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# Comments on minimal interactions in quantum mechanics 

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#### Abstract

We study minimal interactions in quantum systems characterized by position and momentum operators defined as the direct product of a traceless finite matrix and an ordinary canonical coordinate.


## 1. Introduction

There exists a large variety of quantum mechanical systems described by a complete set of compatible (non-interferring) observables which include, as an obvious additional entity, a multiple of a reducible identity matrix. Consider, for instance, the system described by the Jaynes-Cummings model in quantum optics which consists of a two-level atom interacting with a cavity mode [1-3], the well known supersymmetric systems in $(1+1)$ dimensions [4-8], and the familiar ones described by the Pauli wave equation. A common feature of these systems is that their Hilbert spaces are acted upon by an identity matrix of the form $I_{2 \times 2} \otimes I$, where $I$ is an infinite-dimensional identity matrix. There exists a non-trivial way of representing the position and momentum observables of the corresponding systems by making them formally 'traceless' operators. This property permits the introduction of a large number of new minimal interactions into the corresponding free particle wave equations. To be specific, let us consider a general quantum system described by canonical coordinates $Q_{i}$ and $P_{j}$ satisfying the Heisenberg algebra

$$
\begin{equation*}
\left[Q_{i}, P_{j}\right]=\mathrm{i} \hbar \mathbb{I} \delta_{i j} \tag{1.1}
\end{equation*}
$$

where $\mathbb{I} \equiv I_{n \times n} \otimes I$ represents a $n$-block identity matrix such that we may realize these operators in the general form

$$
\begin{equation*}
Q_{i}=\hat{\eta} \otimes q_{i} \quad P_{j}=\hat{\eta} \otimes p_{j} \tag{1.2}
\end{equation*}
$$

where $p_{j}=-\mathrm{i} \hbar \partial / \partial q_{j}$ and $\hat{\eta}$ is a constant $n \times n$ Hermitian matrix operator satisfying $\hat{\eta}^{2}=I_{n \times n}$. From equation (1.2) we can define a label $\Delta$ associated with each representation of the Heisenberg algebra (1.1)

$$
\begin{equation*}
n \geqslant \Delta\left(Q_{i}, P_{j}\right) \equiv|\operatorname{Tr} \hat{\eta}| \geqslant 0 \tag{1.3}
\end{equation*}
$$

Representations satisfying $\Delta=n$ correspond to the usual ones ( $\hat{\eta}=I_{n \times n}$ ) where $Q_{i}, P_{j}$ are reducible operators for $n \geqslant 2$. In this paper we shall be concerned with representations for which $\Delta=0$, i.e. $n$ is an even integer and $Q_{i}, P_{j}$ here are formally 'traceless' operators.

The Hilbert space is abstractly defined as

$$
\begin{equation*}
\mathcal{H}=L^{2}\left(\mathbb{R}^{3}\right) \otimes \mathbb{C}^{n} \tag{1.4}
\end{equation*}
$$

It consists of $n$-component column vectors

$$
\Psi(\boldsymbol{q}, t)=\left(\begin{array}{c}
\psi_{1}(\boldsymbol{q}, t)  \tag{1.5}\\
\vdots \\
\psi_{n}(\boldsymbol{q}, t)
\end{array}\right)
$$

where each component $\psi_{i}$ is a complex valued function of the four-dimensional (flat) spacetime coordinates $\boldsymbol{q}, t$. The scalar product is given by

$$
\begin{equation*}
(\Psi, \Phi) \equiv \int_{\mathbb{V} \subset \mathbb{R}^{3}} \Psi^{\dagger}(\boldsymbol{q}, t) \Phi(\boldsymbol{q}, t) \mathrm{d}^{3} q=\int_{\mathbb{V} \subset \mathbb{R}^{3}} \sum_{i=1}^{n} \psi_{i}^{*}(\boldsymbol{q}, t) \phi_{i}(\boldsymbol{q}, t) \mathrm{d}^{3} q \tag{1.6}
\end{equation*}
$$

The operator $\boldsymbol{Q}$ consists of three self-adjoint operators $Q_{i}$ whose domains are defined as
$\mathcal{D}\left(Q_{i}\right)=\left\{\left.\Psi \in \mathcal{H}\left|\int_{\mathbb{V} \subset \mathbb{R}^{3}}\left(Q_{i} \Psi\right)^{\dagger} Q_{i} \Psi \mathrm{~d}^{3} q=\int_{\mathbb{V} \subset \mathbb{R}^{3}} \sum_{j=1}^{n}\right| q_{i} \psi_{j}\right|^{2} \mathrm{~d}^{3} q<\infty\right\}$.
The momentum operator $P_{j}=-\mathrm{i} \hbar \hat{\eta} \otimes \partial / \partial q_{j}$ can be defined as the Fourier transformation of the position operator $Q_{j}(j=1,2,3)$.

Minimal interactions can now be introduced by means of the prescription $P_{\mu} \rightarrow$ $P_{\mu}-g A_{\mu}$, where $g$ is the coupling constant, $A_{\mu}$ is a gauge field $(\mu=0,1,2,3)$. Note that here $P_{0}=-\mathrm{i} \hbar I_{2 \times 2} \otimes \partial / \partial q_{0}$. This is the basis of the so-called gauge principle whereby the form of the interaction is determined on the basis of local gauge invariance. The covariant derivative $D_{\mu} \equiv(\mathrm{i} / \hbar)\left(P_{\mu}-g A_{\mu}\right)$ turns out to be of fundamental importance to determine the field strength tensor of the theory. It will be the operator which generalizes from electromagnetic-like interactions.

In section 2 we introduce the interaction of a quantum system as above with an electromagnetic-like field specified by $A_{\mu}$ by taking over the procedure of minimal substitution. We then briefly examine the problem of gauge invariance of the theory. In section 3 we consider the one-dimensional problem in the non-relativistic limit. Two cases are discussed: a simple oscillator-like interaction and the interaction of a two-level atom with a two-mode electromagnetic field in a configuration consisting of two counterpropagating travelling waves.

Section 4 deals with the non-relativistic three-dimensional problem for the cases of a spinless and a spin- $\frac{1}{2}$ particle in a central vector potential. Finally, section 5 considers the relativistic problem for a Dirac particle. We study again the case of a spin- $\frac{1}{2}$ particle in a central vector potential taking as an example a Coulomb-like interaction. We also briefly discuss the Zitterbewegung of the free electron.

