Measuring charge based quantum bits by a superconducting single-electron transistor

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Single electron transistors have been proposed to be used as a read-out device for Cooper pair charge qubits. Here we show that a coupled superconducting transistor at a threshold voltage is much more effective in measuring the state of a qubit than a normal-metal transistor at the same voltage range. The effect of the superconducting gap is to completely block the current through the transistor when the qubit is in the logical state 1, compared to the mere diminishment of the current in the normal-metal case. The time evolution of the system is solved when the measuring device is driven out of equilibrium and the setting is analysed numerically for parameters accessible by lithographic aluminium structures.

Nanoscale devices such as Cooper pair boxes or coupled quantum dots have been suggested as scalable and integrable realisations of quantum bits. The two logical states of a qubit are the different charge states, or in the case of a flux qubit, flux states of the system. There are several proposals for quantum gates [1, 2] and for in-states of a qubit are the different charge states, or in the integrable realisations of quantum bits. The two logical states of a Cooper pair charge qubit [8, 9]. Also, a superconducting SET in the Josephson current regime (low biasing voltage) has been experimentally tested [10]. In this work we show that in the regime of high biasing voltages, the superconducting SET [13, 12] leads to a highly efficient quantum nondemolition measurement [13] due to the blocking effect of the gap.

The setting is shown in Fig. 1. In the upper part, the Cooper pair box forms the qubit, its state characterised by the number of excess Cooper pairs in the box, \( n \). In the lower part, the superconducting single-electron transistor is capacitively coupled to the qubit with its state characterised by the excess charge on the island, \( eN \). In addition, the quantum number \( m \) counts the number of charges passing through the SET in left-to-right direction. Without a biasing voltage \( V_{\text{bias}} \) across the SET, there is no dissipative current and no information is received. Moreover, in the absence of a dissipative environment, no dephasing of the composite system will occur and quantum operations on the qubit can be performed.

In order to perform the measurement, a biasing voltage is applied. As different qubit eigenstates correspond to different conductance in the SET, by observing the current one receives information on the state of the qubit. The time needed for the current to give the essential information is called the measurement time. The back action caused by the SET dephases the qubit and eventually destroys also the logical state of the qubit. The corresponding time scales are called the dephasing and the mixing time, respectively. For a good non-demolition measurement of the logical state \( |a\rangle |2 \) and \( |b\rangle |0 \) in \( a|0\rangle + b|1\rangle \), one expects to have a much longer mixing time than a measurement time scale. We show that for a superconducting SET, in a parameter range accessible by aluminium structures, the ratio between the mixing and measurement time scales is in excess of 600.

The total Hamiltonian consists of three parts: the Hamiltonians of the SET, the qubit and the interaction \( H_{\text{set}}, H_{\text{qb}} \) and \( H_{\text{int}} \), respectively. The SET Hamiltonian is defined as

\[
H_{\text{set}} = E_{\text{set}} \left( N - Q_{\text{set}} \right)^2 + H_T + H_L + H_R + H_I, \tag{1}
\]

where \( E_{\text{set}} \) is the charging energy, \( Q_{\text{set}} \) the gate charge of the transistor and

\[
H_T = \sum_{kk'\sigma} T_{kk'\sigma} c_{kk'\sigma}^{\dagger} e^{i\phi} + T_{kk'\sigma}^{\dagger} c_{kk'\sigma} e^{-i\phi} e^{i\Psi} + H.c. \tag{2}
\]

describes quasiparticle tunnelling within the SET. The phase differences \( \phi \) and \( \Psi \) are the conjugate variables of \( N \) and \( m \), respectively. Thus, the operator \( e^{i\phi} (e^{i\Psi}) \) corresponds to quasiparticle tunnelling onto the island (right electrode) of the SET, increasing the quantum number \( N \).
(m) by one. The last three terms in Eq. 2 are defined as $H_r = \sum_{k\sigma} \epsilon_{k\sigma} c_{k\sigma}^\dagger c_{k\sigma}$, $(r = L, R, I)$, where $\epsilon$ labels the transverse channels including the spin and $k$ labels the wave vector within one channel. Describe the non-interacting electrons in the left electrode, the right electrode and the island, respectively. Cooper pair tunnelling is excluded from the SET Hamiltonian due to the high biasing voltage. However, this approximation is not valid for a low voltage, and the contribution from the Cooper pair tunnelling should be added when analysing the effect of the SET on the qubit during logical operations.

