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Non-Landau damping of magnetic excitations in systems with localized and itinerant electrons

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We discuss the form of the damping of magnetic excitations in a metal near a ferromagnetic instability. The paramagnon theory predicts that the damping term should have the form $\gamma(q, \Omega) \propto q^4/(\Omega^2 + \Gamma(q)^2)$ (the Landau damping). However, the experiments on uranium metallic compounds UGe$_2$ and UCoGe showed that $\Gamma(q)$ is essentially independent on $q$. A non-zero $\gamma(q = 0, \Omega)$ is impossible in systems with only one type of carriers (either localized or itinerant) because it would violate the spin conservation. It has been conjectured recently that a near-constant $\Gamma(q)$ in UGe$_2$ and UCoGe may be due to the presence of both localized and itinerant electrons in these materials, with ferromagnetism involving predominantly localized spins. We present the microscopic analysis of the damping of near-critical localized excitations due to interaction with itinerant carriers. We show explicitly how the presence of two types of electrons breaks the cancellation between the contributions to $\Gamma(0, \Omega)$ from the self-energy and vertex correction insertions into the spin polarization bubble. We compare our theory with the available experimental data.

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Introduction.— Recent progress in neutron scattering measurements of spin structure factor in uranium metallic materials UGe$_2$ and UCoGe provided detailed information on the dynamical structure factor of a paramagnetic metal near the transition into an itinerant ferromagnet. The dynamical structure factor $S(q, \Omega)$ at wave-vector $q$ and frequency $\Omega$ is related to the dynamic spin susceptibility $\chi(q, \Omega)$ via the relation $S(q, \Omega) = \chi''(q, \Omega) \coth \frac{\Omega}{2T}$. In a conventional paramagnon theory of a nearly ferromagnetic metal $\chi(q, \Omega) \propto 1/(1-i\Omega/\Gamma(q))$ and $S(q, \Omega) \propto \Gamma(q)/|\Omega^2 + \Gamma(q)^2|$ at $q \ll v_F q$, with $\Gamma(q) \propto q$ for clean systems and $\Gamma(q) \propto q^2$ in the presence of non-magnetic impurities.

This behavior has been observed in three-dimensional (3D) electron itinerant ferromagnets such as Ni, Ni$_3$Al, Fe and MnP. Experiments on UGe$_2$ and UCoGe, however, detected a different behavior – near $T_C$, $\Gamma(q)$ extrapolates to a finite value at $q \rightarrow 0$ both in the paramagnetic and the ferromagnetic state.

In one-component fermionic systems the vanishing of $\chi''(q \rightarrow 0, \omega)$ is the consequence of the conservation of the total fermionic spin $S_z$. In particular, $\chi''(q = 0, \Omega) = 0$ holds when the fermion-fermion interaction is mediated by their own collective spin excitations. The situation changes, however, if there are several different bands crossing the Fermi level, or magnetic impurities and/or strong spin-orbit coupling, or if the system is a two-component one and the neutron scattering measures predominantly the contribution from only one of them in some range of $q$.

This last case is applicable to UGe$_2$ and UCoGe because these systems possess simultaneously itinerant and localized ($5f^2$) electrons. In a recent work, Mineev argued that the spin response at small but finite $q$ is predominantly determined by localized spins, because they predominantly contribute to long-range magnetic order and to static susceptibility above the Curie temperature (the contribution from itinerant fermions to magnetic order is around 1%, Ref. [10]) and $\Gamma(q)$ which only weakly depends on $q$. Indeed, the total spin of the localized and itinerant fermions is a conserved quantity, and strictly at $q = 0$ and finite $\Omega$, $S(0, \Omega)$ it must vanish, i.e., the contributions to $S(0, \Omega)$ from localized and from itinerant fermions must cancel out. However, because the static susceptibility of the localized spins is much larger than the one for itinerant fermions, the contribution to $S(q, \Omega)$ from the localized spins dominates down to $v_F q \ll \Omega$. The experiments on UGe$_2$ are performed well inside this range, at $v_F q \geq \Omega$ (Ref. [12]). The total spin of the localized fermions is not a conserved quantity, and the corresponding $\Gamma(q)$ does not have to vanish at $q \rightarrow 0$.

