On Exotic Systems of Baryons in Chiral Soliton Models

Vladimir Kopeliovich\textsuperscript{a,b,*}
\textsuperscript{a) Institute for Nuclear Research of RAS, Moscow 117312, Russia}
\textsuperscript{b) Nordita}

DOI: 10.1016/0370-2693(91)90821-7

Abstract

The role of zero mode quantum corrections to the energy of baryonic systems with exotic quantum numbers (strangeness) is discussed. A simple expression for the contribution depending on strange inertia is obtained in the $SU(3)$--collective coordinate quantization approach, and it is shown that this correction stabilizes the systems the stronger the greater their baryon number is. Furthermore, systems are considered which could be interpreted in the quark model language as containing additional $q\bar{q}$--pairs. It is argued that a strange skyrmion crystal should have additional binding in comparison with the $SU(2)$--quantized neutral crystal.

1 Introduction

One of the most characteristic predictions of chiral soliton models, such as the Skyrme model, is the possible existence of the $SU(3)$--multiplets of dibaryons, tribaryons, etc., whose strange components could be stable relative to strong interactions \cite{1} - \cite{5}. These predictions followed the paper by Jaffe \cite{6} where the existence of a strange dibaryon, the $H$--particle, was predicted in the MIT bag model. The experimental checking of these predictions would be very important since it would provide evidence for the whole concept of baryons as solitons of some effective lagrangian, or lead to substantial modifications of these models.

The purpose of present note is to clear up some general properties of exotic baryon systems in the framework of chiral soliton models and the collective coordinates quantization picture which have not been well understood in previous investigations. We shall use here the rigid rotator model criticized, e.g. in \cite{7}, so the results obtained should be considered as a starting point for more elaborate investigations. Arguments based on a very simple expression for the quantum correction due to rotation in a "strange" direction show that this correction has the property of stabilizing the baryon systems more strongly than than the quantum corrections due to other zero modes, the value of the binding energy being increased with increasing baryon number $B$ of the system.

Below we shall consider objects which can be interpreted as being built of a minimal number of quarks (valence quarks only) as well as those which contain additional quark-antiquark pairs. The latter may have positive strangeness as well as a ratio of strangeness to baryon number $|S/B| > 3$, for some of possible multiplets. The 'price' to be paid for addition of each $q\bar{q}$-pair is estimated, and it is shown that this price is modest. Finally, arguments are given that in the case of strange skyrmion crystals quantum correction due to zero modes stabilize them more strongly than $SU(2)$--quantized skyrmion crystals (neutron crystals).

\textsuperscript{*}e-mail: kopelio@inr.ru
2 \hspace{1cm} SU(3) - quantum corrections for nonexotic and exotic systems

Our consideration is based on the well known expression for the energy of baryon systems obtained from the SU(2)–solitons by means of rotations in the space SU(3)–collective coordinates [1] - [5]. This energy consists of the classical mass of the soliton $M_{cl}$ and quantum corrections due to zero modes $E_{rot}$ as well as nonzero modes (vibration, breathing, etc) which are difficult to calculate and by this reason are usually omitted:

$$E = M_{cl} + E_{rot} =$$

$$= M_{cl} + \frac{1}{\lambda_s} \left[ C_2(SU(3)) - \frac{3Y'^2}{4} - N(N + 1) \right] + \frac{N(N + 1)}{\lambda_r} + ...$$ \hspace{1cm} (1)

