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2 - this energy scale is called the Kondo temperature, T_K [setting $k_B = 1$]

using $\lambda(E) = \frac{\lambda_0}{1 - \lambda_0 \ln \frac{D_0}{E}}$

we can estimate T_K , using $\lambda(T_K) \approx 1$

$$1 \approx \frac{\lambda_0}{1 - \lambda_0 \ln \frac{D_0}{T_K}} = \frac{1}{\frac{1}{\lambda_0} - \ln \frac{D_0}{T_K}}$$

$$\frac{1}{\lambda_0} \approx \ln \frac{D_0}{T_K}$$

$$T_K \approx D_0 e^{-\frac{1}{\lambda_0}} \quad [\text{up to a factor of } 1(1)]$$

in this rough estimate we could take D_0 to be the original band width or E_F and

λ_0 to be the original "bare" Kondo coupling

$$\sim E_F, \quad T_K \approx E_F \exp\left[-\frac{1}{J \nu}\right]$$

- exponentially small compared to E_F

↳ like 1-QCD - strong interaction coupling

constant also grows with increasing E

- RG improved pert. theory valid for $E \gg T_K$, but at $E \lesssim T_K$ we enter strong-coupling regime where pert. theory completely fails (even when improved by RG)

- it is convenient to write $\lambda(E) \approx \frac{1}{\ln E/T_K}$
for $E \gg T_K$ - elementary bare coupling
we replace it by T_K

- RG improved P.T. can be applied to resistivity as I will now show

- lowest order $R(T) \propto N_{\text{imp}} \lambda_0^2$ where

N_{imp} = impurity concentration - NB - this is

just contribution to $R(T)$ from scattering

off magnetic impurities - must be added

to other contributions

- now if we RG-improve calculation we

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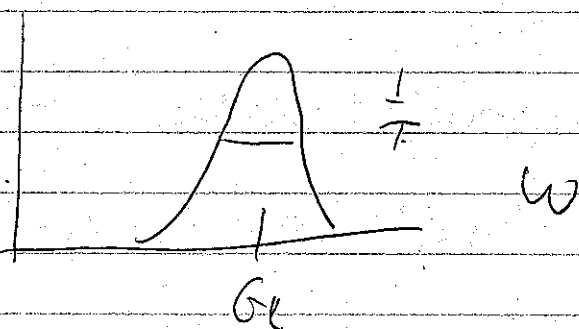
get $R(T) \propto \frac{n_{imp}}{\ln^2 T/T_K}$ for $T \gg T_K$

- this fits experimental data (e numerical calculations of (Kondo model) well at $T \gg T_K$
- N.B. - we cannot use this result as $T \rightarrow T_K$ where it predicts $R \rightarrow \infty$
- I will return to question of what happens at lower T , but first go thru calculation of $R(T)$ to 2nd order in T in some detail - opportunity to learn some other techniques of CMP
- see Hewson, chp 2, Mahan 11.1, AEM chp 13
- the plan is to calculate correction to electron Green's function

$$G_{ret}(\vec{k}, \omega) \propto \frac{1}{\omega - \epsilon_{\vec{k}} + \frac{i}{2T}}$$

- this broader single-electron peak

- In GRET



width = $\frac{1}{T}$, ~~Thus~~ T is time electron spends
in a state of momentum $\vec{k}$ before scattering
(off an impurity)

- we can insert this result in Boltzmann
equation result for resistivity

$$\sigma = \frac{1}{R} = \frac{2e^2}{3} V_F^2 V T$$

where V = density of states at Fermi surface

- Phys 502? - see Ac M

- Boltzmann equation approach makes it
clear that scattering rate at energy

$|G_L| \lesssim T$ contribute to resistivity

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at temperatures T

- using free electron dispersion relation this

reduces to $\frac{1}{R} = \frac{ne^2}{m} T$ - classical Drude formula

- most rigorous derivation of this formula is

from Kubo formula - can express retarded

GF for current operator in terms of

GF for electron operator in this case

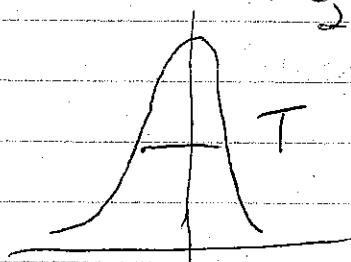
[AEM (13.25)]

$$\sigma^{ab} = e^2 \int \frac{d^3 k}{4\pi^3} T[\mathcal{E}_k] V_k^a V_k^b \left(-\frac{2f}{2\epsilon} \right) \epsilon_k^2$$

