



SpinW factor 2 problem

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1 The problem

There is a well know problem of the different notations of Heisenberg-like Hamiltonians (whether to include factor 1/2 or minus sign, whether to double count or not). The compatibility of the reported parameters relies on the accurate description of the used model. Spin wave theory is often build on top of the Heisenberg-like Hamiltonian. We found separate source of the mismatch of the result, namely magnon dispersion, when using a popular package SpinW [1] and method [10] on which the package is build. SpinW [1] results do not reproduce the test case of 1D ferromagnetic chain, neither 3D ferromagnetic cubic crystal. In this paper we address this mismatch.

The magnon dispersion for such systems plotted with SpinW is twice as small as the results from textbooks [7, 2, 5, 9, 3, 6, 12] (conversion from textbook's notations are discussed in Appendix I). Since there are various notations for the spin hamiltonian, which is the starting point for the magnon dispersion calculation, in this paper all results are presents with respect to the notation of SpinW paper [10]:

$$H = \sum_{mi,nj} \mathbf{S}_{mi}^T \mathbf{J}_{mi,nj} \mathbf{S}_{nj} + \sum_{mi} \mathbf{S}_{mi}^T \mathbf{A}_{mi} \mathbf{S}_{mi} + \mu_B \mathbf{H}^T \sum_{mi} g_i \mathbf{S}_{mi},$$

where double counting is present in the sum and negative J means ferromagnetic alignment. First term describes exchange interaction, second – single ion anisotropy, third – external magnetic field. The indices m, n are indexing the crystallographic unit cells (running from 1 to L), while i and j label the magnetic atoms inside unit cell (running from 1 to N). \mathbf{S}_i is a 3×1 column vector of spin operators $\{S_{mi}^x, S_{mi}^y, S_{mi}^z\}$, $\mathbf{J}_{mi,nj}$ is a matrix of exchange parameters, \mathbf{A}_{mi} - matrix of single ion anisotropy, \mathbf{H} - column vector of external magnetic field.

For the ferromagnetic 3D crystal with one magnetic center in unit cell the solution of SpinW is:

$$E(\mathbf{k}) = \hbar\omega(\mathbf{k}) = SJn \left(\frac{1}{3} (\cos(k_x l) + \cos(k_y l) + \cos(k_z l)) - 1 \right),$$

where l is the length of lattice parameters. While the textbook's result for the same system is:

$$E(\mathbf{k}) = \hbar\omega(\mathbf{k}) = 2SJn \left(\frac{1}{3} (\cos(k_x l) + \cos(k_y l) + \cos(k_z l)) - 1 \right). \quad (1)$$

In the Fig. 1 The magnon dispersion is plotted for both solutions along the k -path specified in [8], $J = -1$, $S = 1$, $n = 6$.

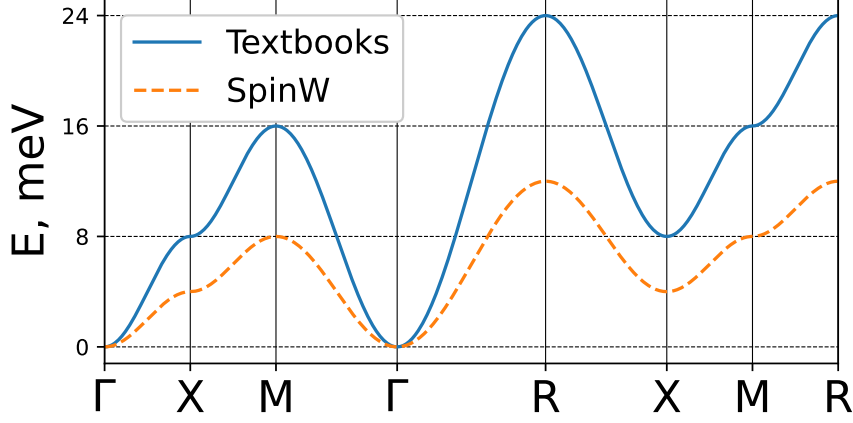


Figure 1: Magnon dispersion comparison between SpinW and textbooks ($J = -1$, $S = 1$).

In the SpinW paper [10] the solution starts by the two consecutive rotations. First one results in the rotation of the exchange matrix $\mathbf{J}'_{mi,nj} = \mathbf{J}_{mi,nj} \mathbf{R}_{n-m}$ and the second rotation defines the vectors \mathbf{u} and \mathbf{v} . These rotations do not affect the following discussion, therefore, we drop the «'» sign in the $\mathbf{J}'_{mi,nj}$ and use the complex valued vectors \mathbf{u} and \mathbf{v} , without recalling their definition. The unique important comment is that for ferromagnetic case (oriented along z axis) the values of the vectors are:

$$\mathbf{u} = (1, i, 0)^T$$

$$\mathbf{v} = (0, 0, 1)^T$$

The single-ion anisotropy and magnetic field can be merged into the exchange term as explained in the SpinW paper [10].

2 The solution

In this section we present the detailed discussion of the correct result and the source of the mismatch. First, we follow the solution of the Heisenberg Hamiltonian presented in SpinW paper [10]. Then we discuss two methods for the solution of the spin wave Hamiltonian in subsections 2.1 and 2.2. At the end we point the source of the mismatch in the original paper [10]. Equation (20) from the SpinW paper [10] is a starting point (where the rotation matrices are absorbed in $\mathbf{J}_{mi,nj}$ and the «'» sign is dropped):

$$H = \sum_{mi,nj} \left\{ \sqrt{\frac{S_i}{2}} \left(\overline{\mathbf{u}}_i^T b_{mi} + \mathbf{u}_i^T b_{mi}^\dagger \right) + \mathbf{v}_i^T (S_i - b_{mi}^\dagger b_{mi}) \right\} \cdot \mathbf{J}_{mi,nj} \cdot \left\{ \sqrt{\frac{S_j}{2}} \left(\overline{\mathbf{u}}_j b_{nj} + \mathbf{u}_j b_{nj}^\dagger \right) + \mathbf{v}_j (S_j - b_{nj}^\dagger b_{nj}) \right\},$$

where b_{mi}^\dagger and b_{mi} are the creation and annihilation operators of the local quantum spin deviations. Overline denotes complex conjugate.

