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Classical ground states of symmetric Heisenberg spin systems

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Abstract

We investigate the ground states of classical Heisenberg spin systems which have point group symmetry. Examples are the regular polygons (spin rings) and the seven quasi-regular polyhedra including the five Platonic solids. For these examples, ground states with special properties, e.g. coplanarity or symmetry, can be completely enumerated using group-theoretical methods. For systems having coplanar (anti-) ground states with vanishing total spin we also calculate the smallest and largest energies of all states having a given total spin *S*. We find that these extremal energies depend quadratically on *S* and prove that, under certain assumptions, this happens only for systems with coplanar S = 0 ground states. For general systems the corresponding parabolas represent lower and upper bounds for the energy values. This provides strong support and clarifies the conditions for the so-called rotational band structure hypothesis which has been numerically established for many quantum spin systems.

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1. Introduction

The study of small interacting spin systems is not only of theoretical interest but also of importance for the experimental investigation of recently synthesized magnetic molecules [1–4]. An exact calculation of the thermal expectation values of the relevant quantum observables and other quantities such as correlation functions is possible only in very few cases. Given this situation, a classical treatment yields a first approximation for individual spins *s* with $s \gg 1$ which is often astonishingly good, as well as bounds for the exact quantum values (cf Berezin/Lieb-inequalities). It is also valuable as a guide for developing approximation schemes for attacking the quantum theoretical problem.

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One of the fundamental characteristics of a system is its ground state(s) and the corresponding ground state energy. The problem of determining the exact classical ground state has been considered for a long time, see e.g. [5, 6], but has only been solved for a few special systems. These include arrays of interacting spins occupying the sites of a regular N-polygon (spin rings) with nearest-neighbour coupling constant J > 0 (antiferromagnetic in our convention). The corresponding ground state energy is

$$E_{\min} = -2JNs^2 \qquad (\text{even } N) \tag{1}$$

corresponding to an alternating configuration of spin orientations in the N-polygon. The analogous result

$$E_{\min} = 2JNs^2 \cos((N-1)\pi/N) \qquad (\text{odd } N) \tag{2}$$

does not appear to be in the literature, although it is easily established, see (77). Another class of systems where the classical ground states are easily determined are the so-called bi-partite systems, which can be divided into two subsets *A* and *B* such that there are only interactions (with a positive coupling constant *J*) between spins $a \in A$ and $b \in B$. In this case any anti-parallel configuration between *A* and *B* will minimize the energy. The regular polygons with even *N* considered above and the cube are simple examples of bi-partite systems.

Recently, the exact classical ground state of a system consisting of N = 30 spins occupying the vertices of an icosidodecahedron [7, 8] has been determined [9]. The method employed there can be applied to any system which can be decomposed into a set of triangles without common edges and which is 3-colourable. Examples are the octahedron (as mentioned in [9]) and the cuboctahedron (this is obtained by joining the midpoints of the edges of a cube with their nearest neighbour, see section 6.4.3). All these systems have the following property: among their ground states are coplanar ground states with S = 0. These can be modified to remain ground states in the presence of a magnetic field \mathcal{H} , the corresponding energy depending quadratically on \mathcal{H} , cf [9]. The construction of the modified ground states is similar to folding of an umbrella and will be referred to as the 'umbrella construction'. This yields a definite prediction for the magnetization versus magnetic field \mathcal{H}_c is applied, and for $\mathcal{H} > \mathcal{H}_c$ the magnetization is saturated.

In the present paper we will derive a further consequence for systems having (among all ground states) coplanar ground states with S = 0. The same formal equation as that characterizing the ground state in the presence of a magnetic field occurs for the ground state subject to the extra constraint that the square of the total spin assumes a given value. Generalizing the solution in [9] we obtain by the umbrella construction a whole family of states assuming the minimal energy $E_{\min}(S)$. It turns out that for these systems the function $E_{\min}(S)$ is quadratic in S of the form

$$E_{\min}(S) = \frac{j - j_{\min}}{N}S^2 + j_{\min}N.$$
 (3)

Here *j* is the row sum $(j = \sum_{\nu} J_{\mu\nu})$ of the adjacency matrix \mathbb{J} defined below. \mathbb{J} is chosen in such a way that *j* does not depend on μ , and j_{\min} is the smallest eigenvalue of \mathbb{J} . Conversely, the existence of a lower parabola (3) implies the coplanarity of some of the S = 0 ground states, see theorem 3 for details. We do not claim that the states constructed from the coplanar S = 0 ground states are unique. Of course, in general there will be many ground states in the presence of a magnetic field or for given *S* which are different from those of the umbrella construction.

Analogous results hold for the maximal eigenvalue $E_{max}(S)$ of systems with S = 0 coplanar anti-ground states. Hence these systems show an exact rotational band structure

(RBS) which has been conjectured to apply for a general class of quantum spin systems [10]. Thus our findings strongly support that conjecture in the sense that the considered class of spin systems has exact RBS-parabolas in the classical limit $s \rightarrow \infty$. On the other hand we have found classes of systems which deviate from the exact RBS-parabolas, especially for small values of *S*, see section 6. Hence our work may be viewed as a first step towards an understanding of the conditions for the occurrence of RBS-parabolas. We will call systems, which exactly attain their lower RBS-parabola (3), 'parabolic'. For general systems the corresponding parabolas represent lower and upper bounds for the energy values, as will be proved in section 3. Sometimes these bounding parabolas will be very good approximations of the boundaries of the energy spectrum even for non-parabolic systems.

Other questions arise in connection with spin systems having coplanar S = 0 ground states. Has every classical spin system coplanar ground states or even coplanar S = 0 ground states? If not, what is the criterion for having coplanar ground states? Are there systems with both coplanar and non-coplanar ground states? We do not have complete answers to these questions, but we can show the following:

- In general the spin triangle with arbitrary coupling constants J_1 , J_2 , J_3 has only coplanar ground states with S > 0.
- A spin system which we call 'pentagonal star' has a non-coplanar ground state with S > 0. The existence of coplanar states with the same energy can be excluded on the basis of numerical calculations.
- For certain systems the existence of coplanar, symmetric ground states can be excluded. By 'symmetric' we mean roughly that the state has the same symmetry as the spin system itself (for details see section 5).
- Other systems, such as the dodecahedron and the icosahedron, have symmetric noncoplanar ground states and apparently no coplanar ones. These ground states can be geometrically visualized in terms of what is known [7] as the 'great stellated dodecahedron' and the 'great icosahedron', respectively.
- The tetrahedron has a symmetric non-coplanar ground state with S = 0 and a variety of other non-symmetric ground states including coplanar ones, all of the same energy.
- The same is true for other systems with full permutational symmetry ('*N*-pantahedron'), except that these systems do not have symmetric ground states for N > 4.
- The cuboctahedron has a lower RBS-parabola but interestingly it has two symmetric ground states with S = 0, a coplanar one and a non-coplanar one. The latter corresponds to a degenerate stellated figure which we have not found in the literature and which is composed of four stars of David, see section 6.4.

Thus we are faced with a surprising variety of possibilities of ground states of different types. The main open problem is to rigorously prove the existence of spin systems without coplanar ground states. It appears that quantitative statements can be made mainly if symmetric states are concerned, since in this case we may use standard methods of group theory, in particular representation theory of the point groups.

The paper is organized as follows: sections 2 and 3 contain the general definitions and results which are independent of symmetry assumptions. The main result of section 3 is that systems with coplanar S = 0 ground states have RBS-parabolas (theorem 2) and its converse (theorem 3). These parabolas represent lower and upper energy bounds even for nonparabolic systems (theorem 1). In section 4 we address questions of extension of ground states to larger systems and 'higher order frustration'. Section 5 introduces symmetry assumptions and the definition and simple properties of symmetric states. This machinery is applied in section 6 to the investigation of particular examples. We consider the general triangle which is generically non-parabolic as well as classes of symmetric spin systems including the quasi-regular polyhedra. We construct symmetric ground states from certain irreducible representations of the respective point groups. The paper closes with a summary.

2. Notation, definitions and general results

We consider as the phase space of the classical spin system the *N*-fold product of the unit sphere S^2 . Instead of using canonical coordinates we will denote a state by a sequence *s* of unit vectors s_{μ} , $\mu = 1, ..., N$ with components s_{μ}^i , i = 1, 2, 3. The total spin is $S \equiv \sum_{\mu} s_{\mu}$ with components S^i , i = 1, 2, 3. If not indicated otherwise, the abstract letters *i*, *j*, ... will always be (upper) indices and there is no danger of confusion with the square of a vector, e.g. S^2 . We will use bracketed indices to denote by, say, $x^{(i)}$ the vector with the components x^i , similarly for matrices. For example, the $n \times m$ matrix $A_{(i)(j)}$ consists of *n* row vectors $A_{i(j)}$ and of *m* column vectors $A_{(i)j}$.

If not mentioned otherwise, the Hamilton function (or 'energy') will be of the form

$$H_0(s) = \sum_{\mu\nu} J_{\mu\nu} s_{\mu} \cdot s_{\nu}. \tag{4}$$

If a magnetic field $\vec{\mathcal{H}}$ is to be included we add a Zeeman term and obtain

$$H_h = H_0 - \mathbf{h} \cdot \mathbf{S} \tag{5}$$

where $h \equiv \sqrt{s(s+1)}g\mu_B \vec{\mathcal{H}}$ contains the common combination of the spectroscopic splitting factor g and the Bohr magneton μ_B and is scaled with the function $\sqrt{s(s+1)}$ of the corresponding spin quantum number s.

Note that the exchange parameters $J_{\mu\nu}$ are not uniquely determined by the Hamiltonian H_0 via (4). Different choices of the $J_{\mu\nu}$ leading to the same H_0 will be referred to as different 'gauges'. We will adopt the following gauges. First, the antisymmetric part of \mathbb{J} does not occur in the Hamiltonian (4). Hence we will follow common practice and choose $J_{\mu\nu} = J_{\nu\mu}$. Thus the $J_{\mu\nu}$ can be considered as the entries of a real symmetric matrix \mathbb{J} . Second, since $\mathbf{s}_{\mu} \cdot \mathbf{s}_{\mu} = 1$ we may choose arbitrary diagonal elements $J_{\mu\mu}$ without changing H_0 , as long as their sum vanishes, $\operatorname{Tr} \mathbb{J} = 0$. The usual gauge chosen throughout the literature is $J_{\mu\mu} = 0$, $\mu = 1, \ldots, N$, which will be called the 'zero gauge'. In this paper, however, we will choose another gauge, to be called the 'homogeneous gauge', which is defined by the condition that the row sums

$$J_{\mu} \equiv \sum_{\nu} J_{\mu\nu} \tag{6}$$

will be independent of μ . Note that the eigenvalues of \mathbb{J} may non-trivially depend on the gauge. We found that the eigenvalues of \mathbb{J} are only relevant for energy estimates if the row sum of \mathbb{J} is constant. This would restrict the applicability of large parts of our theory if we adopt the zero gauge. However, by introducing the homogeneous gauge we can apply our results to a very general class of systems.