$M_{cl}$ is bound, at least for not very large baryon numbers $B$ (the masses of toroidal bound skyrmions with $B = 2, 3, 4$ have been calculated at first in ref. [8]). $C_2(SU(3)) = [p^2 + q^2 + pq + 3(p + q)]/3$ is the SU(3) Casimir operator, $p$ and $q$ being the numbers of upper (lower) indeces in the spinor describing the SU(3) multiplet under consideration, $Y'$ and $N$ are the right hypercharge and isospin (the mistake in the interpretation of $N$ in [8] has been corrected in [5]), $\lambda_s$ and $\lambda_r$ are the moments of inertia characterizing the system, they differ by a factor of two from their standard definition (accepted in classical mechanics, e.g.). The terms depending on the orbital momentum, like $J(J + 1)/\Lambda_r$ will be discussed later. The expression (1) was obtained for systems possessing the generalized axial symmetry [8, 5]. As it was recently established in [9] for systems with baryon numbers $B = 3 - 6$ at least, there exist configurations of more complicated form than the torus-like one, which have even lower energy. However, we expect that some changes in the spatial form of the solitons will not influence the overall structure of the expression (1), so, our our arguments hold also for this case.

Let us consider the SU(3)–multiplets for which the right hypercharge $Y'$ is the largest in the multiplet (we call them the "minimal" multiplets):

$$Y' = \frac{p + 2q}{3}$$ \hspace{1cm} (2)

From the quantization condition [10] it follows that $Y' = N_cB/3$, but here we shall restrict ourselves to the case of the number of colors $N_c = 3$. Since the right isospin $N = p/3$, it is easy to establish that the coefficient of $1/\lambda_s$ in (1) is equal to $(p + 2q)/2 = 3B/2$. So, for the whole family of minimal multiplets with fixed baryon number and $N$ varying from 0 or 1/2 up to $3B/2$ we obtain the universal relation

$$E_{rot} = \frac{3B}{2\lambda_s} + \frac{N(N + 1)}{\lambda_r} + \frac{J(J + 1)}{\Lambda_r} +$$

$$+ J_z^2 \left[ \frac{1}{B^2\lambda_z} - \frac{1}{B^2\lambda_r} - \frac{1}{\Lambda_r} \right].$$ \hspace{1cm} (3)

It was already noted in [1] that the first coefficient in (3) is the same for the octet and decuplet of baryons (hyperons). Therefore, measurements of their masses cannot help in cross-checking the $\lambda_s$. The same holds also for families of baryonic systems with arbitrary baryon numbers. These families consist of $10-$, $27-$, $35-$, $28-$plets for $B = 2$, of $35-$, $64-$, $81-$, $80-$ and $55-$plets for $B = 3$, etc [11].

For toroidal few-baryon systems considered in [4, 8] $\lambda_s(B)$ increases proportional to the baryon number of the system, or even faster. The same holds also for $\lambda_r(B)$. The orbital inertia $\Lambda_r(B)$ is considerably larger than $\lambda_s$ and $\lambda_r$, it grows with increasing $B$ faster than $B^2$. By this reason the depending on $\Lambda_r$ contribution in (3) is smaller than the first two terms, if $J$ is not very large, and does
not play a crucial role in the binding of the system. It follows immediately from (3) that the energy (due to zero modes) per unit baryon number decreases with increasing $B$ and goes to zero for rotations in the "strange" direction. The energy due to isotopic rotations decreases like $1/B^2$ for the multiplets with the smallest value of $N$ (1/2 or 1), but is approximately constant for $N = N_{max} = 3B/2$.

The contribution to the binding energy per unit baryon number for the systems with smallest $N$ increases with increasing $B$ and approaches the value

$$\frac{\epsilon}{B} \simeq \frac{3}{2\lambda_s(B = 1)} + \frac{3}{4\lambda_r(B = 1)}.$$  

(4)

Since usually $\lambda_s$ is a few times smaller than $\lambda_r$, "strange" rotational energy binds baryonic systems more strongly than isotopic rotations \[1, 4, 5\]. The numerical results for the binding energy depend on the values of the model parameters. It follows from (4) that the asymptotic relative binding energy per baryon equals to 0.33 in the Skyrme model \[1, 4\] and to 0.46 in the model with explicit scalar meson \[5\]. These values seem to be too large, but they are smaller in the case of few-baryon systems. The relative binding energy of quantized dibaryons (i.e. divided by the total energy of possible final states) is less sensitive to the type of the model and to the values of the parameters, and in many cases of interest is close to 0.2.