The qubit Hamiltonian is defined as

$$H_{qb} = E_{qb} (n - Q_{qb})^2 - E_J \cos \Theta,$$

(3)

where $E_{qb}$ is the charging energy, $Q_{qb}$ qubit’s gate charge and the last term describes the transfer of Cooper pairs to and from the box ($e^{int}|n\rangle = |n+1\rangle$). The quasiparticle tunnelling in the qubit is suppressed by the Coulomb blockade, and the microscopic degrees of freedom have already been integrated out. The interaction Hamiltonian describes the Coulomb interaction between the qubit and the transistor and is defined as

$$H_{int} = E_{int} n N,$$

(4)

where $E_{int}$ is the charging energy. All charging energies $E_{int}$, $E_{set}$ and $E_{qb}$ are determined by the capacitances of the system, and the gate charges $Q_{qb} = -eV_n/E_{qb}$ and $Q_{set} = -eV_N/2E_{set}$ depend on the effective gate voltages $V_n$ and $V_N$ that, for a symmetric bias, are determined by the gate voltages $V_{qb}$ and $V_{set}$ as given in [3].

Concentrating on the values of $Q_{qb}$ around the degeneracy point $Q_{qb} = \frac{s}{N}$, only the low-energy charge states $n = 0$ and $n = 1$ are relevant. This assumption is required for the Cooper pair box to constitute a quantum bit. In the qubit’s two-state approximation, the diagonalised operator becomes $H_{qb} = -\frac{i}{2} \Delta E \sigma_z$, where

$$\Delta E = \sqrt{[E_{qb}(1 - 2Q_{qb})]^2 + E_J^2},$$

(5)

while the operator $n$ becomes non-diagonal $n = \frac{1}{2} (1 - \cos(\eta) \sigma_z - \sin(\eta) \sigma_x)$, where the mixing angle is given by tan $\eta = E_J / [E_{qb}(1 - 2Q_{qb})]$. Assuming the gate charge $Q_{qb}$ sufficiently different from the degeneracy point 0.5, the mixing angle $\eta$ is small in the charge-qubit approximation $E_J \ll E_{qb}$. By rearranging the operators, the final form for the total Hamiltonian can be written as $H = H_0 + H_T$, where

$$H_0 = H_L + H_R + H_{qb} + E_{set} (N - Q_{set})^2 + H_{int}.$$

(6)

We analyse the measurement process by mastering equation techniques. The master equation for the system reads

$$\frac{\partial \sigma(t)}{\partial t} + \frac{i}{\hbar} [H_0, \sigma(t)] = \text{Tr}_{L,R,1} \int_0^t \Sigma(t - t') \sigma(t') dt',$$

(7)

where the trace is taken over the microscopic degrees of freedom of the transistor’s left and right electrodes and the island. The elements of the transition matrix $\Sigma(t - t') = -\frac{1}{\pi^2} \text{Tr} \{ [V, U(t - t') \{ V, \cdot \} U(t' - t)] \}$ are calculated by using the diagrammatic technique developed in [1] and [3]. However, the rules for converting diagrams into integral equations must be changed to include the superconducting density of states

$$N(x) = \begin{cases} \frac{|x|}{\sqrt{2^{\Delta^2} - x^2}} N_0, & \text{if } |x| > \Delta \\ 0, & \text{otherwise}, \end{cases}$$

(8)

where $\Delta$ is the superconducting gap. In zero-temperature limit, the matrix elements $\Sigma(t - t')$ can be written as

$$\Sigma_{k,l}(t - t') = F(t - t')^2 e^{-i E(t - t')/\hbar},$$

(9)

where $F(t) = \int_0^\infty N(x) e^{-(x+i\varepsilon)^2} dx$, and $\varepsilon$ is a cut-off exponent which ensures the convergence of the integral. With the superconducting density of states $N(x)$, the function $F(t)$ is a so-called Basset function of first order [14] which must be analysed numerically.

By performing the Laplace transform on the master equation (7), the right-hand side becomes $\Sigma(s) \sigma(s)$. Assuming the density matrix $\sigma$ to change slowly in a time scale of $\hbar/E$, the calculations can be restricted to the regime $s \ll E$. In the normal-metal case, $\Sigma(s)$ varies only slowly as a function of small $s$, and therefore the zeroth-order approximation is reasonably good. However, when the energy $E/2$ is close to the gap energy $\Delta$, $\Sigma(s)$ has a strong $s$-dependence which is approximately linear for small $s$, as shown by Fig. 3. Using the linear approximation $\Sigma(s) = a + b \cdot s$ and performing the inverse Laplace transform, the right-hand side becomes $(\Gamma + \Lambda \partial/\partial t) \sigma(t)$. Moving all the derivatives of the density matrix to the left-hand side gives as a master equation

$$(1 - \Lambda) \frac{\partial \sigma(t)}{\partial t} + \frac{i}{\hbar} [\sigma(t), H_0] = \Gamma \sigma(t),$$