After the expansion the Hamiltonian has zero the energy term E_0 , the one-operator terms, and the two-operator term $H^{(2)}$, which is the center of attention in linearised spin-wave theory. We focus on this term, taking into account the property of the exchange matrix $\mathbf{J}_{mi,nj} = \mathbf{J}_{i,j}(d)$,

$$\mathbf{d} = \mathbf{r}_n - \mathbf{r}_m:$$

$$\begin{aligned} H^{(2)} = \frac{\sqrt{S_i S_j}}{2} & \left(\bar{\mathbf{u}}_i^T \mathbf{J}_{i,j}(\mathbf{d}) \bar{\mathbf{u}}_j b_{mi} b_{nj} + \bar{\mathbf{u}}_i^T \mathbf{J}_{i,j}(\mathbf{d}) \mathbf{u}_j b_{mi} b_{nj}^\dagger \right. \\ & + \mathbf{u}_i^T \mathbf{J}_{i,j}(\mathbf{d}) \bar{\mathbf{u}}_j b_{mi}^\dagger b_{nj} + \mathbf{u}_i^T \mathbf{J}_{i,j}(\mathbf{d}) \mathbf{u}_j b_{mi}^\dagger b_{nj}^\dagger \left. \right) \\ & - \mathbf{v}_i^T \mathbf{J}_{i,j}(\mathbf{d}) \mathbf{v}_j \left(S_i b_{nj}^\dagger b_{nj} + S_j b_{mi}^\dagger b_{mi} \right) \end{aligned}$$

The next step of the solution is to apply Fourier transformation in order to move from the creation and annihilation operators of the local quantum spin deviations (b_{mi}^\dagger and b_{mi}) to the creation and annihilation operators of the collective quantum excitations ($b_i^\dagger(\mathbf{k})$ and $b_i(\mathbf{k})$).

$$b_{mi} = \frac{1}{\sqrt{L}} \sum_{\mathbf{k} \in \text{B.Z.}} b_i(\mathbf{k}) e^{i\mathbf{k}\mathbf{r}_m},$$

$$b_{mi}^\dagger = \frac{1}{\sqrt{L}} \sum_{\mathbf{k} \in \text{B.Z.}} b_i^\dagger(\mathbf{k}) e^{-i\mathbf{k}\mathbf{r}_m},$$

After the Fourier transformation the Hamiltonian has the form:

$$\begin{aligned} H^{(2)} = \sum_{ij} \sum_{\mathbf{k}} & \left[\frac{\sqrt{S_i S_j}}{2} \bar{\mathbf{u}}_i^T \mathbf{J}_{i,j}(\mathbf{k}) \bar{\mathbf{u}}_j b_i(\mathbf{k}) b_j(-\mathbf{k}) + \frac{\sqrt{S_i S_j}}{2} \bar{\mathbf{u}}_i^T \mathbf{J}_{i,j}(\mathbf{k}) \mathbf{u}_j b_i(\mathbf{k}) b_j^\dagger(\mathbf{k}) \right. \\ & + \frac{\sqrt{S_i S_j}}{2} \mathbf{u}_i^T \mathbf{J}_{i,j}(-\mathbf{k}) \bar{\mathbf{u}}_j b_i^\dagger(\mathbf{k}) b_j(\mathbf{k}) + \frac{\sqrt{S_i S_j}}{2} \mathbf{u}_i^T \mathbf{J}_{i,j}(-\mathbf{k}) \mathbf{u}_j b_i^\dagger(\mathbf{k}) b_j^\dagger(-\mathbf{k}) \\ & \left. - S_i \mathbf{v}_i^T \mathbf{J}_{i,j}(\mathbf{0}) \mathbf{v}_j b_j^\dagger(\mathbf{k}) b_j(\mathbf{k}) + S_j \mathbf{v}_j^T \mathbf{J}_{i,j}(\mathbf{0}) \mathbf{v}_i b_i^\dagger(\mathbf{k}) b_i(\mathbf{k}) \right] \end{aligned}$$

We follow the definitions from the equation (26) of the SpinW paper [10]:

$$\mathbf{J}_{i,j}(\mathbf{k}) = \sum_{\mathbf{d}} \mathbf{J}_{i,j}(\mathbf{d}) e^{-i\mathbf{k}\mathbf{d}}$$

$$A(\mathbf{k})^{i,j} = \frac{\sqrt{S_i, S_j}}{2} \mathbf{u}_i^T \mathbf{J}_{i,j}(-\mathbf{k}) \bar{\mathbf{u}}_j,$$

$$B(\mathbf{k})^{i,j} = \frac{\sqrt{S_i, S_j}}{2} \mathbf{u}_i^T \mathbf{J}_{i,j}(-\mathbf{k}) \mathbf{u}_j,$$

$$C(\mathbf{k})^{i,j} = C^{i,j} = \delta_{i,j} \sum_l S_l \mathbf{v}_l^T \mathbf{J}_{i,l}(\mathbf{0}) \mathbf{v}_l.$$

Within this notation the Hamiltonian becomes:

$$\begin{aligned} H^{(2)} = \sum_{ij} \sum_{\mathbf{k}} & \left[\overline{B^{i,j}(\mathbf{k})} b_i(\mathbf{k}) b_j(-\mathbf{k}) + \overline{A^{i,j}(\mathbf{k})} b_i(\mathbf{k}) b_j^\dagger(\mathbf{k}) \right. \\ & + A^{i,j}(\mathbf{k}) b_i^\dagger(\mathbf{k}) b_j(\mathbf{k}) + B^{i,j}(\mathbf{k}) b_i^\dagger(\mathbf{k}) b_j^\dagger(-\mathbf{k}) \\ & \left. - 2C^{i,j} b_i^\dagger(\mathbf{k}) b_j(\mathbf{k}) \right] \quad (2) \end{aligned}$$

It is important to note, that the Hamiltonian (2) is the last point before the solution of SpinW paper [10] differs from the solution of Colpa (subsection 2.1) or White and Bayne (subsection 2.2).

Next step is to rewrite the Hamiltonian in the quadratic form:

$$H = \sum_{\mathbf{k}^?} \mathbf{x}^\dagger(\mathbf{k}) h(\mathbf{k}) \mathbf{x}(\mathbf{k}), \quad (3)$$

where

$$\mathbf{x}(\mathbf{k}) = \left[b_1(\mathbf{k}), \dots, b_N(\mathbf{k}), b_1^\dagger(-\mathbf{k}), \dots, b_N^\dagger(-\mathbf{k}) \right]^T \quad (4)$$

and (in the SpinW paper [10])

$$h(\mathbf{k}) = \begin{pmatrix} \mathbf{A}(\mathbf{k}) - \mathbf{C} & \mathbf{B}(\mathbf{k}) \\ \mathbf{B}^\dagger(\mathbf{k}) & \mathbf{A}(-\mathbf{k}) - \mathbf{C} \end{pmatrix} \quad (5)$$

where \dagger means hermitian conjugate.

The writing of the Hamiltonian in the basis (4) remind of the mathematical formalism picture of particle and antiparticle, nevertheless, this interpretation is not physically justified. Mathematically there is a close connection between the introduction of particle-antiparticle picture and negative energy solutions. However, in the case of linearised spin-wave theory (LSWT) there are no negative energy solutions. Therefore, there is no need to utilize particle-antiparticle interpretation.

There is a question mark near the \mathbf{k} under the sum, since that is the place where SpinW solution and what we are going to do next differ. In the article of Colpa [4], in the textbook by Rezende [7] (page 83) the restriction $\mathbf{k} > 0$ is implied, which means that for each \mathbf{k} in the sum $-\mathbf{k}$ is not in the sum (it is not in the set of indexes over which the sum is carried out). Alternatively, in the article of White [11] and in the textbook by White and Bayne [12] the factor 1/2 added in front of the quadratic Hamiltonian (3) with no restriction to \mathbf{k} , which leads to the same result as with the restriction on \mathbf{k} mentioned above. Finally, in the textbook [6] one of these two details has to be implied, since the result is the same as in other sources. However, Jensen and Mackintosh do not discuss it explicitly.