Let us go to the "nonminimal" representations for which the largest possible hypercharge of the system is larger than their baryon numbers. Such multiplets have to contain components with positive strangeness. In the quark model language these systems could contain some number $m$ of quark-antiquark pairs. We have now

$$\frac{p + 2q}{3} = B + m$$

and

$$Y' = \frac{p + 2q}{3} - m,$$

if we assume that in a nonminimal representation the number of indices is increased by $m$ for both $p$ and $q$ in comparison with corresponding minimal representation. Of course, this is not the only possibility. In this case the right isospin $N = (p + m)/2 = N_0 + m$, and an elementary calculation yields

$$C_2(SU(3)) - N(N + 1) - \frac{3B^2}{4} = \frac{3B}{2} + m \left(\frac{3B}{2} + m - N + 1\right).$$

(5)

So, the increase in energy of the system due to the addition of $m$ $q\bar{q}$ pairs contains terms linear in $m$ as well as quadratic ones. At fixed $B$ and $m$ expression (5) decreases with increasing $N$ and is minimal for

$$N_{max} = \frac{3B}{2} + m,$$

the second term in the above formula being equal to $m$. Taking into account the contribution to the energy due to isotopic rotations we obtain for the total increase of the energy due to the addition of $m$ quark-antiquark pairs:

$$\delta E_{rot} = m \left[\frac{3B/2 + 1 + m - N}{\lambda_s} + \frac{2N + 1 - m}{\lambda_r}\right].$$

(6)

For an octet of baryons, e.g., $\delta E_{rot} = 2/\lambda_s + 3/\lambda_r$, at $m = 1$. The contribution of the rotation into the "strange" direction decreases with increasing $N$ down to $m/\lambda_s$, so, when strange quarks are

---

\[1\] Even for a rotating neutron star the energy of the orbital rotation is considerably smaller than the sum of the rotational energies of individual skyrmions (neutrons) if the angular velocity of rotation is not very large (period $T > 10^{-3}$).

\[2\] In this case for starting configuration $(p_0, q_0) = (1, 1)$, $N_0 = 1/2$, and after addition of one $q\bar{q}$-pair we have $(p, q) = (2, 2)$, $N = 3/2$. 

---
"dissolved" in multiplets of higher dimension (in \( p \) or \( N \)) the energy necessary for this "dissolving" decreases. However, the energy of the isotopic rotations increases.

There are also other ways to go to nonminimal representations by means of asymmetric increase of \( p \) and \( q \). E.g., one can increase separately \( p \) or \( q \) by \( 3m \) which will correspond to the addition of \( m \) or \( 2m \) quark-antiquark pairs. In the first case \( N = (m + p)/2 \), in the second one \( N = p/2 + m \), the expression for \( \delta E \) remains the same with the substitution \( m \to 2m \) in the latter case. In the Skyrme model with the parameters \( F_\pi = 108 \text{ MeV}, e = 4.84 \) we have \( \lambda^{-1}_s \simeq 0.3 \text{ GeV} \) and \( \lambda^{-1}_r \simeq 0.1 \text{ GeV} \), and the energy surplus for \( m = 1 \) equals

\[
\delta E_{rot} = \left[ 0.3 \left( \frac{3B}{2} + 2 - N \right) + 0.2N \right].
\]

In the model with an additional scalar meson [5] it is larger by a factor \( \sim 1.5 \). So, \( \delta E \) has the desired order of magnitude because it is expected that such states should be above the threshold for the decay due to strong interactions.

3 Quantum corrections in case of skyrmion crystals

The binding energy of the extended skyrmion crystal consists of two parts: the classical binding energy and the binding energy due to quantum corrections [12] - [17]. The first one depends on the particular symmetry of the crystal. We shall concentrate now on the contribution of the quantum corrections to the binding energy and show that there is a principal distinction between the contribution due to the rotations in the "strange" and "nonstrange" direction.