(10)

where $\Gamma$ and $\Lambda$ are tridiagonal matrices consisting of the zeroth- and first-order terms of the $\Sigma(s)$ matrix, respectively. To guarantee the existence of the inverse of $(1 - \Lambda)$, the elements of the $\Lambda$ coefficient are required to be small. This requirement is fulfilled when the tunnelling rate within the SET is small (that is, $|T^L|, |T^R| \ll 1$). Multiplying equation (10) from the left by $(1 - \Lambda)^{-1}$ the final form of the master equation is obtained as

$$\frac{\partial \sigma(t)}{\partial t} = (1 - \Lambda)^{-1} \left[ -\frac{i}{\hbar} [\sigma(t), H_0] + \Gamma \sigma(t) \right] = \tilde{\Gamma} \sigma(t).$$

(11)

This describes a Markovian evolution of the density matrix and the sequential tunnelling approximation, which is a reasonable assumption when the current in the SET is low.
The elements of the $\Sigma(s)$ matrix are analysed by explicitly writing the Laplace transform

$$\Sigma(s) = \int_0^\infty \Sigma(t)e^{-st}dt = \int_0^\infty F(t)^2e^{(-s+iE)t}dt. \quad (12)$$

The linear approximation is done by calculating $\Sigma(s)$ for two values of $s = 0.0$ and $s = 0.1$ and fitting a line.

Elements of the matrix $\tilde{\Gamma}$ describing the transitions within the qubit are proportional to a small mixing angle $\epsilon$ defined by $\tan \epsilon = \frac{E_{\text{int}} \sin \eta}{(\Delta E + E_{\text{int}} \cos \eta)}$. By approximating these elements by zeroes, the matrix $\tilde{\Gamma}$ separates into four parts: one part describing the system when the qubit is in (diagonal) state 00, one part when the qubit is in state 11 and two parts for the nondiagonal qubit states 01 and 10. The first two parts give rise to the measurement time as they describe the transistor’s evolution for different qubit eigenstates. The coherence of the qubit is described by the strength of the non-diagonal elements, and thus the rate at which the non-diagonal elements vanish gives the dephasing time. For the mixing time, one has to include the small mixing elements proportional to $\epsilon$. To ease the treatment below, a Fourier transformation is performed in $m$-space: $\sigma^k(t) = \sum_m e^{ikm} \sigma^m(t)$.

The measurement time is determined by the first two parts of the matrix $\tilde{\Gamma}$. The equation corresponding to the qubit state 00 is

$$\frac{d}{dt} \left( \frac{\sigma^{N,k}_{00}}{\sigma^{N+1,k}_{00}} \right) = \begin{pmatrix} -\Gamma_{L00} & \Gamma_{R00} e^{i\epsilon k} \\ \Gamma_{L00} & -\Gamma_{R00} \end{pmatrix} \begin{pmatrix} \sigma^{N,k}_{00} \\ \sigma^{N+1,k}_{00} \end{pmatrix}, \quad (13)$$

and the other part is a similar matrix equation for the qubit state 11. These equations describe slowly damping (due to shot noise) conductance peaks that propagate in time. The measurement has been performed once the peaks can be distinguished from each other, and following the treatment in [1] the measurement time $t_{\text{ms}}$ is defined as the time when the width of the peaks is smaller than the distance between their centers yielding

$$t_{\text{ms}} := \left( \frac{\Gamma_{00} - \Gamma_{11}}{\sqrt{2\Gamma_{00}f_0^2 + 2\Gamma_{11}f_1^2}} \right)^2, \quad (14)$$

where the group velocities are $v_i(\epsilon) := \frac{\Gamma_{L\epsilon i} \Gamma_{R\epsilon i}}{\Gamma_{L\epsilon i} + \Gamma_{R\epsilon i}}$, with $i = 0, 1$. The factors $f_i = \frac{(\Gamma_{L\epsilon i} + \Gamma_{R\epsilon i})}{(\Gamma_{L\epsilon i} + \Gamma_{R\epsilon i})}$ are known as Fano factors [3].

The dephasing time is determined by the evolution of the nondiagonal qubit states given by the remaining two parts of the matrix. Choosing $k = 0$ has the effect of tracing out the $m$-variables and simplifying the corresponding submatrices

$$\frac{d}{dt} \left( \frac{\sigma^{N+1}_{01}}{\sigma^{N+1}_{01}} \right) = \begin{pmatrix} i\Delta E^0 - \Gamma_{L01} & \Gamma_{R01} \\ i\Delta E^1 + \Gamma_{L01} & -\Gamma_{R01} \end{pmatrix} \begin{pmatrix} \sigma^{N+1}_{01} \\ \sigma^{N+1}_{01} \end{pmatrix}. \quad (15)$$

Again, these equations have plane wave solutions, and the corresponding eigenvalues determine the dephasing time of the system. The time corresponding to each eigenvalue $\lambda$ is defined as $1/\text{Re}(\lambda)$, and the dephasing time is chosen to be the longer of these two times. The shorter time describes partial dephasing. In the superconducting case these must be evaluated numerically.

