2)-S(1,1) transition (as shown below), then

$$\Delta \varepsilon' = \frac{t_s^2 - (hv - g\mu_B(B + b_0))^2}{hv - g\mu_B(B + b_0)} - \sqrt{(hv)^2 - (2t_s)^2},$$
 (1)

where the first term is the detuning position of the $T^+(1,1)$ -S transition and the second the one of the S(0,2)-S(1,1) transition. The best fits are obtained with t_s = $8.7\pm0.1\,\mu\text{eV}$ and b_0 = $109\pm16\,\text{mT}$. This value for b_0 matches our simulations of the stray field of the μ magnet at the DD location very well (see Methods).

An important question left open so far is whether the red $\Delta m=0$ line involves predominantly transitions from S(0,2) to S(1,1) or to $T^0(1,1)$. The transition to S(1,1) does not require a change in the (total) spin and is thus expected to be excited more strongly than that to $T^0(1,1)$. However, relaxation from S(1,1) back to S(0,2) will be stronger as well, so it is not obvious what steady-state populations will result in either case. Furthermore, given the small energy difference between S(1,1) and $T^0(1,1)$, the two transitions are not resolved in Figure 2. Figure 3a helps to answer this question: The observation that $\Delta \varepsilon^+ > \Delta \varepsilon^-$ indicates that the $\Delta m=0$ line originates from the transition to S(1,1) and not to $T^0(1,1)$. For the former we expect $\Delta \varepsilon^\pm = E_z \pm J$, with

$$J(\varepsilon, t_s) = \frac{t_s^2}{\varepsilon} + O\left[\left(\frac{t_s}{\varepsilon}\right)^4\right]$$

the exchange energy, whereas the latter would result in $\Delta \varepsilon^+ \lesssim \Delta \varepsilon^-$ (in both scenarios, b_0 causes an extra fixed offset in both $\Delta \varepsilon^\pm$, but it does not contribute to their difference; for transitions to both S(1,1) and $T^0(1,1)$, we plotted the expected $\Delta \varepsilon^\pm(B)$ dependence for our DD parameters in the Supplementary Fig. 3a). This interpretation is consistent with the increase of $\Delta \varepsilon^+$ with larger interdot tunnel

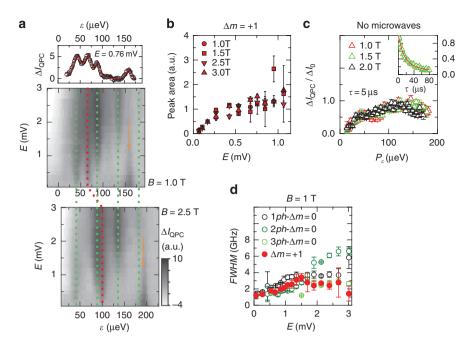


Figure 4 | Power-dependence of the photon-assisted tunnelling spectra and spontaneous relaxation. (a) $\Delta I_{\rm QPC}$ as a function of the double-dot detuning ε and the microwave amplitude E for 11 GHz and B = 1.0 T (middle panel) and B = 2.5 T (lower panel). The excitation scheme from Figure 1d is employed with reference pulse amplitude $P\varepsilon$ = 1.5 mV. The green (red) dashed lines mark the multi-photon Δm = 0 (Δm = +1) PAT transitions. The reference signal stemming from the pulse is marked by orange arrows. The voltage amplitude E is measured at the end of the coaxial lines at room temperature. The uppermost panel displays a linecut measured at 1T. The red line is a least-squares fit by a sum of four lorentzian peaks to the data. (b) The fitted Lorentzian peak area of the Δm = +1 line from $\bf a$ is plotted as a function of microwave amplitude E and magnetic field B. (c) Normalized QPC current averaged over τ =5 μ s immediately after full mixing at the ST^* anticrossing for various magnetic fields at zero microwave power. The spontaneous spin relaxation after mixing is measured at a distance P_{ε} to the ST^* mixing point (see scheme in Fig. 2c). The dependence on the averaged time interval τ is displayed in the inset for B = 1.5 T and P_{ε} = 2 mV together with a least-squares fit (red solid line). (d) The fitted Lorentzian linewidth at half maximum (FWHM) of the Δm = +1 and the lines Δm = 0 from $\bf a$ is plotted as a function of microwave amplitude E for B = 1T. The error bars in $\bf b$, $\bf d$ are determined from the Lorentzian least-squares fit (see the upper panel of Fig. 4a).

coupling, hence larger J (Fig. 3a inset; the slopes are not affected). In Supplementary Note 3 'The $\Delta m=0$ PAT transition', we give more arguments for our interpretation by analysing quantitatively the intercept $\Delta \varepsilon^+(B=0\,\mathrm{T})$ for various microwave frequencies, as plotted in Supplementary Figure S3b.