The quantity

$$Nj \equiv \sum_{\mu\nu} J_{\mu\nu} \tag{7}$$

is gauge independent. If exchange parameters satisfying $\tilde{J}_{\mu\nu} = \tilde{J}_{\nu\mu}$ are given in the zero gauge, the corresponding parameters $J_{\mu\nu}$ in the homogeneous gauge are obtained as follows:

$$J_{\mu\nu} \equiv \widetilde{J}_{\mu\nu} \qquad \text{for} \quad \mu \neq \nu \tag{8}$$

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$$J_{\mu\mu} \equiv j - \widetilde{J}_{\mu}. \tag{9}$$

It follows that

$$j = J_{\mu} = \sum_{\nu} \widetilde{J}_{\mu\nu} + J_{\mu\mu}.$$
 (10)

A spin system is called *antiferromagnetic* (AF) iff all $J_{\mu\nu} \ge 0$ and not all $J_{\mu\nu} = 0$ for $\mu \ne \nu$. A system is called *connected* iff its spin sites cannot be decomposed into two disjoint classes, say \mathcal{A} and \mathcal{B} , such that $J_{\mu\nu} = 0$ if $\mu \in \mathcal{A}$ and $\nu \in \mathcal{B}$.

In the special case where all exchange parameters $J_{\mu\nu}$, $\mu \neq \nu$ are 0 or *J*, the system can be essentially represented by its graph Γ . A graph $\Gamma = (\mathcal{V}, \mathcal{E})$ consists of a set of 'vertices' \mathcal{V} and a set of 'edges' $\mathcal{E} \subset [\mathcal{V}]^2$. Here $[\mathcal{V}]^2$ denotes the set of subsets of \mathcal{V} with exactly two elements. In our case, the vertices are the spin sites, $\mathcal{V} = \{1, \ldots, N\}$, and the edges represent interacting pairs of spins. For spin systems it is appropriate to consider only graphs without loops and multiple edges, as we do in this paper following, e.g., [11].

There are different ways to graphically represent a spin state s. For specific spin systems, e.g. magnetic molecules, the spin sites $\mu \in \mathcal{V}$ are embedded into the physical 3-space and represented by vectors r_{μ} . One way would be to attach the vectors s_{μ} as small arrows to the sites given by r_{μ} , see section 6.4.3 for one example. However, when using this representation it is difficult to visualize the structure of the s_{μ} in spin space. Hence we will mostly employ another method of representation: we draw the *N* vertices in spin space according to the unit vectors s_{μ} and join them with lines according to the edges of the original spin graph Γ . Thus we obtain a graph isomorphic to Γ which contains additional information about the considered spin configuration.

In graph theory the set of edges is often represented by a matrix called the *adjacency matrix*. We will use this name also for the matrix J in the general case of different exchange parameters. There exists an extended literature about the connection between the structure of a graph and the spectrum of J, see, for example [12], where also applications in chemistry and physics are mentioned.

The graph Γ is called *complete* if $\mathcal{E} = [\mathcal{V}]^2$. We will also call the corresponding spin system where any two spins interact with equal strength, a *pantahedron*.

Being symmetrical, \mathbb{J} has a complete set of *N* real, ordered eigenvalues $j_{\min}, \ldots, j_{\max}$ which are counted according to their multiplicity. One of them is the row sum *j* with $\mathbf{1} \equiv \frac{1}{\sqrt{N}}(1, 1, \ldots, 1)$ as the corresponding eigenvector. For connected AF-systems, *j* will be non-degenerate and equals the largest eigenvalue, $j = j_{\max} > 0$, by the theorem of Frobenius–Perron. For general AF-systems we have the following result:

Lemma 1. For AF-systems

$$j_{\min} < 0 < j. \tag{11}$$

Proof. We have

$$Nj = \sum_{\mu,\nu} J_{\mu,\nu} = \sum_{\mu} J_{\mu,\mu} + \sum_{\mu \neq \nu} J_{\mu,\nu}.$$
 (12)

The first sum of the rhs equals $\operatorname{Tr} \mathbb{J}$ and hence vanishes. The second sum is positive due to the definition of AF-systems, hence j > 0. If $j_{\min} \ge 0$, all eigenvalues j_{μ} would be ≥ 0 and hence $\operatorname{Tr} \mathbb{J} = \sum_{\mu} j_{\mu} > 0$, since at least the eigenvalue j is positive. This contradicts $\operatorname{Tr} \mathbb{J} = 0$.

It will be convenient to rearrange the indices such that $j = j_0$ and the remaining eigenvalues j_1, \ldots, j_{N-1} are ordered according to their size. Hence j_{N-1} will always denote

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the largest eigenvalue of \mathbb{J} other than j, counted once. It may happen that $j_{N-1} = j_{\text{max}}$, for example if $j < j_{\text{max}}$ or if $j = j_{\text{max}}$ is at least twofold degenerate. Since the eigenvalue j_{N-1} will often occur in our paper, especially for AF-systems, we will write $j_{N-1} \equiv j_{\text{maxi}}$.

Sums over $\alpha = 1, ..., N - 1$ but excluding $\alpha = 0$ will be denoted by \sum' . We will denote the α th normalized eigenvector of \mathbb{J} by $C_{\alpha}^{(\nu)}$, i.e.

$$\sum_{\nu} J_{\mu\nu} C^{\nu}_{\alpha} = j_{\alpha} C^{\mu}_{\alpha} \qquad \mu, \alpha = 0, \dots, N-1$$
(13)

and

$$\sum_{\mu} \overline{C^{\mu}_{\alpha}} C^{\mu}_{\beta} = \delta_{\alpha\beta} \qquad \alpha, \beta = 0, \dots, N-1$$
(14)

where we also allow for the possibility of choosing complex eigenvectors.

We note in passing that J in the homogeneous gauge is essentially the Hamilton operator of the corresponding quantum spin system, restricted to the subspace of total magnetic quantum number M = Ns - 1. Thus there is an unexpected connection between 'weakly symmetric' classical states to be defined below and quantum eigenstates of the Hamiltonian lying in the mentioned subspace.

The equations of motion resulting from the Hamiltonian (4) are

$$\frac{\mathrm{d}}{\mathrm{d}t}s_{\mu} = \left(\sum_{\nu} J_{\mu\nu}s_{\nu}\right) \times s_{\mu}.$$
(15)

Using $J_{\mu\nu} = J_{\nu\mu}$ one can immediately show that the total spin vector S is a constant of motion. We now will define various kinds of special states. A *stationary* state will be one with $\frac{d}{dt}s_{\mu} = 0$, for all $\mu = 1, ..., N$. According to (15), this is equivalent to

$$\sum_{\nu} J_{\mu\nu} s_{\nu} = \kappa_{\mu} s_{\mu} \tag{16}$$

for some real numbers κ_{μ} , $\mu = 1, ..., N$. Equation (16) can also be viewed as the solution of the problem to seek states with vanishing variation of the quadratic form

$$\sum_{\mu\nu} J_{\mu\nu} s_{\mu} \cdot s_{\nu} \tag{17}$$

subject to the constraints $s_{\mu} \cdot s_{\mu} = 1$, $\mu = 1, \dots, N$. The κ_{μ} then appear as Lagrange parameters of this variational problem.

Geometrically, (16) means that for a stationary state each spin vector is proportional to the 'centre of mass' of its neighbours. As for general Hamiltonian systems, the stationary states are just the critical points of the Hamilton function, i.e. those points where the gradient of H_0 vanishes. In particular, the states *s* with minimal $H_0(s) = E_{\min}$ or maximal $H_0(s) = E_{\max}$ are stationary. The former will be called *ground states*, and the latter *anti-ground states*. For connected AF-systems the anti-ground state is always of the form $s_{\mu} = e$ for all $\mu = 1, \ldots, N$, i.e. all spins are aligned parallel to an arbitrary unit vector *e* (ferromagnetic ordering). In this case $j_{\max} = j$ and $E_{\max} = Nj$.

Moreover, we will consider the *relative ground states* (*relative anti-ground states*), which are defined as the states of minimal (maximal) energy among all states satisfying $(\sum_{\nu} s_{\nu})^2 = S^2$. Let $E_{\min}(S)$ denote the energies of the relative ground states, and $E_{\max}(S)$ that of the relative anti-ground states. Again, the relative ground states and anti-ground states are among the solutions of the variational problem corresponding to (17) with the additional

constraint $(\sum_{\nu} s_{\nu})^2 = S^2 > 0$. If the extra Lagrange parameter is called χ , we obtain the condition, analogous to (16),

$$\sum_{\nu} J_{\mu\nu} s_{\nu} = \kappa_{\mu} s_{\mu} + \chi S \qquad \text{for} \quad S > 0.$$
⁽¹⁸⁾

The equations of motion (15) then imply that for states satisfying (18) we have

$$\frac{\mathrm{d}}{\mathrm{d}t}s_{\mu} = \chi S \times s_{\mu} \tag{19}$$

i.e., for S > 0 these states are no longer stationary but are precessing around the total spin vector. This is completely analogous to the precession of stationary states of H_0 in the presence of a magnetic field.

The condition S > 0 in (18) is necessary in order to apply the method of Lagrange parameters. S = 0 has to be excluded since it would correspond to a critical point of the constraining function $g(s) = (\sum_{v} s_{v})^{2} - S^{2}$, see for example [13]. To cover also the case S = 0 we consider three constraining equations written in the vector form

$$\sum_{\nu} s_{\nu} = S_0 \tag{20}$$

where S_0 is a fixed vector. Then the above-mentioned problem does not occur. We have three Lagrange parameters which will be written as the components of a vector L. The resulting equations are

$$\sum_{\nu} J_{\mu\nu} s_{\nu} = \kappa_{\mu} s_{\mu} + L.$$
⁽²¹⁾

Taking the vector product of (21) $\times s_{\mu}$ and summing over μ yields $\mathbf{o} = S \times L$. If $S \neq \mathbf{o}$ this is equivalent to $L = \chi S$ and we recover (18). For S = 0, however, (18) has to be replaced by (21).

States satisfying (21) will be called *weakly stationary* states, since they are stationary in a rotating frame. Hence relative (anti-) ground states are weakly stationary.

There exists a simple mechanical model for stationary states. Consider a system of N rigid, massless rods of unit length fixed at the same point ('pendula'). These rods may be represented by unit vectors s_{μ} . Between two rods s_{μ} and s_{ν} we consider springs satisfying Hooke's law with spring constants $2J_{\mu\nu}$. Further there is a constant ('gravitational') force F = Fe. By the cosine theorem, $|s_{\mu} - s_{\nu}|^2 = 2(1 - s_{\mu} \cdot s_{\nu})$, hence the potential energy of our mechanical model will be

$$V(s) = \frac{1}{2} \sum_{\mu\nu} J_{\mu\nu} |s_{\mu} - s_{\nu}|^{2} - F \cdot S = Nj - \sum_{\mu\nu} J_{\mu\nu} s_{\mu} \cdot s_{\nu} - F \cdot S.$$
(22)

The equilibrium points of this mechanical system will satisfy the zero-force equation

$$2\sum_{\mu\nu}J_{\mu\nu}s_{\nu}+F=2\kappa_{\mu}s_{\mu}$$
(23)

where $2\kappa_{\mu}s_{\mu}$ are essentially the constraining forces exerted on the rigid rods. Moreover, for equilibrium *S* will be proportional to the constant force, say $F = -2\chi S$, since otherwise there would be a first-order variation of the potential energy when performing infinitesimal uniform rotations of the system. It follows that equations (23) and (18) are formally identical. Summarizing, there is a 1:1 correspondence between the weakly stationary spin states *s* and the equilibrium states *s* of the considered mechanical systems, if the value of the force *F* runs through all real numbers. This mechanical model may help the reader to visualize some of the weakly stationary states to be considered in this paper.