Let us recall the arguments of ref. [12] which have led to the conclusion that there is an important contribution of the quantum corrections to the binding energy of the crystal. The basic assumption is that flavor rotations of the whole crystal can be described by one and the same set of collective coordinates, i.e. the crystal is coherently rotated in flavor space. It follows immediately that the corresponding moment of inertia of the crystal \( \Lambda = n\lambda \), where \( n \) is the number of unit cells in the crystal, and \( \lambda \) is the moment of inertia of the unit cell. The rotational energy of the whole crystal \( E_{rot} = T(T+1)/n\lambda \) should be compared with \( nt(t+1)/\lambda \) where \( t = 1/2 \) is the isospin of the unit cell. For the neutron crystal \( T = nt = n/2 \), and for large \( n \) we have

\[
E_{rot} = nt^2/\lambda < nt(t+1)/\lambda,
\]

i.e. the effect of binding arises. Its physical sources are the quantum fluctuations of the isospin momentum of the free neutron which make \( \vec{t}_z^2 \) three times greater than \( t_z^2 \). These fluctuations are suppressed inside the crystal, and for the whole crystal they are negligible. It was assumed in the above arguments that the inertia of the unit cell does not differ too much from that of the free neutron.

If we consider now the \( SU(3) \) flavor rotations we obtain under the same assumptions for the energy of the crystal:

\[
E_{rot} = \frac{1}{n\lambda_s} \left[ C_2(SU(3))[P, Q] - \frac{3B^2}{4} - N_{tot}^2 (N_{tot}^2 + 1) \right] + \frac{N_{tot}^2 (N_{tot}^2 + 1)}{n\lambda_r},
\]

\( B = n \), which should be compared with the analogous expression for \( n \) unit cells. \( P \) and \( Q \) are the numbers of indeces in the spinor describing the crystal in the \( SU(3) \) space, \( N_{tot} \) is the corresponding right isospin, \( N_{tot}^2 = P/2 \). The simplest and probably most natural possibility is to assume that \( P = np \) (by analogy with the neutron crystal). In this case we obtain

\[
E_{rot} = \frac{3B}{2n\lambda_s} + \frac{n^2}{4n\lambda_r} = \frac{3}{2\lambda_s} + \frac{n}{4\lambda_r},
\]

\( \delta E_{rot} = \left[ 0.3 \left( \frac{3B}{2} + 2 - N \right) + 0.2N \right].\)
This expression is quite similar to the expression for $E_{rot}$ in the case of few-baryon systems. The distinctions are that now $n\lambda$ enters instead of $\lambda(B)$, $n/2$ instead of $N$, and the orbital rotation energy is absent (for the crystal at rest). So, due to the cancellation of the quadratic terms in $P, Q$ we have obtained that the contribution of the first term in (9) to the energy of the crystal is constant and equal to $3/2\lambda_s$. This is the main difference between "strange" and isotopical rotations. As in the case of few-baryon systems, this term is the same for all SU(3) representations $(P, Q)$ which satisfy the relation $P + 2Q = 3n$.

The discussion of the astrophysical aspects of the possible existence of strange matter is beyond the scope of the present paper. Observational results imply that there is a place for dark baryonic matter in the Universe, although arguments were given that the strange matter should have evaporated at the early stages of the evolution of the Universe [18]. However, the possibility of the existence of strange matter does not seem to be excluded completely.

4 Discussion of some general problems

We have shown that in the collective coordinate quantization approach a transparent expression for the quantum correction to the energy of the baryon system can be obtained, illustrating that rotations into "strange" direction stabilize few-baryon systems as well as skyrmion crystals even more than isotopic rotations. The stabilizing property of zero modes quantum corrections seems to be a simple consequence of a natural property of solitons: their geometrical sizes and moments of inertia increase with increasing baryon number.