In contrary, SpinW paper [10] proceeds to cast the Hamiltonian (2) into quadratic form (3) without any restriction on \mathbf{k} , moreover, it is specifically noted under the sum in equation 23 that $\mathbf{k} \in \text{B.Z.}$.

In the SpinW paper the diagonalization of the quadratic form (3) follows the method by Colpa [4]. In the code itself the diagonalization method by White [11] is mentioned. We compare the starting points of Colpa and White with SpinW before diagonalization in the next two subsections. In the Hamiltonian (2) only the part for the \mathbf{k} is written explicitly for each \mathbf{k} under the sum. Therefore, one needs to add terms for $-\mathbf{k}$ in order to construct the form (6). There are two ways to do it:

- To restrict the sum to the $\mathbf{k} > 0$ and rewrite the Hamiltonian. This approach can be interpreted as the separation of the space into $\mathbf{k} > 0$ and $\mathbf{k} < 0$ parts. We focus on this approach in the subsection 2.1, where we follow the solution of Colpa [4].
- To keep the whole set of \mathbf{k} and add $\sum_{-\mathbf{k}}$ to the Hamiltonian. This approach can be interpreted as the artificial doubling of the \mathbf{k} space with the consecutive multiplication by the factor of 1/2, which keeps the Hamiltonian the same. We discuss this approach in the subsection 2.2, where we follow the solution of White and Bayne [12].

The source of the problem with the equation (5) is discussed in the subsection 2.3.

2.1 Colpa [4]

Colpa discusses the diagonalization of the Bogolyubov Hamiltonian of the form:

$$H = \sum_{r',r=1}^m \left(\alpha_{r'}^\dagger \Delta_{1r'r} \alpha_r + \alpha_{r'}^\dagger \Delta_{2r'r} \alpha_{m+r}^\dagger + \alpha_{m+r'} \Delta_{3r'r} \alpha_r + \alpha_{m+r'} \Delta_{4r'r} \alpha_{m+r}^\dagger \right), \quad (6)$$

with the following comment on the nature of the indices r and $m+r$:

The reason why we consider first eq (2.1) is that it often occurs in practice [in solid-state physics e.g. all operators with indices r correspond to the same wave vector \mathbf{k} , those with $m+r$ to $-\mathbf{k}$; m denotes the number of degrees of freedom in the unit cell (or less)]

We have exactly the situation of the solid-state physics. Note, that in the Hamiltonian (6) the sum is carried out over m and not $2m$, which means that the terms with \mathbf{k} and $-\mathbf{k}$ are written explicitly in the sum. First of all, we write the Hamiltonian in a compact form:

$$H^{(2)} = \sum_{ij} \sum_{\mathbf{k}} \left[2(A^{i,j}(\mathbf{k}) - C^{i,j}) b_i^\dagger(\mathbf{k}) b_j(\mathbf{k}) + \overline{B^{i,j}(\mathbf{k})} b_i(\mathbf{k}) b_j(-\mathbf{k}) + B^{i,j}(\mathbf{k}) b_i^\dagger(\mathbf{k}) b_j^\dagger(-\mathbf{k}) \right] \\ + \sum_i \sum_{\mathbf{k}} A^{i,i}(\mathbf{k}) = \sum_{ij} \sum_{\mathbf{k}} H^{i,j}(\mathbf{k}) + const,$$

where we used the fact that $\mathbf{A}(\mathbf{k})$ is Hermitian (see Appendix II) and the commutator $[b_i(\mathbf{k}) b_j^\dagger(\mathbf{k})] = \delta_{i,j}$. In the following we omit the terms of the constant energy shift. Next step is to imply $\mathbf{k} > 0$:

$$H^{(2)} = \sum_{ij} \sum_{\mathbf{k} > 0} (H^{i,j}(\mathbf{k}) + H^{i,j}(-\mathbf{k})) \\ = \sum_{ij} \sum_{\mathbf{k} > 0} \left[2(A^{i,j}(\mathbf{k}) - C^{i,j}) b_i^\dagger(\mathbf{k}) b_j(\mathbf{k}) + \overline{B^{i,j}(\mathbf{k})} b_i(\mathbf{k}) b_j(-\mathbf{k}) + B^{i,j}(\mathbf{k}) b_i^\dagger(\mathbf{k}) b_j^\dagger(-\mathbf{k}) \right. \\ \left. + 2(A^{i,j}(-\mathbf{k}) - C^{i,j}) b_i^\dagger(-\mathbf{k}) b_j(-\mathbf{k}) + \overline{B^{i,j}(-\mathbf{k})} b_i(-\mathbf{k}) b_j(\mathbf{k}) + B^{i,j}(-\mathbf{k}) b_i^\dagger(-\mathbf{k}) b_j^\dagger(\mathbf{k}) \right]$$

We rewrite this Hamiltonian in the form directly comparable with the quadratic form (3) (here the relation $B^{i,j}(\mathbf{k}) + B^{j,i}(-\mathbf{k}) = 2B^{i,j}(\mathbf{k})$, the Hermiticity of $\mathbf{A}(\mathbf{k})$ and $C^{i,j} = C^{j,i}$ is used, see Appendix II):

$$H^{(2)} = \sum_{ij} \sum_{\mathbf{k} > 0} \left[2(A^{i,j}(\mathbf{k}) - C^{i,j}) b_i^\dagger(\mathbf{k}) b_j(\mathbf{k}) \right. \\ \left. + (\overline{B^{j,i}(\mathbf{k})} + \overline{B^{i,j}(-\mathbf{k})}) b_i(\mathbf{k}) b_j(-\mathbf{k}) \right. \\ \left. + (B^{i,j}(\mathbf{k}) + B^{j,i}(-\mathbf{k})) b_i^\dagger(\mathbf{k}) b_j^\dagger(-\mathbf{k}) \right. \\ \left. + 2(A^{j,i}(-\mathbf{k}) - C^{j,i}) b_i(-\mathbf{k}) b_j^\dagger(-\mathbf{k}) \right] + const \\ = \sum_{ij} \sum_{\mathbf{k}} \left[2(A^{i,j}(\mathbf{k}) - C^{i,j}) b_i^\dagger(\mathbf{k}) b_j(\mathbf{k}) \right. \\ \left. + 2\overline{B^{j,i}(\mathbf{k})} b_i(\mathbf{k}) b_j(-\mathbf{k}) \right. \\ \left. + 2B^{i,j}(\mathbf{k}) b_i^\dagger(\mathbf{k}) b_j^\dagger(-\mathbf{k}) \right. \\ \left. + 2(\overline{A^{i,j}(-\mathbf{k})} - C^{i,j}) b_i(-\mathbf{k}) b_j^\dagger(-\mathbf{k}) \right] + const$$