We can easily obtain the following bounds for the energy of stationary states s:

Lemma 2. $Nj_{\min} \leq E_{\min} \leq H_0(s) = \sum_{\mu} \kappa_{\mu} \leq E_{\max} \leq Nj_{\max}$. Moreover, if the total spin of *s* vanishes we have $H_0(s) \leq Nj_{\max}$.

Proof. By definitions, for stationary states we have $H_0(s) = \sum_{\mu\nu} J_{\mu\nu} s_{\nu} \cdot s_{\mu} = \sum_{\mu} \kappa_{\mu} s_{\mu} \cdot s_{\mu} = \sum_{\mu} \kappa_{\mu} \cdot s_{\mu}$. The expectation value of a real, symmetric matrix is bounded by its smallest and largest eigenvalue, hence

$$j_{\min}\sum_{\mu} x_{\mu}^2 \leqslant \sum_{\mu\nu} J_{\mu\nu} x_{\mu} x_{\nu} \leqslant j_{\max} \sum_{\mu} x_{\mu}^2.$$

$$(24)$$

Choosing $x_{\mu} = s_{\mu}^{i}$ and summing over i = 1, 2, 3 yields

$$Nj_{\min} \leqslant \sum_{\mu\nu i} J_{\mu\nu} s^i_{\mu} s^i_{\nu} = H_0(s) \leqslant Nj_{\max}.$$
(25)

By definition we have $E_{\min} \leq H_0(s) \leq E_{\max}$. Since $E_{\min} = H_0(s)$ for any ground state *s*, and every ground state is stationary, (25) implies $Nj_{\min} \leq E_{\min}$. Arguing analogously with an anti-ground state *s* one may show $E_{\max} \leq Nj_{\max}$.

anti-ground state *s* one may show $E_{\max} \leq Nj_{\max}$. Now let a stationary state *s* satisfy $\sum_{\mu} s_{\mu}^{i} = S^{i} = 0$. The latter equation may be viewed to say that the three vectors $s_{(\mu)}^{i}$, i = 1, 2, 3 are orthogonal to the eigenvector **1** of \mathbb{J} corresponding to the eigenvalue $j = j_{0}$. In the subspace orthogonal to **1** the largest eigenvalue of \mathbb{J} is j_{\max} . Hence $H_{0}(s) \leq Nj_{\max}$ in this case.

The respective bounds of lemma 2 are attained if there exist stationary states where all $\kappa_{\mu} = j_{\text{min}}$ or $\kappa_{\mu} = j_{\text{maxi}}$, or $\kappa_{\mu} = j_{\text{maxi}}$ in the AF case for S = 0. At present, we cannot show the existence of such states under general conditions but this can be done for some specific examples. We will call any stationary state with the property that the values of all κ_{μ} of (16) are independent of μ , a *weakly symmetric* state. If we view the set of components s_{μ}^{i} of a state s as an $N \times 3$ -matrix, the definition of 'weakly symmetric' could be rephrased in the way that all rows of s are eigenvectors of \mathbb{J} for the same eigenvalue \tilde{j} . Recall that for stationary states the constant of proportionality will be the same for all spins s_{μ} , which motivates the wording. Obviously, weakly symmetric (anti-)ground states belong to the eigenvalue j_{\min} , (j_{\max}). If one takes any three eigenvectors of \mathbb{J} with the same eigenvalue to form the three rows of a matrix, one will not get automatically a weakly symmetric state, since the N columns need not be unit vectors. In section 5 we will present a construction procedure which, under certain assumptions, yields (weakly) symmetric states.

Since the energy $H_0(s)$ is invariant w.r.t. rotations/reflections in spin space the abovedefined classes of weakly symmetric states, (weakly) stationary states or of (relative) ground states are also invariant under rotations/reflections.

Another simple property of weakly symmetric ground states is the following:

Lemma 3. Every weakly symmetric state has vanishing total spin if the corresponding eigenvalue \tilde{j} is different from the row sum j. In particular, every weakly symmetric ground state of an AF-system has vanishing total spin.

Proof. We obtain $\sum_{\mu\nu} J_{\mu\nu} s_{\nu} = j \sum_{\nu} s_{\nu} = j S$ and $\sum_{\mu\nu} J_{\mu\nu} s_{\nu} = \tilde{j} \sum_{\mu} s_{\mu} = \tilde{j} S$. Hence $\tilde{j} = j$ or $S = \mathbf{0}$. If s is a weakly symmetric ground state of an AF-system, hence $\tilde{j} = j_{\min}$, we have $j_{\min} < j$ by lemma 1 and its total spin must vanish.

We will call a state *s* collinear if all its spin vectors are (anti-)parallel to a given vector *e*, i.e. $s_{\mu} = \pm e$. Due to equation (16), every collinear state is stationary. Further we will call a state *s* coplanar if all its spin vectors lie in a plane i.e., if there exists a non-zero 3-vector *n* such that $n \cdot s_{\mu} = 0$ for all $\mu = 1, ..., N$. In section 6 we will show that systems such as the regular polygon, the pantahedron and the quasi-regular cuboctahedron and icosidodecahedron admit coplanar weakly symmetric S = 0 ground states. These ground states are important since one can construct from them a whole family of relative ground states, which show a rotational band structure. This will be done in the next section.

3. Criteria for parabolicity

Throughout this section we assume the AF case, although some results will have obvious counterparts for ferromagnetic or general systems. First we will show that the *S*-resolved energy spectrum is bounded by two parabolas. The proof is completely analogous to the quantum case, see [15]. We recall that $C_{\alpha}^{(\nu)}$ denotes the α th normalized eigenvector of \mathbb{J} and consider a transformation of the spin vectors analogous to the transformation onto the eigenbasis of \mathbb{J} . Define

Definition 1. $T_{\alpha} \equiv \sum_{\mu} \overline{C_{\alpha}^{\mu}} s_{\mu}$, and $Q_{\alpha} \equiv \overline{T}_{\alpha} \cdot T_{\alpha}$, $\alpha = 0, \dots, N-1$.

The inverse transformation then yields

$$s_{\mu} = \sum_{\alpha} C_{\alpha}^{\mu} T_{\alpha} \qquad \mu = 1, \dots, N.$$
⁽²⁶⁾

Especially, $T_0 = S/\sqrt{N}$ since $\alpha = 0$ corresponds to the eigenvector 1. The following lemma follows directly from the definitions:

Lemma 4.
$$N = \sum_{\mu} (s_{\mu})^2 = \sum_{\alpha} Q_{\alpha} = \frac{1}{N} S^2 + \sum_{\alpha}' Q_{\alpha}.$$

Then we obtain the following bounds.

Theorem 1.

$$\frac{j - j_{\min}}{N} S^2 + j_{\min} N \leqslant H_0 \leqslant \frac{j - j_{\max}}{N} S^2 + j_{\max} N.$$
(27)

Proof. We rewrite the Hamilton function in the following form and conclude:

$$H_{0} = \sum_{\mu\nu\alpha\beta} J_{\mu\nu} \overline{C_{\alpha}^{\mu}} C_{\beta}^{\nu} \overline{T}_{\alpha} \cdot T_{\beta} = \sum_{\beta} j_{\beta} Q_{\beta} = \frac{J}{N} S^{2} + \sum_{\beta} 'j_{\beta} Q_{\beta}$$
$$\geqslant \frac{j}{N} S^{2} + j_{\min} \sum_{\beta} 'Q_{\beta}$$
$$= \frac{j}{N} S^{2} + j_{\min} \left(N - \frac{1}{N} S^{2} \right) = \frac{j - j_{\min}}{N} S^{2} + j_{\min} N$$
(28)

using (26), the positivity of Q_{β} , and lemma 4. The other inequality follows analogously. \Box

We note that the proof does not depend on the dimension of spin space, but the optimal energy bounds may depend on this dimension. The regular pentagon is an example of a system where the above bounding parabolas are assumed by three-dimensional states, but not by two-dimensional, i.e., coplanar states, see section 6.3.1.

Next we will show that the above-defined bounding parabolas are assumed for systems which admit coplanar S = 0 ground states. To this end we define the aforementioned umbrella construction.

Theorem 2.

(i) Let s be a coplanar weakly symmetric state with S = 0 corresponding to an eigenvalue j of \mathbb{J} and perpendicular to a unit vector **n**. Then the following family of states $\hat{s}(S), 0 \leq S \leq N$ has total spin length S:

$$\hat{s}_{\mu}(S) = \sqrt{1 - \frac{S^2}{N^2}} s_{\mu} + \frac{S}{N} n \qquad \mu = 1, \dots, N.$$
 (29)

(ii) If, moreover, s is a ground state then $\hat{s}(S)$ will be a relative ground state for all $0 \leq S \leq N$. In this case.

$$E_{\min}(S) = \frac{j - j_{\min}}{N} S^2 + j_{\min} N.$$
 (30)

(iii) The analogous case for s being a coplanar weakly symmetric relative anti-ground state for S = 0: then $\hat{s}(S)$ will be a relative anti-ground state and

$$E_{\max}(S) = \frac{j - j_{\max}}{N}S^2 + j_{\max}N.$$
(31)

Proof.

- (i) Obviously, $\hat{s}(S)$ is a state with total spin length S.
- (ii) We obtain $\sum_{\nu} J_{\mu\nu} \hat{s}_{\nu} = \sqrt{1 \frac{S^2}{N^2}} j_{\min} s_{\mu} + \frac{S}{N} j n$ and $H_0(\hat{s}) = \frac{j j_{\min}}{N} S^2 + j_{\min} N$, using $\sum_{\mu} s_{\mu} = S = \mathbf{0}$. Since $H_0(\hat{s})$ is identical with the lower bound of theorem 1, \hat{s} must be a relative ground state.
- (iii) This case is largely analogous, but j_{min} has to be replaced by j_{maxi} , since s has S = 0 and cannot be the total anti-ground state 1. \square

Systems which satisfy equation (30) will be called *parabolic* systems. Hence the essential claim of theorem 2 is that systems with a coplanar S = 0 ground state will be parabolic.

We will now prove a theorem which can be viewed as the converse of theorem 2, in so far it shows that coplanar ground states necessarily appear for parabolic systems, if certain technical assumptions are satisfied. We expect that these assumptions hold under fairly general conditions but will not dwell upon this question further.

