

Spin waves

Part I

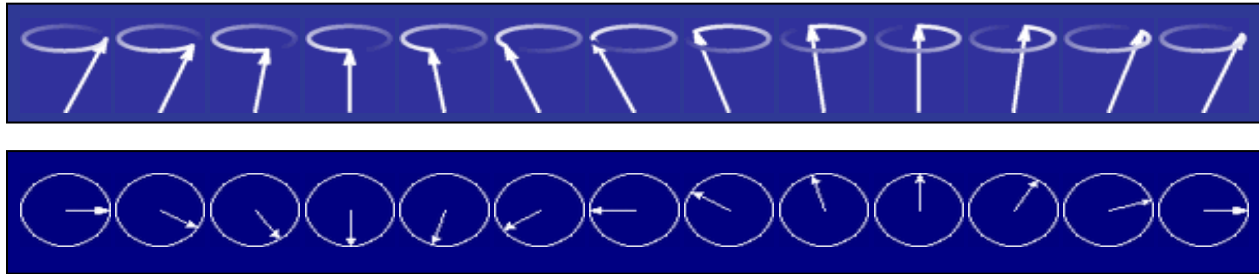
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Why spin waves ?

Time-dependent phenomenon \longrightarrow precession of the spin



Theory developed to describe the excited states of the Heisenberg Hamiltonian

$$\mathcal{H} = \sum_{m,n,i,j} J_{m,i,n,j} \vec{S}_{m,i} \cdot \vec{S}_{n,j}$$

And determine exchange interaction (and anisotropies) via experiments

Why spin waves ?

Bulk systems

$\omega \sim$ A few THz, meV, cm^{-1} in bulk system

$k \sim 0.1 \text{ \AA}^{-1}$

Observed by neutron scattering in (k, ω) space, but also NMR, optical techniques (Raman, $\omega = 0$)

Part I

General considerations

Ferromagnet

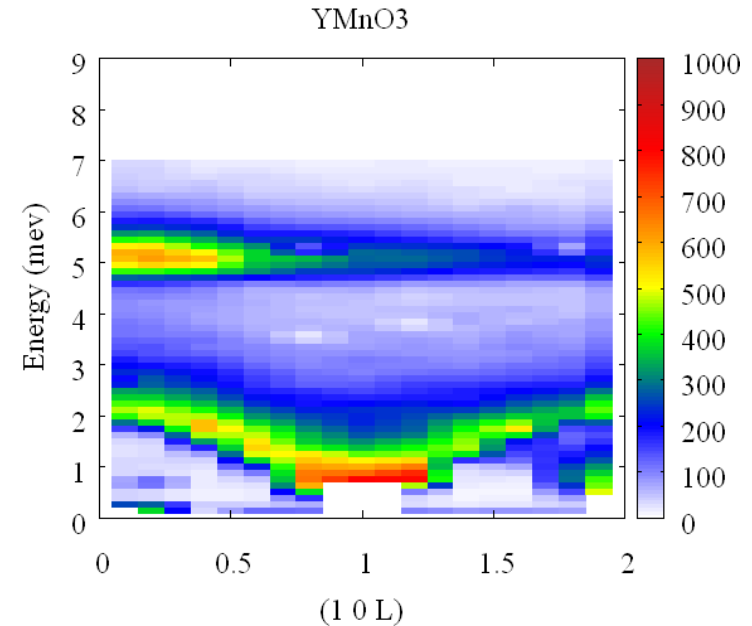
Antiferromagnet

Failure of the theory

Part II

Neutron scattering

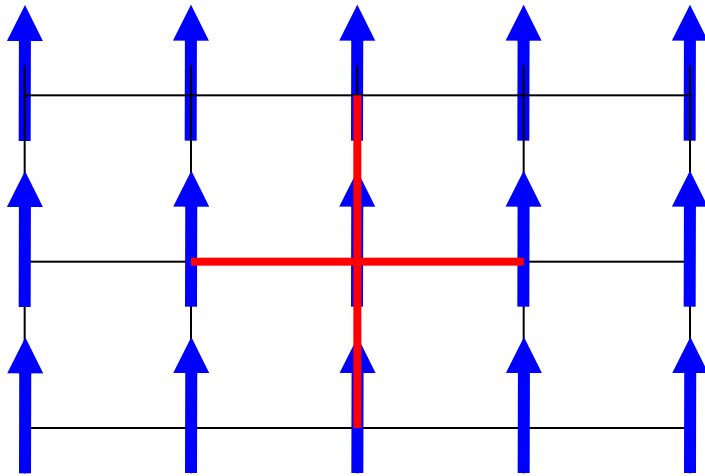
Examples



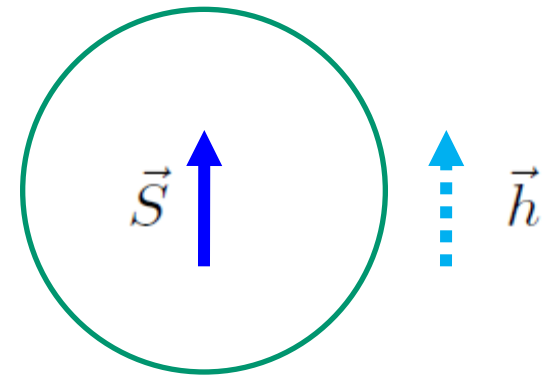
Molecular field

Heisenberg Hamiltonian

$$\mathcal{H} = \sum_{m,n} J_{m,n} \vec{S}_m \cdot \vec{S}_n$$



$J_{m,n}$

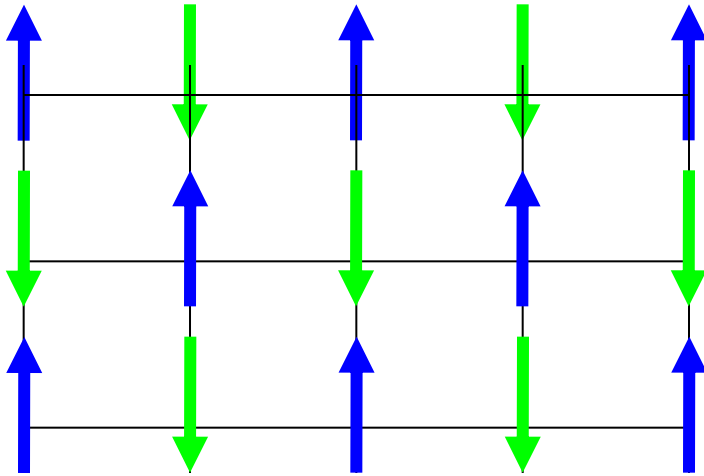


A spin experiences a molecular field due to interaction with its neighbours

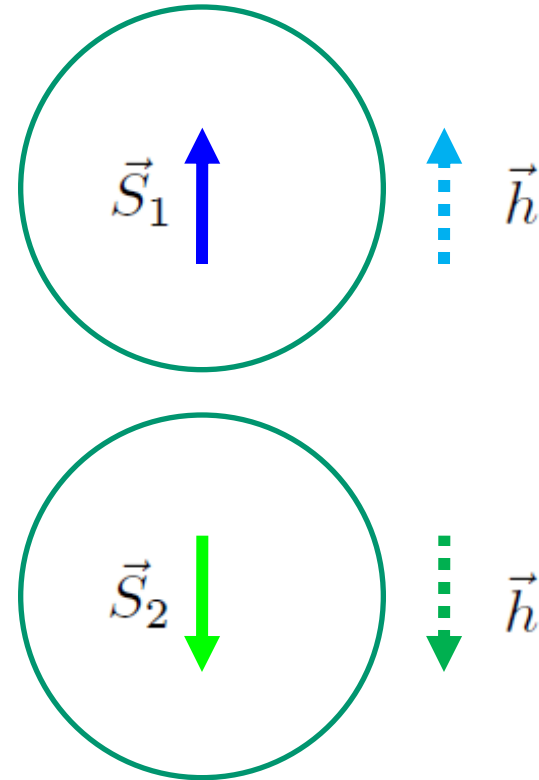
Long range ordering

Molecular field

Depending on interactions, this molecular field can induce a new periodicity

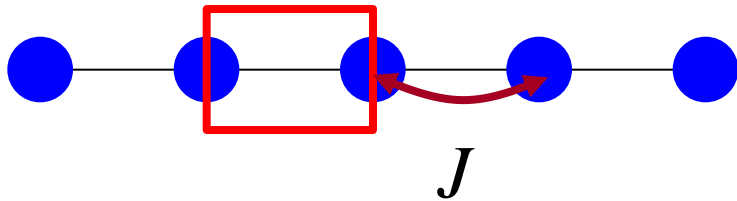


Example : AF ordering

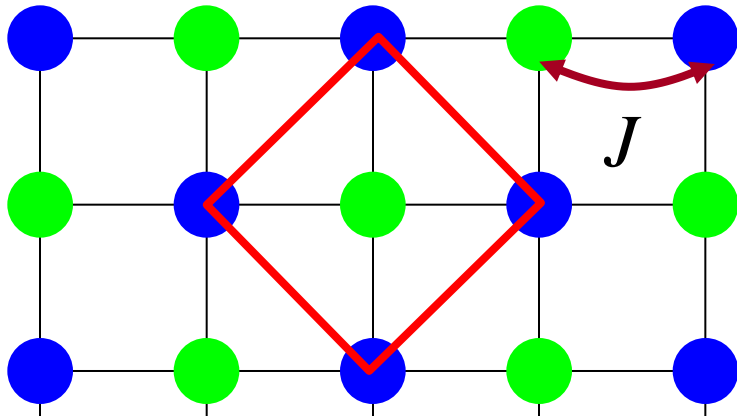


Molecular field

New periodicity, « magnetic unit cell »

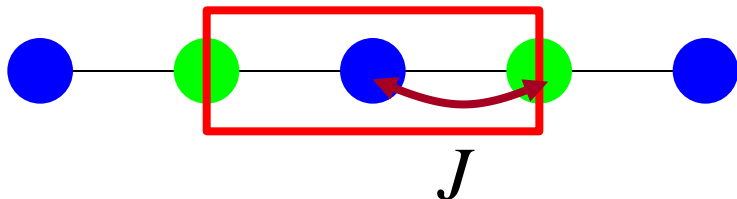


$$\mathcal{H} = \sum_{m,n,i,j} J_{m,i,n,j} \vec{S}_{m,i} \cdot \vec{S}_{n,j}$$



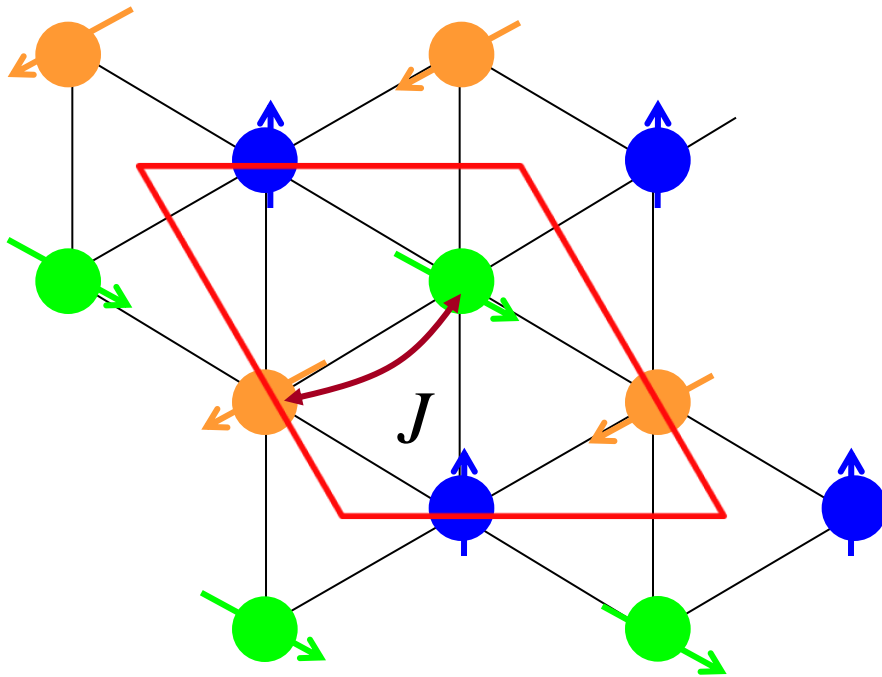
m labels the unit cell

i labels the ion within the unit cell



Molecular field

New periodicity, « magnetic unit cell »