Thus, the matrix $h(\mathbf{k})$ is:

$$h(\mathbf{k}) = \begin{pmatrix} 2(\mathbf{A}(\mathbf{k}) - \mathbf{C}) & 2\mathbf{B}(\mathbf{k}) \\ 2\mathbf{B}^\dagger(\mathbf{k}) & 2(\mathbf{A}(-\mathbf{k}) - \mathbf{C}) \end{pmatrix}, \quad (7)$$

solution of which is the same as in SpinW, but multiplied by the factor 2 and matches with the textbook results. After the diagonalization the matrix is:

$$h'(\mathbf{k}) = \begin{pmatrix} \omega^{(1)}(\mathbf{k}) & 0 \\ 0 & \omega^{(2)}(-\mathbf{k}) \end{pmatrix},$$

where ω is a $N \times N$ diagonal matrix. Diagonalized Hamiltonian looks like (up to a constant term):

$$H^{(2)} = \sum_i \sum_{\mathbf{k}>0} \left(\omega_i^{(1)}(\mathbf{k}) \beta_i^\dagger(\mathbf{k}) \beta_i(\mathbf{k}) + \omega_i^{(2)}(-\mathbf{k}) \beta_i^\dagger(-\mathbf{k}) \beta_i(-\mathbf{k}) \right),$$

with the magnon Hamiltonian being:

$$H_{magnon} = \sum_i \sum_{\mathbf{k}>0} \omega_i^{(1)}(\mathbf{k}) \beta_i^\dagger(\mathbf{k}) \beta_i(\mathbf{k}) + \sum_i \sum_{\mathbf{k}<0} \omega_i^{(2)}(\mathbf{k}) \beta_i^\dagger(\mathbf{k}) \beta_i(\mathbf{k}).$$

Here the magnon dispersion is $E_i(\mathbf{k}) = \hbar\omega_i^{(1)}(\mathbf{k})$ for $\mathbf{k} > 0$ and $E_i(\mathbf{k}) = \hbar\omega_i^{(2)}(\mathbf{k})$ for $\mathbf{k} < 0$. For the ferromagnetic cubic lattice:

$$E(\mathbf{k}) = \hbar\omega_i^{(1)}(\mathbf{k}) = \hbar\omega_i^{(2)}(\mathbf{k}) = 2SJn \left(\frac{1}{3} (\cos(k_x l) + \cos(k_y l) + \cos(k_z l)) - 1 \right),$$

which is the same as textbook's result.

One question remains: the nature of the restriction $\mathbf{k} > 0$. It means that the «positive» and «negative» \mathbf{k} vectors should be separated, but it does not require any particular form of separation. In 3D there are 3 simple non-equivalent separation, which straightforwardly comes to mind: $k_x > 0$, $k_y > 0$ and $k_z > 0$, and one could construct more. The meaning of the separation is that if one think of the $-\mathbf{k}$ as \mathbf{k} and about \mathbf{k} as about $-\mathbf{k}$, then the solution is the same, and $\omega^{(1)}(\mathbf{k})$ describes the spectra with $\mathbf{k} < 0$ and $\omega^{(2)}$ describes the part with $\mathbf{k} > 0$. Which implies $\omega^{(1)}(\mathbf{k}) = \omega^{(2)}(\mathbf{k})$ for any \mathbf{k} and the magnon Hamiltonian could be written as:

$$H^{(2)} = \sum_i \sum_{\mathbf{k}} \omega_i(\mathbf{k}) \beta_i^\dagger(\mathbf{k}) \beta_i(\mathbf{k}),$$

where

$$\omega_i(\mathbf{k}) = \omega_i^{(1)}(\mathbf{k}) = \omega_i^{(2)}(\mathbf{k})$$

2.2 White and Bayne

White and Bayne discuss the diagonalization of the Hamiltonian with dipole-dipole interaction in the book [12]. It includes the same math and ideas as the one we need to use for the solution of the Hamiltonian (2). The quadratic form in this case is (page 246, equation (8.41)):

$$H = \frac{1}{2} \sum_{\mathbf{k}} \mathbf{x}_{\mathbf{k}}^\dagger H_{\mathbf{k}} \mathbf{x}_{\mathbf{k}} \quad (8)$$

And the Hamiltonian, which requires diagonalization (page 246, equation (8.40)) is:

$$H = E_0 + \sum_{\mathbf{k}} \left(A_{\mathbf{k}} a_{\mathbf{k}}^\dagger a_{\mathbf{k}} + B_{\mathbf{k}} a_{\mathbf{k}} a_{-\mathbf{k}} + \overline{B_{\mathbf{k}}} a_{\mathbf{k}}^\dagger a_{-\mathbf{k}}^\dagger \right)$$

From those two equations one can deduct the following. The terms with $a_{\mathbf{k}}a_{-\mathbf{k}}$ and $a_{\mathbf{k}}^\dagger a_{-\mathbf{k}}^\dagger$ are introducing the coupling between $+\mathbf{k}$ and $-\mathbf{k}$, therefore in order to solve the Hamiltonian one has to consider the Hamiltonian for «positive» and «negative» value of **each** \mathbf{k} .

First of all we rewrite the Hamiltonian in a compact form, introducing the notations for this subsection:

$$\begin{aligned} H^{(2)} &= \sum_{ij} \sum_{\mathbf{k}} \left[2(A^{i,j}(\mathbf{k}) - C^{i,j})b_i^\dagger(\mathbf{k})b_j(\mathbf{k}) + \overline{B^{i,j}(\mathbf{k})}b_i(\mathbf{k})b_j(-\mathbf{k}) + B^{i,j}(\mathbf{k})b_i^\dagger(\mathbf{k})b_j^\dagger(-\mathbf{k}) \right] \\ &\quad + \sum_i \sum_{\mathbf{k}} A^{i,i}(\mathbf{k}) = \sum_{ij} \sum_{\mathbf{k}} H^{i,j}(\mathbf{k}) + const \\ &= \sum_{\mathbf{k}} \mathbf{H}(\mathbf{k}) + const \end{aligned}$$

$\mathbf{H}(\mathbf{k})$ describe the \mathbf{k} and its coupling with the $-\mathbf{k}$, thus in order to have the quadratic form, which describes both and their couplings (coupling of \mathbf{k} with $-\mathbf{k}$ and coupling of $-\mathbf{k}$ with \mathbf{k}), one has to construct it in the following way (same as White and Bayne do):

$$H = \frac{1}{2} \sum_{\mathbf{k}} [\mathbf{H}(\mathbf{k}) + \mathbf{H}(-\mathbf{k})],$$

which is effectively leads to the same math as in 2.1 and results in the quadratic form (8) with the matrix $h(\mathbf{k})$ as in (7). And to the same result for the diagonalized Hamiltonian:

$$h'(\mathbf{k}) = \begin{pmatrix} \omega^{(1)}(\mathbf{k}) & 0 \\ 0 & \omega^{(2)}(-\mathbf{k}) \end{pmatrix},$$

$$H^{(2)} = \frac{1}{2} \sum_{i,j} \sum_{\mathbf{k}} \left(\omega_i^{(1)}(\mathbf{k})\beta_i^\dagger(\mathbf{k})\beta_i(\mathbf{k}) + \omega_i^{(2)}(-\mathbf{k})\beta_i^\dagger(-\mathbf{k})\beta_i(-\mathbf{k}) \right)$$