So far only single-photon processes were considered, but at higher microwave power, also multi-photon lines emerge (Fig. 4a), mostly for the S(0,2) - S(1,1) transition (green dashed lines in Fig. 4a). Like the single-photon S(0,2) - S(1,1) line, their position in detuning is B-independent.

Discussion

Having shown the power of spin-flip PAT for detailed molecular spin spectroscopy, we now discuss the mechanisms responsible for this process as confirmed by our simulations. As a first possibility, the transitions from S(0,2) to the triplet (1,1) states can take place through a virtual process involving S(1,1): the state S(0,2) is coupled to *S*(1,1) by the interdot tunnel coupling, and an (effective) magnetic field gradient $\Delta \vec{B} = (\Delta B_x^{\perp}, 0, \Delta B^{\parallel})$ across the DD couples the spin part of all the (1,1) states to each other^{14,17}. Here $\Delta \vec{B}$ has a contribution from the effective nuclear field and from the µ magnet. The transition matrix element from S(0,2) to $T^{\pm}(1,1)$ is $\propto t_s(\Delta B^{\perp}/B)$, assuming $E_z \gg J$. In the following, we use the *B*-dependence of this process as a fingerprint and focus on the red $\Delta m = +1$ line in Figure 2a,b, as we can follow it over the entire magnetic field range. The intensity of this line is constant in B, and even if the microwave amplitude Eis varied, we observe no B-dependence in the area under this peak (Fig. 4b). This does not in itself provide evidence that the transition rate is magnetic field independent.

As stated above, the observed PAT lines reflect the steady-state change in the charge configuration resulting from stimulated photon emission and absorption, and spontaneous relaxation. A field-independent steady state could thus be reached from a field dependence of relaxation and excitation that cancel each other. We therefore verify that the spontaneous relaxation rate is field-independent as well. The measurement of the spontaneous relaxation rate is done as follows: Starting from S(0,2), we populate the $T^+(1,1)$ by 50% through a 200-ns detuning pulse¹⁸ with amplitude $P\varepsilon$ in the absence of microwaves, and monitor the decay back to S(0,2). The relaxation rate Γ_s can be extracted from the time-averaged lock-in signal $\Delta I_{\rm OPC}(\tau) = (\Delta I_0 / \tau) \int_0^{\tau} \exp(-t\Gamma_s) dt$, where τ is the time spent in Pauli blockade between the pulses. $\Delta I_0 = I_{OPC}(0)$, which is independent from *B*, is extracted from the fit in the inset of Figure 4c. To cover various values of the detuning and the magnetic field, we next fix $\tau = 5 \,\mu s$ and record $\Delta I_{\rm OPC}$ as a function of P_{ε} for three different magnetic fields. We observe that $\Delta I_{\rm OPC}(5\,\mu s)$ and thus Γ_s are essentially independent of B (Fig. 4c). This holds true for all P_{ε} and hence for all $T^+(1,1)$ -S(0,2) energy splittings. Note that regardless of B, this energy splitting is given by $P\varepsilon$ alone. This reflects exactly the situation in the PAT experiment, for which the microwave frequency alone sets the energy splitting and, thus, also the required phonon energy for the spontaneous relaxation process. The fieldindependence of relaxation suggests that the coupling mechanism is magnetic field independent, and, thus, virtual processes involving S(1,1) do not give a strong contribution to the transition rates.

More recently, two mechanisms were considered that provide a direct, *B*-independent matrix element between S(0,2) and the (1,1) triplet states: (i) the hyperfine contact Hamiltonian is of the form $\sum_{i} \delta(r-r_{i})\vec{I}_{j} \cdot \vec{S}$. Thus, nuclear spins, \vec{I}_{j} , in the barrier regions,

Table 1 | Microwave transitions in the (1,1) spin manifold.