## 2. Minimal interactions

In what follows we shall omit the symbol ' $\otimes$ ' when its presence is obvious as in equation (1.2). Let us consider a quantum system described by the free-particle Hamiltonian $H_{0}=H_{0}\left(P_{j}\right)$. We can incorporate a minimal interaction into $H_{0}$ by making the substitution

$$
\begin{equation*}
P_{\mu} \rightarrow \Pi_{\mu}=P_{\mu}-g A_{\mu} \tag{2.1}
\end{equation*}
$$

where $c P_{0}=I_{n \times n} \otimes \mathrm{i} \hbar \partial / \partial t$ as usual, $A_{\mu}(\mu=0,1,2,3)$ is a 4-vector potential and $g$ is the corresponding coupling constant. Here

$$
\begin{equation*}
A_{\mu}(q)=\tau_{a} A_{\mu}^{a}(q) \tag{2.2}
\end{equation*}
$$

with the $A_{\mu}^{a}(q)$ Hermitian functions of $q=\left(c t, q_{j}\right)$, and $\tau_{a}=\left(I_{n \times n}, \tau_{r}\right)$, where the $\tau_{r}$ are $n^{2}-1$ independent $n \times n$ traceless Hermitian matrices. Note that in equation (2.2) there
is a summation on the index $a=1, \ldots, n^{2}$. The system will now be described by the Schrödinger wave equation

$$
\begin{equation*}
H_{0}(\boldsymbol{\Pi}) \Psi=\left(\mathrm{i} \hbar \partial_{0}-g \tau_{a} A_{0}^{a}(q)\right) \Psi \tag{2.3}
\end{equation*}
$$

where $\Pi_{i}=P_{i}-g A_{i}(i=1,2,3)$ and the wavefunction $\Psi$ is itself a $n$-component spinor. The minimal replacement (2.1) must also satisfy the (formal) Hermiticity condition

$$
\begin{equation*}
H_{0}^{\dagger}(\boldsymbol{\Pi})=H_{0}(\boldsymbol{\Pi}) \tag{2.4}
\end{equation*}
$$

which in general restricts the form of the vector field (2.2) (see, for instance, section 5).
The field strength is defined by the commutators between the components of the covariant derivative $D_{\mu} \equiv(\mathrm{i} / \hbar) \Pi_{\mu}$. By using equations (2.1) and (2.2) we find
$\frac{\mathrm{i} g}{\hbar} F_{i j} \equiv\left[D_{i}, D_{j}\right]=\frac{\mathrm{i} g}{\hbar}\left\{\hat{\eta}\left(\partial_{j} A_{i}-\partial_{i} A_{j}\right)+\left[A_{j}, \hat{\eta}\right] \partial_{i}-\left[A_{i}, \hat{\eta}\right] \partial_{j}+\frac{\mathrm{i} g}{\hbar}\left[A_{i}, A_{j}\right]\right\}$
$\frac{\mathrm{i} g}{\hbar} F_{i 0} \equiv\left[D_{i}, D_{0}\right]=\frac{\mathrm{i} g}{\hbar}\left\{-\hat{\eta}\left(\partial_{i} A_{0}\right)+\left[A_{0}, \hat{\eta}\right] \partial_{i}-\partial_{0} A_{i}+\frac{\mathrm{i} g}{\hbar}\left[A_{0}, A_{i}\right]\right\}$.
If we want to keep the field strength components $F_{i j}$ and $F_{i 0}$ antisymmetric and symmetric under spatial inversion $\boldsymbol{q} \rightarrow-\boldsymbol{q}$, we must demand the components of the gauge potential $A_{0}$ and $A_{i}$ to be (up to a constant multiple of the identity matrix) symmetric and antisymmetric fields, respectively. Note that equations (2.1)-(2.5) are in direct correspondence to the usual Abelian case (an electromagnetic interaction) and not to a new non-Abelian generalization. The non-Abelian case follows straightforwardly mutatis mutandis as in the ordinary case and, to be brief, we shall not treat it here.

The solution of the wave equation (2.3) describes completely the state of the particle moving under the influence of the potential $A_{\mu}(q)$. This wave equation can be made gauge invariant under the combined local gauge transformation

$$
\begin{align*}
& A_{\mu}(q) \rightarrow A_{\mu}^{\prime}(q)=A_{\mu}(q)+\delta A_{\mu}(q) \\
& \Psi(q) \rightarrow \Psi^{\prime}(q)=\Psi(q)+\delta \Psi(q) \tag{2.6}
\end{align*}
$$

In the above we consider an infinitesimal local phase transformation for $\Psi(q)$ of the form

$$
\begin{equation*}
\Psi^{\prime}(q)=U(q) \Psi(q) \tag{2.7}
\end{equation*}
$$

where

$$
\begin{equation*}
U(q)=\mathbb{I}+\mathrm{i} g \tau_{a} \zeta^{a}(q)+\mathrm{O}\left(\zeta^{2}\right) \tag{2.8}
\end{equation*}
$$

The requirement is that

$$
D_{\mu}^{\prime} \Psi^{\prime}(q)=U(q) D_{\mu} \Psi(q)
$$

with

$$
\begin{equation*}
D_{\mu}^{\prime}=\hat{\eta} \partial_{\mu}-\frac{\mathrm{i} g}{\hbar}\left(A_{\mu}(q)+\delta A_{\mu}(q)\right) \tag{2.9}
\end{equation*}
$$

Equation (2.9) holds true if

$$
\begin{align*}
& \delta \Psi(q)=\mathrm{i} g \tau_{a} \zeta^{a}(q) \Psi(q)  \tag{2.10}\\
& \delta A_{0}(q)=\tau_{a} \partial_{0} \zeta^{a}(q)+\mathrm{i} g\left[\tau_{a} \zeta^{a}(q), A_{0}(q)\right] \\
& \delta A_{k}(q)=\hat{\eta} \tau_{a} \partial_{k} \zeta^{a}(q)+\zeta^{a}(q)\left[\tau_{a}, \hat{\eta}\right] \partial_{k}+\mathrm{i} g\left[\tau_{a} \zeta^{a}(q), A_{k}(q)\right]
\end{align*}
$$

which corresponds to a direct generalization of a usual local Abelian gauge transformation.

## 3. The one-dimensional case

A one-dimensional system is characterized by canonical coordinates $Q$ and $P$ satisfying

$$
\begin{equation*}
[Q, P]=\mathrm{i} \hbar \mathbb{I} . \tag{3.1}
\end{equation*}
$$

For $n=2$ we may represent these operators in the general form

$$
\begin{equation*}
Q=\hat{\eta} q \quad P=\hat{\eta} p \quad \mathbb{I}=I_{2 \times 2} \otimes I \tag{3.2}
\end{equation*}
$$

where $p=-\mathrm{i} \hbar \partial / \partial q, \operatorname{Tr} \hat{\eta}=0$ and $\hat{\eta}^{2}=I_{2 \times 2}$. The free-particle Schrödinger equation can be written as

$$
\begin{equation*}
H_{0} \Psi=\frac{1}{2 m} P^{2} \Psi=\frac{1}{2 m} p^{2} \Psi=\mathrm{i} \hbar \partial_{t} \Psi . \tag{3.3}
\end{equation*}
$$

As one would expect, the Hamiltonian $H$ still describes a Schrödinger 'free particle', but $\Psi$ is a two-component wavefunction now.