Theorem 3. Consider a connected AF-system and let $t \mapsto s(t)$ be a smooth curve for $0 < t < \epsilon$ of relative ground states such that the limits $t \rightarrow 0$ of $\mathbf{s}(t)$ and its derivatives up to second order exist, in particular $\frac{dS}{dt}(0) \neq 0$ for $S^2(t) \equiv \left(\sum_{\mu} \mathbf{s}_{\mu}(t)\right)^2$. Moreover, let $\mathbf{s}(0)$ be a weakly symmetric ground state with S = 0 and assume parabolicity, $E_{\min}(S) = \frac{j - j_{\min}}{N}S^2 + Nj_{\min}$, at least for the interval covered by $0 < t < \epsilon$.

Then $\mathbf{s}(0)$ will be coplanar.

Proof. We will indicate the limit $t \to 0$ of a function of t by skipping the argument, e.g. $s_{\mu} \equiv s_{\mu}(0)$. Differentiation with respect to t will be indicated by an overdot. Since $s_{\mu}(t) \cdot s_{\mu}(t) = 1$, differentiation yields

$$s_{\mu} \cdot \dot{s}_{\mu} = 0 \qquad \dot{s}_{\mu} \cdot \dot{s}_{\mu} + s_{\mu} \cdot \ddot{s}_{\mu} = 0. \tag{32}$$

We may assume that S(t) = S(t)e where e is a constant unit vector. Being a weakly stationary

relative ground state, $\mathbf{s}(t)$ satisfies (18) for t > 0:

$$j_{\mu}(t) \equiv \sum_{\nu} J_{\mu\nu} s_{\nu}(t) = \kappa_{\mu}(t) s_{\mu}(t) + \chi(t) \sum_{\nu} s_{\nu}(t)$$
(33)

with the limit value $S \equiv \sum_{\nu} s_{\nu} = 0$ according to the assumptions of the theorem. But we cannot *a priori* assume that $\chi(t)$ has a finite limit value for $t \to 0$. Thus we solve (33) for $L(t) \equiv S(t)\chi(t)$ and obtain

$$L(t) = \frac{j_{\mu}(t) \cdot e - j_{\mu}(t) \cdot s_{\mu}(t)e \cdot s_{\mu}(t)}{1 - (e \cdot s_{\mu}(t))^{2}}.$$
(34)

If *s* is collinear, the proof is complete. If *s* is not collinear, we find some μ such that $(e \cdot s_{\mu}(t))^2 \neq 1$ for all $0 \leq t < \epsilon$. For this value of μ , (34) defines a smooth function of *t*, independent of μ , with a finite limit value $L \equiv L(0)$. Moreover, we can solve (33) for $\kappa_{\mu}(t), \mu = 1, \ldots, N$ and obtain

$$\kappa_{\mu}(t) = \sum_{\nu} J_{\mu\nu} s_{\nu}(t) \cdot s_{\mu}(t) - L(t) \boldsymbol{e} \cdot s_{\mu}(t).$$
(35)

This shows that also $\kappa_{\mu}(t)$ is a smooth function of t and has a finite limit for $t \to 0$. Since, by assumption, s is weakly symmetric this limit must be $\kappa_{\mu} = j_{\min}$ and hence L = 0. Therefore, the smooth function $\chi(t) = \frac{L(t)}{S(t)}$ for t > 0 has a finite limit $\chi = \lim_{t\to\infty} \frac{\dot{L}(t)}{\dot{S}(t)}$, employing l'Hospital's rule and that $\dot{S}(0) \neq 0$ by assumption.

Differentiating (33) twice and taking the limit $t \rightarrow 0$ yields

$$\sum_{\nu} J_{\mu\nu} \dot{s}_{\nu} = \dot{\kappa}_{\mu} s_{\mu} + j_{\min} \dot{s}_{\mu} + \chi \dot{S}$$
(36)

and

$$\sum_{\nu} J_{\mu\nu} \ddot{s}_{\nu} = \ddot{\kappa}_{\mu} s_{\mu} + 2\dot{\kappa}_{\mu} \dot{s}_{\mu} + j_{\min} \ddot{s}_{\mu} + 2\dot{\chi} \dot{S} + \chi \ddot{S}.$$
(37)

Multipying (33), (36) and (37) with s_{μ} and summing over μ yields

$$H = \sum_{\mu\nu} J_{\mu\nu} s_{\mu} \cdot s_{\nu} = \sum_{\mu} \kappa_{\mu} + \chi S^2 = N j_{\min}$$
(38)

$$\frac{1}{2}\dot{H} = \sum_{\mu\nu} J_{\mu\nu} \dot{s}_{\nu} \cdot s_{\mu} = \sum_{\mu} \dot{\kappa}_{\mu} = \sum_{\nu} j_{\min} s_{\nu} \cdot \dot{s}_{\nu} = 0$$
(39)

$$\frac{1}{2}\ddot{H} = \sum_{\mu\nu} J_{\mu\nu} (\dot{s}_{\nu} \cdot \dot{s}_{\mu} + s_{\mu} \cdot \ddot{s}_{\nu}) \tag{40}$$

$$= j_{\min} \sum_{\mu} \dot{s}_{\mu}^{2} + \chi \dot{S}^{2} + \sum_{\mu} \ddot{\kappa}_{\mu} + j_{\min} \sum_{\mu} \ddot{s}_{\mu} \cdot s_{\mu}$$
(41)

$$=\sum_{\mu}\ddot{\kappa}_{\mu} + \chi \dot{S}^2 \tag{42}$$

using (32) and (36). On the other hand,

$$H(t) = \sum_{\mu} \kappa_{\mu}(t) + \chi(t)S^{2}(t)$$
(43)

$$\dot{H}(t) = \sum_{\mu} \dot{\kappa}_{\mu}(t) + \dot{\chi}(t)S^{2}(t) + 2\chi(t)S(t)\dot{S}(t)$$
(44)

and

$$\ddot{H} = \sum_{\mu} \ddot{\kappa}_{\mu} + 2\chi \dot{S}^2. \tag{45}$$

Comparison with (42) yields $\sum_{\mu} \ddot{\kappa}_{\mu} = 0$ and

$$\dot{H} = 2\chi \dot{S}^2. \tag{46}$$

Hence, using the assumption of parabolicity and l'Hospital's rule,

$$\frac{j - j_{\min}}{N} = \left. \frac{\mathrm{d}H}{\mathrm{d}(S^2)} \right|_{t=0} = \frac{\dot{H}}{\frac{\mathrm{d}^2}{\mathrm{d}t^2}S^2} = \frac{2\chi\dot{S}^2}{2\dot{S}^2} = \chi.$$
(47)

Summing (36) over μ yields

$$\dot{s}\dot{S} = \sum_{\mu} \dot{\kappa}_{\mu} s_{\mu} + j_{\min} \dot{S} + N\chi \dot{S}$$
(48)

whence, by (47) and the assumption $\dot{S} \neq 0$,

$$\sum_{\mu} \dot{\kappa}_{\mu} s_{\mu} = 0. \tag{49}$$

We rewrite (36) in the form

$$\sum_{\nu} (J_{\mu\nu} - j_{\min}\delta_{\mu\nu} - N\chi E_{\mu\nu})\dot{s}_{\nu} = \dot{\kappa}_{\mu}s_{\mu}$$
(50)

where *E* is the projector onto J's eigenvector **1** with constant entries $E_{\mu\nu} = \frac{1}{N}$. We expand both sides of (50) into real eigenvectors of J and obtain

$$\dot{\kappa}_{\mu}s_{\mu} = \sum_{\alpha}{}^{\prime}K_{\alpha}C_{\alpha}^{\mu} \tag{51}$$

$$\dot{s}_{\nu} = \sum_{\beta} \frac{1}{j_{\beta} - j} K_{\beta} C_{\beta}^{\mu} + \frac{1}{N} \dot{S}.$$
(52)

Now $\sum_{\mu} \dot{\kappa}_{\mu} s_{\mu} \cdot \dot{s}_{\mu} = 0$, $\frac{1}{N} \dot{S} \cdot \sum_{\mu} \dot{\kappa}_{\mu} s_{\mu} = 0$ and $\sum_{\mu} C^{\mu}_{\alpha} C^{\mu}_{\beta} = \delta_{\alpha\beta}$ imply

$$0 = \sum_{\mu} \sum_{\alpha} {}^{\prime} K_{\alpha} C^{\mu}_{\alpha} \cdot \sum_{\beta} {}^{\prime} \frac{1}{j_{\beta} - j} K_{\beta} C^{\mu}_{\beta} = \sum_{\alpha} {}^{\prime} \frac{K^{2}_{\alpha}}{j_{\alpha} - j}.$$
(53)

Since $j > j_{\alpha}$ for $\alpha > 0$ and connected AF-systems, this is only possible if all $K_{\alpha} = \mathbf{0}$. Hence, by (52) $\dot{s}_{\nu} = \frac{1}{N}\dot{S}$ and $s_{\nu} \cdot \frac{1}{N}\dot{S} = 0$. Thus *s* is coplanar.

An analogous theorem concerning $E_{\max}(S)$ and coplanar anti-ground states can be proved similarly.

4. Extension of ground states

We consider spin systems with $J_{\mu\nu} = 0$ or J and which can thus essentially be characterized by their graph $\Gamma = (\mathcal{V}, \mathcal{E})$. A sub-graph (representing a sub-system) $\widetilde{\Gamma} = (\widetilde{\mathcal{V}}, \widetilde{\mathcal{E}})$ is defined by the conditions $\widetilde{\mathcal{V}} \subset \mathcal{V}, \widetilde{\mathcal{E}} \subset \mathcal{E}$ and $\widetilde{\mathcal{E}} \subset [\widetilde{\mathcal{V}}]^2$. Let

$$\widetilde{H}_0 = 2J \sum_{(\mu,\nu)\in\widetilde{\mathcal{E}}} s_\mu \cdot s_\nu \tag{54}$$

be the Hamiltonian of the subsystem. Let s be any state of Γ and \tilde{s} be its restriction to $\tilde{\Gamma}$. If s happens to be a ground state of Γ , \tilde{s} need not be a ground state of $\tilde{\Gamma}$, but, of course, $\tilde{H}_0(\tilde{s}) \geq \tilde{E}_{\min}$.

The inverse problem is the extension problem. Given a ground state \tilde{s} of $\tilde{\Gamma}$, can it be extended to a ground state s of Γ (such that \tilde{s} is the restriction of s)? In general, this will

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not be possible. For example, if $\tilde{\Gamma}$ is the graph of a dimer with ground state $\tilde{s} = \uparrow \downarrow$, the extension of \tilde{s} to a global ground state is only possible for bi-partite systems. We have no general theory describing the obstacles against extension (which could be called frustration of higher order), but we can describe the typical situation where extension is possible. For an analogous treatment in the quantum case, see [14].

The situation we have in mind is one where the graph of the spin system Γ_1 is decomposed into smaller graphs which are copies of some graph Γ_2 in such a way that each edge lies in exactly one of the copies. For example, an icosidodecahedron can be decomposed into 20 triangles with disjoint edges. We have local states by considering the ground states of the copies of Γ_2 . But these local states must fit together in order make it possible to construct a global ground state of Γ_1 . For the example of decomposition into triangles the condition of fitting-together of local ground states is equivalent to the 3-colourability of the large system. The icosidodecahedron happens to satisfy this condition and hence possesses a ground state of the energy which is 20 times the ground state energy of the triangle. For general Γ_2 the colourability condition has to be replaced by the existence of a map from Γ_1 to Γ_2 which maps edges onto edges. Such a map may be called a 'graph homomorphism'.