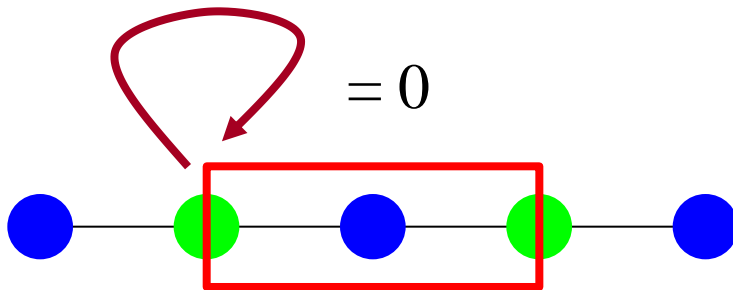
$$\mathcal{H} = \sum_{m,n,i,j} J_{m,i,n,j} \vec{S}_{m,i} \cdot \vec{S}_{n,j}$$

m labels the unit cell

i labels the ion within the unit cell

Molecular field

Define Interactions



$$\mathcal{H} = \sum_{m,n,i,j} J_{m,i,n,j} \vec{S}_{m,i} \cdot \vec{S}_{n,j}$$

$$J_{m,1,m,1}$$

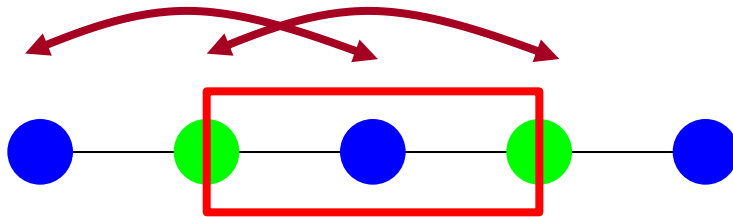
$$J_{m,1,n,1}$$

$$J_{m,1,m,2}$$

$$J_{m,1,n,2}$$

Molecular field

Define Interactions



$$\mathcal{H} = \sum_{m,n,i,j} J_{m,i,n,j} \vec{S}_{m,i} \cdot \vec{S}_{n,j}$$

$$J_{m,1,m,1}$$

$$J_{m,1,n,1}$$

$$J_{m,1,m,2}$$

$$J_{m,1,n,2}$$

Molecular field

Define Interactions

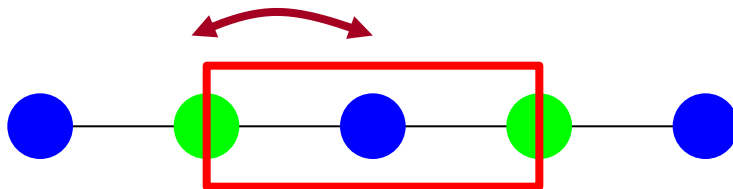
$$\mathcal{H} = \sum_{m,n,i,j} J_{m,i,n,j} \vec{S}_{m,i} \cdot \vec{S}_{n,j}$$

$$J_{m,1,m,1}$$

$$J_{m,1,n,1}$$

$$J_{m,1,m,2}$$

$$J_{m,1,n,2}$$



Molecular field

Define Interactions

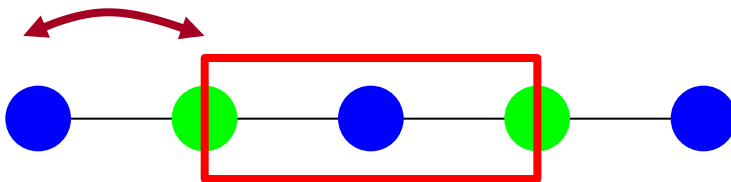
$$\mathcal{H} = \sum_{m,n,i,j} J_{m,i,n,j} \vec{S}_{m,i} \cdot \vec{S}_{n,j}$$

$$J_{m,1,m,1}$$

$$J_{m,1,n,1}$$

$$J_{m,1,m,2}$$

$$J_{m,1,n,2}$$



Molecular field

Mean field approximation

$$\mathcal{H} = \sum_{m,n,i,j} J_{m,i,n,j} \vec{S}_{m,i} \cdot \vec{S}_{n,j}$$



One site Hamiltonian, broken symmetry

$$\mathcal{H} \approx \sum_{m,i} \vec{S}_{m,i} \left(\sum_{n,j} J_{m,i,n,j} \langle \vec{S}_{n,j} \rangle \right)$$

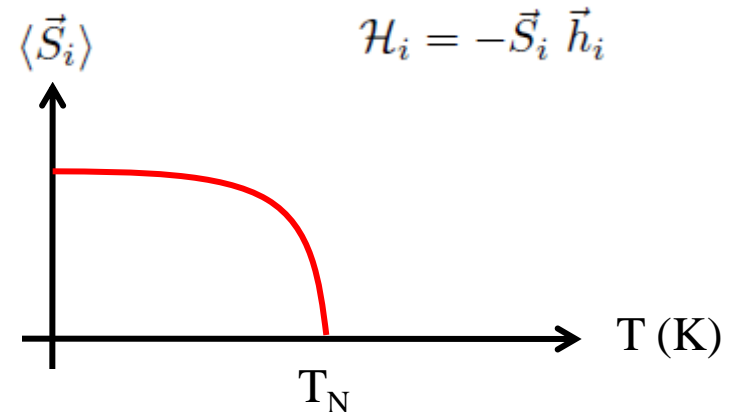


Easy diagonalization

$$\langle \vec{S}_i \rangle = \frac{1}{Z_i} \text{Trace} e^{-\mathcal{H}_i/T} \vec{S}_i$$

$$Z = \sum_{n=-S,S} e^{n h/T}$$

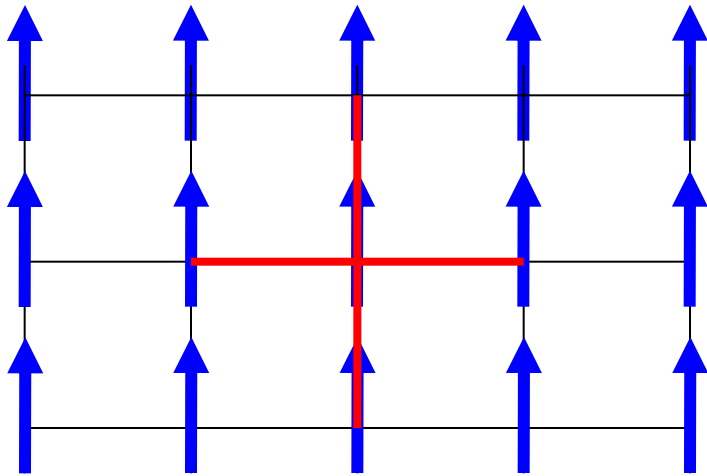
$$S^z |n\rangle = n|n\rangle \quad n = -, -S, \dots, S$$



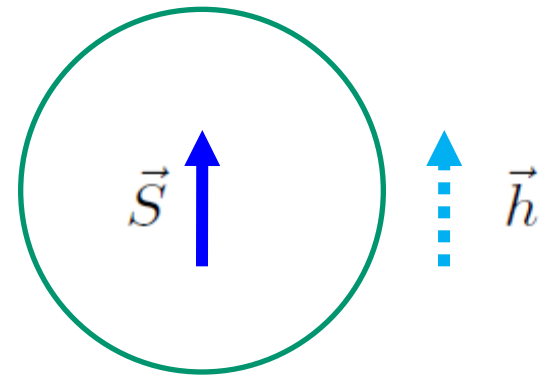
Mermin and Wagner theorem : no spontaneous broken symmetry at finite temperature in 1 and 2 dimension

Spin in a field

$$\mathcal{H} = \sum_{m,n} J_{m,n} \vec{S}_m \cdot \vec{S}_n$$



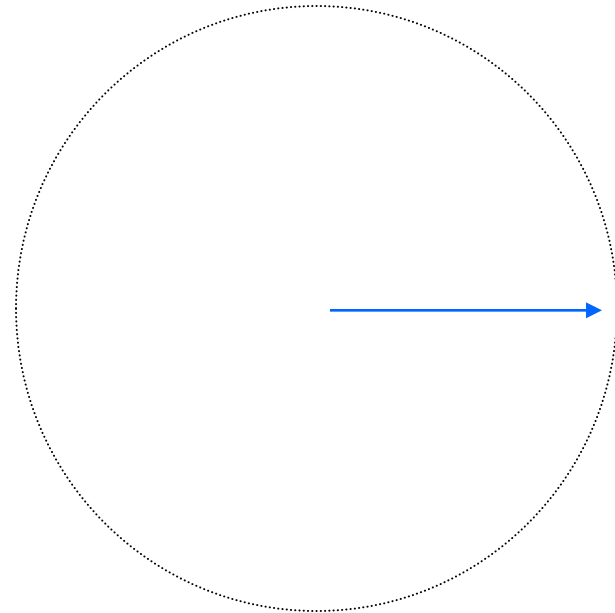
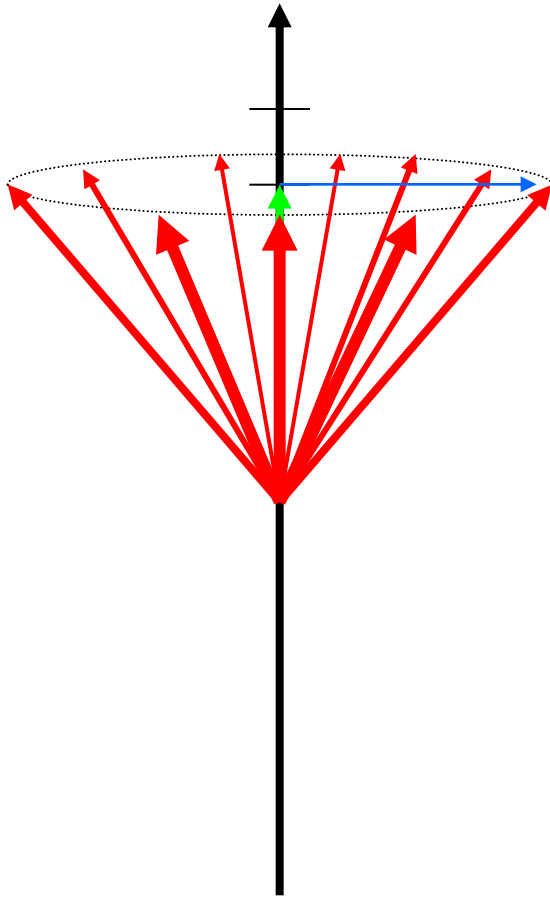
$J_{m,n}$



