1. Introduction

Standard quantum dynamics cannot describe how a particular measured outcome is realized in an experiment—this feature underlies our failure to unify classical physics with its quantum roots. A direct response to this problem is to take the view that standard quantum theory is incomplete and that a complete quantum theory should provide a dynamical description of quantum state reduction. In this way, it is hoped to provide an objective description of quantum measurement without recourse to a classical measuring device.

The great success of quantum theory does not leave much room to manoeuvre when proposing alternatives. Any new theory must agree with the experimentally verified predictions of quantum theory. This roughly implies that the dynamics should reduce to Schrödinger dynamics for micro systems and that state reduction effects should become more important with system size. More specifically, the form of any modifications we make should involve a random factor (at some level of approximation) to reflect the random nature of state reduction, and we would expect some nonlinearity in the dynamics, reflecting a feedback from the state vector to the probability of the outcome [1].
These features can be adequately represented by a stochastic differential equation for the quantum state \([2, 3]\) (for general reviews see \([4, 5]\)), the most well-known model of this type being the continuous spontaneous localization (CSL) model \([6]\). In the CSL model, state reduction occurs onto a basis of smeared position states leading to definite localized positions for systems of macroscopic size. However, there is scope for further model building, principally because we can choose from many other preferred bases for state reduction (e.g. \([7–9]\)). Until we have observational evidence of state reduction dynamics we are free to experiment with different proposals.

Here, we investigate the state reduction process for a simplified form of quantum electrodynamics. We consider a charged two-state quantum system interacting with a quantum electromagnetic field mode. We investigate a model which describes state reduction onto a coherent field state basis. State reduction for the two-state system happens indirectly via standard interaction with the field mode. We find that the two-state system evolves towards a generalized current eigenstate.

The appealing feature of this model is that the stable eigenstates of the system are a natural description of the classical world whereby the field is represented by a coherent state and the two-state system (representative of charged matter) by a current eigenstate. Modifications to Schrödinger’s equation only involve the field mode—it is not necessary to make further modifications in order to independently describe state reduction in the two-state system. Also, it is possible to generalize the model to field theory without generating infinite particle creation \([10]\) (this is a problem which appears, e.g., with the CSL model \([11, 12]\)), although it should be pointed out that this results in a loss of Lorentz invariance.

The idea that dynamical state reduction might be induced from one field to another is not new. It has appeared in previous attempts to describe state reduction in fermionic field theory by invoking a scalar field with state reduction properties to which the fermion is coupled \([11, 12]\). To our knowledge the process has not been previously studied in numerical detail.

In the following section we outline our model, detailing the Hamiltonian for our simplified QED system and the modified stochastic dynamics of the state vector. In section 3, we consider the evolution of a specific initial state vector composed of a superposition of current states. We outline the state reduction process and derive a characteristic time scale. In section 4, we present some numerical results which confirm the analytic results and intuition gained in section 3. We end with some conclusions.

2. The model

2.1. QED Hamiltonian

Our simplified QED Hamiltonian is given by

\[
H = \omega a^\dagger a + \frac{1}{2} \nu \sigma_z + g \sigma_x (a + a^\dagger).
\]

(1)

The \(\sigma_i\) operators are the standard Pauli matrices and \(a, a^\dagger\) are the annihilation and creation operators for a standard harmonic oscillator. The parameters \(\omega, \nu\) and \(g\) are constants which we interpret as the field-mode frequency, the two-state energy gap and the coupling, respectively.

Models of this type are commonly used in quantum optics to describe the interaction of a two-state atom with an electromagnetic field \([13]\). In particular, a further approximation to the interaction term to exclude processes such as those where a photon is created as the atom is exited (the rotating wave approximation) leads to the celebrated Jaynes–Cummings model. Alternatively, this model can be derived from full QED by restricting space to just two points (time remains continuous) and considering only one photon polarization. The result is two
fermion states and one field mode. This simplification is inadequate for a number of reasons. Most importantly, the model can no longer be described as a gauge theory as there is no local symmetry associated with the fermion field. We also lose the concepts of conserved 4-current and relativistic invariance.