In the present work we address the same issue from a microscopic perspective. We show that the analysis of $\Gamma(q)$ at $q \rightarrow 0$ for the systems with spin-spin interaction is rather non-trivial, as the calculation of a spin-polarization bubble for interacting fermions requires the consideration, on equal footing, of the renormalizations coming from (i) the self-energy diagram, (ii) the Maki-Thompson (MT)-type vertex correction diagram, and (iii) the Aslamazov-Larkin (AL)-type diagrams. The latter have formally one extra power of the coupling and are seemingly less important, but, as we will see later, for one-component systems they are actually of the same order as the other terms. The reason is that the extra power of the coupling gets absorbed into the fermionic damp-
The bubble for free fermions $\chi(q, \Omega)$ contains the total spin susceptibility $\chi^{tot}(q, \Omega) = \chi(q, \Omega)/(1 - U\chi(q, \Omega))$. Each side vertex and each interaction vertex contains spin $\sigma$–matrices (not specified). The conservation of total electronic spin $S = 0$ has been emphasized in Ref. 14, where the authors considered a Fermi liquid with the self-energy $\Sigma(\omega) = \lambda\omega$ and demonstrated that the combination of self-energy, MT, and AL corrections preserves spin conservation. Here we demonstrate this for the more general case, when the self-energy also includes thermal damping $\Sigma^{\prime}(T)$. Further we point out the difference between one- and two-component systems. We argue that for two-component systems, the AL diagrams become irrelevant at weak coupling because the balance between the damping and coupling is lost and the extra power of the coupling is not cancelled out. The remaining self-energy and MT insertions into the spin polarization bubble add up and yield a non-zero $\Gamma(q = 0)$, which is proportional to the imaginary part of the single-particle self-energy of a conduction electron $\Sigma^{\prime}(T)$. We compute $\Sigma^{\prime}(T, \Omega)$ due to interaction with near-critical localized carriers and show that it has a peak at $T_c$, in agreement with the experiments.

**One-component itinerant system.**—To set the stage for the analysis of the two-component systems, consider first a one-component system of itinerant electrons, for which the the total electronic spin $S_i$ is conserved. The conservation of $S_i$ implies that its zero-momentum Fourier component does not depend on time $t$, hence both $\chi(q, \Omega)$ and the dynamical structure factor $S(q, \Omega)$ should vanish at $q = 0$ and any non-zero frequency $\Omega$. As an example (and to simplify the formulas), consider a two-dimensional ($2D$) system with isotropic dispersion $k^2/(2m)$. For non-interacting fermions $\chi_0(q, \Omega) = (m/\pi)(1 - \sqrt{\Omega^2 - v_F q^2})$, and $\chi_0(0, \Omega)$ and $S_0(0, \Omega)$ vanish, as they should. At finite $q$ and at $\Omega \ll v_F q$, $\chi_0(q, \Omega) \approx (m/\pi)(1 + \gamma \Omega/T_0(q))$ with $T_0(q) \propto q^2$ at small $q$. This quantity is non-zero, however it vanishes at $q = 0$. In the presence of disorder, $\Gamma_0(q)$ acquires a diffusive form $\Gamma_0(q) = \sqrt{\nu q^2 + \gamma^2 - \gamma}$, where $\gamma$ is the impurity scattering rate, and it still vanishes at $q = 0$ ($\Gamma_0(q) \propto q^2$ at small $q$). We emphasize that the vanishing of $\Gamma_0(q = 0)$ directly follows from spin conservation, because if $\Gamma_0(0)$ was finite, it would immediately imply a nonzero $S_0(0, \Omega)$, in violation of the spin conservation.

Because the four-fermion interaction $U$ is $SU(2)$ spin invariant, $\Gamma(q = 0)$ should vanish in an interacting system as well. The total susceptibility $\chi^{tot}(q, \Omega)$ is proportional to the fully renormalized spin polarization bubble $\chi(q, \Omega)$ (a fully dressed particle-hole bubble with spin $\sigma$–matrices in the vertices). In the Random Phase Approximation (RPA), $\chi^{tot}(q, \Omega) = \chi(q, \Omega)/(1 - \bar{U}\chi(q, \Omega))$, where $\bar{U} = U/2$. Using the free-fermion form of $\chi(q, \Omega)$ one immediately reproduces the paramagnon formula $S(q, \Omega) \propto \Gamma(q)/(\Omega^2 + \Gamma(q)^2)$, with $\Gamma(q) \propto q$.

The conservation of the total spin implies that the fully renormalized $\chi^{tot}(q, \Omega) \propto \chi(q, \Omega)$ should vanish at $q = 0$. For free fermions, this holds, as we demonstrated above. To go beyond free fermions, we need a model for fermion-fermion interaction. With the uranium compounds in mind, we consider a nearly ferromagnetic metal. Following earlier work, we assume that the low-energy physics is described by an effective model in which the bare $U$ is replaced by an effective dynamical interaction $U_{eff}(q, \omega)$ in the spin channel, which is mediated by $\chi^{tot}(q, \omega)$. In the RPA, $U_{eff}(q, \omega) = \bar{U}^2 \chi^{tot}(q, \Omega)/\Omega = \bar{U}^2 \chi(q, \Omega)/(1 - \bar{U}\chi(q, \Omega))$. Using $\bar{U}\chi(0, 0) \approx 1$, valid near a ferromagnetic transition, this can be simplified to $U_{eff}(q, \omega) \approx \bar{U}/(1 - \bar{U}\chi(q, \Omega))$. We will show how the condition $\chi(q = 0, \Omega) = 0$ is satisfied at each order in $\bar{U}$.