A spin experiences a molecular field due to the interaction with its neighbors

Spin in a field

Precession of a spin in a magnetic field



Spin in a field

Classical mechanics

$$\mathcal{H} = -\vec{S} \cdot \vec{h}$$

$$\frac{d}{dt} \vec{S} = \vec{S} \times \vec{h} \quad \text{Equation of motion}$$

$$\frac{d}{dt} \begin{pmatrix} S^x \\ S^y \\ S^z \end{pmatrix} = \begin{pmatrix} S^x \\ S^y \\ S^z \end{pmatrix} \times \begin{pmatrix} 0 \\ 0 \\ h \end{pmatrix} = h \begin{pmatrix} S^y \\ -S^x \\ 0 \end{pmatrix}$$

$$\begin{aligned} S^+ &= S^x + i S^y & S^x &= \frac{S^+ + S^-}{2} \\ S^- &= S^x - i S^y & S^y &= \frac{S^+ - S^-}{2i} \end{aligned}$$

Spin in a field

Classical mechanics

$$\frac{d}{dt} \begin{pmatrix} S^+ \\ S^- \\ S^z \end{pmatrix} = \begin{pmatrix} -i\hbar & & \\ & i\hbar & \\ & & 0 \end{pmatrix} \begin{pmatrix} S^+ \\ S^- \\ S^z \end{pmatrix}$$

$\omega = \hbar$

$$\begin{aligned} S^+(t) &= S^+(t=0) e^{-i\omega t} \\ S^-(t) &= S^-(t=0) e^{+i\omega t} \\ S^z(t) &= S^z(t=0) \end{aligned}$$

The spin precesses around S^z with a frequency ω proportional to \hbar

1 degree of freedom

$$(S^x)^2 + (S^y)^2 = \frac{S^+ S^- + S^- S^+}{2} = \frac{S^+(t=0) S^-(t=0) + S^-(t=0) S^+(t=0)}{2}$$

Spin in a field

Quantum mechanics

Spin operators in the local basis

$$\vec{S} = (S^1, S^2, S^3)$$

$$S^3|\ell\rangle = \ell |\ell\rangle \quad \ell = -S, \dots, S$$

$$S^+|\ell\rangle = \sqrt{S(S+1) - \ell(\ell+1)} |\ell+1\rangle$$

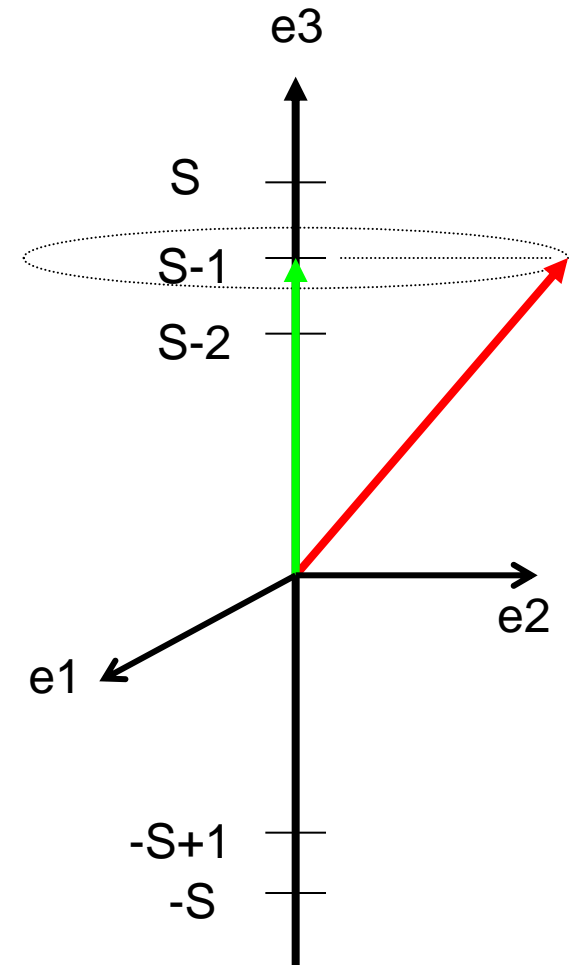
$$S^-|\ell\rangle = \sqrt{S(S+1) - \ell(\ell-1)} |\ell-1\rangle$$

$$[S^+, S^-] = 2 S^3$$

Eigenvalues

$$\mathcal{H} = -\vec{S} \cdot \vec{h} = -h S^3$$

$$E_\ell = h \ell \quad \ell = -S, \dots, S$$



Spin in a field

Quantum mechanics

$$\mathcal{H} = -\vec{S} \cdot \vec{h} = -h S^3$$

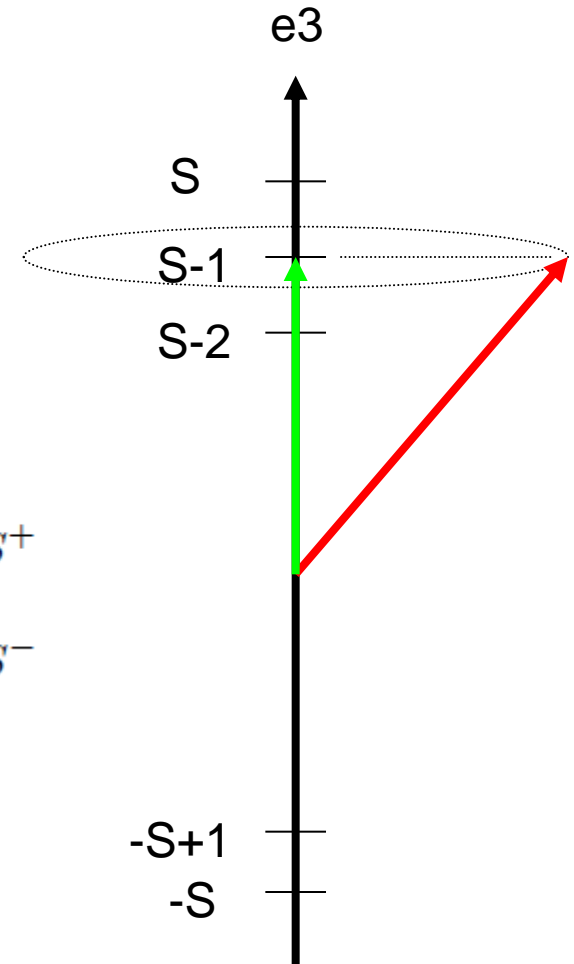
$$\frac{d}{dt} \vec{S} = -i [\vec{S}, \mathcal{H}] \quad \text{Equation of motion}$$

$$\frac{d}{dt} S^3 = -i [S^3, \mathcal{H}] = 0$$

$$\frac{d}{dt} S^+ = -i [S^+, \mathcal{H}] = -i \times -h \times [S^+, S^3] = -ihS^+$$

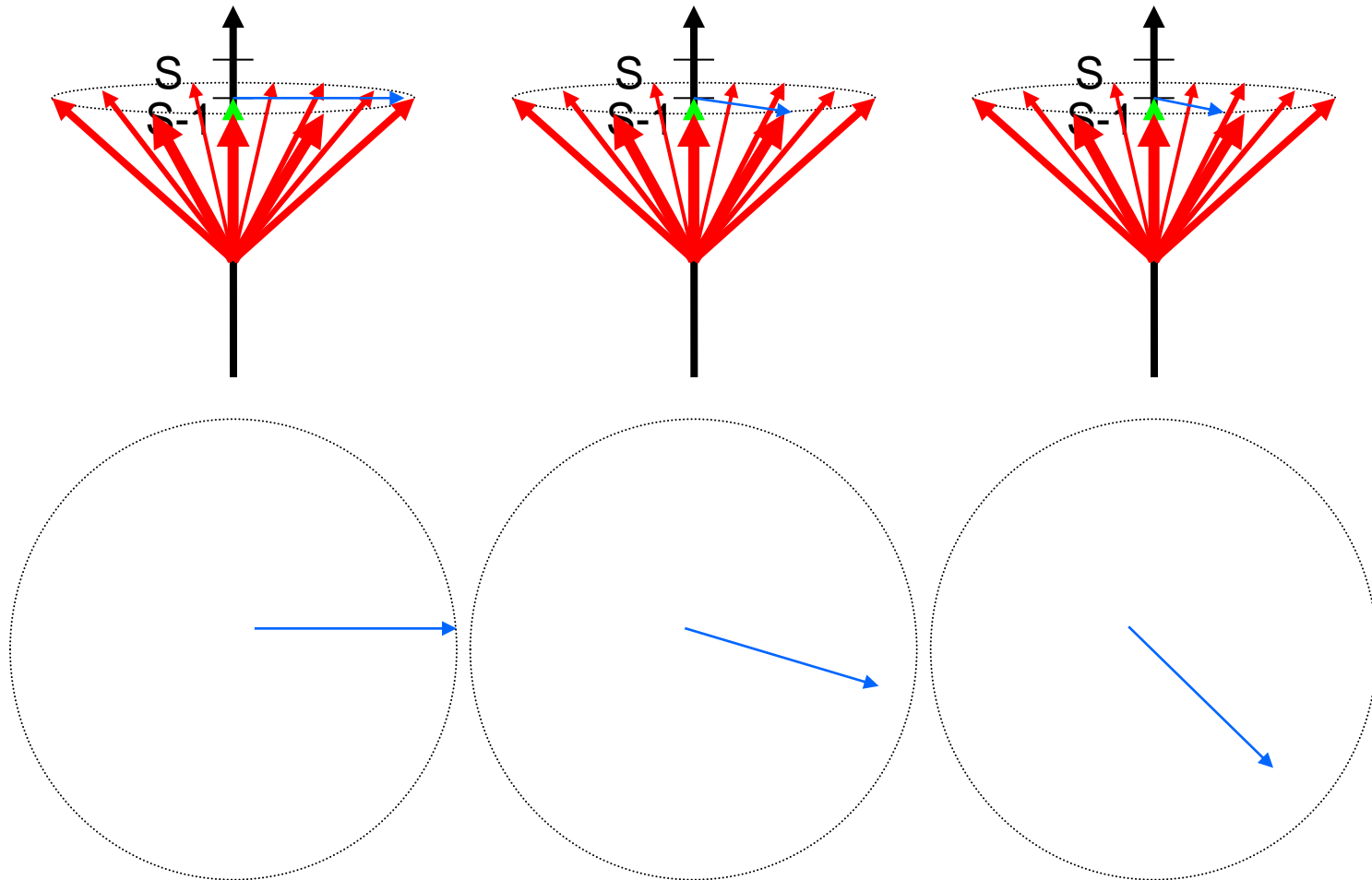
$$\frac{d}{dt} S^- = -i [S^-, \mathcal{H}] = -i \times -h \times [S^-, S^3] = +ihS^-$$