- for isotropic dispersion relation can replace

$$V_k^a V_k^b \rightarrow \frac{1}{3} \delta^{ab} V_k^2 \text{ inside integral}$$

$$(V_k^2 = \nabla \mathcal{E}_k^2)$$



T contribution to σ for $\epsilon_k \lesssim T$
= 2nd half of Phys 502

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- Calculation of Electron GF - $\langle \psi_{\vec{k}} \bar{\psi}_{\vec{k}'} \rangle$

$$\begin{array}{c} \omega, \vec{k} \quad \omega, \vec{k}' \\ \text{---} \textcircled{\text{---}} \text{---} = \text{---} + \text{---} \text{---} \text{---} + \dots \end{array}$$

$\omega, \vec{k} \quad \omega, \vec{k}' \quad \omega, \vec{k}$

- we will average over positions of defects

Random array of impurities - restores translation

- invariance and makes G a function of

energy unless $\vec{k} = \vec{k}'$

- similar to our calculation of χ except

that now we sum over impurity state

and over all momenta in action, not

just in strip $D' < E_{\vec{k}} < D$ as for RGr

- NB - no $O(\hbar)$ contribution due to $\vec{S} = 0$

$\mathcal{H}_0$ - convenient to work in position frequency

space

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$$G_{2\beta}(\vec{r}, \vec{r}'; i\omega_n) = - \langle \Psi_2(\vec{r}, \omega_n) \bar{\Psi}_\beta(\vec{r}', \omega_n) \rangle$$

where $\langle \rangle$ includes trace over impurity spin states

$$\mathcal{H} = - \frac{J^2}{2!} \mathcal{S}$$

$$\mathcal{H}_{WT} = - \frac{J^2}{2!} \int_{-\infty}^{\infty} d\tau \bar{\Psi}(\vec{0}, \tau) \frac{\sigma^a}{2} \Psi(\vec{0}, \tau) \cdot \vec{S}(\tau)$$

$$\mathcal{H}_{2\beta} = - \frac{J^2}{2!} \sum_{ab} \langle \bar{\Psi} \left(\frac{\sigma^a}{2} \Psi \right) \left(\frac{\sigma^b}{2} \Psi \right) \bar{\Psi} \rangle \langle S^a S^b \rangle$$

$$\langle S^a S^b \rangle = \frac{1}{4} S^a S^b \quad \text{no } T\text{-dependence.}$$



diagram corresponds to a correction to ground state energy

only need consider

$$\bar{\Psi} \bar{\Psi} \sigma^a \Psi \bar{\Psi} \sigma^b \Psi \cdot \bar{\Psi} + \text{equivalent diagram}$$

(factor of 2)

- Contraction of electron spin indices gives

$$\frac{\sigma^a}{2} \frac{\sigma^b}{2} = \frac{(\sigma^a \sigma^b)}{4}$$

sum over multiply

$$\sum_{ab} \left(\frac{\sigma^a}{2} \frac{\sigma^b}{2} \right)_{\alpha\beta} \frac{1}{4} f^{ab} = \frac{3}{16} \delta_{\alpha\beta}$$

[using $(\sigma^a)^2 = I$]

using $G^{(0)}(\vec{r} = \vec{r}', i\omega_n) = - \langle \Psi(\vec{r}, \omega_n) \bar{\Psi}(\vec{r}', \omega_n) \rangle$

we obtain

$$\delta G_{\alpha\beta}(\vec{r}, \vec{r}', i\omega_n) = \frac{3}{16} \delta_{\alpha\beta} T^2$$

$$G^{(0)}(\vec{0}, i\omega_n) G^{(0)}(\vec{0}, i\omega_n) G^{(0)}(-\vec{r}', i\omega_n)$$

- frequency is conserved since no T -dependence arises from empty spin operator in this case.

$$G^{(0)}(\vec{0}, i\omega_n) = \frac{1}{V} \sum_{\vec{k}} \frac{1}{i\omega_n - \epsilon_{\vec{k}}}$$

- return to evaluating this later

- convenient to write

$$\delta G(\vec{r}, \vec{r}', i\omega_n) = G^{(0)}(\vec{r}, i\omega_n) T(i\omega_n) G^{(0)}(-\vec{r}', i\omega_n)$$

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(suppressing $\vec{r}, \vec{r}'$)

$$\text{where } T(i\omega_n) = \frac{3}{16} J^2 \frac{1}{V} \sum_{\vec{k}} \frac{1}{i\omega_n - \epsilon_{\vec{k}}}$$

N.B. T is a function of frequency only,
not position (or momentum)