A graph homomorphism $h : \Gamma_1 \longrightarrow \Gamma_2$ is a pair of maps $h = (h_{\mathcal{V}}, h_{\mathcal{E}})$ such that $h_{\mathcal{V}} : \mathcal{V}_1 \longrightarrow \mathcal{V}_2, h_{\mathcal{E}} : \mathcal{E}_1 \longrightarrow \mathcal{E}_2$ and $\{\mu, \nu\} \in \mathcal{E}_1$ implies $h_{\mathcal{E}}\{\mu, \nu\} = \{h_{\mathcal{V}}(\mu), h_{\mathcal{V}}(\nu)\}$. The definition of a graph isomomorphism is analogous.

Proposition 1. Consider two spin systems with graphs Γ_1 and Γ_2 and the following properties:

- (*i*) there exists a graph homomorphism $h : \Gamma_1 \longrightarrow \Gamma_2$,
- (ii) there exists a decomposition into sub-graphs $\Gamma_1 = \bigcup_{\mu=1}^k \Gamma_\mu$, which is a disjoint decomposition with respect to edges,
- (iii) the restriction of h to Γ_{μ} , $h_{\mu} : \Gamma_{\mu} \longrightarrow \Gamma_2$ is a graph isomorphism for each $\mu = 1, ..., k$.

Moreover, let $s^{(2)} : \mathcal{V}_2 \longrightarrow S^2$ be a ground state of Γ_2 . Then $s^{(1)} \equiv h_{\mathcal{V}} \circ s^{(2)}$ will be a ground state of Γ_1 .

Proof. According to the disjoint decomposition of \mathcal{E}_1 , H_1 will be a sum of terms H_μ which are each minimized by the state $s^{(1)}$. Hence $s^{(1)}$ is a ground state of Γ_1 .

The above-mentioned construction of ground states for the icosidodecahedron, the octahedron and the cuboctahedron follows the description given in this proposition. Other examples show that the two given conditions of the proposition are essential. The dodecahedron is 3-colourable but it cannot be decomposed into edge-disjoint triangles. Actually its ground state energy is lower than that of the state obtained by the 3-colouring. A simpler example is the bi-partite 3-chain which can be mapped by a graph homomorphism onto the triangle but is not isomorphic to the triangle. On the other hand it is possible to connect four triangles such that their edges are still disjoint but form a tetrahedron which cannot be 3-coloured. This is then a 'higher order frustrated' system with a ground state energy which is larger than four times the ground state energy of a triangle.

5. Symmetric states

In this section, we will assume suitable symmetry properties of the spin system and correspondingly consider the notion of 'symmetric states'. We make use of some simple concepts and results of group theory, which may be found in many textbooks, e.g. [16].

Let $J_{\mu\nu} = 0$ or J and consider the graph $\Gamma = (\mathcal{V}, \mathcal{E})$ characterizing the spin system. Further let \mathcal{G} denote the symmetry group of the graph Γ . Hence \mathcal{G} consists of all permutations of \mathcal{V} which map edges onto edges. \mathcal{G} has a 'natural' representation by real $N \times N$ -matrices, by permuting the standard basis of \mathbb{R}^N in the same way as the spin sites $\mu \in \mathcal{V}$. We will denote the $N \times N$ -matrix representing any symmetry $g \in \mathcal{G}$ by \mathbf{g} and by \mathbb{G} the set of all $N \times N$ -matrices obtained in this way. Obviously,

Lemma 5. The adjacency matrix \mathbb{J} commutes with all $\mathbf{g} \in \mathbb{G}$.

Hence

Lemma 6. The eigenspaces of \mathbb{J} split into orthogonal direct sums of irreducible subrepresentations of the natural representation of \mathcal{G} .

The irreducible representations of the relevant groups are well known, see e.g. [16], and can easily be associated with the different (subspaces of) eigenspaces of J.

For specific spin systems, e.g. magnetic molecules, the spin sites $\mu \in \mathcal{V}$ are embedded into the physical 3-space and the group \mathcal{G} could also be identified with one of the point groups, i.e. finite subgroups of $\mathcal{O}(3, \mathbb{R})$. There are only a finite number of possibilities, see e.g. [16], 3.1.2. and 3.1.3. The most interesting cases are

- the dihedral group \mathcal{D}_n ,
- the tetrahedral group \mathcal{T} ,
- the octahedral group \mathcal{O} ,
- the icosahedral group \mathcal{Y} ,

as well as the improper point groups attached to them.

In what follows, we make the following general assumption:

Assumption 1. All exchange parameters $J_{\mu\nu}$ are 0 or J, and the group G operates transitively on V.

The latter means that for each pair μ , $\nu \in \mathcal{V}$ there is a permutation $g \in \mathcal{G}$ such that $g(\mu) = \nu$. Hence the above assumption expresses the equivalence of all spin sites with respect to their interaction.

A state *s* can be transformed in two different ways: either the unit vectors s_{μ} can be permuted according to a symmetry $g \in \mathcal{G}$, or the whole configuration *s* can be rotated/reflected in spin space. Mathematically, this means that the two groups \mathcal{G} and $\mathcal{O}(3, \mathbb{R})$ operate independently on the set of states. If two matrices $\mathbf{g} \in \mathbb{G}$ and $R \in \mathcal{O}(3, \mathbb{R})$ are given, we define their action on a state *s* by matrix multiplication with *s* which is again considered as an $N \times 3$ -matrix:

Definition 2. $(R, \mathbf{g}) \bullet s \equiv Rs\mathbf{g}^{-1}$.

The transformations $R \in \mathcal{O}(3, \mathbb{R})$ will also be referred to as 'rotations/reflections in spin space'.

We have the following obvious results:

Lemma 7. Let $\mathbf{g} \in \mathbb{G}$ and $R \in \mathcal{O}(3, \mathbb{R})$ and s be a state. Then

(i) s and Rsg^{-1} have the same total spin length S,

(ii) If s is (weakly) stationary, then Rsg^{-1} is (weakly) stationary,

(iii) If s is weakly symmetric, then Rsg^{-1} is weakly symmetric,

(iv) if s is a (anti-) ground state, then Rsg^{-1} is a (anti-) ground state.

Having defined the action of the product group $\mathcal{O}(3, \mathbb{R}) \times \mathbb{G}$ on states, it is natural to consider the corresponding subgroup leaving a given state invariant:

Definition 3. For any state s let $\mathcal{G}_s \equiv \{(R, \mathbf{g}) \in \mathcal{O}(3, \mathbb{R}) \times \mathbb{G} | Rs\mathbf{g}^{-1} = s\}$. \mathcal{G}_s is called the symmetry group of s.

Definition 4. A state *s* is called symmetric if the projection onto the second factor π_2 : $\mathcal{G}_s \longrightarrow \mathbb{G}$ is surjective, i.e. if for each $\mathbf{g} \in \mathbb{G}$ there exists a rotation/reflection $R \in \mathcal{O}(3, \mathbb{R})$ such that $R^{-1}s\mathbf{g} = s$.

One may thus say that for symmetric states s every permutation $\mathbf{g} \in \mathbb{G}$ of the vectors s_{μ} can be compensated by a suitable rotation/reflection in spin space.

For example, if $(\mathcal{V}, \mathcal{E})$ is the graph of a regular *N*-polygon, the coplanar symmetric states are in 1:1-correspondence to the roots of unity, $z_n^N = 1$, namely

$$s_{\mu} = z_n^{\mu} = \exp(i\mu n2\pi/N)$$
 $n, \mu = 1, \dots, N.$ (55)

Here $\mathcal{G} = \mathcal{D}_N$, the dihedral group of order 2*N*.

We have the following:

Lemma 8. Let $s_{(\mu)}^{(i)}$ be a non-vanishing $N \times 3$ -matrix such that for all $g \in \mathbb{G}$ there exists some $R \in \mathcal{O}(3, \mathbb{R})$ such that sg = Rs. Then, for some suitable $\lambda \in \mathbb{R}$, λs will be a symmetric state.

Proof. It remains to show that $\lambda s_{\mu}^{(i)}$ will be unit vectors for all $\mu = 1, ..., N$. Fix some indices μ, ν . Using assumption 1, we choose a $g \in \mathbb{G}$ such that $(sg)_{\mu} = s_{\nu} = Rs_{\mu}$. Since $R \in \mathcal{O}(3, \mathbb{R})$ the two vectors s_{ν} and s_{μ} have the same length. Because μ, ν were arbitrarily chosen, all vectors s_{μ} have the same length, say λ^{-1} , which completes the proof. \Box

If a state *s* is not collinear, then Rs = s implies R = 1. In this case for each $\mathbf{g} \in \pi_2(\mathcal{G}_s)$ there exists a unique $R \in \mathcal{O}(3, \mathbb{R})$ such that $R^{-1}s\mathbf{g} = s$. We will write $R = \rho_s(\mathbf{g})$. It is easily shown that the map $\mathbf{g} \mapsto \rho_s(\mathbf{g})$ is a linear representation of the subgroup $\pi_2(\mathcal{G}_s) \subset \mathbb{G}$. If *s* is collinear we may uniquely fix *R* by choosing $R \in \{1, -1\}$ and thus obtain a representation ρ_s also in this case.

For the case of symmetric *s* we have $\pi_2(\mathcal{G}_s) = \mathbb{G}$ and hence ρ_s will be an *n*-dimensional representation of \mathbb{G} with $n \leq 3$.

In the case of a symmetric ground state it is thus necessary that the eigenspace of \mathbb{J} corresponding to the smallest eigenvalue j_{\min} contains an irreducible representation of \mathbb{G} of dimension $n \leq 3$. If it does not, we can exclude symmetric ground states.

For practical purposes of constructing (anti-)ground states it would be desirable to invert the process and to reconstruct *s* from a given *n*-dimensional subrepresentation ρ of the natural representation of $\mathcal{G}, n \leq 3$. We will now describe this procedure and consider, for sake of simplicity, only the case n = 3. Let \mathcal{S} be a three-dimensional subspace of \mathbb{R}^N left invariant by all $g \in \mathbb{G}$ and $s_{(\mu)}^i, i = 1, 2, 3$ a basis of \mathcal{S} . It follows that the three vectors $\sum_{\mu} g_{\mu(\nu)} s_{\mu}^i, i = 1, 2, 3$ are linear combinations of the basis vectors, hence

$$\sum_{\mu} g_{\mu\nu} s^i_{\mu} = \sum_j \rho^i_j(g) s^j_{\nu} \tag{56}$$

or, in matrix notation, $\rho(g)s = sg$. Being a real representation of a finite group, ρ is equivalent to an orthogonal representation, i.e. $\tilde{\rho}(g) \equiv T\rho(g)T^{-1}$ is orthogonal with some invertible 3×3 -matrix *T*. It follows that

$$\widetilde{\rho}(g)(Ts) = (Ts)g \tag{57}$$

which means that Ts is a symmetric state with $\rho_{Ts} = \tilde{\rho}$. Note that the orthogonality of the representation is crucial in order to obtain a row $(Ts)^{(i)}_{\mu}\mu = 1, \ldots, N$ of unit vectors, compare lemma 7. Hence we have proved the following.