With the magnon Hamiltonian to be:

$$H_{magnon} = \sum_{i,j} \sum_{\mathbf{k}} \frac{1}{2} (\omega_i^{(1)}(\mathbf{k}) + \omega_i^{(2)}(\mathbf{k}))\beta_i^\dagger(\mathbf{k})\beta_i(\mathbf{k})$$

For the ferromagnetic cubic lattice:

$$\hbar\omega_i^{(1)}(\mathbf{k}) = \hbar\omega_i^{(2)}(\mathbf{k}) = 2SJn \left(\frac{1}{3} (\cos(k_x l) + \cos(k_y l) + \cos(k_z l)) - 1 \right),$$

$$E(\mathbf{k}) = \hbar\frac{1}{2}(\omega_i^{(1)}(\mathbf{k}) + \omega_i^{(2)}(\mathbf{k})) = 2SJn \left(\frac{1}{3} (\cos(k_x l) + \cos(k_y l) + \cos(k_z l)) - 1 \right),$$

which is the same as textbooks result.

2.3 SpinW solution

We suspect that the source of the mistake in SpinW paper [10] lies in the construction of the matrix $h(\mathbf{k})$. The diagonalization of the bosonic Hamiltonian with the terms mixing $\pm\mathbf{k}$ requires to have the term with $-\mathbf{k}$, for each \mathbf{k} , even for the negative ones. It means that for some particular \mathbf{k}_0 one has to add the Hamiltonian for $-\mathbf{k}_0$ and for $-\mathbf{k}_0$ one is required to add

the Hamiltonian for \mathbf{k}_0 . Instead, what we suppose happened with the derivation of (5) from (2) is the following:

$$\begin{aligned}
H^{(2)} &= \sum_{ij} \sum_{\mathbf{k}} \left[\overline{B^{i,j}(\mathbf{k})} b_i(\mathbf{k}) b_j(-\mathbf{k}) + \overline{A^{i,j}(\mathbf{k})} b_i(\mathbf{k}) b_j^\dagger(\mathbf{k}) \right. \\
&\quad + A^{i,j}(\mathbf{k}) b_i^\dagger(\mathbf{k}) b_j(\mathbf{k}) + B^{i,j}(\mathbf{k}) b_i^\dagger(\mathbf{k}) b_j^\dagger(-\mathbf{k}) \\
&\quad \left. - 2C^{i,j} b_i^\dagger(\mathbf{k}) b_j(\mathbf{k}) \right] \\
&= \sum_{ij} \sum_{\mathbf{k}} \left[\overline{B^{i,j}(\mathbf{k})} b_i(\mathbf{k}) b_j(-\mathbf{k}) + \overline{A^{i,j}(-\mathbf{k})} b_i(-\mathbf{k}) b_j^\dagger(-\mathbf{k}) \right. \\
&\quad + A^{i,j}(\mathbf{k}) b_i^\dagger(\mathbf{k}) b_j(\mathbf{k}) + B^{i,j}(\mathbf{k}) b_i^\dagger(\mathbf{k}) b_j^\dagger(-\mathbf{k}) \\
&\quad \left. - C^{i,j} b_i^\dagger(\mathbf{k}) b_j(\mathbf{k}) - C^{i,j} b_i^\dagger(-\mathbf{k}) b_j(-\mathbf{k}) \right],
\end{aligned}$$

which is algebraically correct, since $-\mathbf{k}$ is present in the sum for each \mathbf{k} , however, it effectively takes part of the $\mathbf{H}(\mathbf{k})$ Hamiltonian and substitutes it with the corresponding part of the $\mathbf{H}(-\mathbf{k})$ Hamiltonian, which leads to the underestimation of the resulting matrix $h(\mathbf{k})$ by the factor of 2.

3 Appendix I

3.1 Fundamentals of Magnonics [7]

In «Fundamentals of Magnonics» the derivation of magnon dispersion is done in chapter 3 «Quantum Theory of Spin Waves: Magnons».

The Hamiltonian is defined on page 72, eq. 3.6 as follows:

$$H = -g\mu_B \sum_i H_z S_i^z - J \sum_{i,\delta} \vec{S}_i \cdot \vec{S}_{i+\delta},$$

where \vec{S}_i is spin angular momentum operator as site i , $\langle \dots \rangle$ and $\vec{\delta}$ is the vector connecting site i with its nearest neighbors. $\langle \dots \rangle$ Notice also that the factor 2 in the exchange energy does not appear explicitly because each pair of spins is counted twice in the sum over lattice sites.

The definition of the Hamiltonian differs from SpinW with the sign, the following notation change is necessary at the end:

$$J \rightarrow -J$$

Magnon dispersion is provided on page 78 in eqs. 3.35 and 3.36

$$E_k = A_k = g\mu_B H_z + 2zJS(1 - \gamma_k),$$

where γ_k is the structure factor given by

$$\gamma_k = \frac{1}{z} \sum_{\vec{\delta}} e^{i\vec{k}\cdot\vec{\delta}}$$

where z is the number of neighbors (n in the notation of this paper). γ_k for the cubic system is (it is provided on page 79 in eq. 3.37):

$$\gamma_k = \frac{1}{3}(\cos(k_x a) + \cos(k_y a) + \cos(k_z a))$$

where a is a lattice parameter (l in the notation of this paper). The final equation from the [7] in the notation of SpinW is

$$\hbar\omega(\mathbf{k}) = 2nJS \left(\frac{1}{3} (\cos(k_x l) + \cos(k_y l) + \cos(k_z l)) - 1 \right)$$

3.2 Magnetism in condensed matter [2]

The derivation of magnon dispersion for the ferromagnetic 1D chain is discussed in the section 6.6.6 «Magnons».

The definition of the Hamiltonian is provided on page 122 in eqs. 6.9 and 6.10:

(1) We begin with a semiclassical derivation of the spin wave dispersion. First, recall the Hamiltonian for the Heisenberg model,

$$\hat{\mathcal{H}} = - \sum_{\langle ij \rangle} J \hat{\mathbf{S}}_i \cdot \hat{\mathbf{S}}_j$$

(which is eqn. 6.4) In a one-dimensional chain each spin has two neighbours, so the Hamiltonian reduces to

$$\hat{\mathcal{H}} = -2J \sum_i \hat{\mathbf{S}}_i \cdot \hat{\mathbf{S}}_{i+1}$$

with the comment to the equation (6.4) on the page 116 being:

where the constant J is the exchange integral and the symbol $\langle ij \rangle$ below the \sum denotes a sum over nearest neighbours. The spins \mathbf{S}_i are treated as three-dimensional vectors ...

