THE ROLE OF SPACE AND TIME IN NON-RELATIVISTIC QUANTUM PHYSICS

Rafael-Andrés Alemañ-Berenguer

Dpt. of Material Sciences, Optics and Electronic Tecnology. University *Miguel Hernández*, Avda. Universidad, s/n. Edif. Torrevaillo - 032021 - Elche (Spain) Astronomical Society of Alicante. Group of Gravitation, Cosmology and Celestial Mechanics, Mail-Box Nº 616, 03080-Alicante (España)

agrupacion.astroalicante@gmail.com

Abstract: Since the very begining of quantum theory there started a debate on the proper role of space and time in it. Some authors assumed that space and time have their own algebraic operators. On that basis they supposed that quantum particles had "coordinates of position", even though those coordinates were not possible to determine with infinite precision. Furthermore, time in quantum physics was taken to be on an equal foot, by means of a so-called "Heisenberg's fourth relation of indeterminacy" concerning time and energy. In this paper, the proper role of space and time in the core of non-relativistic quantum physics is analyzed. We will find that, rigorously, that relation for time and energy shows a different root. For the role of space, it will be discussed that there is no "coordinate of position" in the conceptual structure of quantum physics because quantum particles are not point-like objects.

1. Introduction

It could be thought that some of the paradoxes typical in non-relativistic quantum physics perhaps have their origin in the specific role of time as a physical quantity in the quantum theory. In fact this is a very debated matter. It was even pointed out by Von Neumann:

"(...) an essential weakness that is, in fact, the main weakness of the quantum mechanics: its non-relativistic character, which distinguishes time t of the three space coordinates x, y,

z, and presupposes a concept of objective simultaneity. In fact, while all the other quantities (especially those x, y, z closely connected with t by Lorentz transformations) are represented by operators, an ordinary numeric parameter corresponds to the time t, just as in the classic mechanics."

There is no doubt that the elementary quantum theory was elaborated in a non concordant format with the einsteinian relativity, but it is erroneous to suppose that the space coordinates are represented by means of operators. The truth is that t is given the

same treatment as space coordinates (x, y, z); none of those magnitudes possesses a functional operator associated to her. Let us see the reason of this.

2. The birth of a non-mechanical "mechanics"

Quantum theory was developed following the same recipes as hamiltonian mechanics because of two crucial reasons, partly historical and partly logical. Since the victory in the XVIII century of the newtonian natural philosophy, most of the scientists had the conviction that any fundamental theory of the matter –and, generally speaking, of the whole physics- had to be some kind of "mechanics"; that is to say, a group of equations of motion for particles that interact among them obeying some law of forces that could be more or less complex. The field notion introduced later on by Faraday in the XIX century did not change substantially the situation. In fact, the one who mathematically elaborated Faraday's ideas, the famous James Clerk Maxwell, obtained his equations imagining a mechanical ether subjected to Newton laws. It was natural, therefore, that the new theory of quanta were called "quantum mechanics", a not very fortunate name for a parcel of physics born with the purpose of explaining spectroscopic series, distributions of electromagnetic frequencies, specific heats, and a diversity of quantities deeply far from the genuine mechanical magnitudes.

On the other hand, and also in the XIX century, the mathematical methods of Hamilton demonstrated to embrace the mechanical description of particles as much as that of waves too. This supposed an absolute novelty in the traditional procedures of classic mechanics. Many authors have often speculated if with something more than perspicacity -as if Hamilton had lacked it- the brilliant Irishman had been able to take a step more and to discover himself the mathematical framework of the future quantum theory. It does not seem that the state of things were so simple¹:

"(...). It has been said that if Hamilton has advanced a little more, he would have discovered the equation of Schroedinger. It is not this way; he lacked experimental authority to give that jump. In the days of Hamilton it was considered that the classical Mechanics was rigorously certain and justifications based on the experience to consider did not exist that went an approach to a broader theory. (...)."