The spin rotates around S^z with a frequency ω proportional to h

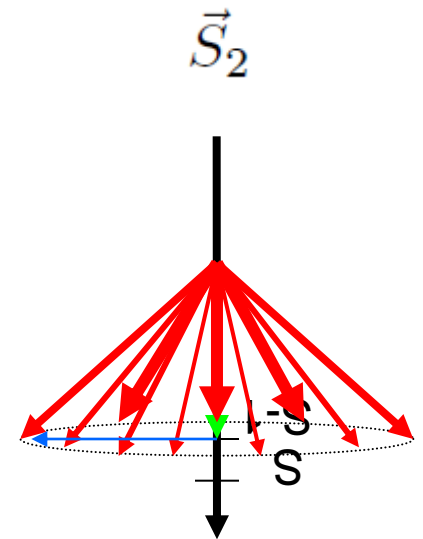
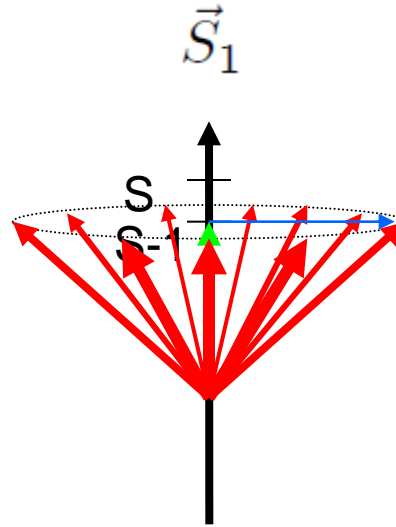
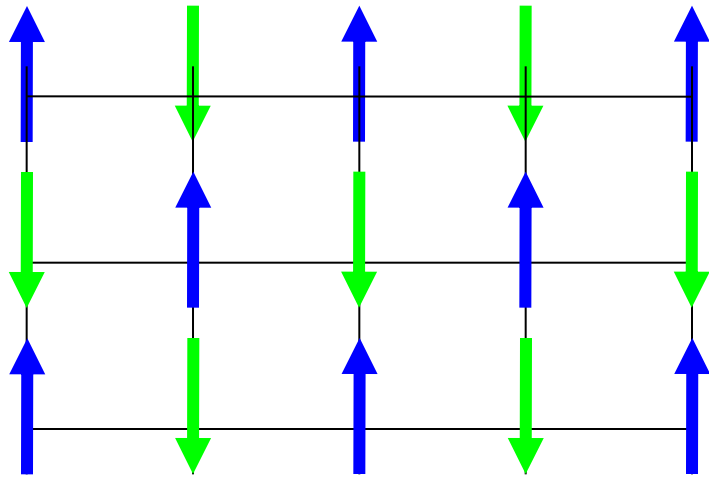


Coupled spins

Back to the problem of coupled spins ...



Transformation to local basis



Cartesian coordinates

$$\vec{S} = \begin{pmatrix} S^x \\ S^y \\ S^z \end{pmatrix}$$

Local coordinates

$$\vec{\sigma} = \begin{pmatrix} \sigma^+ \\ \sigma^- \\ \sigma^3 \end{pmatrix}$$

$$S_i^\alpha = \sum_{\mu=1,2,3} R_i^{\alpha,\mu} \sigma_i^\mu$$

Equation of motion

Classical mechanics

$$\mathcal{H} = \sum_{m,n,i,j} J_{m,i,n,j} \vec{S}_{m,i} \cdot \vec{S}_{n,j}$$

Equation of motion :

N coupled ...

$$\frac{d}{dt} \vec{S}_{m,i} = \vec{S}_{m,i} \times \left(- \sum_{n,j} J_{m,i,n,j} \vec{S}_{n,j} \right)$$

... non linear equations



Equation of motion

Classical mechanics

1. Molecular field : small deviations around the direction of the ordered moment

$$\vec{S} = \langle \vec{S} \rangle + \delta \vec{S} \quad \longrightarrow \quad \vec{S}_{m,i} = \langle \vec{S}_{m,i} \rangle + \delta \vec{S}_{m,i}$$

2. Take advantage of the new periodicity (Fourier transform) : reduce the number of coupled equations



Magnetic unit cell

$$\begin{aligned} \vec{S}_{k,i} &= \frac{1}{\sqrt{N}} \sum_m e^{ikR_m} \vec{S}_{m,i} \\ \vec{S}_{m,i} &= \frac{1}{\sqrt{N}} \sum_k e^{-ikR_m} \vec{S}_{k,i} \end{aligned} \quad \longrightarrow \quad \vec{S}_{k,i} = \langle \vec{S}_i \rangle \delta_{k=0} + \delta \vec{S}_{k,i}$$

3. Exchange $J_{m,i,n,j} = J_{m,i,m+\Delta,j} = \sum_q e^{iq\Delta} J_{q,i,j}$

Equation of motion

Classical mechanics

1. Fourier transform



$$\frac{d}{dt} \vec{S}_{m,i} = \vec{S}_{m,i} \times \left(- \sum_{n,j} J_{m,i,n,j} \vec{S}_{n,j} \right)$$

$$\frac{d}{dt} \vec{S}_{k,i} = - \sum_{k'} \vec{S}_{k-k',i} \times \sum_j J_{k',i,j} \vec{S}_{k',j}$$

2. Ordered moment + small deviations : linearization $\vec{S}_{k,i} = \langle \vec{S}_i \rangle \delta_{k=0} + \delta \vec{S}_{k,i}$

$$\frac{d}{dt} \delta \vec{S}_{k,i} \approx - \left(\sum_{k'} \delta \vec{S}_{k-k',i} \times \sum_j J_{k',i,j} \langle \vec{S}_j \rangle \delta_{k'=0} + \sum_{k'} \langle \vec{S}_i \rangle \delta_{k-k'=0} \times \sum_j J_{k',i,j} \delta \vec{S}_{k',j} \right)$$

$$\approx \sum_j \left(\underbrace{\delta_{i,j} \sum_{\ell} J_{0,i,\ell} \langle \vec{S}_{\ell} \rangle \times - \langle \vec{S}_i \rangle \times J_{k,i,j}}_{\vec{h}_{k,i,j}} \right) \delta \vec{S}_{k,j}$$



$\vec{h}_{k,i,j}$ effective magnetic field acting on $\delta \vec{S}_{k,j}$

Equation of motion

Classical mechanics

Effective magnetic field acting on $\delta\vec{S}_{k,j}$

$$\frac{d}{dt} \begin{pmatrix} \delta S_{k,i}^x \\ \delta S_{k,i}^y \\ \delta S_{k,i}^z \end{pmatrix} = \sum_j \vec{h}_{k,i,j} \times \begin{pmatrix} \delta S_{k,j}^x \\ \delta S_{k,j}^y \\ \delta S_{k,j}^z \end{pmatrix}$$

Local coordinates (use local transformation)

$$\frac{d}{dt} \begin{pmatrix} \delta\sigma_{k,i}^+ \\ \delta\sigma_{k,i}^- \\ \delta\sigma_{k,i}^3 \end{pmatrix} = \sum_j \left(R_i^T \vec{h}_{k,i,j} \times R_j \right) \begin{pmatrix} \delta\sigma_{k,j}^+ \\ \delta\sigma_{k,j}^- \\ \delta\sigma_{k,j}^3 \end{pmatrix}$$

L magnetic ions per magnetic unit cell : **L coupled linear equations**

Approximations :

1. Ordered phase
2. Small deviations around the ordered moment : « linear spin wave theory » (large S, low T)

Ferromagnet

From the general equations of motion

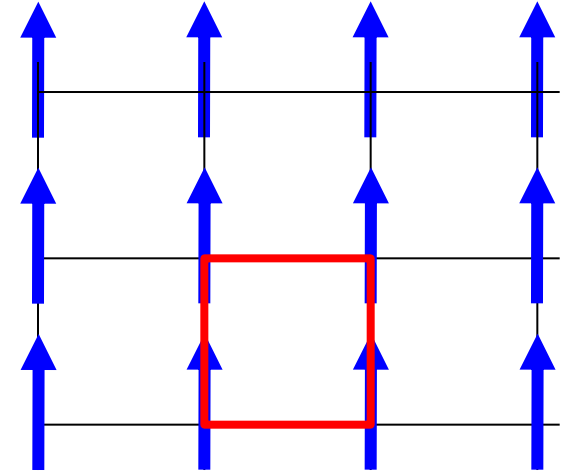
$$\frac{d}{dt} \delta \vec{S}_{k,i} \approx \sum_j \left(\delta_{i,j} \sum_\ell J_{0,i,\ell} \langle \vec{S}_\ell \rangle \times - \langle \vec{S}_i \rangle \times J_{k,i,j} \right) \delta \vec{S}_{k,j}$$

back to the simple ferromagnetic case :

$$\frac{d}{dt} \delta \vec{S}_k \approx (J_0 - J_k) \langle \vec{S} \rangle \times \delta \vec{S}_k$$

$$\frac{d}{dt} \begin{pmatrix} \delta S_k^x \\ \delta S_k^y \\ \delta S_k^z \end{pmatrix} \approx (J_0 - J_k) \begin{pmatrix} 0 \\ 0 \\ \langle S \rangle \end{pmatrix} \times \begin{pmatrix} \delta S_k^x \\ \delta S_k^y \\ \delta S_k^z \end{pmatrix} = (J_0 - J_k) \langle S \rangle \begin{pmatrix} -\delta S_k^y \\ \delta S_k^x \\ 0 \end{pmatrix}$$

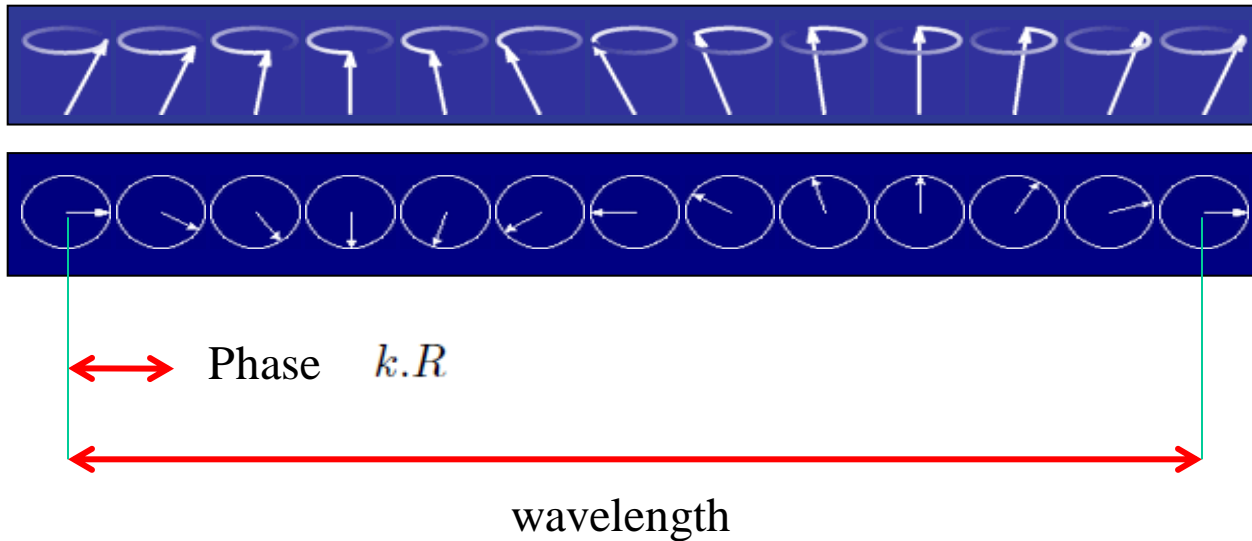
$$\frac{d}{dt} \begin{pmatrix} \delta S_k^+ \\ \delta S_k^- \\ \delta S_k^z \end{pmatrix} = \begin{pmatrix} -i\omega_k & & \\ & i\omega_k & \\ & & 0 \end{pmatrix} \begin{pmatrix} \delta S_k^+ \\ \delta S_k^- \\ \delta S_k^z \end{pmatrix}$$



Ferromagnet

$$\begin{aligned}\delta S_k^+(t) &= \delta S_k^+(t=0) e^{-i\omega_k t} \\ \delta S_k^-(t) &= \delta S_k^-(t=0) e^{+i\omega_k t} \\ \delta S_k^z(t) &= \delta S_k^z(t=0)\end{aligned}$$