The diagrams for the $\chi(q = 0, \Omega)$ to order $\bar{U}$ are shown in Fig. 1IT1. The first two diagrams contain the self-energy insertions, the third is the MT vertex correction diagram. Each of the three diagrams contain the product of four Green’s function and one dynamical $U_{eff}(q, \omega) \propto \bar{U}$. In explicit form $\chi(q = 0, \Omega) = I_{se}(\Omega) + I_{MT}(\Omega)$ is

\begin{align}
I_{se}(\Omega) &= -6 \int \int G^2(k, \omega)G(k, +q, \omega + \Omega')(G(k, \omega - \Omega) + G(k, \omega + \Omega))U_{eff}(q, \Omega') \\
I_{MT}(\Omega) &= 2 \int \int G(k, \omega)G(k, \omega - \Omega)G(k, +q, \omega + \Omega')G(k + q, \omega - \Omega + \Omega')U_{eff}(q, \Omega')
\end{align}
where the prefactors are due to the summation over spin indices, \(G(k, \omega) = (\omega - \varepsilon_k + i\delta \sigma q\omega)^{-1}\), and we adopt the notation \(\int \equiv \int \frac{d^3k \omega}{(2\pi)^3} \int \frac{d^dq d\Omega'}{(2\pi)^d+} \) for a \(d\)-dimensional system. Applying several times the identity:

\[
G(k, \omega)G(k, \omega - \Omega) = (G(k, \omega - \Omega) - G(k, \omega))/\Omega,
\]

we explicitly re-write the two terms as

\[
I_{se}(\Omega) = 6A(\Omega), \quad I_{MT}(\Omega) = 2A(\Omega)
\]

where

\[
A(\Omega) = \frac{1}{\Omega^2} \int \frac{d^3k \omega}{(2\pi)^3} \int \frac{d^3k \omega_1}{(2\pi)^3} \int \frac{d^3k \omega_2}{(2\pi)^3} \int \frac{d^dq d\Omega'}{(2\pi)^d+} G(k + q, \omega + \Omega') U_{eff}(q, \Omega')
\]

We see that the self-energy and MT diagrams are of the same sign and add up: \(I_{se}(\Omega) + I_{MT}(\Omega) = (4/3)I_{se}(\Omega)\).


The frequency-dependent part of \(I_{se}(\Omega)\) can be evaluated exactly at small \(\Omega\). We obtain \(I_{se}(\Omega) = I_{se}(0) + 3iU m\Omega/(2\pi\Sigma' T)\), where \(\Sigma' T\) is the imaginary part of the fermionic self-energy at zero frequency. If we stop here and associate \((4/3)I_{se}(\Omega)\) with the fully renormalized \(\chi(0, \Omega)\) to order \(U\), we would conclude that \(\Gamma(q = 0)\) becomes finite, in apparent violation of the spin conservation principle. It turns out, however, that there are two other contributions to \(\chi(0, \Omega)\) to order \(U\). They come from the two AL diagrams shown in Fig. 1c. By power counting, these diagrams are of order \(U^2\), but we will see that one power of \(U\) actually cancels out.

The two AL diagrams are equivalent and add up, such that one can consider one of them and multiply the result by 2. In explicit form we have:

\[
I_{AL}(\Omega) = 16 \int \frac{d^3k d\omega}{(2\pi)^3} \int \frac{d^3k_1 d\omega_1}{(2\pi)^3} \int \frac{d^3k_2 d\omega_2}{(2\pi)^3} \int \frac{d^dq d\Omega'}{(2\pi)^d+} U_{eff}(q, \Omega') G(k_1, \omega_1) G(k_2, \omega_2 - \Omega) G(k_2 + q', \omega_1 + \Omega') G(k_2 + q', \omega_2 + \Omega')
\]

where, we recall, \(U_{eff}^{-1} = 1/\bar{U} - \chi(q, \Omega)\) and \(\chi(q', \Omega') = 2 \int G(k, \omega)G(k + q', \omega + \Omega')\). We see that the r.h.s. of Eq. 4 is of order \(U\), not \(U^2\), as one could assume by looking at the l.h.s. of this equation. This cancellation of one power of \(U\) is the natural consequence of the fact that the dispersion of the effective interaction \(U_{eff}(q, \Omega)\) is due to the interaction with the same fermions whose susceptibility we consider.