Neverthless, that experimental arguments did exist in the begining of the XXth century. So, it was almost unavoidable to appeals to the hamiltonian formalism in the nascent quantum physics, as long as its physical referents -the quantum objectsmanifested as much a corpuscular as a wave-like behaviour. And it was made this way, decorating the analytic mechanics of Hamilton with the algebra of operators of Von

¹ Goldstein (1990), p. 596.

Neumann on the functional space of Hilbert. This cocktail of famous names often darkens the relative simplicity of the situation. In the analytic mechanics of Hamilton, a system of *n* particles is described by means of 3n couples of conjugated dynamic variables that are usually represented as q_k and p_k in its canonical form. These variables obey the relationships expressed in the brackets of Poisson:

$$\{q_k, p_l\} = \delta_{kl}; \{q_k, q_l\} = \{p_k, p_l\} = 0.$$

We define with it a 6*n*-dimensional point in the space of phases of the system. And evolution in time is characterized by means of the hamiltonian functional of these dynamic variables, $H = H(q_k, p_l)$ that are assumed to be not explicitly dependent of *t*:

$$dq_k/dt = \{q_k, H\}, \ dp_l/dt = \{p_l, H\}$$

In all the physical theories, except in General Relativity, it is supposed that space and time (or the space-time if we speak about Special Relativity) constitute a merely passive stage in which the natural processes occur, a sort of indispensable inert background to describe the physical phenomena. This makes necessary to distinguish among the space-time coordinates (t, x, y, z) and the dynamical variables q_n and p_n corresponding to the space of phases. The first ones on their own are good to label mathematically the different points of space and time; this is the reason why they could be denominated "field variables", although they were only acting with respect to a metric field that defined distances among these points. On the contrary, q_n and p_n are dynamical variables associated, for example, to the position and the impulse of a certain physical object. They do not label in a generic way all the points of a continuous manifold (as space-time) used as basic framework to formalize our theories; they only refer to the points that the physical object in question occupies in fact. Otherwise, q_n and p_n specify the states of specific material systems, while coordinates t, x, y, z characterize the continuous space-time adopted as our background manifold where we can immerse this specific material systems.

Now we find easier to distinguish, firstly, among the variable of position of a point-like particle, q_x , and the space coordinate of the point that this particle occupies in a certain instant, x. It is true that we have the algebraic relationship $q_x = x$ (and similarly for the rest of coordinates), but we must make a difference between the point-like particle (as a physical entity endowed with mass-energy, position, speed and acceleration) and the geometric coordinate x of a fixed point in a preexistent empty space.

We cannot forget the essential role carried out by symmetries in our understanding of physical laws. This laws are not modified when we change the position of the origin of our reference system (space-temporary symmetry of displacements), neither when we rotate their axes a certain angle (space symmetry of rotations). The Lorentz transformations add the equivalence of systems in relative inertial motion, what is put forwards in the symmetry of space-time rotations. But again we must underline an outstanding distinction: the mandatory fulfilment of certain symmetry requirements in nature, only concerns to the physical laws (that is to say, to the formal representation of the entirety of phenomena and physically permissible processes), not necessarily to the individual and concrete physical systems. A lot of situations will show material systems that, because of the asymmetry of their configuration, for instance, will not be rotationally symmetrical. And it does not mean that the rotational symmetry of the natural laws has been infringed.

The symmetries of spatial displacements are generated by means of the total impulse, P, time symetries depend on total energy², H, and the generator of rotational symmetries is the total angular momentum, J. If all these symmetries are on an equal foot, we may wonder about the priority usualli given to hamiltonian functional, H, representing the time evolution of the physical systems. In our description of nature, what is the priority of time displacements upon the spatial ones owed to? The answer must be sought in the historical tradition of analytical mechanics, mostly engaged to the study of point-like masses and rigid bodies, all which trivially transform under spatial displacements. However, the case of classical fields (electromagnetic, distribution of speeds or densities in a fluid, etc.) is very different, because those displacements in space are anything but trivial. In such situations, P and H get the same importance; so, in Special Relativity energy and linear momentum constitute the components of a tetravector in space-time.