The definition of the Heisenberg model is found for the first time in the section 4.2.1 on the page 76 in eqs. 4.7 and 4.8:

This motivates the Hamiltonian of the Heisenberg model:

$$\hat{\mathcal{H}} = - \sum_{ij} J_{ij} \mathbf{S}_i \cdot \mathbf{S}_j,$$

where J_{ij} is the exchange constant between the i^{th} and j^{th} spins. The factor of 2 is omitted because the summation includes each pair of spins twice. Another way of writing eqn 4.7 is

$$\hat{\mathcal{H}} = -2 \sum_{i>j} J_{ij} \mathbf{S}_i \cdot \mathbf{S}_j,$$

where the $i > j$ avoids the «double-counting» and hence the factor of two returns. Often it is possible to take J_{ij} to be equal to a constant J for nearest neighbours spins and to be 0 otherwise.

The eq. 6.9 corresponds to the definition in eq. 4.7 and the eq. 6.10 corresponds to the definition in eq. 4.8. The definition in eq. 4.7 differs from SpinW with the sign, the following notation change is necessary at the end:

$$J \rightarrow -J$$

The Hamiltonian is solved specifically for the ferromagnetic 1D chain and not for the 3D cubic system with the final result (equation 6.20 on page 123 and equation 6.25 on page 124)

$$\begin{aligned} \hbar\omega &= 4JS(1 - \cos(qa)), \\ E(q) &= -2NS^2J + 4JS(1 - \cos(qa)), \end{aligned}$$

Magnon dispersion from eq. 6.20 is plotted in the book on page 123 in figure 6.12 (Fig. 2). Path from 0 to π/a is the same as the Γ -X path in Fig. 1. If the parameters $J = 1$, $S = 1$ are substituted into the eq. 6.20 then those two graphs are exactly the same.

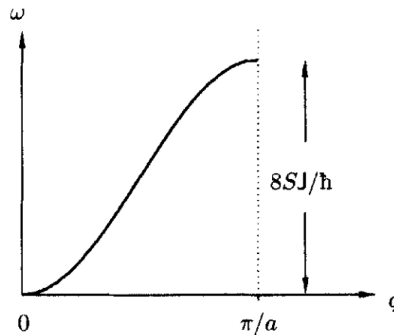


Fig. 6.12 The spin wave dispersion relation for a one-dimensional chain of spins.

Figure 2: Magnon dispersion plot from «Magnetism in condensed matter».

For the cubic system eq. 6.20 in SpinW notation looks like:

$$\hbar\omega(\mathbf{k}) = 2nJS\left(\frac{1}{3}(\cos(q_x a) + \cos(q_y a) + \cos(q_z a)) - 1\right)$$

3.3 Magnetisation oscillations and waves [5]

The derivation of magnon dispersion for the ferromagnet is discussed in the section 7.4 «Elements of microscopic spin-wave theory».

The definition of the Hamiltonian is provided on page 205 in eq. 7.82:

$$\hat{\mathcal{H}} = \gamma\hbar \sum_f \hat{S}_f^z - \sum_f \sum_{f' \neq f} I_{ff'} \mathbf{S}_f \mathbf{S}_{f'}$$

$$\text{where } \mathbf{S}_f \mathbf{S}_{f'} = \hat{S}_f^x \hat{S}_{f'}^x + \hat{S}_f^y \hat{S}_{f'}^y + \hat{S}_f^z \hat{S}_{f'}^z.$$

The double counting is present in this Hamiltonian, thus the definition of the Hamiltonian differs from SpinW with the sign, the following notation change is necessary at the end:

$$J \rightarrow -J$$

The dispersion law is provided in eq. 7.99 on page 209:

where $r_g = r_f - r_{f'}$, $I_g \equiv I_{ff'}$, and the last sum is over all lattice points except one, the initial. The Hamiltonian (7.98) has the desired form of (7.84), and

$$\varepsilon_k(k) = \gamma\hbar H + 2S \sum_g [1 - \exp(i\mathbf{k}\mathbf{r}_g)] I_g.$$

For the cubic ferromagnet the textbook provides the figure 7.13 (Fig. 3)

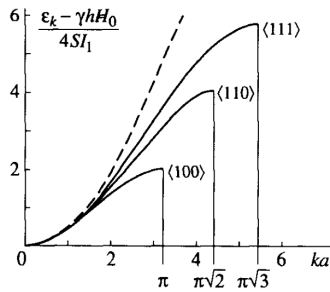
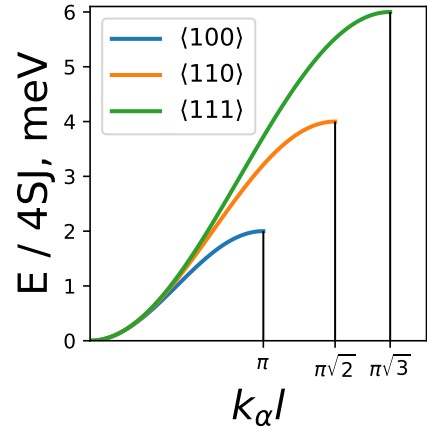


FIGURE 7.13

Dispersion characteristics of spin waves in a ferromagnet with simple cubic spin lattice for different directions of propagation calculated (solid curves) by formula (7.101), i.e., in the nearest-neighbor approximation and without allowance for dipole-dipole interaction. Dashed curve corresponds to the continuum dispersion law.

(a) Original plot



(b) Same plot with use of eq. (1)

Figure 3: Magnon dispersion plot from «Magnetisation oscillations and waves».

In this picture curve $\langle 100 \rangle$ (from 0 to π) corresponds to the path Γ -X, curve $\langle 110 \rangle$ (from 0 to $\pi\sqrt{2}$) to the path Γ -M and curve $\langle 111 \rangle$ (from 0 to $\pi\sqrt{3}$) to the path Γ -R in the Fig. 1. In Fig. 3b the same graph is plotted by using the equation for magnon dispersion from this paper. The picture is produced with the script «codes/dispersion.py» using «custom_moaw» function.

The dispersion law from eq. 7.99 for the cubic system in the notation of SpinW is:

$$\hbar\omega(\mathbf{k}) = 2SI_n \left(\frac{1}{3} (\cos(k_x r_x) + \cos(k_y r_y) + \cos(k_z r_z)) - 1 \right)$$

3.4 The Oxford Solid State Basics [9]

The derivation of magnon dispersion for the ferromagnet is discussed in the exercise 20.3 for the Chapter 20 «Spontaneous Magnetic Order: Ferro-, Antiferro-, and Ferri-Magnetism». The definition of the Hamiltonian is provided on page 229 in eqs. 20.6 and 20.2:

Consider the Heisenberg Hamiltonian

$$\hat{\mathcal{H}} = -\frac{1}{2} \sum_{\langle i,j \rangle} J \mathbf{S}_i \cdot \mathbf{S}_j + \sum_i g \mu_B \mathbf{B} \cdot \mathbf{S}_i$$

and for this exercise set $\mathbf{B} = 0$.