Substituting Eq. 4 into Eq. 4 we find after some algebra that \(I_{AL}(\Omega)\) has the same form as self-energy and MT contributions and is given by

\[
I_{AL}(\Omega) = -8A(\Omega)
\]

As a result, \(I_{tot}(\Omega) = I_{se}(\Omega) + I_{MT}(\Omega) + I_{AL}(\Omega) = 0\), i.e., \(\Gamma(q)\) vanishes at \(q = 0\), as it indeed should for consistency with the spin conservation principle. For completeness, we analyzed other diagrams which contain \(U^2\) in the prefactor but found that they remain of order \(U^2\), up to logarithmic corrections. Only in the two AL diagrams one power of \(U\) is cancelled out.

Two-component systems.— We now analyze how this result changes when we consider a two-component system consisting of localized and itinerant electrons. As we discussed in the Introduction, we focus on the range \(v_F q \sim \Omega\) where the full dynamical susceptibility almost coincides with the one for localized spins, \(\chi_L(q, \Omega)\). By itself (i.e., with no itinerant fermions present) \(\chi_L(0, \Omega)\) vanishes. We consider how it is modified due to the interaction with the itinerant electrons.

We use the same model as before with the effective spin-fermion coupling

\[
H = \Phi \sum_q s_q \cdot S_{-q}
\]

where \(s_q = \sum_{k, \alpha, \beta} c_{k, \alpha}^\dag c_{k, \beta}^\dag \sigma_{\alpha, \beta} c_{k+q, \alpha, \beta}\) with \(c^\dag_{k, \alpha}, c_{k, \alpha}\) being the creation and annihilation operators for itinerant electrons, \(\sigma_{\alpha, \beta}\) are Pauli matrices, and \(S_q\) describes the localized spins. This interaction gives rise to the correction to the susceptibility of the localized spins \(\chi_L(0, \Omega) = \chi_L(0, \Omega) + I_{tot}(\Omega)\), where \(I_{tot}(q, \Omega)\) is the fully renormalized spin-polarization bubble of itinerant fermions. Simultaneously, Eq. 7, taken to second order, gives rise to effective interaction between itinerant carriers, mediated by the localized spins: \(\Phi_{eff}(q, \Omega) \sim \Phi^2 \chi_L(q, \Omega)\), like in Fig. 1a-c. The crucial difference with the previous case is that now the localized spins have their own dynamics even in the absence of the interaction with itinerant carriers. This dynamics is consistent.
with the conservation of the total spin of localized carriers, e.g., in a paramagnetic state it is spin diffusion: 
\[ \chi_L(q, \Omega) = \chi_0/(q^2 + \xi^2 - i\Omega/(Dq^2)). \]
However, when we include the spin-spin interaction between localized and itinerant carriers, we find that the contributions from AL diagrams no longer cancel out the contribution from self-energy and MT terms because the difference 
\[ \chi_L^{-1}(q', \Omega') - \chi_L^{-1}(q', \Omega + \Omega) \]
in the analogue of Eq (9) is non-zero even when \( \Phi = 0 \). As the consequence, the extra \( \Phi^2 \) in the AL diagrams does not cancel out, and the AL contribution \( I_{AL}(\Omega) \) becomes small compared to self-energy and MT terms. The sum of these two is, according to Eq. (2), \( (4/3)I_{se}(\Omega) = A + iB\Omega/\Sigma''(T) \), where \( A \) and \( B \) are constants. Then, to leading order in \( \Phi \),
\[ \chi_L^{\text{tot}}(q \to 0, \Omega) = \frac{4}{3}I_{se}(\Omega) = A + iB\frac{\Omega}{\Sigma''(T)} \]  
(8)
i.e., the system has a non-zero \( \Gamma_L(q \to 0) \propto \Sigma''(T) \). This result holds for both 2D and 3D systems.

At small but finite \( q \) and at \( \Omega < Dq^2 \), the static part of \( \chi_L(0, \Omega) \) can be approximated by \( \chi_0\xi^2 \), and once the damping due to interaction with itinerant carriers exceeds the diffusion term, the susceptibility and the structure factor become
\[ \chi_L^{\text{tot}}(q, \Omega) \approx \frac{\chi_0\xi^2}{1 - \frac{\Omega}{\Gamma_L(q)}}; \quad S(q, \Omega) \sim \frac{\chi_0\xi^2\Gamma_L(q)}{\Omega^2 + \Gamma_L^2(q)} \]  
(9)
where \( \Gamma_L(q) \approx \Gamma_L(q = 0) \propto \Phi^2\Sigma''(T)\xi^{-2} \).