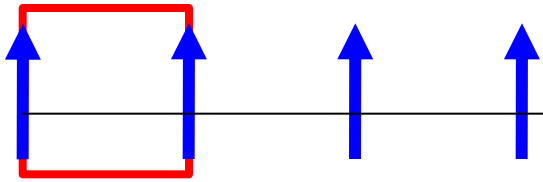
Coupled precessions of the spins around the ordered moment; propagate through the lattice



The dispersion relation connects the wavevector and the frequency ω (energy)

$$\omega_k = - (J_0 - J_k) \langle S \rangle$$

Ferromagnet

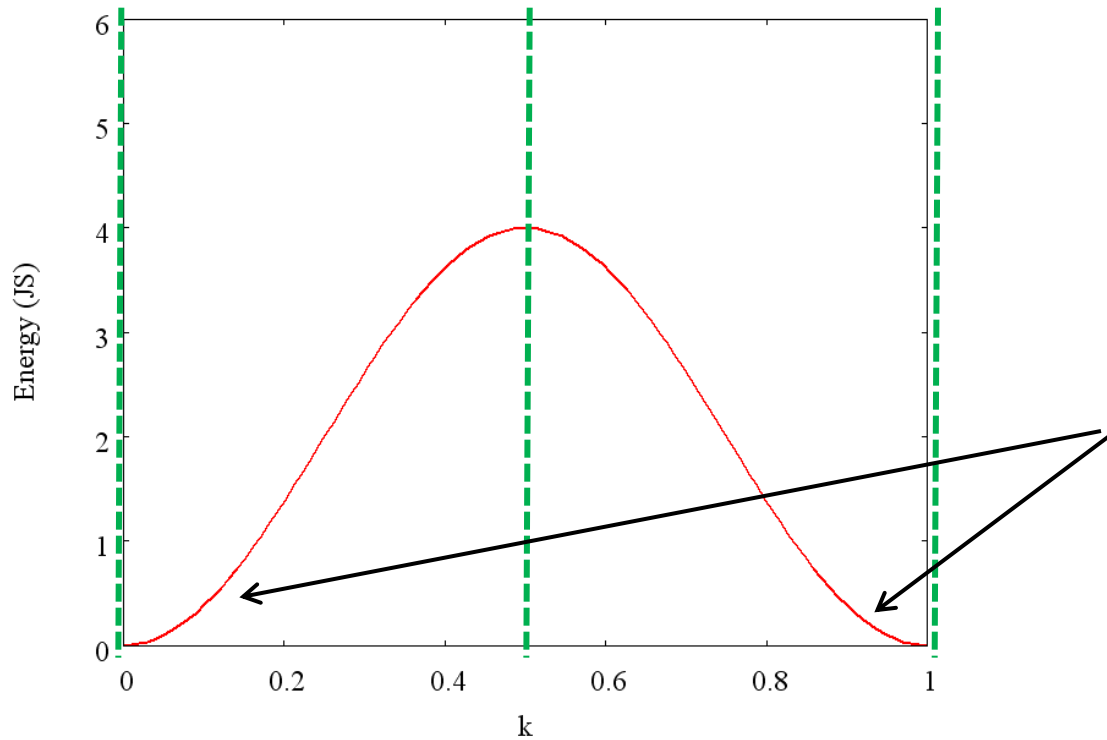


$$\omega_k = 2|J| \langle S \rangle (1 - \cos ka)$$

Zone center

Zone boundary

Zone center of the
next Brillouin zone



$$\omega_k \sim 2 |J| \langle S \rangle \frac{k^2 a^2}{2}$$

Parabolic dispersion

Ferromagnet

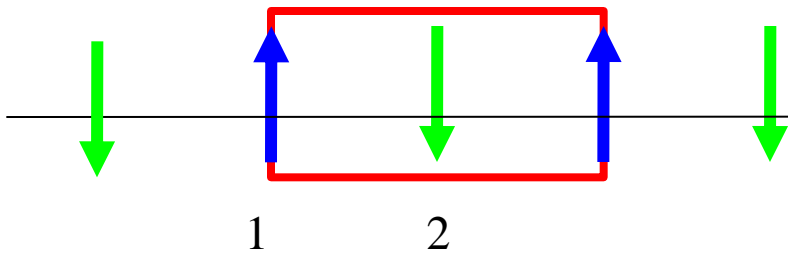
Back to quantum mechanics : spin waves are (quasi) independent Bose modes

$$\mathcal{H} = \sum_k \omega_k b_k^\dagger b_k \quad \langle E \rangle = \sum_k \omega_k n_B(\omega_k) \quad n_B(E) = \frac{1}{e^{E/k_B T} - 1}$$

Check the approximations (correction to the magnetization)

$$\left. \begin{aligned} \langle S \rangle &\approx S - \sum_k n_B(\omega_k) \\ \sum_k &\longrightarrow \int dk^d = \int dk \frac{k^{d-1}}{(2\pi)^d} \\ n_B(E) &\longrightarrow \frac{k_B T}{E} \\ \omega_k &\sim 2 |J| \langle S \rangle \frac{k^2 a^2}{2} \end{aligned} \right\} \begin{aligned} \sum_k n_B(\omega_k) &\longrightarrow \int dk \frac{k^{d-1}}{(2\pi)^d} \frac{T}{k^2} \\ \text{The thermal fluctuations prevents long range} \\ \text{ordering for} \\ &d \leq 2 \\ \text{Breakdown of the spin wave theory is} \\ \text{consistent with Mermin and Wagner theorem} \end{aligned}$$

Antiferromagnet



$$J_{k,1,2} = J_k = J(1 + e^{-ika})$$

$$J_{k,2,1} = J_k^* = J(1 + e^{ika})$$

$$J_{k,1,1} = J_{k,2,2} = 0$$

$$\langle \vec{S}_1 \rangle = -\langle \vec{S}_2 \rangle = \langle \vec{S} \rangle$$

From the general equations of motion

$$\frac{d}{dt} \delta \vec{S}_{k,i} \approx \sum_j \left(\delta_{i,j} \sum_\ell J_{0,i,\ell} \langle \vec{S}_\ell \rangle \times - \langle \vec{S}_i \rangle \times J_{k,i,j} \right) \delta \vec{S}_{k,j}$$

$$\frac{d}{dt} \delta \vec{S}_{k,1} = J_0 \langle \vec{S}_2 \rangle \times \delta \vec{S}_{k,1} - J_k \langle \vec{S}_1 \rangle \times \delta \vec{S}_{k,2}$$

$$\frac{d}{dt} \delta \vec{S}_{k,2} = J_0 \langle \vec{S}_1 \rangle \times \delta \vec{S}_{k,2} - J_k \langle \vec{S}_2 \rangle \times \delta \vec{S}_{k,1}$$

Antiferromagnet

Local coordinates (use local transformation)

4 : There are 2 degrees of freedom

1: Sublattices 1 and 2 are still coupled

$$\frac{d}{dt} \begin{pmatrix} \delta\sigma_{k,1}^+ \\ \delta\sigma_{k,2}^- \end{pmatrix} = -i \langle S \rangle \begin{pmatrix} J_0 & -J_k \\ J_k & -J_0 \end{pmatrix} \begin{pmatrix} \delta\sigma_{k,1}^+ \\ \delta\sigma_{k,2}^- \end{pmatrix}$$

$$\frac{d}{dt} \begin{pmatrix} \delta\sigma_{k,2}^+ \\ \delta\sigma_{k,1}^- \end{pmatrix} = -i \langle S \rangle \begin{pmatrix} J_0 & -J_k \\ J_k & -J_0 \end{pmatrix} \begin{pmatrix} \delta\sigma_{k,2}^+ \\ \delta\sigma_{k,1}^- \end{pmatrix}$$

$$\frac{d}{dt} \delta\sigma_{k,1}^3 = \frac{d}{dt} \delta\sigma_{k,2}^3 = 0$$

3 : Exchange the role of sublattices 1 and 2 (degenerate modes)

2 : Projection on e3 is constant

Antiferromagnet

Additional transformation to decouple sublattice 1 and 2

$$\delta\sigma_{k,1}^+ = u_k \delta A_k + v_k \delta B_k$$

$$\delta\sigma_{k,2}^- = v_k \delta A_k + u_k \delta B_k$$

$$\delta\sigma_{k,2}^+ = u_k \delta A_k + v_k \delta B_k$$

$$\delta\sigma_{k,1}^- = v_k \delta A_k + u_k \delta B_k$$

$$\begin{pmatrix} J_0 & -J_k \\ J_k & -J_0 \end{pmatrix}$$

Details of the transformation :

$$u_k^2 - v_k^2 = 1$$

Spin wave energies

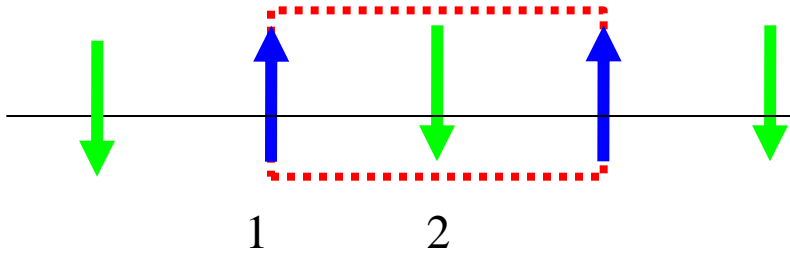
$$\omega_k = \pm \langle S \rangle \sqrt{J_0 - J_k^2}$$

$$u_k^2 = \frac{1}{2} \left(+1 + \frac{J_0 \langle S \rangle}{\omega_k} \right)$$

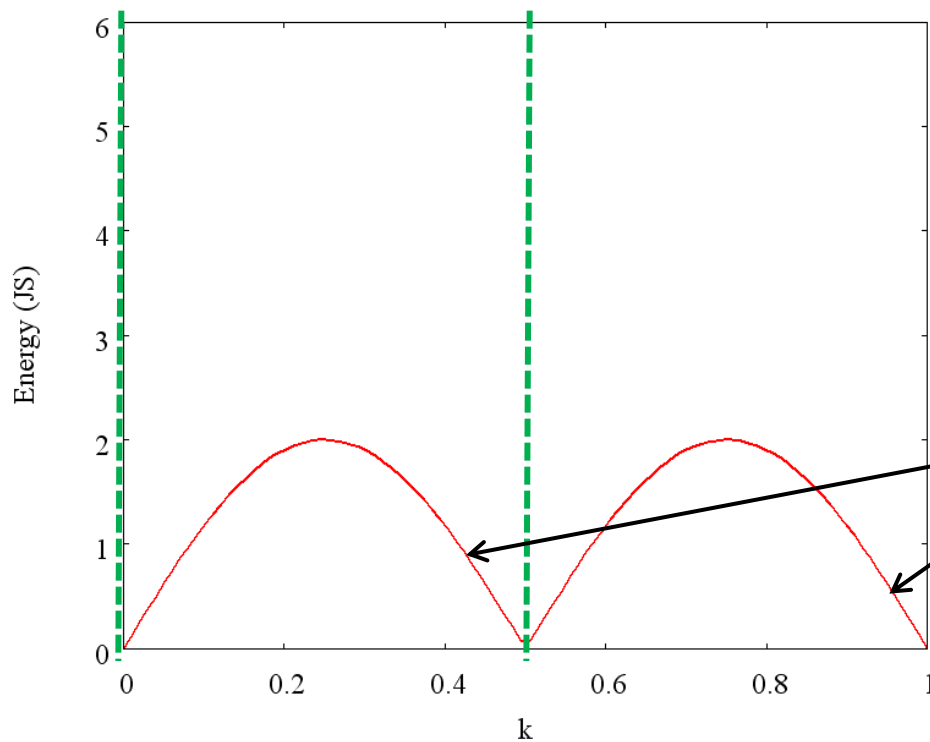
$$v_k^2 = \frac{1}{2} \left(-1 + \frac{J_0 \langle S \rangle}{\omega_k} \right)$$