In the presence of an electromagnetic coupling, equation (2.3) reads ( $g \equiv e / c, e<0$ )

$$
\begin{equation*}
H \Psi=\left(\frac{1}{2 m}\left(P-\frac{e}{c} A_{1}\right)^{2}+e A_{0}\right) \Psi=\mathrm{i} \hbar \partial_{t} \Psi \tag{3.4}
\end{equation*}
$$

where

$$
\begin{equation*}
A_{\mu}=\tau_{a} A_{\mu}^{a}(q) \tag{3.5}
\end{equation*}
$$

with $\mu=0,1$. Here we require the representation to have as low a rank as possible. We choose

$$
\begin{equation*}
\hat{\eta} \equiv \tau_{1} \quad \tau_{a}=\left(I_{2 \times 2}, \tau_{i}\right) \tag{3.6}
\end{equation*}
$$

where the $\tau_{i}$ are the Pauli matrices. Thus the Schrödinger equation becomes

$$
\begin{gather*}
H \Psi=\left(\frac{1}{2 m} p^{2}-\frac{e}{2 m c}\left(-\mathrm{i} \hbar \hat{\eta} \frac{\mathrm{~d} A_{1}(q)}{\mathrm{d} q}+\left\{\hat{\eta}, A_{1}(q)\right\} p\right)+\frac{e^{2}}{2 m c^{2}} A_{1}^{2}(q)+e A_{0}\right) \Psi \\
=\mathrm{i} \hbar \partial_{t} \Psi \tag{3.7}
\end{gather*}
$$

### 3.1. The harmonic oscillator-like case

Let us choose a particular representation satisfying

$$
\begin{equation*}
A_{0}=0 \quad\left\{\tau_{3}, \Pi\right\}=0 \tag{3.8}
\end{equation*}
$$

where $\{$,$\} denotes the anticommutator and \Pi=P-(e / c) A_{1}$. In this case the most general gauge potential will have the form

$$
\begin{equation*}
A_{1}(q)=U_{1}(q) \tau_{1}+U_{2}(q) \tau_{2} \tag{3.9}
\end{equation*}
$$

where $U_{1}, U_{2}$ are general differentiable functions of $q$. Thus equation (3.7) reduces to

$$
\begin{gather*}
H \Psi=\left(\frac{1}{2 m} p^{2}+\mathrm{i} \hbar \frac{e}{2 m c}\left(\frac{\mathrm{~d} U_{1}(q)}{\mathrm{d} q}+\mathrm{i} \tau_{3} \frac{\mathrm{~d} U_{2}(q)}{\mathrm{d} q}\right)-\frac{e}{m c} U_{1}(q) p\right. \\
\left.+\frac{e^{2}}{2 m c^{2}}\left(U_{1}^{2}(q)+U_{2}^{2}(q)\right)\right) \Psi=\mathrm{i} \hbar \partial_{t} \Psi \tag{3.10}
\end{gather*}
$$

Note that in this problem the set $\left\{\mathcal{H}, H, \tau_{3}, \Pi\right\}$ defines a supersymmetric system [7]. Here $i(\Pi)=\operatorname{Tr}\left[\left(\tau_{3} \otimes I\right) \exp (-\beta H)\right]$ (well defined independently of $\beta>0$ if $\exp (-\beta H)$ is of trace class) is an index which measures supersymmetry breaking [7].

To be specific, we set $A_{0}=0, U_{2}(q)=\lambda q$ ( $\lambda$ a constant), and $U_{1}(q)$ a general function of $q$. Equation (3.10) becomes

$$
\begin{equation*}
H \Psi=\left(\frac{1}{2 m}\left(p-\frac{e}{c} U_{1}(q)\right)^{2}-\frac{e \hbar \lambda}{2 m c} \tau_{3}+\frac{e^{2} \lambda^{2}}{2 m c^{2}} q^{2}\right) \Psi=\mathrm{i} \hbar \partial_{t} \Psi \tag{3.11}
\end{equation*}
$$

which corresponds to a harmonic oscillator system. By defining

$$
\begin{equation*}
a \equiv \frac{1}{\sqrt{2 m \hbar \lambda}}\left(\lambda q+\mathrm{i}\left(p-\frac{e}{c} U_{1}(q)\right)\right) \tag{3.12}
\end{equation*}
$$

with

$$
\begin{equation*}
\left[a, a^{\dagger}\right]=I \tag{3.13}
\end{equation*}
$$

we get

$$
H=\left(\begin{array}{cc}
H_{-} & 0  \tag{3.14}\\
0 & H_{+}
\end{array}\right)
$$

where the supersymmetric partner Hamiltonians are given by

$$
\begin{equation*}
H_{\mp}=\frac{e \hbar \lambda}{m c}\left(a^{\dagger} a+\frac{1}{2}\right) \mp \frac{e \hbar \lambda}{2 m c} \tag{3.15}
\end{equation*}
$$

The energy spectrum is then

$$
\begin{equation*}
E_{n}^{+}=\frac{e \hbar \lambda}{m c}(n+1) \quad E_{n}^{-}=\frac{e \hbar \lambda}{m c} n \tag{3.16}
\end{equation*}
$$

for $n=0,1, \ldots$ Thus the ground state $E_{n=0}^{-}$has zero energy and is non-degenerate. All excited states are doubly degenerate. This shows that supersymmetry in this system is unbroken, see also [7].

### 3.2. Atom-field interaction in a cavity

As a second instance, let us consider the interaction of a two-level atom in a configuration consisting of two counterpropagating travelling waves, for instance in a ring configuration $[9,10]$. If the atom is assumed to propagate with momentum $p_{z}$ in the $z$ direction perpendicular to the light field, the atom-field Hamiltonian before the atom gets into the cavity is

$$
\begin{align*}
H_{0} & =H_{K}^{(\text {Atom })}+V^{\text {(Atom) }}+H_{K}^{(\text {Field })} \\
& =\frac{1}{2 m} P^{2}+\frac{\hbar \omega_{0}}{2} \tau_{3}+\hbar \omega\left(a_{1}^{\dagger} a_{1}+a_{2}^{\dagger} a_{2}+\mathbb{I}\right) \tag{3.17}
\end{align*}
$$

where we have chosen $P=-\mathrm{i} \hbar \hat{\eta} \partial / \partial q$, with $q$ the direction of the light field, $\hat{\eta}=\tau_{3}$ and we have absorbed the kinetic energy of the atom associated with the $z$ direction in the definition of $H_{0}$. In the above $\omega_{0}, \omega$ are the atomic transition and field frequencies, respectively. In what follows we assume that the two frequencies are close to resonance, $\omega \sim \omega_{0}$. Once the atom is inside the cavity, the interaction can be incorporated through the minimal replacement

$$
\begin{equation*}
P_{\mu} \rightarrow \Pi_{\mu}=P_{\mu}-\frac{e}{c} A_{\mu} \tag{3.18}
\end{equation*}
$$

with (effective) gauge potential
$A_{0}=0$
$A_{1}=\left(\frac{\hbar c^{2}}{V \omega}\right)^{1 / 2}\left\{\left(a_{2} \tau_{+}-a_{1}^{\dagger} \tau_{-}\right) \exp (\mathrm{i} k q)+\left(a_{2}^{\dagger} \tau_{-}-a_{1} \tau_{+}\right) \exp (-\mathrm{i} k q)\right\}$
where $\tau_{-}, \tau_{+}$are the pseudo-spin lowering and raising operators, respectively. From equations (3.17)-(3.19) we get