The physical reason for a non-zero \( \Gamma_L(q = 0) \) is the fact that, when itinerant and localized electrons interact via spin-spin coupling \( S \cdot S \), only the total combined spin is conserved, while the localized spin \( S \) can change its \( z \)-component and transfer the difference to the spin \( S \) of an itinerant electron. This reasoning parallels the one presented by Mineev \cite{1}. He also found \( \Gamma_L(q = 0) \propto \xi^{-2} \).

Our microscopic consideration expresses \( \Gamma_L(q = 0) \) in terms of the coupling \( \Phi \) and the temperature dependent fermionic self-energy of the itinerant electrons \( \Sigma''(T) \). This allows us to proceed further with the analysis of the temperature dependence of \( \Gamma(0) \) and the comparison with the experimental data.

Fermionic self-energy.—The diagram for the fermionic self-energy is presented in Fig. 1d. The wavy line is the propagator of localized spins \( \chi_L \), which we assume to have a diffusive form. Using the spectral representation of \( G(k, \omega) \) and \( \chi_L(k + q, \omega) \), we obtain
\[ \Sigma''(T) \propto \Phi^2 \int \frac{d \Omega \Omega}{\sinh \frac{\Omega}{2T}} \int d^{d-1}q \frac{Dq^2}{D^2q^4(\xi^{-2} + q^2)^2 + \Omega^2} \]  
(10)
At some distance away from the transition point, \( \Sigma''(T) \propto T^{3/2} \) in 2D and \( (T^2/E_F) \log E_F/T \) in 3D. This dependence holds both in the paramagnetic phase (where Eq. (10) is valid), and in the ferromagnetic state. Right at the transition point, \( T = T_c \), \( \xi = \infty \), \( \Sigma''(T) \) is enhanced: a self-consistent solution yields \( \Sigma''(T) \propto T \log T \) in 3D and \( \Sigma''(T) \propto T^{4/5} \) in 2D. As the result, \( \Gamma_L(T)\xi^2 \propto \Sigma''(T) \) scales as some power of \( T \) below and above the transition, but get an enhancement very near the transition point. In Fig. 2 we compare our theory with the experimental data for \( \Gamma_l(q, T)\xi^2 \) as a function of \( T \) in \( UGe_2 \) (Ref. 1). We recall that we identify \( \Gamma_L(q, T) \) with the measured \( \Gamma_l(q, T) \) over the range \( v_F q \geq \Omega \) probed by the experiments. At asymptotically small \( v_F q \ll \Omega \), \( \Gamma_L(q, T) \) should indeed vanish. The data are consistent with our result that there is a smooth increase of \( \Gamma_l(T)\xi^2 \) with \( T \) in both the ferromagnetic and paramagnetic state, on top of which there is a peak at \( T_c \).

Conclusion.—To conclude, in this paper we presented the microscopic study of the damping term of spin excitations in a metal near a ferromagnetic instability. We demonstrated that in an one-component systems, the spin scattering rate \( \Gamma_l(q) \) vanishes at \( q = 0 \) as the consequence of the spin conservation. We argued that to see this in a loop expansion, one needs to invoke AL scattering processes. We then considered a two-component model with localized and itinerant fermions, which was argued to describe ferromagnetic uranium compounds such as \( UGe_2 \) and \( UCoGe \). Localized spins mostly contribute to long-range order and to Curie susceptibility and for \( v_F q \geq \Omega \), probed by the experiments, the measured damping rate \( \Gamma_l(q, T) \) almost coincides with \( \Gamma_L(q, T) \) for localized spins. We showed that for \( \Gamma_L(q, T) \) the AL diagrams are relatively small. Without the AL contribution, the spin scattering rate \( \Gamma_L(q = 0) \) becomes finite and scales with the temperature dependent part of fermionic self-energy \( \Sigma''(T) \) for itinerant fermions. We found that \( \Gamma_L(0, T)\xi^2 \) has a peak at \( T_c \). This is consistent with the data on \( UGe_2 \).

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12. For charge interaction $I_{se}(\Omega) + I_{MT}(\Omega) = 0$ and the two AL diagrams also cancel each other.
SUPPLEMENTAL MATERIAL.

Cancellation of the diagrams of the spin polarization bubble at zero momentum and a finite frequency

A. Charge channel

We show how self-energy and Maki-Thompson (MT) vertex correction insertions into the particle-hole bubble cancel each other in the charge channel. The diagrams are shown in Fig. 1 in the text. For charge channel, side vertices are spin delta-functions rather than $\sigma$-matrices. For convenience, we perform calculations in Matsubara frequencies.