For the first time Heisenberg Hamiltonian is defined on pages 225 – 226 in eq. 20.2:

Note that we have included a factor of 1/2 out front to avoid overcounting, since the sum actually counts both J_{ij} and J_{ji} (which are equal to each other).

$\langle \dots \rangle$

One can use brackets $\langle i, j \rangle$ to indicate that i and j are neighbors:

$$\hat{\mathcal{H}} = -\frac{1}{2} \sum_{\langle i,j \rangle} J_{ij} \mathbf{S}_i \cdot \mathbf{S}_j$$

In a uniform system where each spin is coupled to its neighbors with the same strength, we can drop the indices from $J_{i,j}$ (since they all have the same value) and obtain the so-called *Heisenberg Hamiltonian*

$$\hat{\mathcal{H}} = -\frac{1}{2} \sum_{\langle i,j \rangle} J \mathbf{S}_i \cdot \mathbf{S}_j$$

The double counting is present in this Hamiltonian as well as the additional factor of 1/2, thus the definition of the Hamiltonian differs from SpinW with the sign and factor 1/2, the following notation change is necessary at the end:

$$J \rightarrow -2J$$

The dispersion law for the cubic system is provided on page 230:

▷ Show that the dispersion curve for «spin-waves» of a ferromagnet is given by $\hbar\omega = |F(\mathbf{k})|$ where

$$F(\mathbf{k}) = g\mu_b|B| + JS(6 - 2(\cos(k_x a) + \cos(k_y a) + \cos(k_z a)))$$

where we assume a cubic lattice

In the notation of SpinW the dispersion law becomes:

$$\hbar\omega(\mathbf{k}) = 2JSn \left(\frac{1}{3} (\cos(k_x l) + \cos(k_y l) + \cos(k_z l)) - 1 \right)$$

3.5 Magnetism and magnetic materials [3]

The derivation of magnon dispersion for the ferromagnet is discussed in the section 5.4.1 «Spin waves».

The definition of the Hamiltonian is provided on page 137 in eq. 5.24:

When there is a lattice, the Hamiltonian¹ is generalized to a sum over all pairs of atoms on lattice sites i, j :

$$\hat{\mathcal{H}} = -2 \sum_{i>j} J_{ij} \mathbf{S}_i \cdot \mathbf{S}_j$$

In this definition there is no double counting ($i > j$), but there is a factor of 2 present, thus the definition of the Hamiltonian differs from SpinW with the sign, the following notation change is necessary at the end:

$$J \rightarrow -J$$

The dispersion law for the cubic system is provided on page 163:

The generalization to a three-dimensional cubic lattice with nearest-neighbour interactions is

$$\hbar\omega_q = 2JS \left[Z - \sum_{\delta} \cos \mathbf{q} \cdot \delta \right],$$

where the sum is over the Z vectors δ connecting the central atom to its nearest neighbours.

In case of the cubic system $Z = 6$ and there are 6 nearest-neighbours with the vectors

$$\begin{aligned} &(l, 0, 0), \quad (0, l, 0), \quad (0, 0, l), \\ &(-l, 0, 0), \quad (0, -l, 0), \quad (0, 0, -l), \end{aligned}$$

And the dispersion in notation of SpinW law becomes:

$$\hbar\omega_q = 2JSZ \left(\frac{1}{3} (\cos(q_x l) + \cos(q_y l) + \cos(q_z l)) - 1 \right),$$

3.6 Rare earth magnetism [6]

The derivation of magnon dispersion for the ferromagnet is discussed in the chapter 5 «Spin waves in the ferromagnetic heavy rare earths».

The definition of the Hamiltonian is provided on page 186 in eq. 5.2.1:

$$\hat{\mathcal{H}} = \sum_i \left[\sum_{l=2,4,6} B_l^0 Q_l^0(\mathbf{J}_i) + B_6^6 Q_6^6(\mathbf{J}_i) - g\mu_B \mathbf{J}_i \cdot \mathbf{H} \right] - \frac{1}{2} \sum_{i \neq j} \mathcal{J}(ij) \mathbf{J}_i \cdot \mathbf{J}_j$$

In the eq. 5.2.1 crystal field and magnetic field are considered in the first sum, while the second term represents Heisenberg Hamiltonian. There is a double counting in the sum as well as the factor 1/2, thus the definition of the Hamiltonian differs from SpinW with the sign and factor 1/2, the following notation change is necessary at the end:

$$J \rightarrow -2J$$

The spin-wave spectra is defined on the page 190 in the eq. 5.2.22:

The energy parameters are

$$U_1 = \frac{1}{2} \sum_{\mathbf{q}} (E_{\mathbf{q}} - A_{\mathbf{q}}); \quad E_{\mathbf{q}} = \sqrt{A_{\mathbf{q}}^2 - B^2}.$$

where $A_{\mathbf{q}}$, A and B are defined in the eqs. 5.2.18 and 5.2.15:

$$A = \frac{1}{J} \{3B_2^0 J^{(2)} - 21B_6^6 J^{(6)} \cos 6\phi + g\mu_B JH \cos(\phi - \phi_H)\}$$

$$B = \frac{1}{J} \{3B_2^0 J^{(2)} + 15B_6^6 J^{(6)} \cos 6\phi\}.$$

$\langle \dots \rangle$

$$A_{\mathbf{q}} = A + J \{\mathcal{J}(\mathbf{0}) - \mathcal{J}(\mathbf{q})\}$$

$A = 0$ and $B = 0$ if no magnetic field nor anisotropic effects are considered. In the case of this paper $L = 0$, thus $J = S$ and the equation for the dispersion law is

$$E_{\mathbf{q}} = S (\mathcal{J}(\mathbf{0}) - \mathcal{J}(\mathbf{q}))$$

$\mathcal{J}(\mathbf{q})$ is defined in the eq. 5.1.1a:

$$\mathcal{J}_{sst}(\mathbf{q}) = \sum_{j \in s' - subl.} \mathcal{J}(ij) e^{-i\mathbf{q} \cdot (\mathbf{R}_i - \mathbf{R}_j)}; \quad i \in s - sublattice,$$

For the cubic lattice it becomes:

$$\mathcal{J}(\mathbf{q}) = J (e^{-iq_x l} + e^{iq_x l} + e^{-iq_y l} + e^{iq_y l} + e^{-iq_z l} + e^{iq_z l}) = 2J (\cos(q_x l) + \cos(q_y l) + \cos(q_z l))$$

The dispersion in the notation of SpinW law becomes:

$$E_{\mathbf{q}} = 2nSJ \left(\frac{1}{3} (\cos(q_x l) + \cos(q_y l) + \cos(q_z l)) - 1 \right)$$

3.7 Quantum theory of magnetism [12]

The derivation of magnon dispersion is discussed in the section 8.2.1 «Spin-waves theory». The definition of the Hamiltonian is provided on page 238 in eq. 8.2:

Let us begin by considering a lattice of spins whose interactions may be described by the Heisenberg exchange interaction (2.89). Suppose we apply a uniform static field which serves to define a z -axis. We now wish to determine how this system responds to the time- and space-dependent field $\mathbf{H}_1 \cos(\mathbf{q} \cdot \mathbf{r}) \cos(\omega t)$. If this field is in the x direction, the total Hamiltonian becomes

$$H = - \sum_i \sum_{j \neq i} J_{ij} S_i \cdot S_j + g\mu_B H_0 \sum_i S_i^z + g\mu_B H_1 \sum_i S_i^x \cos(\mathbf{q} \cdot \mathbf{r}) \cos(\omega t).$$

There is a double counting in the sum, thus the definition of the Hamiltonian differs from SpinW with the sign, the following notation change is necessary at the end:

$$J \rightarrow -J$$

The spin-wave spectra is defined on the page 239 in the eq. 8.10:

$$\omega(\mathbf{k}) = \gamma H_0 + \frac{2NS}{\hbar} [J(0) - J(\mathbf{q})].$$

where $J(\mathbf{k})$ is defined on page 134 in the eq. 4.6:

$$J(-\mathbf{q}') \equiv \frac{1}{N} \sum_{i \neq j} J(\mathbf{R}_i - \mathbf{R}_j) e^{i\mathbf{q}' \cdot (\mathbf{r}_i - \mathbf{r}_j)}$$

And the dispersion in the notation of SpinW law becomes:

$$\omega(\mathbf{q}) = \frac{2nSJ}{\hbar} \left(\frac{1}{3} (\cos(q_x l) + \cos(q_y l) + \cos(q_z l)) - 1 \right)$$

4 Appendix II

First of all, we recall the definitions:

$$\begin{aligned}\mathbf{J}_{i,j}(\mathbf{k}) &= \sum_{\mathbf{d}} \mathbf{J}_{i,j}(\mathbf{d}) e^{-i\mathbf{k}\mathbf{d}} \\ A(\mathbf{k})^{i,j} &= \frac{\sqrt{S_i, S_j}}{2} \mathbf{u}_i^T \mathbf{J}_{i,j}(-\mathbf{k}) \bar{\mathbf{u}}_j, \\ B(\mathbf{k})^{i,j} &= \frac{\sqrt{S_i, S_j}}{2} \mathbf{u}_i^T \mathbf{J}_{i,j}(-\mathbf{k}) \mathbf{u}_j, \\ C(\mathbf{k})^{i,j} &= C^{i,j} = \delta_{i,j} \sum_l S_l \mathbf{v}_l^T \mathbf{J}_{i,l}(\mathbf{0}) \mathbf{v}_l.\end{aligned}$$

There are three statements, that requires proofs:

$$\begin{aligned}A^{i,j}(\mathbf{k}) &= \overline{A^{j,i}(\mathbf{k})} \\ B^{i,j}(\mathbf{k}) &= B^{j,i}(-\mathbf{k}) \\ C^{i,j} &= C^{j,i}\end{aligned}$$

Symmetry of the exchange between two sites (note, that transposition and the switch of indices i, j are two different operations):

$$\mathbf{J}_{i,j}(\mathbf{d}) = \mathbf{J}_{j,i}^T(-\mathbf{d})$$

which leads to the:

$$\mathbf{J}_{i,j}(\mathbf{k}) = \mathbf{J}_{j,i}^T(-\mathbf{k})$$

Then:

$$\overline{A^{j,i}(\mathbf{k})} = \frac{\sqrt{S_i, S_j}}{2} \bar{\mathbf{u}}_j^T \mathbf{J}_{j,i}(\mathbf{k}) \mathbf{u}_i = \frac{\sqrt{S_i, S_j}}{2} \bar{\mathbf{u}}_j^T \mathbf{J}_{i,j}^T(-\mathbf{k}) \mathbf{u}_i$$

since $\overline{A^{j,i}(\mathbf{k})}$ is a complex number, we can transpose it, without modification of the result:

$$\overline{A^{j,i}(\mathbf{k})} = \left(\frac{\sqrt{S_i, S_j}}{2} \bar{\mathbf{u}}_j^T \mathbf{J}_{i,j}^T(-\mathbf{k}) \mathbf{u}_i \right)^T = \frac{\sqrt{S_i, S_j}}{2} \mathbf{u}_i^T \mathbf{J}_{i,j}(-\mathbf{k}) \bar{\mathbf{u}}_j = A(\mathbf{k})^{i,j}$$

Second one:

$$B^{j,i}(-\mathbf{k}) = \frac{\sqrt{S_i, S_j}}{2} \mathbf{u}_j^T \mathbf{J}_{j,i}(\mathbf{k}) \mathbf{u}_i = \frac{\sqrt{S_i, S_j}}{2} \mathbf{u}_j^T \mathbf{J}_{i,j}^T(-\mathbf{k}) \mathbf{u}_i$$

since $B^{j,i}(-\mathbf{k})$ is a complex number, we can transpose it, without modification of the result:

$$B^{j,i}(-\mathbf{k}) = \left(\frac{\sqrt{S_i, S_j}}{2} \mathbf{u}_j^T \mathbf{J}_{i,j}^T(-\mathbf{k}) \mathbf{u}_i \right)^T = \frac{\sqrt{S_i, S_j}}{2} \mathbf{u}_i^T \mathbf{J}_{i,j}(-\mathbf{k}) \mathbf{u}_j = B(\mathbf{k})^{i,j}$$

$C^{i,j} = C^{j,i}$ because of the kronecker delta in the definition.

References

- [1] Spinw. <https://www.spinw.org/>.
- [2] Stephen Blundell. *Magnetism in condensed matter*, 2003.
- [3] John MD Coey. *Magnetism and magnetic materials*. Cambridge university press, 2010.
- [4] JHP Colpa. Diagonalization of the quadratic boson hamiltonian. *Physica A: Statistical Mechanics and its Applications*, 93(3-4):327–353, 1978.
- [5] Alexander G Gurevich and Gennadii A Melkov. *Magnetization oscillations and waves*. CRC press, 1996.
- [6] Jens Jensen and Allan R Mackintosh. *Rare earth magnetism*. Clarendon Press Oxford, 1991.
- [7] Sergio M Rezende. *Fundamentals of magnonics*, volume 969. Springer, 2020.
- [8] Wahyu Setyawan and Stefano Curtarolo. High-throughput electronic band structure calculations: Challenges and tools. *Computational materials science*, 49(2):299–312, 2010.
- [9] Steven H Simon. *The Oxford solid state basics*. OUP Oxford, 2013.
- [10] S Toth and B Lake. Linear spin wave theory for single-q incommensurate magnetic structures. *Journal of Physics: Condensed Matter*, 27(16):166002, 2015.
- [11] RM White, M Sparks, and I Ortenburger. Diagonalization of the antiferromagnetic magnon-phonon interaction. *Physical Review*, 139(2A):A450, 1965.
- [12] Robert M White and Bradford Bayne. *Quantum theory of magnetism*, volume 1. Springer, 1983.