On the other hand, our simplified Hamiltonian has the structure of the QED Hamiltonian: it is composed of free matter and field parts and an interaction between a current operator $\sigma_x$ and the field. (That $\sigma_x$ corresponds to the current can be seen as we restrict full QED to a space of two points.) These features suffice in order to make a tentative attempt to understand quantum state reduction in QED.

Since this model can be applied to quantum optical systems, it is worth pointing out that stochastic adjustments to Schrödinger dynamics have been used in the past to describe the effect of an environment on optical and atomic behaviour [14–16].

2.2. Stochastic Schrödinger equation

We shall consider a modified Schrödinger equation of the form

$$\frac{d}{dt}|\psi_t\rangle = \left\{ \left(-\frac{i}{\hbar} H - \frac{1}{2}\lambda^2 a^\dagger a + \frac{1}{2}\lambda^2 a^\dagger a\right) dt + \lambda \left(a - \frac{1}{2}a^\dagger\right) dB_t + \frac{1}{2}\lambda a^\dagger dB^*_t \right\} |\psi_t\rangle,$$

where $\lambda$ is a constant coupling parameter with units of $[\text{time}]^{-1/2}$ and

$$a_t = \langle \psi_t | a | \psi_t \rangle$$

is the conditional expectation of $a$. The complex Wiener process $B_t$ satisfies $dB_t dB^*_t = 2 dt$, $dB_t dB^*_t = dB^*_t dB_t = 0$, and $E[dB_t] = 0$, where $E[\cdot]$ denotes unconditional expectation [17]. Note that the Schrödinger equation is recovered by setting $\lambda = 0$ and that the additional $\lambda$-dependent terms only involve field operators.

Equation (2) is a stochastic differential equation for the state. It can be derived by looking for a state diffusion process describing stochastic fluctuations in the quantum probability amplitudes of possible $a$-eigenstates [4]. The probabilities for the stochastic fluctuations must be given by the quantum probabilities at each point in time. This ensures that stochastic outcomes match with the initial quantum predictions (and leads to nonlinearity). The equation has the property that general states evolve towards $a$-eigenstates on a time scale approximately inversely proportional to the variance in $\Delta a$ (see below).

Using the rules of Itô calculus we can show that

$$d\langle \psi_t | \psi_t \rangle = 0,$$

i.e. the norm of the state is preserved under the dynamics (2). This allows us to maintain the quantum interpretation of the state as a probability amplitude.

The process for the conditional expectation of $a$ is found to be

$$da_t = \left\{ -i\omega a_t - g\sigma_z a_t - \frac{1}{2}\lambda^2 a_t^2 \right\} dt + \lambda (\langle \Delta a \rangle^2) dB + \lambda \langle \Delta a \rangle dB^*,$$

where $\Delta a = a - a_t$, and we have used the convenient notation $\langle \cdot \rangle_t = \langle \psi_t | \cdot | \psi_t \rangle$. When the state $|\psi_t\rangle$ is an $a$-eigenstate, the process $a_t$ is nonstochastic. In this case, the effect of the modifications to Schrödinger’s equation is a decay in the magnitude of $a_t$ at a rate given by $\lambda^2$. This can be interpreted as being due to the fact that the classical stochastic process $B_t$ in equation (2) is only coupled to photon annihilation events. This results in a gradual loss of energy. Appropriate tuning of $\lambda$ is required to ensure that field energy loss is negligible whilst state reduction time scales are appropriately small for strong fields [10].

In order to demonstrate the state reducing properties of equation (2) we consider the conditional variance in $a$ defined by

$$\text{Var}_t(a) = \langle (\Delta a)^2 \rangle_t,$$

where
The conditional variance process is found to be

$$
\text{d} \text{Var}_t(a) = i\{[\sigma_x(a - a^\dagger)], -\{[\sigma_x], (a - a^\dagger)]\} \text{d}t
- 2\lambda^2 \text{Var}_t(a) \text{d}t - 2\lambda^2 \left\{ \left| \left\langle (\Delta a)^2 \right\rangle \right|^2 + \text{Var}_t^2(a) \right\} \text{d}t
+ \lambda \left\langle |\Delta a|^2 \Delta a \right\rangle, \text{d}B_t + \lambda \left\langle |\Delta a^\dagger|^2 \right\rangle, \text{d}B_t^\dagger.
$$

(7)

The first line on the right side is the effect on \( \text{Var}_t(a) \) due to the interaction between the two-state system and field. We see that \( \text{Var}_t(a) \) increases if there is a negative conditional covariance between \( \sigma_x \) and \( -i(a - a^\dagger) \). As \( \text{Var}_t(a) \) increases, other terms on the right side of equation (7) become more important.