Transition	Ω (MHz)	γ (μs ⁻¹)
$ T^+\rangle \rightarrow S(1,1)\rangle$	5.3	1.9
$ T^+\rangle \rightarrow T^0\rangle$	1.5	0.16
$ T^+\rangle \rightarrow T^-\rangle$	1.0	0.04
$ S(1,1)\rangle \rightarrow T^0\rangle$	27	2.0
$ S(1,1)\rangle \rightarrow T^{-}\rangle$	18	1.9
$ T^0\rangle \rightarrow T^-\rangle$	5.1	0.20

The Rabi frequencies Ω and dephasing rates γ , for transitions between different pairs of states of the (1,1) spin manifold, are calculated with the model explained in Methods (detuning ε = $-40\,\mu$ eV, B_{ext} =1T and 10 μ eV microwave amplitude). The dephasing rates include only phonon decay and co-tunnel effects. Further spin dephasing due to, for example, hyperfine coupling is neglected.

where the left-dot and right-dot orbitals overlap, can flip-flop with the electron spin, \vec{S} , simultaneously with charge tunnelling¹⁹. (ii) The SO Hamiltonian is of the form $p_{x,y}S_{x,y}$ and can directly couple states that differ in both orbital and spin^{20,21}. (When the orbital part of the initial and final state are the same, the SO Hamiltonian does not provide a direct matrix element, and the transition rate becomes B-dependent. 9,22-25) The ratio of the SO-mediated rate and the hyperfine mediated rate can be estimated as $[(\Delta\sqrt{N}/4A)(a/\lambda_{SO})]^2$, which is a few thousand in the experiment (see Supplementary Note 4 'Spin flip tunnelling mechanism' for the derivation). Here Δ is the singledot level spacing, N the number of nuclei in contact with one dot, A the hyperfine coupling strength, a the interdot distance and λ_{so} the spin-orbit length. We, therefore, believe that SO interaction is the dominant spin-flip mechanism for the observed PAT transitions. The presence of a magnetic-field-independent matrix element between S(0,2) and the (1,1) triplets is confirmed by the observation that the intensity of the reference signal shows no field dependence.

Finally, we extract from Figure 4a the fitted linewidth as a function of driving power (Fig. 4d). For small E, we find a width $\gtrsim 1\,\mathrm{GHz}$, similar to that observed earlier for spin-conserving PAT processes⁷. For stronger driving, both the $\Delta m = 0$ and the $\Delta m = +1$ lines are further broadened, up to $E \sim 1.5\,\mathrm{mV}$. If these lines were power broadened, their width would imply transition rates in excess of 1 GHz. However, we do not believe that this is the case, because, in measurements with short microwave bursts, the populations saturated only on a long (10- μ s) timescale (data not shown). Presumably, charge or gate-voltage noise is responsible for the broadening instead.

We have shown that in our DD system, all (1,1) spin states have at least weakly allowed electric-dipole transitions to S(0,2). In materials with high SO interaction like InAs, the effect of the non-spin conserving PAT will be even stronger. In materials with weak SO interaction, a strong gradient magnetic field can be used to facilitate spin-flip tunnelling transitions. In all cases, all (1,1) to S(0,2) transitions can be driven. This opens up the possibility of realizing full quantum control of the two spins in the (1,1) manifold through off-resonant (microwave) -stimulated Raman transitions through the excited S(0,2) state, at small negative detuning and in the presence of a finite ΔB^{\parallel} . Furthermore, we find from our analysis that the vicinity in energy of the S(0,2) state also allows direct (single-frequency) transitions to be induced between any two–spin states within the (1,1) subspace. Direct and stimulated Raman transitions are governed by the same matrix elements (see Methods).