$$u_k v_k = \frac{J_k \langle S \rangle}{2 \omega_k}$$

Antiferromagnet



$$\omega_k = 2|J| \langle S \rangle |\sin ka|$$

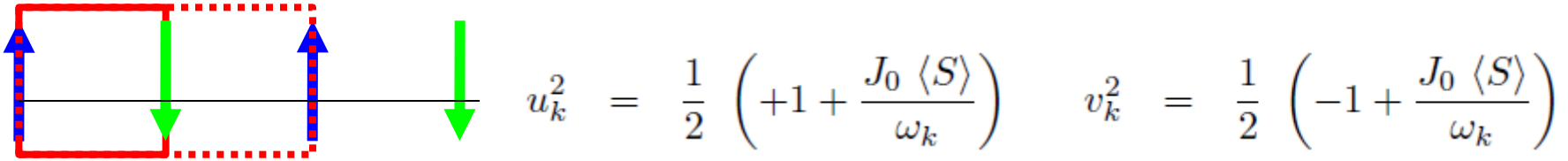


Two degenerate modes

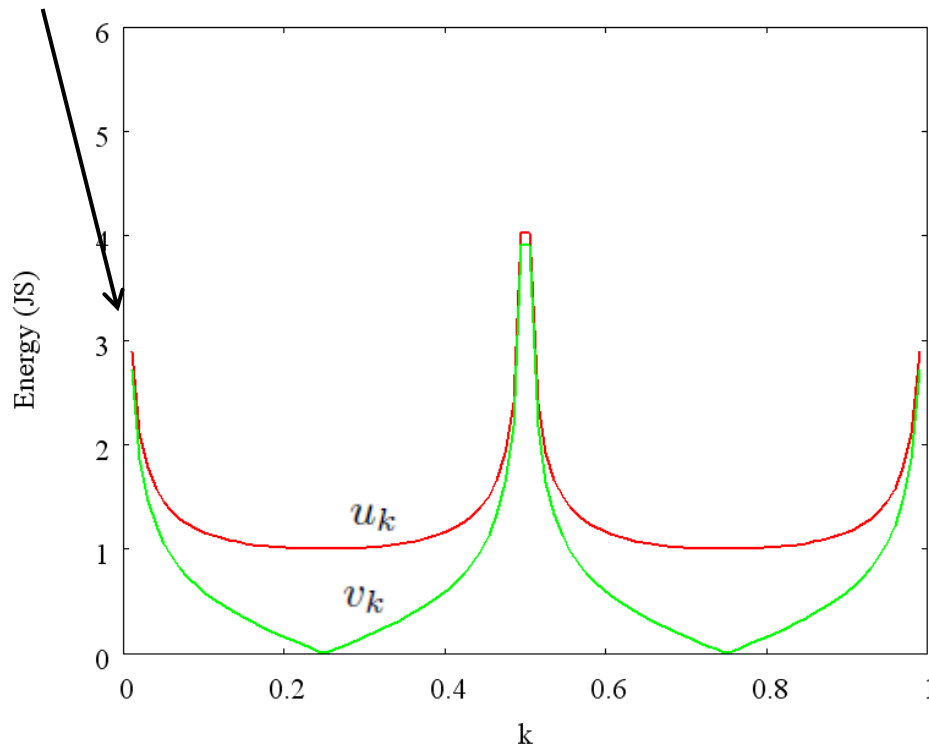
$$\omega_k \sim 2|J| \langle S \rangle ka$$

Linear dispersion

Antiferromagnet



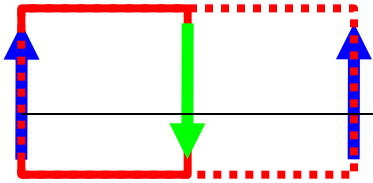
Zone center



$$v_k \sim u_k$$

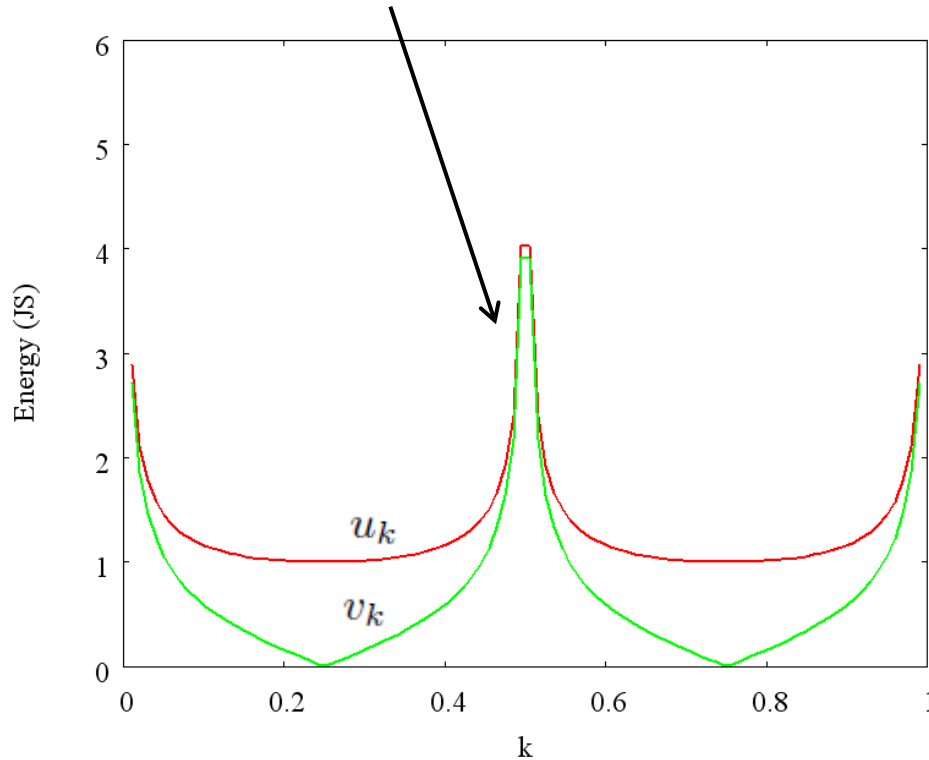


Antiferromagnet



$$u_k^2 = \frac{1}{2} \left(+1 + \frac{J_0 \langle S \rangle}{\omega_k} \right) \quad v_k^2 = \frac{1}{2} \left(-1 + \frac{J_0 \langle S \rangle}{\omega_k} \right)$$

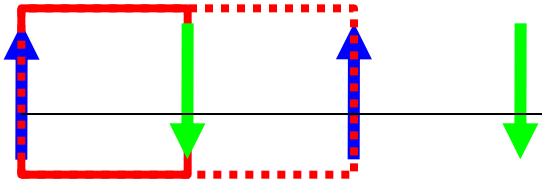
Zone center of the magnetic unit cell,
(Zone boundary of the lattice unit cell)



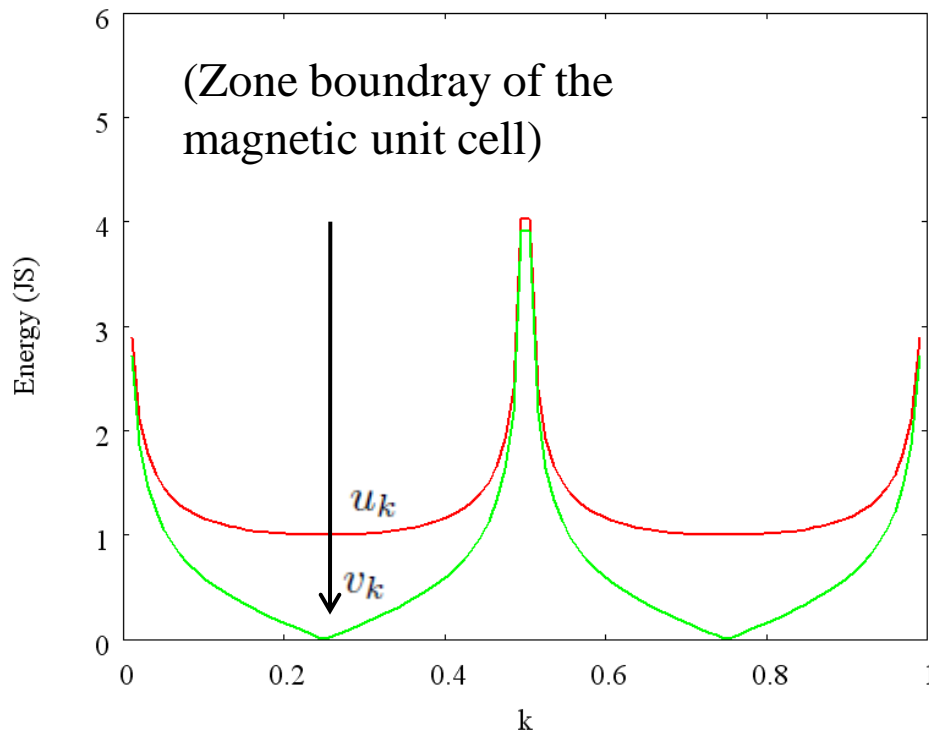
$$v_k \sim -u_k$$



Antiferromagnet



$$u_k^2 = \frac{1}{2} \left(+1 + \frac{J_0 \langle S \rangle}{\omega_k} \right) \quad v_k^2 = \frac{1}{2} \left(-1 + \frac{J_0 \langle S \rangle}{\omega_k} \right)$$



$$v_k = 0, u_k = 1$$



Antiferromagnet

Check the approximations (correction to the magnetization)

$$\langle S \rangle \approx S - \sum_k v_k^2 + (u_k^2 + v_k^2) n_B(\omega_k)$$

$$\langle S \rangle \approx S + \frac{1}{2} - \sum_k \frac{J_0 \langle S \rangle}{\omega_k} \left(n_B(\omega_k) + \frac{1}{2} \right)$$

Thermal fluctuations

$$d \leq 2$$

Quantum fluctuations

$$d \leq 1$$

$$\sum_k \frac{J_0 \langle S \rangle}{\omega_k} \left(n_B(\omega_k) + \frac{1}{2} \right) \longrightarrow \int dk \frac{k^{d-1}}{(2\pi)^d} \frac{1}{k} \left(\frac{T}{k} + \frac{1}{2} \right)$$