$$
\begin{align*}
H=\frac{1}{2 m} \Pi_{1}^{2}+ & \frac{1}{2} \hbar \omega_{0} \tau_{3}+\hbar \omega\left(a_{1}^{\dagger} a_{1}+a_{2}^{\dagger} a_{2}+\mathbb{I}\right) \\
= & \frac{1}{2 m} P^{2}+\hbar \omega^{\prime}\left(a_{1}^{\dagger} a_{1}+a_{2}^{\dagger} a_{2}+\mathbb{I}\right)+\frac{1}{2} \hbar \omega_{0}^{\prime} \tau_{3} \\
& +\hbar \Omega\left\{\left(a_{1}^{\dagger} \tau_{-}+a_{2} \tau_{+}\right) \exp (\mathrm{i} k q)+\left(a_{1} \tau_{+}+a_{2}^{\dagger} \tau_{-}\right) \exp (-\mathrm{i} k q)\right\} \\
& -\hbar \delta \omega\left\{a_{1}^{\dagger} a_{2} \exp (-2 \mathrm{i} k q)+a_{1} a_{2}^{\dagger} \exp (2 \mathrm{i} k q)\right\} \tag{3.20}
\end{align*}
$$

where

$$
\begin{equation*}
\omega^{\prime}=\omega+\delta \omega \quad \omega_{0}^{\prime}=\omega_{0}+2 \delta \omega \quad \delta \omega=|e \mu| \frac{c}{\hbar V \omega} \tag{3.21}
\end{equation*}
$$

and

$$
\begin{equation*}
\Omega=|\mu|\left(\frac{\omega}{\hbar V}\right)^{1 / 2} \tag{3.22}
\end{equation*}
$$

with $\mu=\hbar e / 2 m c$. Note that in equation (3.22) we have obtained the correct value for the coupling constant $\Omega$ being a factor $\sqrt{2}$ larger than the one given for standing waves [11]. The expression for the Hamiltonian (3.20) differs from the one given by Shore, Meystre and Stenholm (SMS) [10]: the gauge potential (3.19) introduces a shift $\delta \omega$ in the field frequency, and the detuning becomes $\omega^{\prime}-\omega_{0}^{\prime}=\omega-\omega_{0}-\delta \omega$. Furthermore, apart from the usual one-photon exchange interaction with coupling constant $\hbar \Omega$, there is a new (momentum conserving) exchange mode contribution in $H$ with coupling constant $\hbar \delta \omega$. Note that for high enough field frequencies (visible optical frequencies) $\delta \omega$ becomes small. For instance, by choosing $\omega \sim 10^{11} \mathrm{~Hz}$ and $V=1 \mathrm{~cm}^{3}$, one gets $\delta \omega / \Omega \sim 0.01$. Thus, in the limit of high field frequencies, the SMS result is re-obtained.

Using the fact that

$$
\begin{equation*}
\hat{M} \equiv a_{1}^{\dagger} a_{1}+a_{2}^{\dagger} a_{2}+\mathbb{I}+\frac{1}{2} \tau_{3} \quad \hat{N} \equiv-\frac{\mathrm{i}}{k} \frac{\mathrm{~d}}{\mathrm{~d} q}+a_{1}^{\dagger} a_{1}-a_{2}^{\dagger} a_{2} \tag{3.23}
\end{equation*}
$$

commute with each other and with the total Hamiltonian $H$, the basis states can be labelled as

$$
\begin{equation*}
\left.\mid M, N ; m^{(\tau)}, \tau\right) \equiv \exp \left(\frac{\mathrm{i}}{\hbar} p q\right)\left|m_{1}^{(\tau)}\right\rangle\left|m_{2}^{(\tau)}\right\rangle|\tau\rangle \tag{3.24}
\end{equation*}
$$

with

$$
\begin{array}{ll}
M=m_{1}^{(\tau)}+m_{2}^{(\tau)}+1+\frac{1}{2} \tau & \tau= \pm 1 \\
N=p+\hbar k\left(m_{1}^{(\tau)}-m_{2}^{(\tau)}\right) & \tag{3.25}
\end{array}
$$

where we have chosen $m^{(\tau)} \equiv m_{1}^{(\tau)}=0,1, \ldots, M+1-\tau / 2$. Note that $H$ involves generators of the superalgebra $u(2 / 1)$ generated by the bilinear products
$S_{i}^{(+)}=\tau_{-} a_{i}^{\dagger} \quad G_{i}^{j}=a_{i} a_{j}^{\dagger} \quad S_{i}^{(-)}=\tau_{+} a_{i} \quad \tau_{ \pm} \tau_{\mp}=\frac{1}{2}\left(I_{2 \times 2} \pm \tau_{3}\right)$.
Here the exponential factors $\exp ( \pm \mathrm{ikq})$ are implicit in the field operators. In fact $\hat{M}$ displays an $s u(2)$ symmetry (subalgebra) whose generators are

$$
\begin{equation*}
J_{+}=a_{1}^{\dagger} a_{2}+\tau_{+} \quad J_{-}=a_{2}^{\dagger} a_{1}+\tau_{-} \quad J_{3}=a_{1}^{\dagger} a_{1}-a_{2}^{\dagger} a_{2}+\frac{1}{2} \tau_{3} \tag{3.27}
\end{equation*}
$$

with

$$
\begin{equation*}
\left[J_{i}, J_{j}\right]=\mathrm{i} \epsilon_{i j k} J_{k} \tag{3.28}
\end{equation*}
$$

We find that

$$
\begin{equation*}
\left[J_{i}, \hat{M}\right]=\left[J_{i}, \hat{N}\right]=0 \tag{3.29}
\end{equation*}
$$

being (3.24) a basis of representation for the elements of this algebra.
To determine the energy spectrum we note that $\Pi_{1}=P-(e / c) A_{1}$ shares the same symmetry as $H$ :

$$
\begin{equation*}
\left[\Pi_{1}, \hat{M}\right]=\left[\Pi_{1}, \hat{N}\right]=0 \tag{3.30}
\end{equation*}
$$

By using this fact we calculate the matrix elements of $\Pi_{1}$ in the basis (3.24) for given values of $M, N$. After squaring the corresponding $2 M \times 2 M$ diagonal matrix, we replace it in equation (3.20) yielding the exact spectrum of $H$. With the definition (3.24) we find

$$
\begin{align*}
\left(M, N ; m_{a}^{\left(\tau_{a}\right)},\right. & \left.\tau_{a}\left|\Pi_{1}\right| M, N ; m_{b}^{\left(\tau_{b}\right)}, \tau_{b}\right) \\
= & \hbar \tau\left(M+N-2 m_{b}^{\left(\tau_{b}\right)}-\frac{1}{2}\left(\tau_{b}+2\right)\right) k \delta_{m_{a}^{\left(\tau_{a}\right)}, m_{b}^{\left(\tau_{b}\right)}} \delta_{\tau_{a}, \tau_{b}} \\
& +\frac{e f}{c}\left(\sqrt{m_{b}^{\left(\tau_{b}\right)}} \delta_{m_{a}^{\left(\tau_{a}\right)}, m_{b}^{\left(\tau_{b}\right)}-1} \delta_{\tau_{a}, \tau_{b}+2}\right. \\
& -\sqrt{M-m_{b}^{\left(\tau_{b}\right)}-\frac{1}{2}\left(\tau_{b}+2\right)} \delta_{m_{a}^{\left(\tau_{a}\right)}, m_{b}^{\left(\tau_{b}\right)}} \delta_{\tau_{a}, \tau_{b}+2} \\
& +\sqrt{m_{b}^{\left(\tau_{b}\right)}+1} \delta_{m_{a}^{\left(\tau_{a}\right)}, m_{b}^{\left(\tau_{b}\right)}+1} \delta_{\tau_{a}, \tau_{b}-2} \\
& \left.-\sqrt{M-m_{b}^{\left(\tau_{b}\right)}-\frac{1}{2} \tau_{b}} \delta_{m_{a}^{\left(\tau_{a}\right)}, m_{b}^{\left(\tau_{b}\right)}} \delta_{\tau_{a}, \tau_{b}-2}\right) \tag{3.31}
\end{align*}
$$