The accented formal similarity among the q_x behavior and of x under time displacements and spatial rotations, has notably darkened the background differences between both magnitudes. And the use of the notation x for the particles position (equally for the other coordinates) has still carried bigger confusion to such an extent that we will hardly find many textbooks where the distinction among both variables is explicitly pointed out. Even more, the bold efforts of some authors to include the time coordinate, t, as conjugated canonical variable of H, were bound to be a failure (provided we stay inside the orthodox hamiltonian scheme). The hamiltonian functional H depends on the original canonical variables, q_n and p_n (and sometimes also on t); therefore time cannot be itself an independent canonical variable. The mistake is consequence, again, of confusing space-time coordinates (a mathematical label assigned to the points of the space-time) with dynamical variables (estates that characterize the physical systems located in space-time).

 $^{^{2}}$ We should not assume that *H* always stands for the total energy of a physical system.

3. Space and time in quantum physics

In the elementary quantum theory the situation is completely similar: the existence of a space-time background, continuous and inert, is presupposed, the points of which are specified by means of space-time coordinates that are classical variables without dispersion ("*c*-numbers" of Dirac). The symmetries and space-time transformations are expressed in terms of such coordinates. The dynamical variables, on the other hand, are indeed quantized, due to which they are substituted by self-adjoint operators in a Hilbert space. All the hamiltonian formulae conserve their validity by only replacing the Poisson brackets for quantum commutators, according to the rule $\{A, G\} \rightarrow (i\hbar)^{-1} [\hat{A}, \hat{G}]$. In particular, the canonical variables are substituted by operators that obey the following commutation relationships:

$$\left[\hat{q}_{k},\hat{p}_{l}\right]=i\hbar\delta_{kl};\left[\hat{q}_{k},\hat{q}_{l}\right]=\left[\hat{p}_{k},\hat{p}_{l}\right]=0$$

We arrive now at one of the keys of this controversy: the substitution of dynamical variables for operators and Poisson brackets for quantum commutators, expose the inherent limitations of representing by means of typical point-particles magnitudes, quantum estates that are in no way associated to point-like objects. In fact, we usually preserve the notation q_i for the cartesian component of the position of the quantum particle, considered as a material point, and similarly for the momentum components, p_j . But it happens to be that p_j has its corresponding differential operator, unlike the variable q_x , which is replaced by the so-called "multiplication operator", $x \cdot ($). This last one is not a genuine operator because of its lack of true self-functions. The Dirac deltas are not even authentic functions in a rigorous mathematical sense, for which no quantum state can be developed as linear superposition of self-functions of the operator position.

The operator linear momentum does not suffer from the precedent complications because the notion of speed, or the linear momentum, is compatible as much with ideally point-like objects (the point-like mass of the classical mechanics) as with extensive entities (an ideal plane wave). However, the dynamical variable q_i only applies to classical objects ideally reducible to a point, which is impossible for quantum ones. For that reason we speak about the propagation speed of a plane wave, but we do not speak about its point-like position; as we do know in geometrical optics, the non-wave limit of a plane wave is a ray, not a point.

A big amount of texts on elementary quantum physics open the explanations considering only a mass-point, what implicates an error from the very begining. We know that quantum objects are spatially extensive entities³, the time-dependent wave function is not written $\psi(q_x, t)$, as we could expect, but $\psi(x, t)$, where x denotes not the

³ That is why the usual name is "electronic field" -insted of "electrons"- or "material fields" in general.

instantaneous position of a pointlike corpuscle, but a geometric coordinate that embraces the whole. And it is natural because a free quantum object is represented by means of an infinite plane wave. The function of quantum state, in fact, is a magnitude still located in a higher level of abstraction, whose formal characterization is given in a functional space of Hilbert with an specific algebra. Anyway, the usual notation, in which x and t appear in an equal foot as arguments of the function ψ , incites us to wonder why t is not an operator like x. The answer, obviously, comes on remembering that neither t nor x are true operators.