- it is called the T -matrix and is a generalization
of the T -matrix in ordinary potential
scattering ($S = I + i T$)

- suppose impurity was not at origin, but
at arbitrary location $\vec{R}$

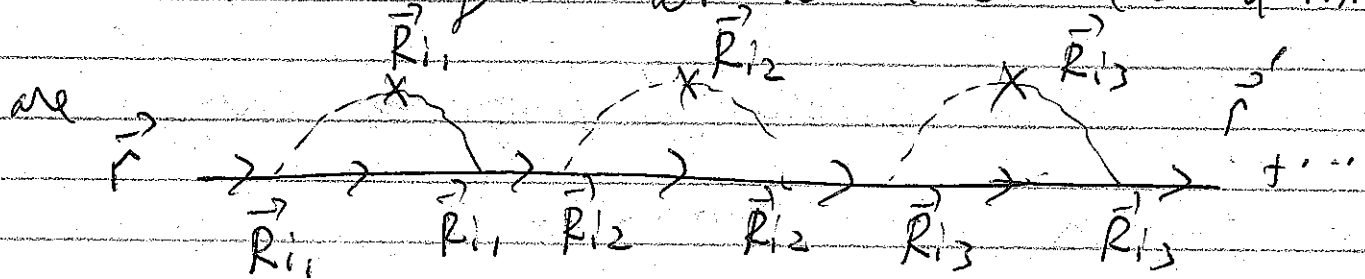
$$\text{then } S_M(\vec{r}, \vec{r}', i\omega_n) = G^0(\vec{r} - \vec{R}, i\omega_n) T(i\omega_n)$$

$$\text{due to } \overbrace{\psi(\vec{r}) \bar{\psi}(\vec{R})}^{G^0(\vec{r} - \vec{R}, i\omega_n)} \overbrace{\psi(\vec{R}) \bar{\psi}(\vec{r}')}^{\text{}} \quad \overbrace{\psi(\vec{R}) \bar{\psi}(\vec{r}')}^{\text{}}$$

- now suppose we have many impurities
at locations $\vec{R}_1, \vec{R}_2, \dots$

$$H_{int} = T \sum_{i=1}^{N_{imp}} \psi^\dagger(\vec{R}_i) \frac{\sigma}{2} \psi(\vec{R}_i) \cdot \vec{J}_i$$

- dominant diagrams at small T and small N_{imp}



$$= g^{(0)}(\vec{r} - \vec{R}_{i_1}) T g^{(0)}(\vec{R}_{i_1} - \vec{R}_{i_2}) T g^{(0)}(\vec{R}_{i_2} - \vec{R}_{i_3}) T \dots T g^{(0)}(\vec{R}_{i_n} - \vec{r}')$$

where $n = 0, 1, 2, 3, \dots, \infty$

- all $g^{(0)}$'s and T 's at same frequency, $i\omega_n$

- now sum over $i_1, i_2, i_3, \dots$ in from 1 to N_{imp}

- in dilute limit we can ignore diagrams where same impurity occurs more than once

in a diagram

$$\sum_{i_p=1}^{N_{imp}} \rightarrow N_{imp} \int d^3 \vec{R}_p$$

since impurity locations are random and since we will also be integrating over $\vec{r}$ and $\vec{r}'$ to get resistivity

n^{th} order correction is

$$[n_{\text{imp}} T(i\omega_n)]^n \int d^3 \vec{r}_1 \dots d^3 \vec{r}_n$$

$$J^{(0)}(\vec{r} - \vec{r}_1) J^{(0)}(\vec{r}_1 - \vec{r}_2) \dots J^{(0)}(\vec{r}_n - \vec{r}') J^{(0)}(\vec{r}_n - \vec{r}')$$

- NB - translation invariance is restored by averaging over impurity locations

- in k -space

$$S(\vec{k}, i\omega_n) = (n_{\text{imp}} T(i\omega_n))^n [J^{(0)}(\vec{k}, i\omega_n)]^{n+1}$$