In figure 1 , we show the energy spectra of the interacting Hamiltonian $H$ for a rubidium atom $\left(m=1.4192 \times 10^{-22} \mathrm{~g}\right)$ in units of $\hbar \omega, \omega=21506.5 \mathrm{MHz}\left(63 \mathrm{p}_{3 / 2}-61 \mathrm{~d}_{3 / 2}\right.$ transition), for the particular case $N=0$. On the left we display the levels of $\hat{M}$ which are $2 M$ degenerate. The energy levels of $H_{0}$ depicted on the right by full lines are, excluding the lowest one within each subspace, doubly degenerate. After the interaction is switched on, only the excited states within each subspace are shifted as shown in the same figure.

## 4. Three dimensions

### 4.1. A spinless non-relativistic particle in a Coulomb field

We first consider the Hamiltonian $H$ for a spinless particle moving in an electromagnetic field $A_{\mu}$

$$
\begin{equation*}
H=\frac{1}{2 m}\left(\boldsymbol{P}-\frac{e}{c} \boldsymbol{A}\right)^{2}+e A_{0} \tag{4.1}
\end{equation*}
$$

where $P_{i}=-\mathrm{i} \hbar \hat{\eta} \partial / \partial q_{i}$ and $A_{i}=\tau_{a} A_{i}^{a}$, with $\tau_{a}$ Pauli matrices and $A_{i}^{0}, A_{i}^{1}, A_{i}^{3}, A_{0}$ vanishing fields. Here $\hat{\eta} \equiv \tau_{1}$, while $A_{i}^{2}$ is a central vector field of the form

$$
\begin{equation*}
A_{i}^{2}(q)=U(q) \frac{q^{i}}{q} \tag{4.2}
\end{equation*}
$$

with $U(q)$ a real function of $q=|\boldsymbol{q}|$. We obtain

$$
\begin{equation*}
H=-\frac{\hbar^{2}}{2 m} \nabla^{2}-\tau_{3} \frac{\hbar e}{2 m c q^{2}} \frac{\mathrm{~d}}{\mathrm{~d} q}\left(q^{2} U(q)\right)+\frac{e^{2}}{2 m c^{2}} U^{2}(q) \tag{4.3}
\end{equation*}
$$

It is clear that this Hamiltonian is not supersymmetric since the term proportional to $\mathrm{d} / \mathrm{d} q()$ involves a function which is different from $U$.


Figure 1. Schematic representation of the $2 M$ degenerate energy levels for $N=0$ in units of $\hbar \omega$ and their splitting after the atom-field interaction is turned on. Notice that, except for the lowest energy levels, there is a residual double degeneracy in the excited levels within each subspace.

By choosing $U=m Z e / \hbar$, the Schrödinger equation for stationary states leads to the following two equations:

$$
\begin{equation*}
\left\{-\frac{\hbar^{2}}{2 m} \frac{\mathrm{~d}^{2}}{\mathrm{~d} q^{2}}+\frac{\hbar^{2}}{2 m} l(l+1) \frac{1}{q^{2}} \mp \frac{Z e^{2}}{q}+\frac{Z^{2} e^{4} m}{2 \hbar^{2}}\right\} \psi_{\mp}=E_{\mp} \psi_{\mp} . \tag{4.4}
\end{equation*}
$$

Thus the Hilbert space $\left\{\Psi=\left(\Psi_{+}, \Psi_{-}\right)\right\}$with

$$
\begin{equation*}
\Psi_{+}=\binom{\psi_{+}}{0} \quad \Psi_{-}=\binom{0}{\psi_{-}} \tag{4.5}
\end{equation*}
$$

consists of the complete set of bound and scattering states of the problem. Note that there is a double degeneracy for the continuum solutions: the eigenvalues $E\left(E>Z^{2} e^{4} m / 2 \hbar^{2}\right)$ of (4.4) are those satisfying $E=E_{+}=E_{-}$, with $\Psi_{ \pm}$the corresponding well known scattering solutions for the scalar potentials $\pm Z e^{2} / q$. The discrete solutions are described by $\Psi_{+}$ with energy eigenvalues

$$
\begin{equation*}
E_{n}=\frac{Z^{2} e^{4} m}{2 \hbar^{2}}\left(1-\frac{1}{n^{2}}\right) \tag{4.6}
\end{equation*}
$$

with $n=1,2, \ldots$. Because $H$ is proportional to the square of $\boldsymbol{P}-(e / c) \boldsymbol{A}$ we see that $E_{n} \geqslant 0$.

### 4.2. A spin- $\frac{1}{2}$ non-relativistic particle

The squared momentum of a spin- $\frac{1}{2}$ free particle may be written in the form

$$
\begin{equation*}
\boldsymbol{P}^{2}=\left(\sum_{k=1}^{3} P_{k} \otimes \sigma_{k}\right)^{2} \tag{4.7}
\end{equation*}
$$

Thus the free-particle wave equation reads

$$
\begin{equation*}
H \Psi=\frac{1}{2 m} \boldsymbol{P}^{2} \Psi=\frac{1}{2 m}\left(\sum_{k} P_{k} \otimes \sigma_{k}\right)^{2} \Psi=\mathrm{i} \hbar \partial_{t} \Psi \tag{4.8}
\end{equation*}
$$

with $P_{k}=\hat{\eta} p_{k}$. Under minimal interaction the Hamiltonian becomes

$$
\begin{align*}
H & =\frac{1}{2 m}\left(\left(P_{k}-\frac{e}{c} A_{k}\right) \otimes \sigma_{k}\right)^{2}+e A_{0} \\
& =\frac{1}{2 m}\left(P_{k}-\frac{e}{c} A_{k}\right)^{2}-\mathrm{i} \frac{e}{m c} \varepsilon_{i j k}\left[P_{i}, A_{j}\right] \otimes \frac{1}{2} \sigma_{k}+e A_{0} \tag{4.9}
\end{align*}
$$

where there is summation over the indices $i, j, k$. In equation (4.9) we have chosen the same two-dimensional matrix structure for $P_{\mu}, A_{\mu}$ as in the previous case. From equation (2.5) the magnetic field is defined by

$$
\begin{equation*}
B_{k} \equiv \frac{1}{2} \varepsilon_{i j k} F_{i j}=\frac{\mathrm{i} c}{2 \hbar e} \varepsilon_{i j k}\left[\Pi_{i}, \Pi_{j}\right]=-\frac{\mathrm{i}}{\hbar} \varepsilon_{i j k}\left[P_{i}, A_{j}\right] \tag{4.10}
\end{equation*}
$$