Proposition 2. Let ρ be a three-dimensional subrepresentation of the natural representation of G. Then there exists a three-dimensional symmetrical state s such that ρ_s is equivalent to ρ .

Further there holds

Proposition 3. Each stationary symmetric state *s* is weakly symmetric.

Proof. We will use a matrix notation and write $\kappa \equiv \operatorname{diag}(\kappa_1, \ldots, \kappa_N)$. Then condition (16) for stationarity may be written as $s\mathbb{J} = s\kappa$. For $\mathbf{g} \in \mathbb{G}$ we have $\widetilde{\kappa} \equiv \mathbf{g}^{-1}\kappa\mathbf{g} = \operatorname{diag}(\kappa_{\mathbf{g}(1)}, \ldots, \kappa_{\mathbf{g}(N)})$. We conclude $s\kappa\mathbf{g} = s\mathbb{J}\mathbf{g} = s\mathbf{g}\mathbb{J} = \rho_s(\mathbf{g})s\mathbb{J} = \rho_s(\mathbf{g})s\kappa = s\mathbf{g}(\mathbf{g}^{-1}\kappa\mathbf{g}) = \rho_s(\mathbf{g})s\tilde{\kappa}$. Hence $s\kappa = s\tilde{\kappa}$, which is in components $\kappa_{\mu}s_{\mu} = \kappa_{\mathbf{g}(\mu)}s_{\mu}$ whence κ_{μ} is independent of μ (using assumption 1).

The converse of proposition 3 is not true. There are ground states of the icosidodecahedron which are weakly symmetric, but not symmetric, see section 6.5. The condition of stationarity in proposition 3 can be replaced by another condition.

Definition 5. Let \mathcal{G}_{μ} denote the subgroup of \mathcal{G} leaving $\mu \in \mathcal{V}$ fixed and $\mathbb{G}(\mu)$ the matrix group generated by its natural representation. A symmetric state *s* will be called isotropic if $\rho_s(\mathbb{G}(\mu))$ contains at least one rotation with an angle $\alpha \neq 0, \pi$.

Lemma 9. If s is isotropic and $\rho_s = \rho_{s'}$ then $s_{\mu} = \pm s'_{\mu}$.

Proof. s_{μ} is invariant under all rotations/reflections $\rho_s(\mathbf{g}), \mathbf{g} \in \mathbb{G}(\mu)$. Since some of these is a rotation with an angle $\alpha \neq 0, \pi$, the axis of rotation will be unique and $\rho_s = \rho_{s'}$ implies $s_{\mu} = \pm s'_{\mu}$.

Proposition 4. Each isotropic state is stationary, hence weakly symmetric.

Proof. Using the above matrix notation we obtain $s \mathbb{J}\mathbf{g} = s\mathbf{g}\mathbb{J} = \rho_s(\mathbf{g})s\mathbb{J}$ hence $s\mathbb{J}$ is a symmetric (not normalized) state with $\rho_{s\mathbb{J}} = \rho_s$. Using that *s* is isotropic and lemma 9 we conclude $(s\mathbb{J})_{\mu} = \pm \lambda s_{\mu}$, hence *s* is stationary, and, by proposition 3, weakly symmetric. \Box

6. Examples

We will mainly consider AF-systems and set J > 0 throughout this section if not mentioned otherwise.

6.1. The general spin triangle

We consider spin systems with N = 3 and general coupling coefficients J_1 , J_2 and J_3 . The special symmetric case $J_1 = J_2 = J_3$ is atypical and its properties are discussed later. Another special case is the 3-chain with $J_1 = 0$, $J_2 = J_3$, which is probably the simplest example of a non-parabolic system. Its investigation will be left for the reader.

Let s_{ν} , $\nu = 1, 2, 3$, be any spin configuration and consider the 3×3 -matrix S with coefficients

$$S_{\mu\nu} = \mathbf{s}_{\mu} \cdot \mathbf{s}_{\nu}. \tag{58}$$

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Figure 1. The three-dimensional convex set \mathcal{P} defined in (59) as seen from the view point (4, -2, -2). The shape of its projection onto the plane perpendicular to (4, -2, -2) is essentially identical with that shown in figure 2.

It is positive and its diagonal consists of 1's. Conversely, any 3×3 -matrix with these properties is of the form (58). This follows from the spectral theorem. These matrices can be written in the form

$$S = \begin{pmatrix} 1 & w & v \\ w & 1 & u \\ v & u & 1 \end{pmatrix} \equiv [u, v, w]$$
(59)

where u, v, w are real numbers subject to the constraint

det
$$S = 1 - (u^2 + v^2 + w^2) + 2uvw \ge 0$$
 and $u^2, v^2, w^2 \le 1$. (60)

Hence the set \mathcal{P} of the matrices of the form (58) can be considered as a compact convex subset of \mathbb{R}^3 . It has the form of an 'inflated tetrahedron', see figure 1. Since the map

$$\pi: \mathcal{P} \longrightarrow \mathbb{R}^2 \tag{61}$$

$$\pi[u, v, w] \equiv (3 + 2(u + v + w), J_1u + J_2v + J_3w)$$
(62)

$$= \left(S^2(\mathbf{s}), \frac{1}{2}H_0(\mathbf{s})\right) \tag{63}$$

is affine, the set $\mathcal{P}_2 \equiv \pi[\mathcal{P}]$ of all possible values of the energy versus square of the total spin will be a compact convex set too. To simplify the geometry we will assume $J_1 + J_2 + J_3 = 0$. Then \mathcal{P}_2 can be considered as the orthogonal projection of \mathcal{P} onto a suitable plane. If we introduce new coordinates in the u, v, w-space by

$$\sigma = u + v + w \tag{64}$$

$$\epsilon = J_1 u + J_2 v + J_3 w \tag{65}$$

$$\tau = (J_2 - J_3)u + (J_3 - J_1)v + (J_1 - J_2)w$$
(66)

the orthogonal projection π is essentially the projection onto the σ , ϵ -plane. Hence the



Figure 2. Energy ϵ versus square of total spin, S^2 , for the triangle with $J_1 = 0$, $J_2 = 1$, $J_3 = -1$ according to (70). The broken lines represent the bounding parabolas of theorem 1.

boundary of \mathcal{P}_2 can be obtained by solving the two equations

$$\det \mathcal{S} = 0 \tag{67}$$

$$\nabla \det S \cdot \begin{pmatrix} J_2 - J_3 \\ J_3 - J_1 \\ J_1 - J_2 \end{pmatrix} = 0.$$
 (68)

The first equation (67) defines the boundary of \mathcal{P} , the second one (68) is satisfied by those points of the boundary of \mathcal{P} which are orthogonally projected onto the boundary of \mathcal{P}_2 . By using computer algebra software it is straightforward to express (67) and (68) as equations for σ , ϵ , τ and to eliminate τ between (67) and (68). Yet the result for the general case is too complex to be reproduced here. We choose the special case

$$J_1 = 0 \qquad J_2 = 1 \qquad J_3 = -1 \tag{69}$$

and obtain the following equation for the boundary of \mathcal{P}_2

$$0 = S^{2}(9 - 10S^{2} + S^{4})^{2} + (-27 + 288S^{2} + 18S^{4} - 24S^{6} + S^{8})\epsilon^{2} - 8(3 + S^{2})^{2}\epsilon^{4} + 16\epsilon^{6}$$
(70)

where we have re-substituted $\frac{S^2-3}{2}$ for σ . The actual boundary of \mathcal{P}_2 is only a part of the family of curves defined by (70) and is shown in figure 2. Here we have an example where the curves $E_{\min}(S)$ and $E_{\max}(S)$ can be explicitly calculated and are not parabolas, although they are close to their bounding parabolas which are represented as broken lines in figure 2. This is not in contradiction to theorem 2 since the ground state of this system is coplanar but has S > 0.

We expect that similar figures also appear for the general system of N spins. Parabolic systems are probably rare cases, some of which will be considered in the remaining subsections.

6.2. S_N -symmetric systems

Recall that the *N*-pantahedron was defined as a system of *N* spins where each pair (μ, ν) interacts with equal strength J > 0. These systems serve mainly as exactly solvable model examples for the sake of illustration, but the dimer, the triangle and the tetrahedron are also of experimental relevance [17–19]. The symmetry group of the *N*-pantahedron is obviously the group S_N of all permutations of *N* spin sites.

By completing squares the Hamiltonian can be written as

$$H_0 = J(S^2 - N). (71)$$

Since the energy depends only on S, each state is a relative (anti-)ground state with

$$E_{\min}(S) = E_{\max}(S) = J(S^2 - N).$$
(72)

For S = 0 there exist symmetrical ground states only if $N \leq 4$. These are the obvious dimer, triangle and tetrahedron configurations in spin space. The dimension *dim* of the irreducible representations of S_N is given by the number of standard Young tableaus (see [16]) and hence dim = 1 or $dim \ge N - 1$. The one-dimensional representations are either the trivial one corresponding to the total anti-ground state or the one attaching to each permutation its sign, which is not generated by a symmetric state. Thus there are no symmetrical ground states for N > 4. For example, a cyclic permutation of three spins leaving two other spins invariant cannot be compensated by a reflection/rotation in spin space.

The coplanar *N*-polygon configuration in spin space does however define a weakly symmetric ground state with $\kappa_{\mu} = -1$. Hence, the construction of theorem 1 yields symmetrical relative ground states $\hat{s}(S)$. We will check our equation (30) for this case. Let \mathbb{E} denote the $N \times N$ -matrix with $E_{\mu\nu} = 1/N$ for all μ , ν . It is the matrix of the projector onto the one-dimensional subspace spanned by the vector $\mathbf{1} = \frac{1}{\sqrt{N}}(1, 1, \dots, 1)$. The adjacency matrix for the *N*-pantahedron is $\mathbb{J} = J(N\mathbb{E}-\mathbf{1})$. Hence its eigenvalues are j = J(N-1) with eigenvector $\mathbf{1}$ and $j_{\min} = -J$ with the (N-1)-dimensional eigenspace of vectors orthogonal to $\mathbf{1}$. Thus $j_{\min} = j_{\max i}$ and the two functions $E_{\min}(S)$ and $E_{\max}(S)$ coincide in accordance with what has been said above.