By the way, just as they have been defined the operators associated to q_i and p_j are not enclosed, and their spectrum of allowed values extends to the whole real line. When periodic contour conditions are imposed to the variable position, self-values of the linear momentum operator become discreet. And if the wave function must annul in the ends of a finite space interval (the ordinary example of the "particle in a box"), not even exists then a self-adjoint operator for the linear momentum. The insistence of considering t as if it were a genuine operator, would take us to expect that it should obey the relationship $[t, H] = i\hbar$. Being this way, t should possess a continuous spectrum of self-values from $-\infty$ until $+\infty$, in so far it embraces all the moments in time. In consequence, the same behavior should be shown by the hamiltonian H, against the obvious evidence that there exist systems with discreet self-values for the energy.

This reasoning convinced to not few authors about the impossibility of building an "operator time", while the presumed existence of an operator position accentuated the asymmetry between space and time moving away still more the quantum theory from a relativistic spirit. Whereas there is not an authentic problem; neither *x* nor *t* are operators, and the formal symmetry among both coordinates stays. An alternative, certainly, consists on formally defining an operator of time evolution (not an operator "time") that provides the transition from a particular state $\psi_0(x)$ in an instant t_0 until another later state $\psi_t(x)$ in an instant *t*.

We would have this way:

$$\Psi_t(x) = \hat{U}(t)\Psi_0(x)$$

where the operator $\hat{U}(t)$ is equal to an exponential function $\exp_e[-iHt]$. It is easy to prove that the time-evolution operator, although linear, is not hermitian; its self-values, $\exp_e[-iE_nt]$, are not real. For that reason, $\hat{U}(t)$ cannot be considered but a purely formal artifice concieved to express the transition from an initial state to another final one by

means of a linear operator that only depends on H and t, in a mathematical language comparable to that of other authentic physical operators⁴.

The situation is still more delicate when incorporating the Special Relativity in the elementary quantum theory, because then we are even deprived of the position pseudo-operator managed until that moment⁵. In 1949, T.D. Newton and E.P. Wigner published a well-known paper⁶ in which they showed an almost univoc characterization of an operator called "of position" by means of its behavior under displacements and spatial rotations. However, the operator defined this way turns out to be non-covariant in the relativistic sense. Even more, due to the positive sign of the energy in the ordinary physical systems, if in a certain moment we have a self-state of this operator (a "located state", in Newton-Wigner terminology), after an interval of infinitely brief time the subsequent state is extended over all the space. So unpleasant behavior has been a source of plentiful literature around the discussion about the meaning and real utility of the concept of "localizable particle" in the framework of a consistent quantum-relativistic theory⁷.

The truth is that in the usual relativistic versions of quantum physics, neither position nor duration are counted among the basic notions. The main role in this context is played by the operator of quantum field parametrized by means of the space-time coordinates considered as classic magnitudes without dispersion (again, those "*c*-numbers" of Dirac).

4. Heisenberg inequalities

Another source of confusion is what we can denominate the "Schroedinger approach": the idea that elementary particles are no more than ultramicroscopic wave packets (which is equivalent to accept a wave ontology as the ultimate one for the quantum realm and, in general, for the whole physical reality). If a wave packet is a localized disturbance that results from the sum of many different wave forms, the more strongly localized the packet is, the more frequencies (or linear momenta) are needed to allow the constructive superposition in the region of localization and destructive

⁴The time evolution operator plays an outstanding role in higher study of quantum systems depending on time. But it does not change its non-physical condition.

⁵In a relativistic quantum theory the notion of a particle with a "definite location" –and, besides, the notion of wave function as a carrier of a probability density– is still more controversial than in a non-relativistic situation. A good discussion could be found in Malament (1996), or Halvorson & Clifton (2002).

⁶ Newton & Wigner (1949).

⁷ For the Dirac spinors of quantum objects with spin = $\pm \frac{1}{2}$, the Newton-Wigner operator is equivalent to the Foldy-Wouthuysen "position average" operator. More details in Foldy & Wouthuysen (1950).