Thus the Hamiltonian can be written as

$$
\begin{align*}
H & =\frac{1}{2 m}\left(\left(P_{k}-\frac{e}{c} A_{k}\right) \otimes \sigma_{k}\right)^{2}+e A_{0} \\
& =\frac{1}{2 m}\left(P_{k}-\frac{e}{c} A_{k}\right)^{2}+\frac{\boldsymbol{g e}}{2 m c} B_{k} \otimes S_{k}+e A_{0} \tag{4.11}
\end{align*}
$$

where $S_{k} \equiv \frac{1}{2} \sigma_{k}$, and $\boldsymbol{g}=2$ is the gyromagnetic constant. Hence this Hamiltonian represents the interaction of a magnetic moment $(\hbar e / 2 m c) \sigma$ with the magnetic field $\boldsymbol{B}$ felt in the rest frame of the particle.

As an example let us consider again a central potential of the form $A_{\mu}=\left(0, \tau_{2} U(q) \hat{\boldsymbol{q}}\right)$. The Schrödinger equation for the system becomes

$$
\begin{align*}
H \Psi=(- & \left.\frac{\hbar^{2}}{2 m} \nabla^{2}+\frac{e^{2}}{2 m c^{2}} U^{2}(q)-\frac{\hbar e}{2 m c} \tau_{3} \otimes\left(\frac{\mathrm{~d} U(q)}{\mathrm{d} q}+\frac{2 U(q)}{q}\left(\mathbb{I}+\frac{1}{\hbar} \boldsymbol{L} \cdot \sigma\right)\right)\right) \Psi \\
& =\mathrm{i} \hbar \partial_{t} \Psi \tag{4.12}
\end{align*}
$$

where now the four-component spinor wavefunction $\Psi$ has the form

$$
\begin{equation*}
\Psi=\binom{\Phi_{+}}{\Phi_{-}} \tag{4.13}
\end{equation*}
$$

Note that the Hamiltonian $H$ in equation (4.12) is supersymmetric [12]. Hence, under a particular (vector-like) minimal interaction, $H$ turns out to be supersymmetric only if $\boldsymbol{g}=2$. The supersymmetric partners $V_{ \pm}$are in this case

$$
\begin{equation*}
V_{ \pm}(q)=\frac{e^{2}}{2 m} U^{2}(q) \pm \frac{\hbar e}{2 m}\left(\frac{\mathrm{~d} U(q)}{\mathrm{d} q}+\frac{2 U(q)}{q}\left(\mathbb{I}+\frac{1}{\hbar} \boldsymbol{L} \cdot \sigma\right)\right) \tag{4.14}
\end{equation*}
$$

According to equation (2.5) the magnetic field is defined by

$$
\begin{align*}
B_{k} & \equiv F_{i j}=\left[A_{j}, \hat{\eta}\right] \partial_{i}-\left[A_{i}, \hat{\eta}\right] \partial_{j} \\
& =-\frac{2 U(q)}{q} \tau_{3} L_{k} \tag{4.15}
\end{align*}
$$

( $i, j, k$ cyclic), i.e. it is proportional to the angular momentum of the particle.
The standard spin-orbit coupling in quantum mechanics due to an electric potential $V(q)$ takes the form

$$
\begin{equation*}
H_{Q M}^{(L S)}=\frac{\hbar e}{4 m^{2} c^{2} q} \frac{\mathrm{~d} V(q)}{\mathrm{d} q} \boldsymbol{L} \cdot \sigma \tag{4.16}
\end{equation*}
$$

In contrast with this, the spin-orbit coupling required by (4.12) is

$$
\begin{equation*}
H^{(L S)}=\frac{e}{m} \frac{U(q)}{q} \boldsymbol{L} \cdot \sigma \tag{4.17}
\end{equation*}
$$

If we now look for a direct identification

$$
\begin{equation*}
U(q) \rightarrow \frac{\hbar}{4 m c^{2}} \frac{\mathrm{~d} V(q)}{\mathrm{d} q} \quad U^{2}(q) \rightarrow \frac{2 m c^{2}}{e} V(q) \tag{4.18}
\end{equation*}
$$

we get $U(q)=\left(4 m^{2} c^{3} / \hbar e\right) q$. This system corresponds to a harmonic oscillator with frequency $\omega=4 m c^{2} / \hbar$. This result can be interpreted by saying that the harmonic oscillator would be supersymmetric if the ground-state energy is sufficient to create a pair [12, 13].

Notice that the spin-orbit coupling introduces a correction to the result given in (4.4) for the Coulomb field,

$$
\begin{equation*}
-\frac{Z e^{2}}{q} \rightarrow-\frac{\hbar e}{2 m}\left(\frac{2 m Z e / \hbar}{q}\left(\mathbb{I}+\frac{1}{\hbar} \boldsymbol{L} \cdot \sigma\right)\right)=-\frac{Z e^{2}}{q}\left(\mathbb{I}+\frac{1}{\hbar} \boldsymbol{L} \cdot \sigma\right) \tag{4.19}
\end{equation*}
$$

associated to the $\Phi_{+}$wavefunction component. Thus the resulting effective potential will strongly depend on the values of the orbital and total angular momenta $j, l$ of the state system (see section 5 for the solution of the corresponding relativistic problem).

## 5. The relativistic case for a spin- $\frac{1}{2}$ particle

Our considerations can also be extended to relativistic quantum mechanics. In the case of a Dirac particle we already count with a $4 \times 4$ (formally) traceless Hamiltonian. Here it seems natural to also look for appropriate $4 \times 4$ (formally) traceless momentum and coordinate operators. To this end we notice that the Dirac wave equation can be written as

$$
\begin{equation*}
H \Psi=\left(c \boldsymbol{\Sigma} \cdot \boldsymbol{P}+m c^{2} \beta\right) \Psi=\mathrm{i} \hbar \partial_{t} \Psi \tag{5.1}
\end{equation*}
$$

where

$$
P_{k} \equiv-\gamma_{5} p_{k}=\mathrm{i} \hbar \gamma_{5} \nabla_{k} \quad \Sigma_{k}=-\gamma_{5} \alpha_{k}=\left(\begin{array}{cc}
\sigma_{k} & 0  \tag{5.2}\\
0 & \sigma_{k}
\end{array}\right)
$$

with

$$
\gamma_{5}=-\left(\begin{array}{cc}
0 & I_{2 \times 2}  \tag{5.3}\\
I_{2 \times 2} & 0
\end{array}\right)
$$