6.3. D_N -symmetric systems

6.3.1. The N-polygon. The N-polygon has the dihedral group D_N as symmetry group. We have already mentioned the set of S = 0, stationary, coplanar, symmetric states s given by

$$s_{\mu} = z_n^{\mu} \qquad \mu = 0, \dots, N-1$$
 (73)

where $z_n = \exp(in2\pi/N)$, n = 1, ..., N, is a root of unity, $z_n^N = 1$. Indeed, these states are the complex eigenvectors of the corresponding adjacency matrix

$$\mathbb{J} = J \begin{pmatrix} 0 & 1 & 0 & \dots & \dots & 0 & 1 \\ 1 & 0 & 1 & 0 & \dots & \dots & 0 \\ 0 & 1 & 0 & 1 & 0 & \dots & 0 \\ \vdots & & & & & & \\ 0 & \dots & \dots & 0 & 1 & 0 & 1 \\ 1 & 0 & \dots & \dots & 0 & 1 & 0 \end{pmatrix}.$$
(74)

The eigenvalues are obtained by

$$(\mathbb{J}s)_{\mu} = J(z_n^{\mu+1} + z_n^{\mu-1}) = 2J\cos(2\pi n/N)s_{\mu}$$
(75)

and the corresponding energies are

$$E_n = 2JN\cos(2\pi n/N)$$
 $n = 0, ..., N - 1.$ (76)

It follows that

$$E_{\min} = N j_{\min} = \begin{cases} E_{N/2} = -2JN & \text{for } N \text{ even} \\ E_{(N-1)/2} = 2JN \cos(\pi (N-1)/N) & \text{for } N \text{ odd} \end{cases}$$
(77)

and

$$E_{\max} = E_0 = Nj = 2NJ.$$
 (78)

This follows since the bounds of lemma 1 are attained by these states. For N = 5 the ground state of the pentagon may be visualized as a pentagram in spin space and analogously for other odd N.



Figure 3. The shaded region represents the spectrum E versus S^2 of coplanar states of the pentagon as determined by numerical methods. The straight lines are the boundaries of the full spectrum. Also a part of the boundary of the coplanar spectrum which can be analytically calculated is displayed.

Moreover,

$$E_{\max}(0) = E_1 = N j_{\max} = 2N J \cos(2\pi/N).$$
(79)

The relative ground states of theorem 1 have energies

$$E_{\min}(S) = \begin{cases} 2J(2\frac{S^2}{N} - N) & \text{for } N \text{ even} \\ 2J((1 - \cos(\pi \frac{N-1}{N}))\frac{S^2}{N} + N\cos(\pi \frac{N-1}{N})) & \text{for } N \text{ odd} \end{cases}$$
(80)

and

$$E_{\max}(S) = 2J\left((1 - \cos(2\pi/N)\frac{S^2}{N} + N\cos(2\pi/N)\right).$$
(81)

Both parabolas meet in the anti-ground state with $E_{\text{max}} = 2JN$ for S = N.

6.3.2. Coplanar states of the pentagon. The regular pentagon as a special case dealt with in the preceding subsection assumes its bounding parabolas. However, the extremal energies $E_{\min}(S)$ and $E_{\max}(S)$ can only be realized by non-coplanar states, except for S = 0 and S = 5. This has not been rigorously proved but shown by numerical evidence, see figure 3. The spectrum *E* versus S^2 realized by coplanar states is a subset of the full spectrum with concave boundaries $E_{\min}^{coplanar}(S)$ and $E_{\max}^{coplanar}(S)$.

The permutation of spin sites $(0 \leftrightarrow 0, 1 \rightarrow 2 \rightarrow 4 \rightarrow 3 \rightarrow 1)$ leaves the Hamiltonian invariant and maps relative ground states onto anti-ground states, hence $E_{\min}^{coplanar}(S) = S^2 - 5 - E_{\max}^{coplanar}(S)$. The (anti-) ground state with S = 1 and E = -7 (resp. E = 3) is especially remarkable: its state vectors s_{μ} occupy the vertices of a regular hexagon leaving one vertex free. The anti-ground state $(E_{\max}^{coplanar}(1) = 3)$ satisfies equation (18) with $\chi = 0$, hence it is a stationary state, not only weakly stationary as other relative anti-ground states. Actually, $E_{\max}^{coplanar}$ has a local minimum at S = 1, see figure 3. For axisymmetric weakly stationary states in the neighbourhood of the hexagonal state, equation (18) can be solved analytically by using computer algebra software and hence E and S^2 can be expressed as functions of a common parameter, although of forbidding complexity. The resulting curve turns out to be a part of the upper boundary of the coplanar spectrum and is displayed in figure 3, as well as the corresponding lower boundary part.

6.3.3. The pentagonal star. We obtain the pentagonal star (N = 6) by joining the five vertices of the pentagon with its midpoint, see figure 4. This system has D_5 symmetry, but it



Figure 4. Pentagonal star as an example of a spin system without coplanar ground states.



Figure 5. Three-dimensional ground state of the pentagonal star: The vertices correspond to six unit vectors in spin space and the edges correspond to the ten bonds of the pentagonal star of figure 4, see the remarks after (10).

is not \mathcal{D}_5 -symmetric in our sense, since \mathcal{D}_5 does not operate transitively on the six spin sites. Nevertheless, we will discuss this example since it has the interesting property that its ground states are not coplanar and $E_{\min}(S)$, $E_{\max}(S)$ are only piecewise parabolic.

If we add to the ten edges of the pentagonal star the five edges of the corresponding pentagram we obtain the 15 edges of the 6-pantahedron. Hence its Hamiltonian reads

$$H = S^2 - 6 - H_5 \tag{82}$$

where H_5 is the Hamiltonian of the pentagram, which is the same as that for the pentagon, up to a suitable permutation of the spin sites. This shows that the configurations minimizing H for a given S are exactly those which maximize H_5 , analogously for $E_{\max}(S)$. The maximal values for H_5 are given by the parabola, compare (81),

$$E_{\max}^{(5)}(S_5) = \frac{2(1 - \cos(2\pi/5))}{5}S_5^2 + 10\cos(2\pi/5)$$
(83)

where S_5 is the length of the total spin of the five vertices of the pentagon. Here S_5 has to be chosen maximal for given S, which means that $S_5 = S + 1$ for $0 \le S \le 4$ and $S_5 = 5$ for $4 \le S \le 6$. This yields

$$E_{\min}(S) = \begin{cases} \frac{5+\sqrt{5}}{10}(S-4)(S+1+\sqrt{5}) & \text{if } 0 \leqslant S \leqslant 4\\ S^2 - 16 & \text{if } 4 \leqslant S \leqslant 6. \end{cases}$$
(84)

Hence the total ground state is attained not for S = 0, but for $S = \frac{3-\sqrt{5}}{2} = 0.381966...$ with $E_{\min} = -5 - 2\sqrt{5} = -9.47214...$ The corresponding spin configuration is shown in figure 5.

It remains to show that no other, coplanar ground state exists. Employing the numerical result of the preceding subsection for the pentagon, $E_{\max}^{\text{coplanar}}(S) < E_{\max}(S)$ for 0 < S < 5, it is easily shown that all coplanar states with $S = \frac{3-\sqrt{5}}{2}$ have an energy larger than E_{\min} . The minimal energy for all coplanar states appears to be $E_{\min}^{\text{coplanar}} = -9$, realized by a hexagonal ground state with S = 0, but we do not yet have a rigorous proof for this claim.

Analogously one can show that

$$E_{\max}(S) = \begin{cases} \frac{5-\sqrt{5}}{10}(S+4)(S-1+\sqrt{5}) & \text{if } 1 \leq S \leq 6\\ S^2 - 16 + \frac{5(1+\sqrt{5})}{2} & \text{if } 0 \leq S \leq 1. \end{cases}$$
(85)

6.4. O-symmetric systems

6.4.1. The cube. As already mentioned the cube allows a bi-partition and hence possesses a ground state of the form

$$s_{\mu} = (-1)^{\mu} e \qquad \mu = 1, \dots, 8$$
 (86)

if the vertices are suitably labelled. From this we obtain relative ground states with

$$E_{\min}(S) = J\left(\frac{3}{4}S^2 - 24\right) \tag{87}$$

since j = 3J, $j_{min} = -3J$. Apart from j and j_{min} the other eigenvalues of \mathbb{J} are $\pm J$ with threefold degeneracy, respectively. The corresponding eigenspaces carry two inequivalent irreducible representations of \mathcal{O} , called F_1 , F_2 , see [16]. According to the considerations in section 5 we may conjecture that these two representations are generated by stationary, symmetric, non-coplanar states.

Indeed, this can be directly verified for the following states: let r_{μ} , $\mu = 1, ..., 8$ denote the unit vectors pointing to the vertices of the cube. The states

$$s_{\mu}^{(1)} \equiv r_{\mu} \qquad \mu = 1, \dots, 8$$
 (88)

and

$$s_{\mu}^{(-1)} \equiv (-1)^{\mu} r_{\mu} \qquad \mu = 1, \dots, 8$$
(89)

have the desired properties. They are easily visualized: $s^{(1)}$ is just the 'cube in spin space' and $s^{(-1)}$ the tetrahedron where each of the four distinct spin vectors is attached to pairs of vertices of the cube connected by space diagonals. $s^{(1)}$ is the anti-ground state for S = 0 with energy $E_{\text{max}}(0) = 8J$. Since it is not coplanar, theorem 1 is not directly applicable. However, it is possible to project the cube $s^{(1)}$ onto a square and, in a second step, to enlarge the square to a square of unit vectors. Since these are linear transformations, the resulting coplanar state $s^{(1)'}$ is weakly symmetric and has the same energy as before, namely $8J = E_{\text{max}}(0)$. Now theorem 1 yields $E_{\text{max}}(S) = J(\frac{1}{4}S^2 + 8)$.

6.4.2. The octahedron. As noted in [9], the octahedron can be decomposed into four triangles with disjoint edges and it is 3-colourable. Hence it has coplanar ground states with S = 0 which are obtained by extensions of the local ground states of the triangles. Since j = 4J and $j_{min} = -2J$ we again obtain the RBS-parabola

$$E_{\min}(S) = J(S^2 - 12). \tag{90}$$

The eigenspace of \mathbb{J} corresponding to j_{\min} is two dimensional and carries a real, irreducible representation of \mathcal{O} . This corresponds to the one-dimensional complex eigenspace of \mathbb{J}



Figure 6. The cuboctahedron (shaded figure) is obtained by joining the midpoints of the cube's edges with their nearest neighbours.

spanned by the vector

$$s_{\mu} = z_{n(\mu)}$$
 $\mu = 1, \dots, 6$ (91)

where the

$$z_n \equiv \exp(in2\pi/3)$$
 $n = 0, 1, 2$ (92)

are the third roots of unity and $\mu \mapsto n(\mu)$ denotes any 3-colouring of the octahedron. It follows that the state *s* is symmetrical.

The remaining three-dimensional eigenspace of \mathbb{J} with eigenvalue $j_{\text{maxi}} = 0$ carries the three-dimensional self-representation of \mathcal{O} and hence corresponds to the symmetric antiground state with S = 0 which can be visualized as a octahedron in spin space. Its energy is $E_{\text{max}} = 0$. Again, as in the case of the cube, there is a weakly symmetric coplanar state with the same energy and theorem 1 yields $E_{\text{max}}(S) = \frac{2J}{3}S^2$. This state has the form of a (not necessarily regular) hexagon such that opposing vertices of the octahedron in real space are mapped onto opposing vertices of the hexagon in spin space.