- sum series

$$J(\vec{k}, i\omega_n) = \sum_{n=0}^{\infty} [J^{(0)}(\vec{k}, i\omega_n)]^{n+1} [n_{\text{imp}} T(i\omega_n)]^n$$

$$= \frac{J^{(0)}}{1 - J^{(0)} n_{\text{imp}} T} = \frac{1}{J^{(0)-1} - n_{\text{imp}} T}$$

$$= \frac{1}{i\omega_n - E_{\vec{k}} - n_{\text{imp}} T(i\omega_n)}$$

- this has the form of a self-energy

$$G = \frac{1}{i\omega_n - \epsilon_F - \Sigma(i\omega_n, \vec{k})}$$

where self-energy depends on $i\omega_n$ only, not $\vec{k}$

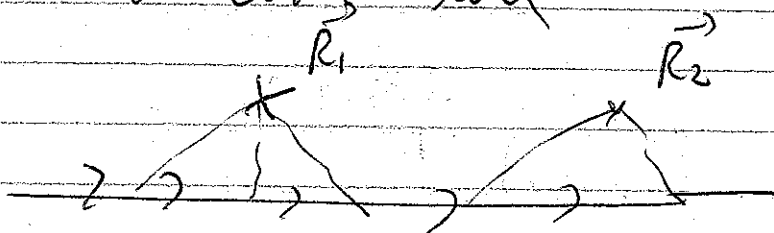
$$\Sigma(\vec{k}, i\omega_n) = N_{\text{imp}} T(i\omega_n)$$

- we actually want retarded G^R &

define $T(\omega)$ which appears in σ resistivity

- other diagrams?

- correcting like

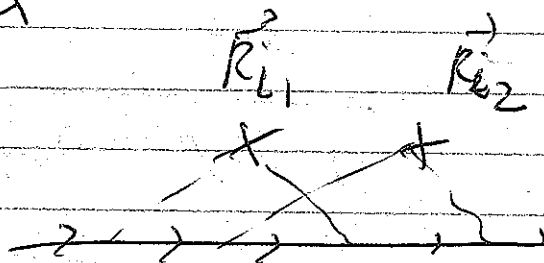


still give $\Sigma = N_{\text{imp}} T(i\omega_n)$ but

give corrections to $T(i\omega_n)$ of higher

order in λ

diagram like



mix different impurities - now ~~drop~~

sum over $i_1, i_2 = 1 \dots N_{\text{imp}}$ can't be done

independently - give corrections to Σ of

higher order in N_{imp} [vertex expansion]

- we ignore them in dilute limit

- finally let's evaluate $T(\omega)$ to $O(J^2)$

$$T(\omega) = \frac{3}{16} \frac{J^2}{V} \sum_{\vec{k}} \frac{1}{\omega - \epsilon_{\vec{k}} + i\eta} \quad (\text{retarded GF})$$

$\text{Re } T$ just shifts Fermi energy a bit [$O(N_{\text{imp}})$]

and leads to other unimportant corrections

$\text{Im } T$ gives scattering rate

$$\text{Im } T = - \frac{3}{16} J^2 \pi \int \frac{d^3 k}{(2\pi)^3} \delta[\omega - \epsilon_{\vec{k}}]$$

$$= - \frac{3}{16} J^2 \pi \nu(\omega)$$

$$\approx - \frac{3}{16} J^2 \pi \nu \quad \text{at } \omega \rightarrow 0$$

$$\text{Im } \Sigma = - \frac{3}{16} \nu J^2 \pi N_{\text{imp}} = - \frac{1}{2\tau}$$

- NB - has correct sign - pole of G_{ret} in negative half-plane

$$\frac{I}{T} = \frac{3}{2} \pi T^2 V \cdot \Pi_{imp}$$

- plug into Kubo formula [or Boltzmann equation result in AEM]

$$\rho = \frac{3}{2e^2 V_F^2 V} \times \frac{I}{T} = \frac{9\pi}{16e^2 V_F^2} T^2 \Pi_{imp}$$

- explicit calculation is $O(T^3)$ gives

$$\rho = \frac{9\pi}{16e^2 V_F^2 V^2} \left[\lambda_0^2 + 2\lambda_0^3 \ln \frac{R_0}{T} + \dots \right]$$

$$\approx \frac{9\pi}{16e^2 V_F^2 V^2} \left[\lambda(T)^2 + \dots \right]$$

- i.e. we get RG-improved result

as we expect - T provides scale at

which to evaluate effective coupling

- correction is for $\frac{R_1}{2}$

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- an ∞ series of diagrams correspond to RG result

$$\Rightarrow f \propto \frac{\Pi_{mp}}{\ln^2 T/T_K} \text{ for } T \gg T_K$$

- we must go beyond RG improved pert. theory
to study $T \lesssim T_K$

- but first return to another pressing
Lecture next Wed 1-2:30 $\approx$ 12:30-2:00 ??