Thus we define

$$
\begin{array}{ll}
Q_{0} \equiv I_{4 \times 4} \otimes q_{0} & P_{0} \equiv I_{4 \times 4} \otimes p_{0} \\
Q_{i} \equiv-\gamma_{5} \otimes q_{i} & P_{j} \equiv-\gamma_{5} \otimes p_{j} \tag{5.4}
\end{array}
$$

where $p_{\mu}=\mathrm{i} \hbar \partial / \partial q^{\mu}$. The operators $Q_{\mu}, P_{\nu}$ satisfy the canonical commutation relations

$$
\begin{equation*}
\left[Q_{\mu}, Q_{\nu}\right]=\left[P_{\mu}, P_{\nu}\right]=0 \quad\left[Q_{\mu}, P_{\nu}\right]=\mathrm{i} \hbar g_{\mu \nu} \mathbb{I} \tag{5.5}
\end{equation*}
$$

where $\mathbb{I}=I_{4 \times 4} \otimes I$ and $g=\operatorname{diag}(-1,1,1,1)$.
Next we introduce the minimal replacement

$$
\begin{equation*}
P_{\mu} \rightarrow P_{\mu}-\frac{e}{c} A_{\mu} \tag{5.6}
\end{equation*}
$$

with

$$
\begin{equation*}
A_{\mu}(q)=\Gamma_{a} A_{\mu}^{a}(q) \quad A_{\mu}^{\dagger}=A_{\mu} \tag{5.7}
\end{equation*}
$$

where the matrices $\Gamma_{a}$ belong to the complete set of 16 Gamma matrices

$$
\begin{equation*}
I_{4 \times 4} \quad \mathrm{i} \gamma_{\mu} \quad \mathrm{i} \gamma_{\mu} \gamma_{v}(v \neq \mu) \quad \gamma_{5} \quad \mathrm{i} \gamma_{5} \gamma_{\mu} . \tag{5.8}
\end{equation*}
$$

By replacing equations (5.6) and (5.7) in equation (5.1) we get

$$
\begin{equation*}
H \Psi=\left(c \Sigma_{k}\left(P_{k}-\frac{e}{c} A_{k}(q)\right)+m c^{2} \beta\right) \Psi=\left(\mathrm{i} \hbar \partial_{t}-e A_{0}(q)\right) \Psi . \tag{5.9}
\end{equation*}
$$

Note that if we demand (formally) $H$ to be an Hermitian operator we find that

$$
\begin{equation*}
\left[\Gamma_{a}, \Sigma_{k}\right] A_{k}^{a}(q)=0 \tag{5.10}
\end{equation*}
$$

which is a constraint on the form of the gauge potential $A_{k}$.
Equation (5.9) may be re-written as

$$
\begin{equation*}
\left(\Upsilon^{\mu} D_{\mu}+\frac{m c}{\hbar}\right) \Psi=0 \tag{5.11}
\end{equation*}
$$

where

$$
\begin{equation*}
\Upsilon^{0}=\beta \quad \Upsilon^{k}=\beta \Sigma_{k} \tag{5.12}
\end{equation*}
$$

Note also that, apart from the basic condition (2.9), gauge invariance will require

$$
\begin{equation*}
\left.\left[\Upsilon^{\mu}, U(q)\right)\right]=0 \tag{5.13}
\end{equation*}
$$

which is a strong restriction on the space of allowed gauge transformations for a Dirac particle. From (5.13) we find that the most general gauge transformation $U$ has the form

$$
\begin{equation*}
U(q)=\operatorname{expi} g\left(\zeta^{0}(q)+\zeta^{1}(q) \beta\right) \tag{5.14}
\end{equation*}
$$

On the basis of the equations (5.6)-(5.10) a variety of problems can be considered. As an example, let us choose

$$
\begin{equation*}
A_{0}(q)=0 \quad A_{k}(q)=\mathrm{i} \beta \gamma_{5} \frac{m Z e}{\hbar} \frac{q_{k}}{q} \tag{5.15}
\end{equation*}
$$

By writing

$$
\begin{equation*}
\Psi=\binom{\Phi_{+}}{\Phi_{-}} \tag{5.16}
\end{equation*}
$$

we see that for steady states equation (5.9) consists of two coupled equations

$$
\begin{equation*}
c \boldsymbol{\sigma} \cdot\left(\boldsymbol{p} \mp \mathrm{i} \frac{m Z e^{2}}{\hbar} \frac{\boldsymbol{q}}{q}\right) \Phi_{ \pm}=\left(E \pm m c^{2}\right) \Phi_{\mp} \tag{5.17}
\end{equation*}
$$

After decoupling them we find
$\left(\frac{1}{2 m} \boldsymbol{p}^{2} \mp \frac{Z e^{2}}{q}\left(I+\frac{1}{\hbar} \boldsymbol{\sigma} \cdot \boldsymbol{L}\right)+\frac{Z^{2} e^{4} m}{2 \hbar^{2}}\right) \Phi_{ \pm}=\left(2 m c^{2}\right)^{-1}\left(E^{2}-m^{2} c^{4}\right) \Phi_{ \pm}$.
This may as well be written in the compact form

$$
\begin{equation*}
\left(\frac{1}{2 m} \boldsymbol{p}^{2}-\frac{Z e^{2}}{q} K\right) \Psi=\left(\epsilon-\epsilon_{0}\right) \Psi \tag{5.19}
\end{equation*}
$$

with

$$
\begin{equation*}
\epsilon=\frac{E^{2}}{2 m c^{2}} \quad \epsilon_{0}=\frac{1}{2} m c^{2}\left(1+Z^{2} \alpha^{2}\right) \tag{5.20}
\end{equation*}
$$

and

$$
\begin{equation*}
K=\beta\left(I+\frac{1}{\hbar} \boldsymbol{\Sigma} \cdot \boldsymbol{L}\right)=\frac{1}{\hbar} \beta \boldsymbol{\Sigma} \cdot \boldsymbol{J}-\beta / 2 \quad \boldsymbol{J} \equiv \boldsymbol{L}+\frac{\hbar}{2} \boldsymbol{\Sigma} \tag{5.21}
\end{equation*}
$$

Consider the non-relativistic limit. For positive energy solutions the norm of $\Phi_{-}$is small in this limit and the wave equation reduces to (5.18) for $\Phi_{+}$. Then $\epsilon-\epsilon_{0}$ is the eigenvalue of the steady states of (4.12) with the corresponding potential given in (4.19).

By using the fact that

$$
\begin{equation*}
\boldsymbol{L} \times \boldsymbol{q}+\boldsymbol{q} \times \boldsymbol{L}=2 \mathrm{i} \hbar \boldsymbol{q} \tag{5.22}
\end{equation*}
$$

we find that $K$ is a constant of motion

$$
\begin{equation*}
[H, K]=0 . \tag{5.23}
\end{equation*}
$$

In addition to this we have that

$$
\begin{equation*}
[J, K]=0 \tag{5.24}
\end{equation*}
$$

Thus we can construct simultaneous eigenfunctions of $H, J^{2}, K, J_{3}$ just as in the case of the relativistic hydrogen atom [14]. From equation (5.21) we get

$$
\begin{equation*}
K^{2}=J^{2}+\frac{1}{4} \tag{5.25}
\end{equation*}
$$

so the eigenvalues of $K$ are

$$
\begin{equation*}
\kappa= \pm\left(j+\frac{1}{2}\right) \tag{5.26}
\end{equation*}
$$