The above RBS energy bounds can also be obtained more simply: since $H_0 = S^2 - (S_{16}^2 + S_{25}^2 + S_{34}^2)$, where $S_{ij} \equiv S_i + S_j$ and J = 1, the minimal energy $E_{\min}(S)$ is obtained for $S_{ij} = 2$ as $E_{\min}(S) = S^2 - 12$. Similarly the energy is maximal for $S_{ij} = \frac{1}{3}S$ which yields $E_{\max}(S) = S^2(1 - \frac{3}{9}) = \frac{2}{3}S^2$.

6.4.3. The cuboctahedron. The cuboctahedron is the quasi-regular polyhedron obtained by joining the midpoints of the cube's edges with their nearest neighbours, see [7, 8], and figure 6.

It can be decomposed into eight triangles with disjoint edges and it is 3-colourable. Hence there is a coplanar ground state with S = 0 giving rise to an RBS-parabola

$$E_{\min}(S) = J\left(\frac{1}{2}S^2 - 24\right) \tag{93}$$

since j = 4 and $j_{\min} = -2$. The eigenvalues of \mathbb{J} together with their degeneracy and irreducible representations of \mathcal{O} are summarized in table 1.

This has been calculated by using the formula connecting the characters with the multiplicity of irreducible representations, c.f. [16], 4.2.31b. From this table it is obvious that the trivial irreducible representation A_1 corresponds to the total anti-ground state E is



Figure 7. A stationary state of the cuboctahedron with energy E = 0. The spin vectors s_{μ} are shown as small arrows attached to vertices of the cuboctahedron.

Table 1. Eigenvalues j_n of the adjacency matrix \mathbb{J} of the cuboctahedron (first column) together with their degeneracy (second column). In the third column the irreducible representations of \mathcal{O} are indicated which occur in the corresponding eigenspaces. The nomenclature A_1, A_2, E, F_1, F_2 follows [16].

j_n/J	Degeneracy	Irreducible representations of \mathcal{O}
-2	5	$E \oplus F_2$
0	3	F_2
2	3	F_1
4	1	A_1

spanned by the complex eigenvector of the coplanar ground state and the self-representation F_1 corresponds to the cuboctahedron in spin space with the energy $E_{\text{max}}(0) = 12 j_{\text{maxi}} = 24 J$. It remains to identify the symmetrical states which generate the two three-dimensional irreducible representations F_2 corresponding to the eigenvalues 0 and -2 of J.

The vertex vectors r_{μ} , $\mu = 1, ..., 12$ of the cuboctahedron may be represented, up to normalization, by integer 3-vectors with components -1, 0, 1 and exactly one 0-component. Then a state vector s_{μ} may be defined by the following rule: invert the first component of r_{μ} after 0 and set the other two components to 0. Here 'after' is understood cyclically, e.g. $(1, 1, 0) \mapsto (-1, 0, 0)$. Thus we obtain a state, see figure 7 where the spins s_{μ} of adjacent vertices are orthogonal. Indeed, this state corresponds to the three-dimensional eigenspace of \mathbb{J} with eigenvalue 0 which is transformed under \mathcal{O} according to F_2 .

A second state s' is obtained by a permutation of the r_{μ} . Each r_{μ} is mapped onto that vertex where the above-defined state vector s_{μ} points to

$$s'_{\mu} \equiv 2s_{\mu} - r_{\mu}$$
 (not normalized). (94)

Alternatively, s'_{μ} is obtained from r_{μ} by the following rule: invert the first component of r_{μ} after 0 and leave the other unchanged. For example, $(1, 1, 0) \mapsto (-1, 1, 0)$. Also this state generates a three-dimensional subspace of the eigenspace of \mathbb{J} with $j_{\min} = -2J$ which transforms under \mathcal{O} according to F_2 .



Figure 8. A ground state of the cuboctahedron formed by four stars of David.

Table 2. Eigenvalues j_n of the adjacency matrix \mathbb{J} of the icosahedron (first column) together with their degeneracy (second column). In the third column the irreducible representations of \mathcal{Y} are indicated which occur in the corresponding eigenspaces. The nomenclature A, F_1 , F_2 , G, H follows [16].

j_n/J	Degeneracy	Irreducible representations of \mathcal{Y}
$-\sqrt{5}$	3	F_2
-1	5	Н
$\sqrt{5}$	3	F_1
5	1	Α

Geometrically, s' is a figure in spin space formed by four stars of David, see figure 8. Thus the cuboctahedron has a coplanar ground state with S = 0 as well as a non-coplanar one with the same energy.

6.5. *Y-symmetrical systems*

We will consider the two remaining Platonic solids as well as the quasi-regular icosidodecahedron and again calculate the decomposition of the eigenspaces of the adjacency matrix \mathbb{J} into \mathcal{Y} -irreducible subspaces for these three cases. As in the case of \mathcal{O} -symmetric systems, the eigenvalue j_{maxi} always corresponds to the three-dimensional self-representation of \mathcal{Y} called F_1 . However, we have not found coplanar states realizing j_{maxi} and hence the question whether $E_{\text{max}}(S)$ will be a parabola remains open.

The situation for ground states is different from the previous examples: for the icosahedron and the dodecahedron we find symmetric ground states with S = 0 corresponding to the threedimensional irreducible representation F_2 of \mathcal{Y} , but no coplanar ones (see tables 2 and 3). Moreover, numerical calculations yield the results

$$E_{\min}^{\text{coplanar}} \approx -43.0614...J > E_{\min} = -20\sqrt{5J}$$
 for the dodecahedron (95)

$$E_{\min}^{\text{coplanar}} \approx -24J > E_{\min} = -12\sqrt{5}J$$
 for the icosahedron. (96)

Hence we conjecture that there are no coplanar ground states and hence, according to theorem 3, $E_{\min}(S)$ will not be an exact parabola. Of course, even overwhelming numerical evidence cannot be considered as a rigorous proof. We can only strictly exclude symmetric

Table 3. Eigenvalues j_n of the adjacency matrix \mathbb{J} of the dodecahedron (first column) together with their degeneracy (second column). In the third column the irreducible representations of \mathcal{Y} are indicated which occur in the corresponding eigenspaces. The nomenclature A, F_1 , F_2 , G, H follows [16].

j_n/J	Degeneracy	Irreducible representations of $\mathcal Y$
$-\sqrt{5}$ -2 0	3 4 4	F ₂ G G
$\sqrt{5}$	5 3 1	H F_1 A

Table 4. Eigenvalues j_n of the adjacency matrix \mathbb{J} of the icosidodecahedron (first column) together with their degeneracy (second column). In the third column the irreducible representations of \mathcal{Y} are indicated which occur in the corresponding eigenspaces. The nomenclature A, F_1 , F_2 , G, H follows [16].

j_n/J	Degeneracy	Irreducible representations of \mathcal{Y}
-2	10	$H \oplus H$
$1 - \sqrt{5}$	3	F_2
-1	4	G
1	4	G
2	5	Н
$1 + \sqrt{5}$	3	F_1
4	1	Α

coplanar ground states, since there are no two-dimensional irreducible representations of \mathcal{Y} . This is in accordance with proposition 4 and the fact that the recently discovered S = 0 ground states of the icosidodecahedron [9] are coplanar and weakly symmetric, but not symmetric (see table 4).

For the two \mathcal{Y} -symmetrical Platonic solids we will investigate more closely the threedimensional geometry of the ground states. By numerical simulation of a heat bath at zero temperature, Schröder [20] has found the angles between adjacent spins to be $\alpha_I \approx 116.6^{\circ}$ for the icosahedron and $\alpha_D \approx 138.2^{\circ}$ for the dodecahedron. From the knowledge of the corresponding irreducible subrepresentation of the natural representation of \mathcal{Y} in \mathbb{R}^N we can now exactly determine the symmetric ground state, see proposition 2. It turns out that these states are well-known stellated geometrical structures, called 'great icosahedron' for the ground state of the icosahedron and 'great stellated dodecahedron' for the ground state of the dodecahedron, see [7, 8]. Thus the angles between adjacent spins are just the angles between neighbouring vertices of these stellated structures and hence have the exact values

$$\alpha_I = \arccos\left(-\frac{\sqrt{5}}{5}\right) \qquad \alpha_D = \arccos\left(-\frac{\sqrt{5}}{3}\right)$$
(97)

in agreement with the numerical findings of [20].

7. Summary

We have tried to give a systematic survey of the properties of (relative) ground states of classical Heisenberg spin systems with particular emphasis on symmetric systems. Further

Example	Section	Ground state characteristics	Parabolicity
General triangle	6.1	Coplanar, $S > 0$	No
N-Pantahedron	6.2	Any $S = 0$ state	Yes
N-polygon	6.3.1	Symmetric, coplanar, $S = 0$	Yes
Pentagonal star	6.3.3	Non-coplanar, $S > 0$	No
Cube	6.4.1	Collinear, $S = 0$	Yes
Octahedron	6.4.2	Coplanar, $S = 0$	Yes
Cuboctahedron	6.4.3	Coplanar, non-coplanar, both $S = 0$	Yes
Icosahedron	6.5	Symmetric, non-coplanar, $S = 0$	Probably no
Dodecahedron	6.5	Symmetric, non-coplanar, $S = 0$	Probably no
Icosidodecahedron	6.5	Coplanar, $S = 0$	Yes

Table 5. Summary of some properties of ground states for the considered examples. Generally, we cannot exclude the existence of other types of ground states not listed in the table.

we have devised various methods of ground state construction, e.g. extension of local ground states, construction of symmetric ground states from irreducible representations of the system's symmetry group, and the construction of relative (anti-) ground states from coplanar ground states with S = 0. The latter procedure yields upper and lower parabolas as the boundaries of the *S*-resolved energy spectrum. Thus we have a sufficient condition for a system to be parabolic. Moreover, under certain assumptions, this condition can also be proved to be necessary. Some properties of the considered examples are summarized in table 5.

The above issues are also relevant for the quantum theory of Heisenberg spin systems. It has been shown for various cases [10, 21] that the shape of the *S*-resolved energy spectrum of the quantum system is very well approximated by the curves $E_{\min}(S)$ and $E_{\max}(S)$ of the corresponding classical system, if only the individual spin quantum number *s* exceeds some moderate value, say s = 2. Hence one can predict semi-quantitative features of the quantum spectrum if one knows the classical spectrum and $s \gtrsim 2$.

For some quantum systems, such as the icosidodecahedron with s = 5/2, where the diagonalization of Heisenberg Hamiltonian is totally impractical, approximation methods such as the density matrix renormalization group method (DMRG) are currently able to provide estimates for the lower boundary of the eigenvalue spectrum. In fact, for the icosidodecahedron the estimates [22] are in very good quantitative agreement with the lower boundary of the parabolic spectrum of the classical system. This agreement suggests that the DMRG results are also close to those for the quantum system with s = 5/2. For other classical systems with the same symmetry group \mathcal{Y} , namely the dodecahedron and the icosahedron we expect non-parabolicity, although the corresponding quantum systems may exhibit approximate rotational bands.

In this sense our results and case studies are to a great extent relevant also for real, quantum spin systems.

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