superposition outside that region. So, representing an arbitrary wave as a superposition of plane waves:

$$\psi(x,0) = \int_{\infty} d^{3}(k) A(k) \exp(ix \cdot k)$$

The amplitudes A(k) can be expressed in turn as a function of $\psi(x,0)$ evaluated at t = 0 by inverting the Fourier transform above:

$$A(k) = (2\pi)^{-3} \int_{\infty} d^3(x) \psi(x,0) \exp(-ix \cdot k)$$

Taking into account that $p = \hbar k$, many authors deduce the Heisenberg relation of indeterminacy $\Delta x \cdot \Delta p_x \ge \hbar$. The same procedure applied to variables as time and angular frequency leads to similar conclusions. If we regard a wave as a function of time, we can write:

$$f(t) = \int_{\infty} d\omega \cdot g(\omega) \cdot \exp(-i\omega t)$$

whre $g(\omega)$ is given by

$$g(\omega) = (2\pi)^{-1} \int_{\infty} dt \cdot f(t) \cdot \exp(i\omega t)$$

Correspondingly, from the features of a wave packet it could be deduced that $\Delta t \cdot \Delta \omega \ge 1$. And if we remember that $E = \hbar \omega$, we arrive at the "fourth relation of indeterminacy" $\Delta t \cdot \Delta E \ge \hbar$.

There are, however, many reason to suspect of the previous results. The main one is tht quantum objects are not wave packets. If it were that way, each individual quantum particle would produce a whole intereference pattern in the double-slit experiment. And this is not what we observe.

From the relationships of commutation, $p_nq_m - q_np_m = -i\hbar\delta_{nm}$, that are postulated by the theory, the definition of quantum average, the definition of standard deviation a (coming from the mathematical statistics⁸), and the inequality of Schwartz⁹ (taken from mathematical analysis), it is obtained without difficulty¹⁰

⁸ For individual quantum objects , statistical dispersions cannot be regarded as those applied to sets of quantum particles. See Uffink & Hilgevoord (1985), or Hilgevoord & Uffink (1988).

⁹ Schwartz (1950-51).

¹⁰ Kennard (1927), Robertson (1929) and Schroedinger (1930).

$\Delta_{\Psi} p_n \Delta_{\Psi} q_m \ge (\hbar/2) \delta_{nm}$

for the component p_n of the momentum and the component q_m of the position of the micro-object represented by ψ .

It is possible, nevertheless, to define a "time of evolution", $\delta_{\psi}t_A$, for any dynamical variable *A* in a state ψ as the interval of necessary time so that the mean value change of *A* would be appreciable compared with their intrinsic dispersion $\Delta_{\psi}A$; algebraically:

$$\delta_{\Psi} t_A \equiv \Delta_{\Psi} A / \{ d \langle A \rangle / dt \}$$

Then we have $\delta_{\psi} t_A \cdot \Delta_{\psi} A \ge \hbar$, that would get to the discussed case when A is the energy¹¹.

5. Conclusions

Most of the debates engendered about the role of space and the time in quantum physics, could have been dissipated distinguishing among the space-time coordinates (that are *c*-numbers) and the dynamical variables (an inheritance of the analytical mechanics by the hamiltonian formalism) that characterize the behavior of the physical systems in space-time. Since the quantum objects are not reducible -not even ideally- to point-like corpuscles, a real "position operator" does not exist in quantum theory, what balances the situation, because neither is there an "operator time." The opposite belief, so common as it is, happens to be founded in a double mistake: on the one hand, to confuse the dynamical variables of position, typical of the particles, with the coordinates of points in space; and on the other, to assign the dynamical variables of position to physical entities, as quantum ones, for which they are essentially inappropriate. The indeterminacy relations are not the same when applied to *x* or *t*. As *t* has no dispersion, the physical meaning of this "fourth Heisenberg's inequality" is different.

When we try to submerge the quantum theory in a relativistic formulation, the covariance requirements for space-time transformations become so demanding that we are even prevented from appealing to a so-called "position operator": the concept of pointlike object gets lost ab initio, still in a much more transparent way that in the non-relativistic quantum theory, and the entirety of the controversy becomes an obsolete one. Finally we would arrive at the domain of the quantum field theory, conceived as the royal road to insert the relativistic covariance in the quantum world. That is, at least,

¹¹ Gillespie (1976), p. 74.

the general consent of the scientific community; a consent that-later on it will be seen-it has their own –and in no way negligible– inconveniences. But this is another story.

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