The bound states are characterized by $\kappa>0$. In this case the wavefunctions $\Phi_{ \pm}$are eigenfunctions of $\boldsymbol{L}^{2}$ with eigenvalues $l_{ \pm}\left(l_{ \pm}+1\right), l_{ \pm}=j \pm \frac{1}{2}$. The solution of (5.18) is then given by
$\Phi_{+}(\boldsymbol{q})=\left\langle\boldsymbol{q} \mid+, n^{\prime} j j_{3}\right\rangle=\sum_{m_{+}, s} c\left(j j_{3} ; l_{+} m_{+}, \frac{1}{2} s\right) q_{n^{\prime}, l_{+}}(q) Y_{l_{+} m_{+}}(\theta, \varphi) \chi_{s}$
$\Phi_{-}(\boldsymbol{q})=\left\langle\boldsymbol{q} \mid-, n^{\prime} j j_{3}\right\rangle=\sum_{m_{-}, s} c\left(j j_{3} ; l_{-} m_{-}, \frac{1}{2} s\right) q_{n^{\prime}-1, l_{-}}(q) Y_{l_{-} m_{-}}(\theta, \varphi) \chi_{s}$.
Here $c(\ldots)$ are Clebsch-Gordan coefficients, $q_{n^{\prime}, l_{ \pm}}$are radial wavefunctions of the nonrelativistic hydrogen atom, $Y_{l_{ \pm} m_{ \pm}}(\theta, \varphi)$ are spherical harmonics and $\chi_{s}$ correspond to the spinors

$$
\begin{equation*}
\chi_{1 / 2}=\binom{1}{0} \quad \chi_{-1 / 2}=\binom{0}{1} \tag{5.28}
\end{equation*}
$$

From (5.19) the energy eigenvalues are given by
$E_{n^{\prime}, \kappa}=m c^{2}\left(1+(Z \alpha)^{2}\left(1-\frac{|\kappa|^{2}}{\left(n^{\prime}+|\kappa|\right)^{2}}\right)\right)^{1 / 2} \quad n^{\prime}=0,1,2 \ldots$
The degeneracy of the energy levels is infinity: the levels $E_{n_{i}^{\prime}, \kappa_{i}}, i=1,2,3, \ldots$, satisfying

$$
\begin{equation*}
\frac{\left|\kappa_{1}\right|^{2}}{\left(n_{1}^{\prime}+\left|\kappa_{1}\right|\right)^{2}}=\frac{\left|\kappa_{2}\right|^{2}}{\left(n_{2}^{\prime}+\left|\kappa_{2}\right|\right)^{2}}=\cdots \tag{5.30}
\end{equation*}
$$

have the same energy.
It is worth mentioning here that the Dirac oscillator [15] can also be re-obtained through a minimal interaction of the form (5.6)-(5.7) in the Dirac equation. This is done by choosing

$$
\begin{equation*}
A_{k}(q)=\mathrm{i} \beta \gamma_{5} \frac{m \omega}{e} q_{k} \tag{5.31}
\end{equation*}
$$

where $\omega$ is the frequency for this oscillator. This gauge field gives rise to a harmonic oscillator with a strong spin-orbit coupling which introduces, as in the previous case, an infinite degeneracy. This oscillator has a hidden supersymmetry, responsible for the special properties of its spectrum [16]. In fact we can easily see that by setting $U(q)=(m c \omega / e) q$ in equation (4.12) we re-obtain, up to a constant term in the Hamiltonian, the Dirac oscillator in the non-relativistic limit. It is interesting to note that the vector field $A_{k}$ given in equation (5.31), is a Hermitian operator. This feature is absent in Moshinski's approach [15].

### 5.1. The electron Zitterbewegung

We briefly discuss the Zitterbewegung (trembling motion) of the Dirac electron in the representation (5.4). To this end let us consider a free Dirac particle described by equation (5.1). In the Heisenberg picture, the time derivative of an operator, say $T$, is given by

$$
\begin{equation*}
\frac{\mathrm{d} T}{\mathrm{~d} t}=\frac{\mathrm{i}}{\hbar}[H, T] . \tag{5.32}
\end{equation*}
$$

We can easily see that both the coordinate and the momentum operators are not constants of the motion despite the fact that the particle is free. From (5.1), (5.4) and (5.32), the equations of motion for $Q_{k}$ and $P_{k}$ are

$$
\begin{equation*}
\frac{\mathrm{d} Q_{k}}{\mathrm{~d} t}=\frac{2 \mathrm{i} m c^{2}}{\hbar} \beta Q_{k}+c \Sigma_{k} \quad \frac{\mathrm{~d} P_{k}}{\mathrm{~d} t}=\frac{2 \mathrm{i} m c^{2}}{\hbar} \beta P_{k} \tag{5.33}
\end{equation*}
$$

while from equations (5.32) and (5.33) we get

$$
\begin{equation*}
\frac{\mathrm{d}^{2} Q_{k}}{\mathrm{~d} t^{2}}=\frac{2 \mathrm{i}}{\hbar}\left(H \frac{\mathrm{~d} Q_{k}}{\mathrm{~d} t}-c^{2} P_{k}\right) \tag{5.34}
\end{equation*}
$$

Let us define [17], for an observable $G$ and for energy $E$

$$
\begin{equation*}
G_{A} \equiv \frac{1}{2}\left(G+H E^{-1} G H E^{-1}\right) \tag{5.35}
\end{equation*}
$$

as the observable relative to which the Zitterbewegung takes place. By using (5.4) and (5.35) we get

$$
\begin{align*}
& \boldsymbol{\xi}_{Q} \equiv \boldsymbol{Q}-\boldsymbol{Q}_{A}=m c^{2} H^{-1} \beta \boldsymbol{Q}-\frac{\mathrm{i} \hbar c}{2} H^{-1}\left(\boldsymbol{\Sigma}-c \boldsymbol{P} H^{-1}\right) \\
& \boldsymbol{\xi}_{\boldsymbol{P}} \equiv \boldsymbol{P}-\boldsymbol{P}_{A}=m c^{2} H^{-1} \beta \boldsymbol{P} \tag{5.36}
\end{align*}
$$

with

$$
\begin{equation*}
T_{r}\left(\xi_{Q}\right)=T_{r}\left(\xi_{P}\right)=\mathbf{0} \tag{5.37}
\end{equation*}
$$

This result strongly differs from the usual expressions [17]. In particular the (odd) position and momentum operators $\boldsymbol{Q}, \boldsymbol{P}$ are themselves present in the corresponding Zitterbewegung coordinates.

## 6. Final comments

In this paper, we have studied minimal interactions in a wide class of quantum systems characterized by position and momentum operators defined as the direct product of a (Hermitian and unitary) finite traceless matrix and an ordinary canonical coordinate. This approach allows us to obtain in a simple fashion supersymmetric systems in quantum mechanics. However, we are not restricted only to this class of systems as was shown in the examples given in sections 3.2 and 4.1. In fact, equations (2.1) and (2.2) also include ordinary minimal interactions. In section 2 we worked out in detail the bound states of a two-level atom interacting with a two-mode electromagnetic field in a particular familiar configuration. This example suggests a useful gauge approach to some problems in cavity QED. A promising natural development of the present work is its extension to relativistic quantum field theory. We hope to report on such an extension elsewhere.

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