The fact that the quantum corrections due to "strange" rotations decrease with increasing baryon number of the system may also mean that the whole approach of the SU(3)– collective coordinates quantization becomes more selfconsistent with increasing $B$. Indeed, several attempts to describe the $B=1$ hyperon spectrum [10, 19, 20] have met problems since the strange inertia $\lambda_s$ is especially small for $B=1$, and therefore the quantum correction is especially large in this case. As shown above, the relative value of this correction drops as $1/B$, so the problems in the $B=1$ sector disappear for systems with large $B$. However, if this argument is correct, the same result should be obtained in other versions of quantization, e.g. in that of Callan and Klebanov. In other words, it may turn out that systems with large $B$ are in some sense simpler objects than system with $B=1$. There remain, however, some effects which role has not been clarified till now.

First of all, the vibrational corrections remain to be calculated. They should increase the energy of the system, as well as of the whole crystal. For the system with $B=1$ they have been analyzed in [21, 22]. An elegant method for estimating them has been proposed in [23], and for the case of the skyrmion crystal in [15].

The other problem is the applicability of the soliton approach with increasing baryon number. It is clear that under usual conditions the soliton (skyrmion) cannot be too large, e.g. with baryon number of the order of the Avogadro number. But it is unclear what parameter defines the maximal baryon number of the soliton.

We have not included in our consideration the meson mass terms in the effective lagrangian of the model. According to our previous experience [3] - [5] the mass terms do not play a crucial role in the binding of quantized skyrmions, although they define the mass splittings inside the multiplets, which happened to be too small in comparison with experimental data, especially in Syracuse model

---

3The first term in Eq. (9) should be compared with corresponding contribution for $B = n$ separate cells (baryons), equal to $3n/2\lambda_s$
The mass terms induce also the mass splittings of the quantized skyrmion crystals.

The author is greatly indebted to Andreas Wirzba for many useful discussions and reading of the manuscript, and to Andy Jackson for enlightening discussions of some points of principle.

References

1. A.P. Balachandran, F. Lizzi, V.G. Rogers and A. Stern, Nucl. Phys. B256 (1985) 525
2. J. Kunz and P.J. Mulders, Phys. Lett. B215 (1988) 449
3. V.B. Kopeliovich, Yad. Fiz. 51 (1990) 241
4. V.B. Kopeliovich and B.E. Stern, NORDITA preprint 89/34 (1989)
5. V.B. Kopeliovich, B. Schwesinger and B.E. Stern, Phys. Lett. B242 (1990) 145
6. R.L. Jaffe, Phys. Rev. Lett. 38 (1977) 196
7. G. Holzwarth and B. Schwesinger, Rep. Prog. Phys. 49 (1986) 825
8. V.B. Kopeliovich and B.E. Stern, JETP Lett. 45 (1987) 203
9. E. Braaten, L. Carson and S. Townsend, Phys. Lett. B235 (1989) 147
10. E. Guadagnini, Nucl. Phys. B236 (1984) 35
11. N.W. Park, J. Schechter and H. Weigel, Phys. Lett. B224 (1989) 171
12. I. Klebanov, Nucl. Phys. B262 (1985) 133
13. E. Wüst, G.E. Brown and A.D. Jackson, Nucl. Phys. A468 (1988) 419
14. A.S. Goldhaber and N.S. Manton, Phys. Lett. B198 (1987) 231
15. T.S. Walhout, Nucl. Phys. A484 (1988) 392
16. A.D. Jackson and J.J.M. Verbaarschot, Nucl. Phys. A484 (1988) 419
17. L. Castillejo, et al, Nucl. Phys. A501 (1989) 801
18. C. Alcock and A. Olinto, Phys. Rev. D39 (1989) 1233
19. M. Praszalowicz, Phys. Lett. B158 (1985) 264
20. M. Chemtob, Nucl. Phys. B256 (1985) 600
21. I. Zahed, A. Wirzba and U. Meissner, Phys. Rev. D33 (1986) 830
22. M.I. Eremeev, R.V. Konoplich, A.E. Kudryavtsev, B.V. Martenyanov and S.G. Rubin, ITEP, Moscow, preprint 77 (1987)
23. T.S. Walhout, Phys. Lett. B227 (1989) 10