If we ignore the interaction with the two-state system the variance process is a super-martingale (the drift of the process is negative-definite). Integrating and taking the unconditional expectation of equation (7) we then find

$$
\mathbb{E}[\text{Var}_t(a)] = \text{Var}_0(a) - 2\lambda^2 \int_0^t \text{d}u \mathbb{E}[\text{Var}_u(a)] - 2\lambda^2 \int_0^t \text{d}u \mathbb{E}\left\{ \left| \left\langle (\Delta a)^2 \right\rangle \right|^2 + \text{Var}_u^2(a) \right\}.
$$

(8)

This demonstrates that the expected variance will decrease in time and therefore that the field approaches an \( a \)-eigenstate.

We can estimate the time scale \( \tau_a \) for state reduction to an \( a \)-eigenstate from the following:

$$
\frac{\mathbb{E}[\text{Var}_t(a)] - \text{Var}_0(a)}{\text{Var}_0(a)} \sim \frac{t}{\tau_a}.
$$

(9)

By freezing the stochastic terms on the right side of equation (8) at time \( t = 0 \) we find

$$
\tau_a \sim \frac{\text{Var}_0(a)}{\lambda^2 \left\{ \left| \left\langle (\Delta a)^2 \right\rangle \right|^2 + \text{Var}_0^2(a) \right\}}.
$$

(10)

Even though \( \lambda \) must be very small to prevent energy loss, the reduction time scale \( \tau_a \) can be small if the variance in \( a \) is large. This demonstrates the specific way in which the effectiveness of state reduction is dependent on the size of the system. A superposition of greatly differing \( a \)-eigenstates has large variance and state reduction will occur rapidly; if \( \text{Var}_0(a) \sim \left| \left\langle (\Delta a)^2 \right\rangle \right| \sim \mathcal{O}(1) \) then state reduction takes an extremely long time due to the small value of \( \lambda \). In principle, the dependence of the reduction time on the \( a \)-variance is a testable result. An experiment designed to test this might look for diminishing quantum interference as a signature of the state reduction [18]. The technical difficulty would be to eliminate any environmental influence.

In order to see that stochastic probabilities for outcomes match with the initial quantum estimates we consider an initial superposition of coherent states

$$
|\psi_0\rangle = \sum_i c_i |\alpha_i\rangle,
$$

(11)

where it is supposed that the coherent states \( |\alpha_i\rangle \), where \( a|\alpha_i\rangle = \alpha_i |\alpha_i\rangle \), are sufficiently separated in phase space such that \( \langle \alpha_i | \alpha_j \rangle \simeq \delta_{ij} \) (this requires that \( |\alpha_i - \alpha_j| \gg 1 \) for \( i \neq j \)). We assume for simplicity that \( H = 0 \) since we wish to focus on the reduction dynamics.

Note that the coherent states are eigenstates of all terms in the evolution equation (2) except the term \( -\lambda^2 a^\dagger a \). This means that most terms in the evolution equation simply modify the coefficients \( c_i \), leaving the set of states \( \{|\alpha_i\rangle\} \) intact. The term \( -\lambda^2 a^\dagger a \) is responsible for diffusing the initial set of coherent states. In order to analyse its effects, consider a state evolution involving just this term:

$$
\frac{\text{d}}{\text{d}t} |\psi_t\rangle = -\lambda^2 a^\dagger a |\psi_t\rangle.
$$

(12)
This equation has the solution
\[ |\psi_t\rangle = \sum_i c_i \exp \left\{ \frac{1}{2} \left( |\alpha_i e^{-i\lambda t}|^2 - |\alpha_i|^2 \right) \right\} |\alpha_i e^{-i\lambda t}\rangle. \] (13)

This indicates that the individual coherent state contributions to the overall superposition will decay to the ground state on a time scale \(\lambda^{-2}\). For times \(t \ll \lambda^{-2}\) we can ignore this effect and assume that the set of coherent states contributing to the superposition is fixed as \(|\alpha_i\rangle\).