The expressions for the two self-energy diagrams in Fig. 1b are:

$$I_{se}^A(\Omega) = 3 \int \frac{d^dq d\Omega'}{(2\pi)^{d+1}} G^2(k, \omega) G(k + q, \omega + \Omega') G(k + q, \omega - \Omega) U_{\text{eff}}(q, \Omega')$$ (1)

$$I_{se}^B(\Omega) = 3 \int \frac{d^dq d\Omega'}{(2\pi)^{d+1}} G^2(k, \omega) G(k + q, \omega + \Omega') G(k + q, \omega + \Omega) U_{\text{eff}}(q, \Omega')$$ (2)

where the factor of 3 is due to summation over spin indices, the Green’s functions are $G(\vec{k}, \omega) = \frac{1}{\omega - \epsilon(\vec{k})}$, and the expression for $U_{\text{eff}}(q, \Omega)$ in given in the main text: $U_{\text{eff}}(q, \omega) \approx \tilde{U} / (1 - \tilde{U} \chi(q, \Omega))$. The $\chi(q, \Omega)$ in this expression is the fully renormalized spin-polarization bubble, but to lowest order in the coupling $\tilde{U}$ it can be replaced by the spin susceptibility of free fermions.

By using several times the known identity:

$$G_1 G_2 = (G_1 - G_2) \times \frac{1}{G_2^{-1} - G_1^{-1}}$$ (3)

we obtain:

$$I_{se}^A(\Omega) = \frac{3}{i\Omega} \int \frac{d^dq d\Omega'}{(2\pi)^{d+1}} G(k + q, \omega + \Omega') G(k, \omega) U_{\text{eff}}(q, \Omega') [G(k, \omega) - G(k, \omega - \Omega)] \frac{1}{(-i\Omega)}$$

$$= -\frac{3}{i\Omega} \int \frac{d^dq d\Omega'}{(2\pi)^{d+1}} G^2(k, \omega) G(k + q, \omega + \Omega') U_{\text{eff}}(q, \Omega')$$

$$+ \frac{3}{\Omega^2} \int \frac{d^dq d\Omega'}{(2\pi)^{d+1}} G(k + q, \omega + \Omega') U_{\text{eff}}(q, \Omega') [G(k, \omega) - G(k, \omega - \Omega)]$$ (4)

$$I_{se}^B(\Omega) = \frac{3}{i\Omega} \int \frac{d^dq d\Omega'}{(2\pi)^{d+1}} G(k + q, \omega + \Omega') G(k, \omega) U_{\text{eff}}(q, \Omega') [G(k, \omega) - G(k, \omega + \Omega)] \frac{1}{(-i\Omega)}$$

$$= \frac{3}{i\Omega} \int \frac{d^dq d\Omega'}{(2\pi)^{d+1}} G^2(k, \omega) G(k + q, \omega + \Omega') U_{\text{eff}}(q, \Omega')$$

$$+ \frac{3}{\Omega^2} \int \frac{d^dq d\Omega'}{(2\pi)^{d+1}} G(k + q, \omega + \Omega') U_{\text{eff}}(q, \Omega') [G(k, \omega) - G(k, \omega + \Omega)]$$ (5)

Therefore the total contribution from the two diagrams is:

$$I_{se}(\Omega) = I_{se}^A(\Omega) + I_{se}^B(\Omega) = \frac{3}{\Omega^2} \int \frac{d^dq d\Omega'}{(2\pi)^{d+1}} [2G(k, \omega) - G(k, \omega + \Omega) - G(k, \omega - \Omega)] G(k + q, \omega + \Omega') U_{\text{eff}}(q, \Omega')$$ (6)

Similarly, the vertex-correction MT diagram reads:

$$I_{MT} = 3 \int \frac{d^dq d\Omega'}{(2\pi)^{d+1}} G(k, \omega) G(k + q, \omega + \Omega') G(k + q, \omega - \Omega + \Omega') U_{\text{eff}}(q, \Omega')$$
Using (3) we rewrite it as
\[ I_{MT}(\Omega) = -\frac{3}{\Omega^2} \int \frac{d^d q d\Omega'}{(2\pi)^{d+1}}[G(k, \omega) - G(k, \omega - \Omega)][G(k + q, \omega + \Omega') - G(k + q, \omega - \Omega + \Omega')] U_{eff}(q, \Omega') \]
\[ = -\frac{3}{\Omega^2} \int \frac{d^d q d\Omega'}{(2\pi)^{d+1}}[G(k, \omega) - G(k, \omega - \Omega)] G(k + q, \omega + \Omega') U_{eff}(q, \Omega') \]
\[ + \frac{3}{\Omega^2} \int \frac{d^d q d\Omega'}{(2\pi)^{d+1}}[G(k, \omega) - G(k, \omega - \Omega)] G(k + q, \omega - \Omega + \Omega') U_{eff}(q, \Omega') \]
\[ (7) \]
Shifting the frequency: \( \omega' = \omega - \Omega \) in the last line and then adding the two terms, we finally obtain:
\[ I_{MT} = -\frac{3}{\Omega^2} \int \frac{d^d q d\Omega'}{(2\pi)^{d+1}} \left[ 2G(k, \omega) - G(k, \omega + \Omega) - G(k, \omega - \Omega) \right] G(k + q, \omega + \Omega') U_{eff}(q, \Omega') \]
\[ (8) \]
Adding \( I_{se} \) and \( I_{MT} \) together, we obtain in the charge channel:
\[ I_{se}(\Omega) + I_{MT}(\Omega) = 0 \]
\[ (9) \]
There are also two Aslamazov-Larkin (AL) diagrams (Fig. 1c). They are identical, up to sign. For charge channel (side vertices with spin \( \delta \)-functions) signs are different, and the the two AL diagram cancel each other. Then the total \( I_{se}(\Omega) + I_{MT}(\Omega) + I_{AL}(\Omega) = 0 \), as it is required by charge conservation.