In order to analyse the mixing time of the system, also the mixing coefficients must be included in the $\Gamma$ matrix. Once again, the SET degrees of freedom are irrelevant and $k$ is chosen to be 0. Now the matrix has eight eigenvalues: four of the eigenvalues describe the dephasing, two describe the measurement and one eigenvalue is zero (describing the trace preserving symmetry). The remaining real eigenvalue $\lambda_{\text{mix}}$ gives the mixing time as the $t_{\text{mix}} := 1/\lambda_{\text{mix}}$, which must be analysed numerically.

If the mixing time is very large compared to the measurement time, the measuring device disturbs the probability amplitudes of the qubit $|a|e^{i\theta}$ or $|b|e^{i\theta}$ only a little. The uncertainties of the charge $\Delta Q$ and its conjugate variable (phase or flux) $\Delta \Phi$ are linked to each other by the uncertainty principle $\Delta Q \cdot \Delta \Phi \geq \hbar/2$. According to this principle, if the precision of the charge measurement is very high, the phase becomes completely undetermined. The information that can be gathered from the SET contains no information of the phase and thus the precision of measuring the charge within the qubit $|a|^2$ or $|b|^2$ can be very high. Furthermore, since charge is not disturbed by the measurement for $t_{\text{mix}} \gg t_{\text{meas}}$, we are essentially doing the classic quantum non-demolition measurement.

The time scales are calculated for a specific set of parameters. With aluminium structures in mind, the superconducting gap of the SET is chosen to be $\Delta = 2.3 K$. 

Figure 2: The complex transition coefficient $\Sigma(s+iE)$ plotted as a function of $s$ with fixed $E = 10.0$ for different values of the gap $\Delta$. The real components are drawn in solid curves and the imaginary parts in dotted ones. The corresponding gap values are in the order from the topmost curve to the lowest curve: $\Delta = 0.0, \Delta = 3.0, \Delta = 4.5, \Delta = 4.9, \Delta = 5.1$ and $\Delta = 5.3$. For small values of $s$, the real components are nearly linear.
These fix the charging energies \( E \) to the state 1. The two peaks are clearly distinguishable from the capacitances normal-metal SET.

The curve in the box shows the corresponding evolution for a in the text. The graph shows the separation of the two peaks, parameters are for aluminium structures and they are given measurement times is very high, i.e. \( t \) gives the measurement time \( t \). The mixing effects are small, as the peaks respond to the qubit's state \( n = 1 \) is blocked by the presence of the gap. The mixing time is much larger than the measurement time \( t_{\text{mix}}/t_{\text{ms}} \gg 1 \). for parameter ranges accessible by aluminium structures.

Figure 3: The probability \( P(m, t) \) that \( m \) electrons have tunnelled in a superconducting SET plotted as a function of time \( t \) for initial amplitudes \( \sqrt{0.75} (n = 0) \) and \( \sqrt{0.25} (n = 1) \). The parameters are for aluminium structures and they are given in the text. The graph shows the separation of the two peaks, the faster corresponding to the qubit's state 0 and the slower to the state 1. The two peaks are clearly distinguishable from each other, and thus the measurement time is very small. The curve in the box shows the corresponding evolution for a normal-metal SET.

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[14] H. Schoeller and G. Schönh, Phys. Rev. B 50, 18436 (1994).
[15] J. Spanier and K.B. Oldham, An Atlas of Functions (Springer-Verlag, Berlin, 1987).
[16] The Basset function of integer order \( K_n \) is generally defined as an elliptic integral \[ K_n(z) = \frac{\sqrt{z}}{2\pi \Gamma(1/2)(2n+1)} \int_0^\infty \frac{e^{-z} z^{n+1/2}}{(z+1)^{n+3/2}}dz. \] The numerical analysis of the Basset function for small values of the argument \( z \) is based on presenting it as a Neumann series of Bessel functions. For a large argument, a good algorithm can be written by applying the Padé technique on the asymptotic approximation for large \( z \)
\[ K_n(z) \sim \sqrt{\frac{\pi}{2z}} z^{-n} \left[ 1 + \sum_{j=1}^{\infty} \frac{(y-1/2)\cdots(y+j-1)}{2^j j!} \right], \]
where \( \langle x \rangle_j = x(x+1)(x+2)\cdots(x+j-1) \) is the Fochhammer polynomial.