The calculated Rabi frequencies Ω_{ij} for the direct transitions are shown in Table 1. For each pair of states, the dephasing rate γ_{ij} from phonon decay and co-tunnelling effects is calculated as well. In the absence of spin dephasing, rotations will be limited by the achievable Ω_{ij}/γ_{ij} , with the overall amplitude damping in time t going as $t\gamma_{ij}/2$, leading to an infidelity of a π -pulse of order $\pi\gamma_{ij}/(2\Omega_{ij})$. Considering dephasing due to the random, quasi-static nuclear field with characteristic dephasing time $T_2^* \sim 10-30$ ns, we find for our

device parameters that $\gamma_{ij} \ll 1/T_2^*$. Therefore, phonon mediated decay does not increase dephasing beyond the hyperfine contribution. Importantly, a quasi-static dephasing contribution (here from the hyperfine field) affects Rabi oscillations less strongly than rapidly fluctuating fields (here from phonons); it permits Rabi oscillations to be observed with periods much longer than $T_2^{*\ 27,28}$.

Controllability of the (1,1) manifold can be best achieved by using only the three couplings between the triplet manifold and the singlet, corresponding to the largest Rabi frequencies. With envelope amplitudes of $\hbar\Omega\!\lesssim\!10\,\mu\text{eV}$, just within the perturbative limits of our theory, Rabi frequencies of order 10 MHz are possible, similar to the frequencies obtained in similar quantum dots for single-spin rotations. Thus, we find that the complete (1,1) subspace is controllable using direct transitions near the S(0,2) crossing, with many rotations possible $^{3.27,28}$, enabling microwave-induced entangling gates.

In addition, our observations suggest new measurement techniques that do not rely on Pauli spin blockade²⁶. An example is a measurement that distinguishes parallel from anti-parallel spins while acting non-destructively on the $S(1,1)-T^0(1,1)$ subspace, which can be achieved by coupling resonantly the $T^+(1,1)$ and $T^-(1,1)$ states to S(0,2), followed by charge readout. This constitutes a partial Bell measurement and leads to a new method for producing and purifying entangled spin states that enables universal measurement-based quantum computation²⁹⁻³¹.

Methods

Sample fabrication. 30-nm thick TiAu gates are fabricated on a 90-nm deep (Al_{0.3}, Ga_{0.7})As/GaAs two-dimensional electron gas (2DEG) by means of e-beam lithography. The double-dot axis is aligned along the [110] GaAs crystal direction (z-direction), which is parallel to the external magnetic field direction \vec{B} (Fig. 1a). The 2DEG is Si δ-doped (40 nm away from the hetero-interface), exhibits an electron density of 2.05×10¹¹ cm⁻² and a mobility of 2.06×10⁶ cm² per Vs at 1 K in the dark. The grounded, 275 nm thick, 2-μm wide and 10-μm long Co μ magnet is evaporated on top of an 80-nm thick dielectric layer, aligned along \vec{B} (magnetic easy axis) and placed ~400 nm away from the closest dot centre. We calculate³² that, at the double-dot position, the μ magnet adds b_0 ~110 mT to B_z and generates a magnetic field gradient of $\Delta B^{\parallel} \approx 6$ mT per 50 nm and a transverse gradient of $\Delta B^{\perp} \sim -6$ mT per 50 nm at saturation ($B_z \approx 2$ T).

Measurement. The sample is mounted in an Oxford KelvinOx 300 dilution refrigerator at 30 mK. Left and right side gate voltages, $V_{\rm L}$ and $V_{\rm R}$, are set by low-pass filtered dc lines and ~60 dB attenuated coaxial lines combined with bias-tees with a cutoff frequency of 30 Hz. The pre-amplified current through the quantum-point contact is read out by a lock-in amplifier locked to the 880 Hz on-off modulation of the microwaves. The bias across the double dot is set to 0 μ V. Voltage pulses to the left and right side gates are generated with a Sony Textronix AWG520. The microwaves are generated with a HP83650A and combined with the pulses to the right side gate. Microwave bursts and detuning pulses are synchronized to ensure that the microwave excitation is switched off during the detuning pulses that generate the reference signal (Fig. 1d).

The gates constitute an open-ended termination of the transmission line and, therefore, the microwaves predominantly generate an AC electric field giving rise to purely electric-dipole PAT transitions. The AC magnetic field contribution from the displacement current is negligible at the dot position. The highest contribution is the displacement current from the right side gate to the 2DEG. The separation between them is 90 nm and the shortest distance between the right dot and the displacement current is more than 100 nm. The relevant area of the displacement current is the area of the right side gate closest to the dot, which is roughly 40 nm by 200 nm. At an AC voltage of 1 mV and at the highest frequency of 20 GHz, the displacement current generates an AC magnetic field of less than 2 nT at the right dot, corresponding to a Rabi frequency of 11 Hz. This is three orders of magnitude slower than the spin relaxation time (Fig. 4) and therefore negligible (see also the reasoning in ref. 9).