**B. Spin Channel**

For spin channel, the side vertices contain spin \( \sigma \)-matrices \( \sigma_{\alpha\beta} \). The overall sign in \( I_{se} \) is now 6, while the overall sign in \( I_{MT} \) is (-2). Then \( I_{MT}(\Omega) = 1/3 I_{se}(\Omega) \) and \( I_{se}(\Omega) + I_{MT}(\Omega) = (4/3) I_{se}(\Omega) \). In explicit form
\[ \frac{4}{3} I_{se} = -\frac{8}{\Omega^2} \int \frac{d^d q d\Omega'}{(2\pi)^{d+1}} \left[ 2G(k, \omega) - G(k, \omega + \Omega) - G(k, \omega - \Omega) \right] G(k + q, \omega + \Omega') U_{eff}(q, \Omega') \]
\[ (10) \]
Simultaneously, however, the two AL diagrams add up instead of canceling out. The total contribution from the two AL diagrams is:
\[ I_{AL}(\Omega) = 16 \int \frac{d^d k d\omega}{(2\pi)^{d+1}} \int \frac{d^d q d\Omega'}{(2\pi)^3} \int \frac{d^d p d\omega'}{(2\pi)^{d+1}} G(k, \omega) G(k, \omega - \Omega) G(p, \omega') G(k + q, \omega + \Omega') \]
\[ \times \left[ G(p, \omega' + \Omega) G(p - q, \omega' - \Omega') + G(p, \omega' - \Omega) G(p + q, \omega + \omega' + \Omega') \right] U_{eff}(q, \Omega') U_{eff}(q, \Omega + \Omega') \]
\[ (11) \]
The product of the two effective interactions can be rewritten in the form:
\[ U_{eff}(q', \Omega') U_{eff}(q', \Omega' + \Omega) = \frac{U_{eff}(q', \Omega') - U_{eff}(q', \Omega' + \Omega)}{\chi(q', \Omega' + \Omega) - \chi(q', \Omega')} \]
\[ U_{eff}(q', \Omega' + \Omega) = \frac{U_{eff}(q', \Omega') - U_{eff}(q', \Omega')}{{\chi(q', \Omega' + \Omega) - \chi(q', \Omega')}} \]
\[ (12) \]
where, we recall, \( \chi(q', \Omega') = 2 \int G(k, \omega) G(k + q', \omega + \Omega') \). Note that the resulting expression is of first order in \( U_{eff} \). Physically, this is the consequence of the fact the dynamics of collective excitations comes solely from the interaction with fermions (i.e., that there is no damping of a collective boson if \( \tilde{U} = 0 \)).
Using (12) for \( U_{\text{eff}}(\mathbf{q}', \Omega')U_{\text{eff}}(\mathbf{q}', \Omega' + \Omega) \) and using the identity Eq. (3) several times, we obtain:

\[
I_{AL}(\Omega) = 16 \int \int \int G(\mathbf{k}, \omega)G(\mathbf{k}, \omega - \Omega)G(\mathbf{p}, \omega')G(\mathbf{p}, \omega' - \Omega)G(\mathbf{k} + \mathbf{q}, \omega + \Omega')G(\mathbf{p} + \mathbf{q}, \omega + \Omega') \\
\times [U_{\text{eff}}(\mathbf{q}, \Omega') - U_{\text{eff}}(\mathbf{q}, \Omega' + \Omega)] \frac{1}{2u \left[ \Pi(\mathbf{q}, \Omega' + \Omega) - \Pi(\mathbf{q}, \Omega') \right]} \\
= -\frac{16}{\Omega^2} \int \int \int [G(\mathbf{k}, \omega) - G(\mathbf{k}, \omega - \Omega)] [G(\mathbf{p}, \omega') - G(\mathbf{p}, \omega' - \Omega)] G(\mathbf{k} + \mathbf{q}, \omega + \Omega')G(\mathbf{p} + \mathbf{q}, \omega + \Omega') \\
\times [U_{\text{eff}}(\mathbf{q}, \Omega') - U_{\text{eff}}(\mathbf{q}, \Omega' + \Omega)] \frac{1}{2u \left[ \Pi(\mathbf{q}, \Omega' + \Omega) - \Pi(\mathbf{q}, \Omega') \right]} \\
-\frac{16}{\Omega^2} \int \int \int [G(\mathbf{k}, \omega)G(\mathbf{p}, \omega') - G(\mathbf{k}, \omega - \Omega)G(\mathbf{p}, \omega') - G(\mathbf{k}, \omega)G(\mathbf{p}, \omega' - \Omega) + G(\mathbf{k}, \omega - \Omega)G(\mathbf{p}, \omega' - \Omega)] \\
\times G(\mathbf{k} + \mathbf{q}, \omega + \Omega')G(\mathbf{p} + \mathbf{q}, \omega' + \Omega') [U_{\text{eff}}(\mathbf{q}, \Omega') - U_{\text{eff}}(\mathbf{q}, \Omega' + \Omega)] \frac{1}{2u \left[ \Pi(\mathbf{q}, \Omega' + \Omega) - \Pi(\mathbf{q}, \Omega') \right]}
\]

where we used short-hand notation \( \int \int \equiv \int \frac{d^d k d\omega}{(2\pi)^{d+1}} \int \frac{d^d q d\Omega'}{(2\pi)^{d+1}} \int \frac{d^d p d\omega'}{(2\pi)^{d+1}} \).

At this stage we can separate the integration \( \int \frac{d^d p d\omega'}{(2\pi)^{d+1}} \) of a product of two Green’s functions in each term and rewrite the last equation as

\[
I_{AL}(\Omega) = -\frac{8}{\Omega^2} \int \int \frac{[U_{\text{eff}}(\mathbf{q}, \Omega') - U_{\text{eff}}(\mathbf{q}, \Omega' + \Omega)]}{2u \left[ \Pi(\mathbf{q}, \Omega' + \Omega) - \Pi(\mathbf{q}, \Omega') \right]} [G(\mathbf{k}, \omega)G(\mathbf{k} + \mathbf{q}, \omega + \Omega')\Pi(\mathbf{q}, \Omega') + G(\mathbf{k}, \omega - \Omega)G(\mathbf{k} + \mathbf{q}, \omega + \Omega') \\
\times \Pi(\mathbf{q}, \Omega' + \Omega) - G(\mathbf{k}, \omega - \Omega)G(\mathbf{k} + \mathbf{q}, \omega + \Omega')\Pi(\mathbf{q}, \Omega') - G(\mathbf{k}, \omega)G(\mathbf{k} + \mathbf{q}, \omega + \Omega')\Pi(\mathbf{q}, \Omega' + \Omega)] \\
= \frac{8}{\Omega^2} \int \int \frac{[U_{\text{eff}}(\mathbf{q}, \Omega') - U_{\text{eff}}(\mathbf{q}, \Omega' + \Omega)]}{2u} [G(\mathbf{k}, \omega)G(\mathbf{k} + \mathbf{q}, \omega + \Omega') - G(\mathbf{k}, \omega - \Omega)G(\mathbf{k} + \mathbf{q}, \omega + \Omega')] \\
= \frac{8u^2}{\Omega^2} \int \int U_{\text{eff}}(\mathbf{q}, \Omega')G(\mathbf{k} + \mathbf{q}, \omega + \Omega') [2G(\mathbf{k}, \omega) - G(\mathbf{k} + \mathbf{q}, \omega + \Omega') - G(\mathbf{k}, \omega - \Omega')]
\]

(13)

where now \( \int \int \equiv \int \frac{d^d k d\omega}{(2\pi)^{d+1}} \int \frac{d^d q d\Omega'}{(2\pi)^{d+1}} \), and the last expression is obtained by shifting the frequency variable \( \Omega' + \Omega \) to \( \Omega' \) in \( U_{\text{eff}}(\mathbf{q}, \Omega' + \Omega) \) and in the corresponding Green’s functions. Comparing this expression with Eq. (10) we immediately see that

\[
I_{AL}(\Omega) = -\frac{4}{3} I_{se}(\Omega),
\]

(14)

such that the total contribution to the spin polarization bubble

\[
I_{se}(\Omega) + I_{MK}(\Omega) + I_{AL}(\Omega) = 0,
\]

(15)
as it should be due to spin conservation.