Breakdown of the spin wave theory is consistent with Mermin and Wagner theorem

Summary

Spin waves : excited states of the Heisenberg Hamiltonian

L ions per magnetic unit cell : L branches

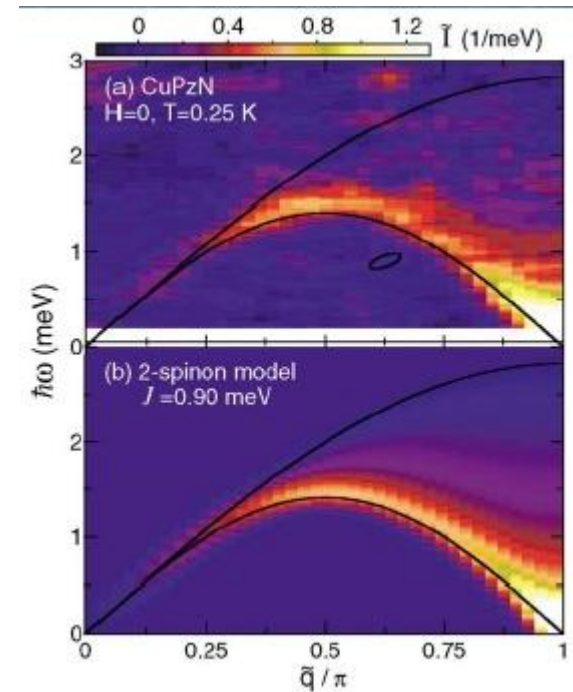
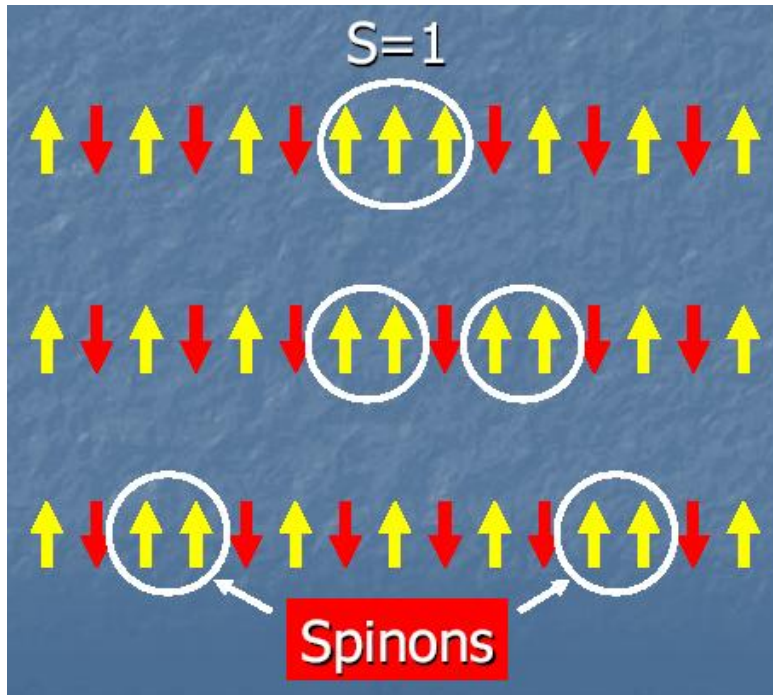
Approximations

1) Ordered phase

2) Small deviations around the ordered moment : large S, low T

Quasi independent modes (bosons) and important role of quantum fluctuations (low dimension)

Beyond spin wave theory



Spin $\frac{1}{2}$: no long range order, no spin waves

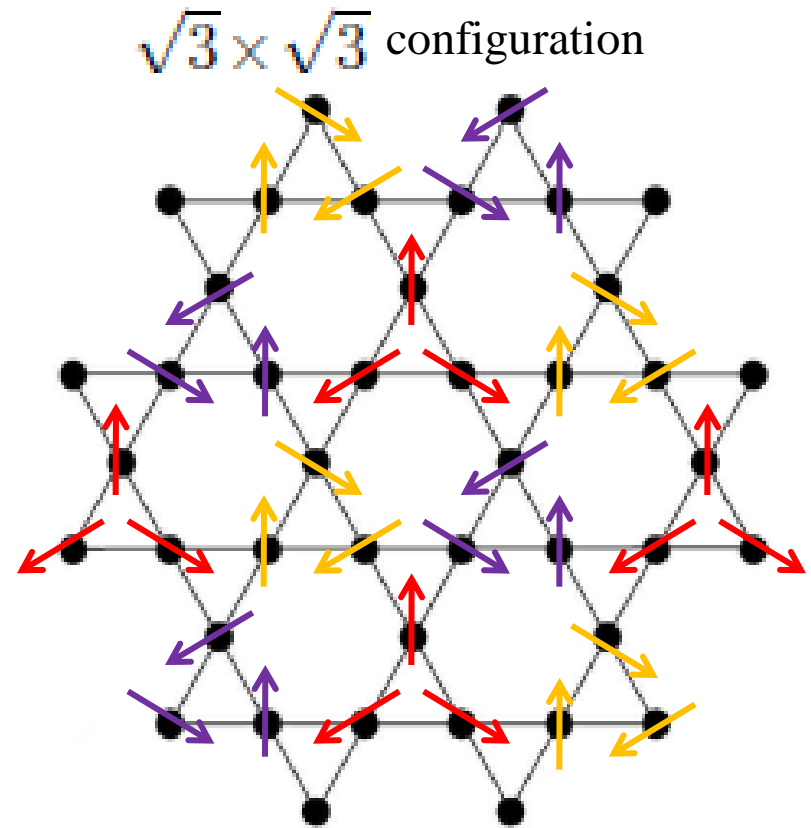
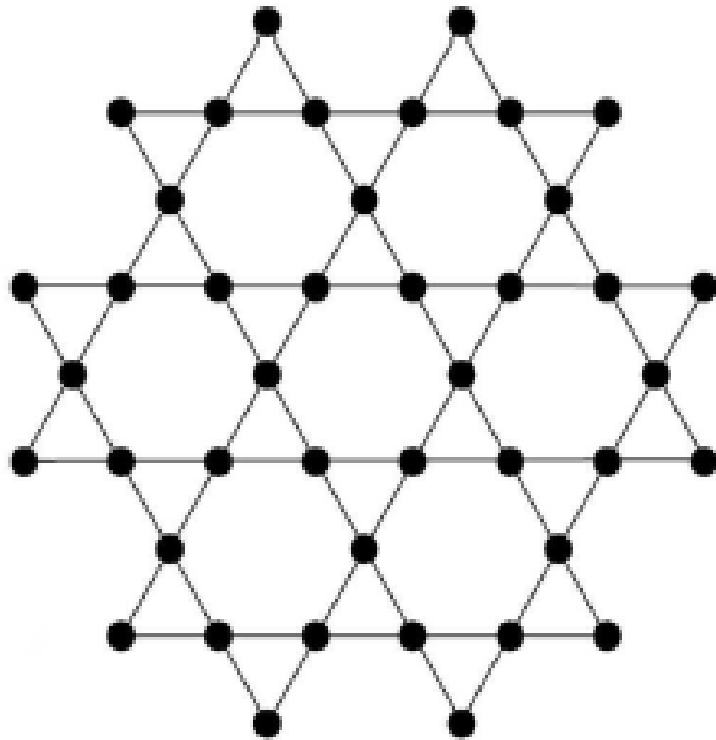
A spin 1 excitation = 2 spinons : continuum and no dispersion relation

Beyond spin wave theory

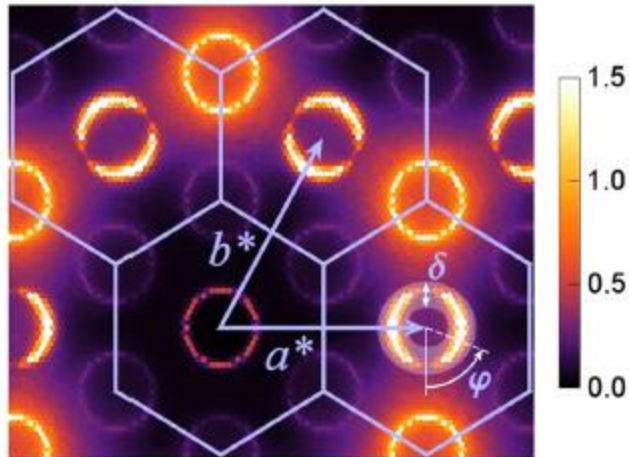
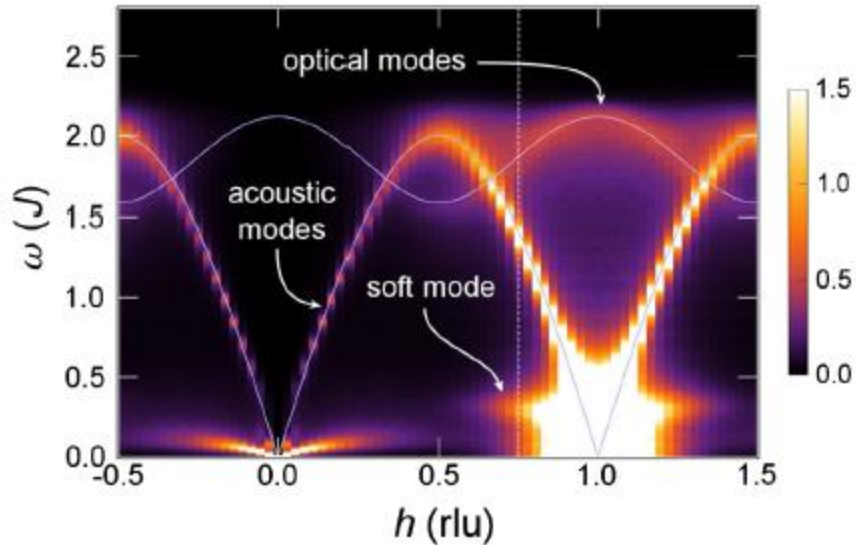
Kagome Lattice

Degenerate ground state : no long range order

The system keeps fluctuating : liquid and co-planar regimes (order by disorder)

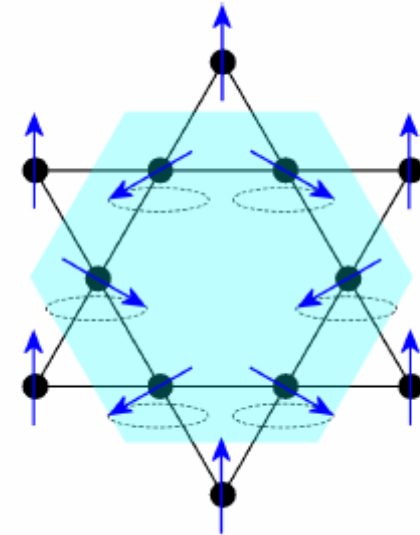


Beyond spin wave theory



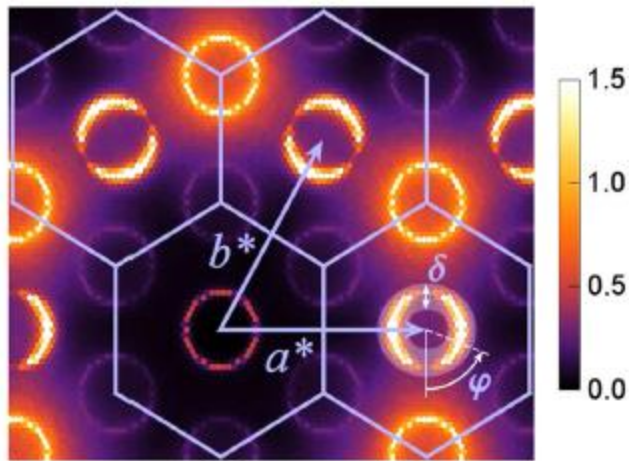
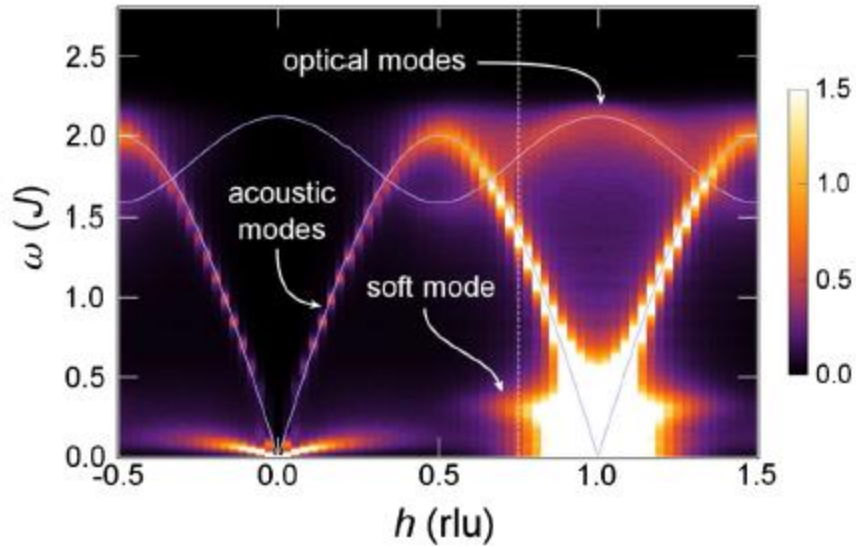
Beyond spin wave theory : calculate the equation of motion for each spin (\sim molecular dynamics) in classical mechanics (no approximation):