Simulation. The. Hamiltonian describing the two-spin system near the (1,1)-(0,2) transition is taken to be a five-state system, with four (1,1) spin states and a (0,2) spin singlet¹⁷ in the presence of an external magnetic field B and a magnetic field gradient $\Delta \vec{B}$, which includes both the quasistatic nuclear field and the field from the μ magnet. This is given by $H = g\mu_B P_{11}(\vec{S}_1 \cdot (\vec{B} + \Delta \vec{B}) + \vec{S}_2 \cdot (\vec{B} - \Delta \vec{B})) P_{11} - \varepsilon | S(0,2) | S(0,2) | + H_t$, were P_{11} is the projector onto the (1,1) subspace, ε is the detuning due to the difference in gate potentials from the left and right gates and the tunnel coupling $H_t = t_s(|S(1,1)) | S(0,2) | + t_{SO}(|T^*(1,1)) | S(0,2) | + t_{SO}(|T^*($

To find the signal, we expect theoretically from the experiment, we add a weak, rapidly oscillating term to the Hamiltonian: $\varepsilon \to \varepsilon_0 + \Omega \cos(\nu t)$. We diagonalize H with $\Omega = 0$, then make a rotating frame transformation in which levels are grouped into bands n (defined by a projector P_n) where the states in a band n are much closer in energy than hv, while the energy difference between states in band n and n+1 are within 2/3rds of hv. Each band rotates at a rate nv, and we can then make a rotating wave approximation, keeping terms due to δ that couple adjacent bands, that is, our perturbation in the rotating frame and rotating wave approximation is $V = \Omega \sum_n P_n \mid (0,2)S \rangle (0,2)S \mid P_{n+1} + h.c.$

Next, we add dissipation and dephasing by including relaxation due to coupling of the electron charge to piezoelectric phonons in a two-orbital (Heitler–London-like) model. We thereby neglect deformation phonons as the energy scales examined in the experiment (7–22 GHz) are much smaller than the characteristic frequency scale of a phonon on the length scale of the dot $c_{ph}/l_{dor}\sim60$ –120 GHz. To determine the coupling, we take as an ansatz for the electronic wavefunctions the Fock–Darwin states, given by Gaussians, and calculate the coupling after orthogonalizing the states with the perturbation $V_{ph}=\Sigma_{e,k}f(k_z)\sqrt{(\hbar/\rho\omega_{e,k})}e^{ik\cdot\vec{r}}\beta_e(a_{e,k}-a_{e,-k}^\dagger)^{33}, \text{ where } f(k_z)\sim1 \text{ for the energy scales we are working with. We then use Fermi's golden rule to calculate excitation and relaxation from thermal and spontaneous emission of phonons. Finally, we numerically solve the superoperator for the steady state and compare the expectation value of <math display="inline">|S(0,2)\rangle\langle S(0,2)|$ with and without the excitation Ω , mimicking the effect of the lock-in detection.

Coherent transitions within the (1,1) manifold. We conclude with a discussion of the controllability of the (1,1) charge manifold using modulated microwaves on ε . To evaluate this, we consider the high-magnetic field, large detuning limit, where $-\varepsilon$, $g\mu_B \mid B_z \mid \gg t_s \gg g\mu_B \mid \Delta B \mid$, where t_s is the characteristic tunnelling energy scale. In this limit, we can use a Schrieffer–Wolff transformation for H_0+V into the (1,1) subspace, where

$$H_0 = \begin{pmatrix} b & 0 & 0 & 0 & 0 \\ 0 & -db & 0 & 0 & 0 \\ 0 & 0 & db & 0 & 0 \\ 0 & 0 & 0 & -b & 0 \\ 0 & 0 & 0 & 0 & -\varepsilon \end{pmatrix}$$
 (2)