We can define an approximate coherent state projection operator by
\[ P_{\alpha_k} = |\alpha_k\rangle \langle \alpha_k| . \]

For \(\tau_a \ll t \ll \lambda^{-2}\) the conditional expectation \(\langle \psi_t | P_{\alpha_k} | \psi_t \rangle\) is approximately either 1 or 0 depending on whether the state is reduced to \(|\alpha_k\rangle\) or not. From equation (2) we find
\[ d\langle \psi_t | P_{\alpha_k} | \psi_t \rangle = \lambda|c_k(t)|^2 (\alpha_k - a_i) dB_t + \lambda|c_k(t)|^2 (\alpha_k^* - a_i^*) dB_T^*, \] (14)
from which we conclude that
\[ \langle \psi_0 | P_{\alpha_k} | \psi_0 \rangle = \mathbb{E}\left[ \langle \psi_t | P_{\alpha_k} | \psi_t \rangle \right] = \mathbb{E}\left[ 1_{|\psi_t\rangle = |\alpha_k\rangle} \right], \] (15)
valid for \(\tau_a \ll t \ll \lambda^{-2}\). The left-hand equality gives the initial quantum probability for the outcome \(|\psi_t\rangle = |\alpha_k\rangle\); the right-hand equality gives the stochastic probability for the same outcome.

3. An experiment

Consider an initial situation where the field is in a coherent state \(|\alpha\rangle\) and the two-state system is in a superposition of current \((\sigma_x)\) eigenstates:
\[ |\psi(0)\rangle = \left\{ c_1 \begin{pmatrix} 1 \\ 1 \end{pmatrix} + c_2 \begin{pmatrix} 1 \\ -1 \end{pmatrix} \right\} |\alpha\rangle. \] (16)

In order to understand the behaviour of this model without explicitly solving it we consider some possible scenarios depending on the relative sizes of the model parameters.

Since the field state is initially coherent, the variance in \(a\) is initially small. As this is our measure of system size, we can assume that equation (2) initially reduces to the Schrödinger equation.

3.1. \(\omega, \nu\) domination

First consider the case where the coupling \(g\) tends to zero. We then have
\[ |\psi(t)\rangle_{g=0} = \exp \left\{ -it \left( \omega a^\dagger a + \frac{1}{2} \nu \sigma_z \right) \right\} |\psi(0)\rangle = \left\{ c_1(t) \begin{pmatrix} 1 \\ 1 \end{pmatrix} + c_2(t) \begin{pmatrix} 1 \\ -1 \end{pmatrix} \right\} |e^{-it\omega}\alpha\rangle, \] (17)
where
\[ c_1(t) = c_1 \cos \frac{1}{2} t \nu - ic_2 \sin \frac{1}{2} t \nu, \quad c_2(t) = c_2 \cos \frac{1}{2} t \nu - ic_1 \sin \frac{1}{2} t \nu. \] (18)
Here we find independent state-space rotations for the two-state system and the field.
3.2. \( g \) domination

Next consider the case where the field frequency \( \omega \) and the two-state energy gap \( \nu \) tend to zero. In this limit we find the following solution to the Schrödinger equation:

\[
|\psi(t)\rangle_{\omega=\nu=0} = e^{i\phi_1} \left( \begin{array}{c} 1 \\ i \end{array} \right) |(\pi -igt)\rangle + c_2 e^{i\phi_2} \left( \begin{array}{c} 1 \\ -1 \end{array} \right) |(\pi +igt)\rangle,
\]

where \( \phi_i \) are real-valued phase factors. The interaction between two-state system and field leads to a correlation between current eigenstates and field coherent states.

These two limits provide an overall picture of two competing processes. When the interaction \( g \) is dominant, different current eigenstates become associated with different coherent field states. When the interaction is negligible we find that current states rotate into one another.