Propagative modes as well as soft modes

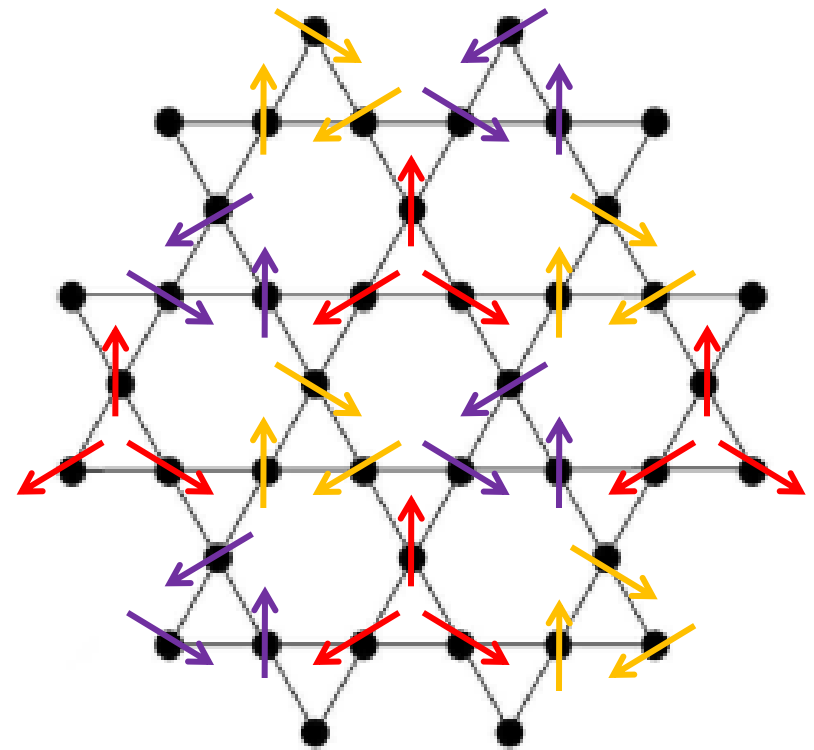


Robert et al, PRL 101, 117207 (2008)

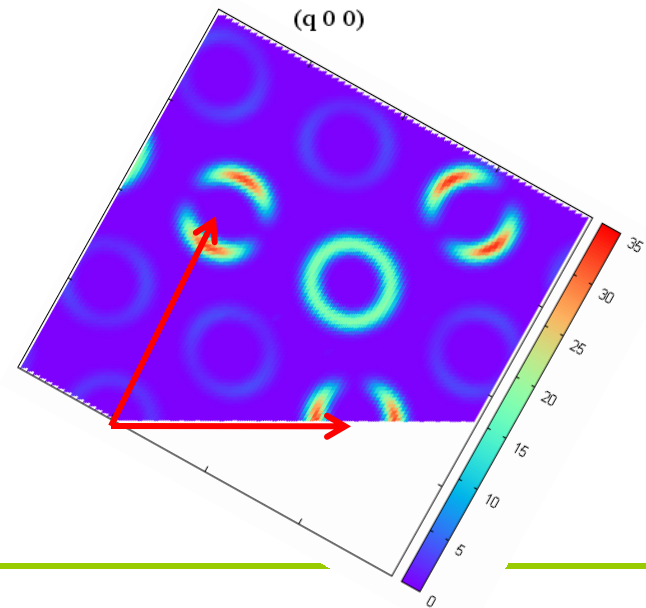
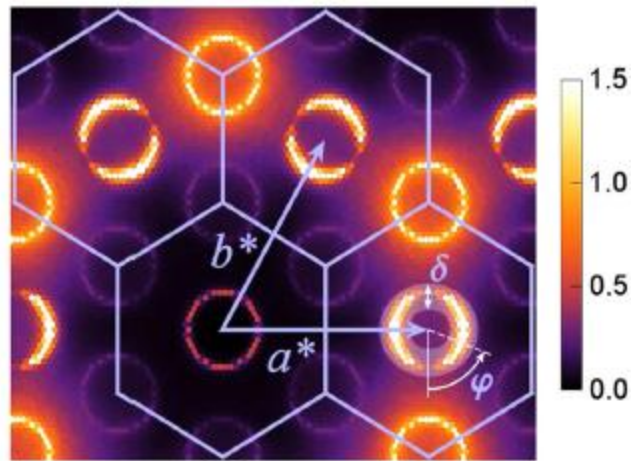
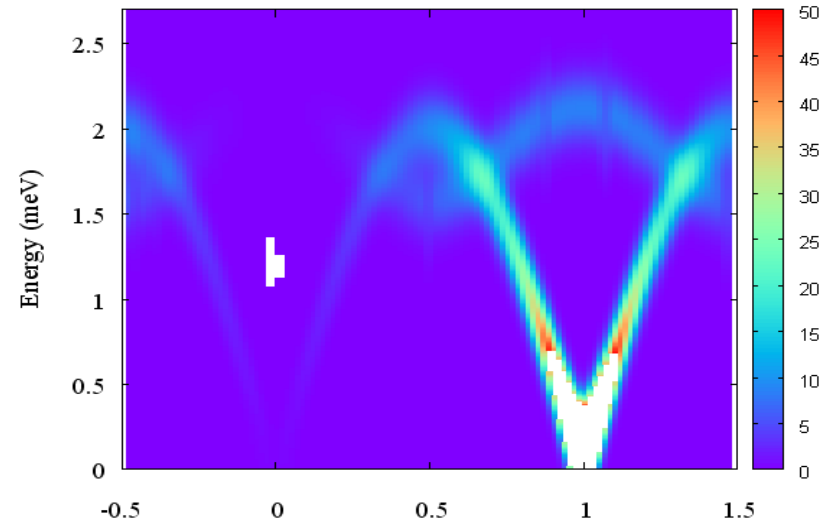
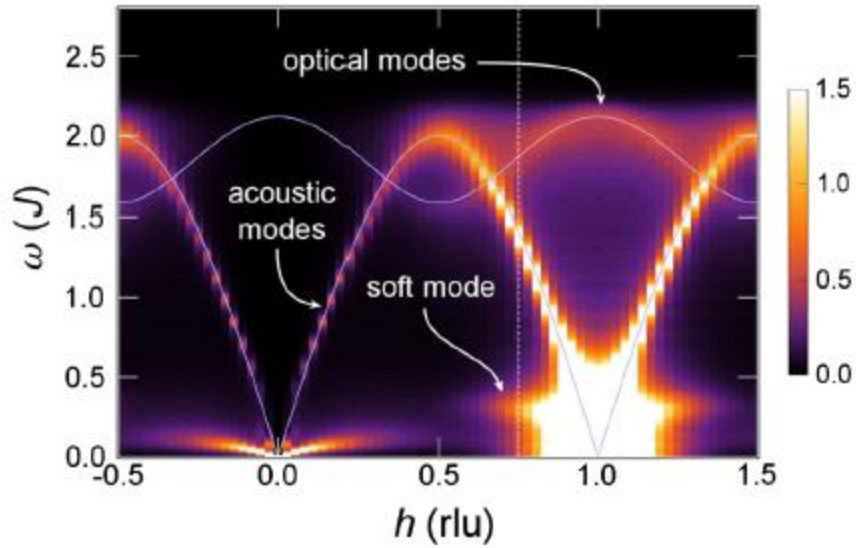
Beyond spin wave theory



$\sqrt{3} \times \sqrt{3}$ Configuration



Beyond spin wave theory



Thanks for your attention

Questions

Practical

To be continued ...Part II : how to observe spin waves ?



References

- [1] P.W. Anderson, Phys. Rev. 83, 1260 (1951)
- [2] R. Kubo, Phys. Rev. 87, 568 (1952)
- [3] T. Oguchi, Phys. Rev 117, 117 (1960)
- [4] D.C. Mattis, *Theory of Magnetism I*, Springer Verlag, 1988
- [5] R.M. White, *Quantum Theory of Magnetism*, Springer Verlag, 1987
- [6] A. Auerbach, *Interacting electrons and Quantum Magnetism*, Springer Verlag, 1994.

Quantum mechanics

Holstein-Primakov representation of the spin « seen from the classical picture » :

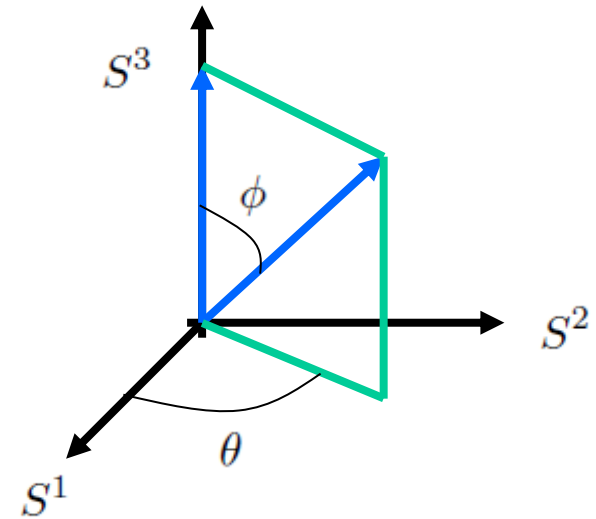
$$\begin{aligned} S^1 &= S \cos \theta \sin \phi & S^+ &= S \sin \phi e^{+i\theta} \\ S^2 &= S \sin \theta \sin \phi & S^- &= S \sin \phi e^{-i\theta} \\ S^3 &= S \cos \phi \end{aligned}$$

New variable : deviation D

$$\boxed{S^3 = S - D} \quad \cos \phi = 1 - \frac{D}{S}$$



$$\begin{aligned} S^+ &= \sqrt{2S} \sqrt{1 - \frac{D}{2S}} \sqrt{D} e^{+i\theta} \\ S^- &= \sqrt{2S} \sqrt{1 - \frac{D}{2S}} \sqrt{D} e^{-i\theta} \\ S^3 &= S - D \end{aligned}$$



Quantum mechanics

Deviation $D \xrightarrow{\text{red arrow}} n_b$

Boson field : $n_b = b^\dagger b = 0, 1, 2, \dots, \infty$ $[b, b^\dagger] = 1$

$$S^+ = \sqrt{2S} \sqrt{1 - \frac{n_b}{2S}} b$$

$$S^- = \sqrt{2S} b^\dagger \sqrt{1 - \frac{n_b}{2S}}$$

$$S^3 = S - n_b$$

$$S^+ \approx \sqrt{2S} b$$

$$S^- \approx \sqrt{2S} b^\dagger$$

$$S^3 = S - n_b$$