in the basis $(T^-(1,1),|\downarrow\uparrow\rangle,|\uparrow\downarrow\rangle,T^+(1,1),S^-(0,2))$ with $b=g\mu_{\rm B}B_{\nu}db=g\mu_{\rm B}\Delta B_{\parallel}$. Defining the un-normalized coupling vector

$$|v\rangle = u|T^-\rangle + t|\downarrow\uparrow\rangle - t^*|\uparrow\downarrow\rangle + u|T^+\rangle$$
, the tunnel coupling is represented by $V = |v\rangle\langle S(0,2)| + |S(0,2)\rangle\langle v|$, with $u = t_{SO,y}, t = (-it_{SO,z} - t_s)/\sqrt{2}$. We seek *S*

such that $\exp(\lambda S)(H_0 + \lambda V)\exp(-\lambda S)$ is block diagonal (has no terms coupling (1,1) to (1,1)) to order λ^2 . Here λ is a perturbation theory parameter that we will take to be unity after the analysis. S is given by solution of $P([S,H_0] + V)Q = 0$, where P projects onto the (1,1) charge space and Q projects onto the (0,2) configuration. It takes the form $S = |R| \langle S(0,2)| - |S(0,2) \rangle \langle R|$ with

takes the form
$$S = |B\rangle \langle S(0,2)| - |S(0,2)\rangle \langle B|$$
, with
$$|B\rangle = \frac{u}{\varepsilon + b} |T^{+}\rangle + \frac{t}{\varepsilon - db} |\downarrow\uparrow\rangle - \frac{t}{\varepsilon + db} |\uparrow\downarrow\rangle + \frac{u}{\varepsilon - b} |T^{-}\rangle$$

Finally, for the (1,1) manifold, we get the effective Hamiltonian $H_{\mathrm{eff}} = H_0 + \frac{1}{2} (|B\rangle \langle v| + |v\rangle \langle B|)$ to order λ^2 of $H_0 + 1/2[S, V]$.

Within this framework, we can now evaluate the effects to order λ^2 expected from a time-dependent perturbation of the form $V_1 = \Omega(t)\cos(vt)|S(0,2)| \cdot \langle S(0,2)|$. We get coupling between (1,1) and (0,2) at order λ , and coupling between (1,1) states at order λ^2 . The charge-changing coupling becomes (taking $\lambda \rightarrow 1$):

$$e^{\lambda S} |S(0,2)| \langle S(0,2) | e^{-\lambda S} = |S(0,2)| \langle S(0,2) | + \lambda [S, |S(0,2)| \langle S(0,2) |] + \frac{\lambda^2}{2} [S, [S, |S(0,2)| \langle S(0,2) |]]$$

$$= |B| \langle S(0,2) | + |S(0,2)| \langle B | + |B| \rangle B | + \frac{(1 - \langle B |B|) |S(0,2)| \langle S(0,2) |]}{2}$$
(3)

From this expression, we see that microwave transitions are possible through time-dependent driving of ε , if $|B\rangle$ has a non-zero matrix element for the desired state. We also see immediately that Raman coupling through $|S(0,2)\rangle$ has the same matrix elements as a direct coupling between (1,1) states through the $|B\rangle\langle B|$ term.

Thus, full controllability of the (1,1) manifold through microwaves can be arrived at by two methods. One is to use stimulated Raman transitions through the $|S(0,2)\rangle$ state, the other is to use direct transitions at order λ^2 . Both yield similar coupling strengths so that the stimulated Raman process does not beat the direct process in this case. However, we remark that in the absence of a magnetic field gradient, controllability of the $|\uparrow\downarrow\rangle$, $|\downarrow\uparrow\rangle$ manifold, without coupling to the $|T^\pm\rangle$

states, can not be achieved this way, as there is zero coupling through microwaves within this theory to the pure $|T^0\rangle$ state. However, with the magnetic-field gradients in the experiment, controllability is recovered with the constraint that operation pulses cannot be faster than db/\hbar .

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Author contributions

L.R.S., F.R.B., V.C. and T.M. performed the experiment; W.W. grew the heterostructure; T.M. fabricated the sample; L.R.S., J.D., J.M.T. and L.M.K.V. developed the theory; J.M.T. did the simulations; all authors contributed to the interpretation of the data and commented on the manuscript; and L.R.S., J.D., J.M.T. and L.M.K.V. wrote the manuscript.

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