3.3. \( \lambda \) domination

Consider the scenario \( \omega \sim \nu \ll g \). From equation (19) we can estimate the growth of variance in \( a \):

\[
\text{Var}_a(a) \simeq 4g^2t^2|c_1|^2|c_2|^2 \simeq |(\Delta a)^2|_t,
\]

and the conditional covariance between \( \sigma_x \) and \(-i(a - a^\dagger)\):

\[
\text{Cov}_t(\sigma_x, -i(a - a^\dagger)) = -i(\sigma_x(a - a^\dagger))_t + i(\sigma_x)_t(\langle a - a^\dagger \rangle)_t
\]

\[
\simeq -8gt|c_1|^2|c_2|^2
\]

(21)

(the approximation results from the fact that we have assumed that different coherent states are nearly orthogonal). As the variance grows, \( \lambda \)-effects come to dominate the evolution. This can be seen from equations (20) and (21) by noting that the \( \text{Var}_a(a) \) and \( |(\Delta a)^2|_t \) terms in equation (7) will soon outgrow the \( \text{Cov}_t(\sigma_x, -i(a - a^\dagger)) \) term. Let us define an approximate time \( s \) at which a transition from \( g \) to \( \lambda \)-dominated behaviour happens. Once in the \( \lambda \)-dominated phase, the \( a \)-variance will tend to zero on a time scale given by equation (10).

The stochastic evolution effectively chooses between the possible states \((1, 1)\rangle |\pi -igt\rangle\) and \((1, -1)\rangle |\pi +igt\rangle\) leading towards a definite field coherent state and a definite current state.

Using (10) and (20) we find that a total state reduction time scale for the current state is given by

\[
\tau_{\sigma_x} \sim s + \frac{1}{\lambda^2 g^2 |c_1|^2 |c_2|^2}.
\]

In order to determine \( s \) we find the value which minimizes \( \tau_{\sigma_x} \). Setting

\[
\frac{d}{ds} \tau_{\sigma_x} = 0 \quad \implies \quad 1 \sim \frac{1}{\lambda^2 g^2 s^3 |c_1|^2 |c_2|^2},
\]

therefore

\[
\tau_{\sigma_x} \sim \frac{1}{\lambda^2 g^2 s^3 |c_1|^2 |c_2|^2}.
\]

This is the characteristic time scale for state reduction of a superposition of current eigenstates in a two-state system coupled to a field mode.
Figure 1. Growth and subsequent decay of $E[\text{Var}_a(t)]$ (filled line). A typical example of a random sample evolution of $\text{Var}_a(t)$ is also shown (dashed line). The initial state involves an equal superposition of two different current states. The coupling between current and field mode leads at first to a growth in $\text{Var}_a(t)$ via standard Hamiltonian dynamics. When $\text{Var}_a(t)$ is sufficiently large, the stochastic dynamics cause a decay in variance as the $a$-state collapses. Parameters are $\omega = \nu = 0.5$, $g = 4.0$, and $\lambda = 0.2$.

4. Numerical results

In order to test these results we have implemented a numerical simulator. Starting from an initial state of the form given in equation (16) with coefficients $c_1 = c_2 = 1/2$ and with $\alpha = 4$, we have generated sample state evolutions according to equation (2). We have chosen $\omega = \nu = 0.5$, $g = 4.0$ and $\lambda = 0.2$. This choice of parameters are such that $g$ dominates the initial evolution. The parameter $\lambda$ is chosen to be large enough to observe state reduction for the modestly sized numerical system, but small enough that energy loss is not too significant.

Figure 1 shows the Monte Carlo estimate for the process $E[\text{Var}_a(t)]$ (filled line). We used 1000 random sample paths to estimate unconditional expectation values at each point in time. The unconditional expectation of the $a$-variance is seen to increase in proportion to $t^2$ (see equation (20)). After a certain time the growth begins to slow. It then decreases in accordance with equation (8). When the unconditional expectation of the $a$-variance is zero the state must be a coherent state.

Also shown in figure 1 is an example of a random sample evolution of $\text{Var}_a(t)$ (dashed line) demonstrating the stochastic behaviour. This process indicates that the initial phase of the evolution is non-stochastic. After a certain time the evolution rapidly becomes dominated by stochastic movements with a notable downward drift. The conditional $a$-variance finally becomes stable again as its value tends to zero (at around $t = 2$).

Figure 2 shows the Monte Carlo estimate for $E[\text{Var}_a(\sigma_x)]$ (filled line) where
\[
\text{Var}_a(\sigma_x) = 1 - \langle \sigma_x \rangle^2.
\]

This provides a measure of whether the system is in a current eigenstate. Again we used 1000 random sample paths to estimate unconditional expectation values. One sample evolution for
Figure 2. Evolution of the unconditional expectation of variance in the current operator $\langle \text{Var}(\sigma_x) \rangle$ (filled line). A random sample path for $\text{Var}(\sigma_x)$ is also shown (dashed line). An initial (equal) superposition of two different current states ($\langle \sigma_x \rangle_0 = 0$) undergoes state reduction to a definite current state ($\langle \sigma_x \rangle_t = 1$).

Figure 3. Expected current reduction time (defined as $E[\tau_{\sigma_x}] = E[\inf \{t > 0; |\langle \sigma_x \rangle_t| > 0.99\}]$) as a function of the current-field coupling $g$. Other parameters are $\omega = \nu = 0, \lambda = 0.2$. The filled line shows a fit to the curve $g^{-2/3}$. The error bars show the one standard deviation region for the reduction time estimated from the simulated data.

$\text{Var}(\sigma_x)$ is shown explicitly (dashed line). We see that the $\sigma_x$-variance is slow to change at first since the induced state reduction requires an initial development of correlation between current
Figure 4. Histogram of current reduction times (defined as \( \tau_{\sigma_x} = \inf \{ t > 0 ; |\langle \sigma_x \rangle_t | > 0.99 \} \)) for \( g = 3 (\omega = v = 0, \lambda = 0.2) \). The smoothed line is the Gaussian kernel density estimate.

eigenstates and coherent states. Eventually (at around \( t = 2 \)) \( \mathbb{E}[\text{Var}_t(\sigma_x)] \to 0 \) indicating that the two-state system must end up in a current eigenstate.

We define the reduction time in our numerics as the stopping time at which \( |\langle \sigma_x \rangle_t | > 0.99 \), i.e.

\[
\tau_{\sigma_x} = \inf \{ t > 0 ; |\langle \sigma_x \rangle_t | > 0.99 \}.
\]

Figure 3 shows numerical estimates of the expectation of this stopping time for a range of \( g \) values. Other parameters are chosen to be \( \lambda = 0.2 \) and \( \omega = v = 0 \). Expectations are determined using 100 sample paths for each point. The estimates are seen to almost perfectly fit the curve \( kg^{-2/3} \) with \( k = 3.593 \). This confirms the form of the characteristic state reduction time given in equation (24). The error bars indicate a range of one standard deviation for \( \tau_{\sigma_x} \).

We observe that the variance in the stopping time decreases as \( g \) increases.

Figure 4 shows a histogram of \( \tau_{\sigma_x} \) results for 100 sample paths with \( g = 3 \). Also shown is the Gaussian kernel density estimate [19]. The distribution of stopping times is approximately normal with a slight skew. The important result is that the density of events with large deviation from the mean stopping time falls away rapidly.

We find that occurrences of \( |\langle \sigma_x \rangle_t | \to 1 \) and \( |\langle \sigma_x \rangle_t | \to -1 \) are even to within statistical error. This agrees with standard quantum predictions for a measurement of \( \sigma_x \) given the initial conditions.

5. Conclusions

We have investigated the state reduction dynamics described by a modified Schrödinger equation for a simplified QED system. Our simplified QED system is composed of a charged two-state system interacting with an electromagnetic field mode. The modified dynamics result in state reduction to an overcomplete coherent field state basis. At the same time, the electromagnetic interaction leads to the development of correlation between current states and field coherent states. This leads to an induced state reduction to a current state basis for the
two-state system. We have considered evolution from a specific initial condition composed of an equal superposition of the two possible current eigenstates with a coherent field state. We have implemented a numerical simulator to generate sample state evolutions. These have confirmed our analytic predictions for the general behaviour and the time scales involved.

The time scale for state reduction of the two-state system is found to behave as $g^{-2/3} \text{Var}^{-1/3}(\sigma_x)$. We might speculatively suggest that the reduction time scale for more general charged matter systems might scale as $\text{Var}^{-1/3}(j_\mu)$ for 4-current $j_\mu$.

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