1/3 Relevant, Irrelevant & Marginal Def'n of RG

Relevant, Irrelevant & Marginal operators

- so far I've done "RG light" & haven't
justified dropping terms generated by

RG transformation such as

$$\int d\tau \sum_{\vec{k}, \vec{k}'} \bar{\Psi}_{\vec{k}}(\tau) \vec{\sigma}_{\vec{k}, \vec{k}'} \frac{d}{d\tau} \Psi_{\vec{k}'}(\tau) \cdot \vec{S}(\tau)$$

- I now want to justify this &

give full def'n of RG, using Kondo model
as an example

- I will first streamline notation and establish connection of Kohn model with relativistic Q.F.T.

- Note that Kohn interaction only involves

S-wave component of $\psi(\vec{r}) \propto \int d^2\hat{r} \psi(\hat{r}, r) \propto \psi_{s,0}$

$$\psi_{s,0} \propto \int d^2R \psi \hat{R} R$$

- in terms of S-wave component, passing to

$\propto$ V limit & normalizing S-wave fermion

$$\text{field} \sim \{ \psi(\vec{r}), \psi^\dagger(\vec{r}') \} = 2\pi \delta(\vec{r} - \vec{r}')$$

(Note unusual 2π , as in Shankar)

$$H_0 = VF \int_{-\Lambda}^{\Lambda} \frac{dk}{2\pi} k \psi^\dagger(k) \psi(k)$$

- here I assumed cut-off $D_0 \ll$ band-width

& linearized dispersion relation $E = VF(k - k_F)$,

shifted $k - k_F \rightarrow k$ and let $\Lambda = \frac{D_0}{VF}$

$$S_0 = \int_{-\infty}^{\infty} \frac{d\omega}{2\pi} \int_{-\infty}^{\infty} \frac{dk}{2\pi} (i\omega - VFk) \bar{\Psi}(\omega, k) \Psi(\omega, k) \quad (137)$$

$$S_{INT} = -\lambda F \int_{-\infty}^{\infty} \frac{dk}{2\pi} \int_{-\infty}^{\infty} \frac{dk'}{2\pi} \bar{\Psi} \int_{-\infty}^{\infty} \frac{d\omega}{2\pi} \int_{-\infty}^{\infty} \frac{d\omega'}{2\pi}$$

$$\bar{\Psi}(\omega', k) \sigma_z \Psi(\omega - \omega', k') \cdot \tilde{J}(\omega)$$

where I introduced Fourier transform of impurity spin operator

- S_{INT} involves frequency, not momentum

- if we Fourier transform back to

time & space, with a 1D Co-ordinate x

$$S_0 = - \int_{-\infty}^{\infty} dT \int_{-\infty}^{\infty} dx \bar{\Psi}(x, T) \left[\frac{d}{dT} - iV \frac{d}{dx} \right] \Psi(x, T)$$

- corresponds to a relativistic (Dirac,

left-moving only) fermion, with $(- \rightarrow) VF$
 [obtained by imposing $x \in \mathbb{R}^+$]
 $S_{INT} = -\lambda VF \int_{-\infty}^{\infty} dT \bar{\Psi}(0, T) \sigma_z \Psi(0, T) \cdot \tilde{J}(T)$

- this destroys Lorentz invariance, & even spatial translation invariance

R G step reduces $\Lambda \rightarrow \frac{\Lambda}{s}$ [$s > 1$]

- reducing Λ in infinitesimal steps $s \approx 1$

- after R G step N_s

$$S_0 = \int_{-\infty}^{\infty} \frac{d\omega}{2\pi} \int_{-\frac{\Lambda}{s}}^{\frac{\Lambda}{s}} \frac{dk}{2\pi} (i\omega - v_F k) \bar{\Psi}(\omega, k) \Psi(\omega, k)$$

- it is convenient & conventional to rescale

k ~~to~~ after reducing cut-off $\tilde{k} = k \cdot s$

to make S_0 have same form as before

$$\int_{-\frac{\Lambda}{s}}^{\frac{\Lambda}{s}} dk = \frac{1}{s} \int_{-\Lambda}^{\Lambda} d\tilde{k}$$

[will eventually drop the $\sim$]

- also rescale $\tilde{\omega} = \omega s$ so $(\omega - v_F k)$

factor remain unchanged

$$S_0 \Rightarrow \frac{1}{s^3} \int_{-\infty}^{\infty} \frac{d\tilde{\omega}}{2\pi} \int_{-\Lambda}^{\Lambda} \frac{d\tilde{k}}{2\pi} (i\tilde{\omega} - v_F \tilde{k}) \bar{\Psi}(\tilde{\omega}, \tilde{k}) \Psi(\tilde{\omega}, \tilde{k})$$

- finally we rescale fields to leave S_0

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○ invariant $\Psi' = S^{3/2} \Psi$, $\bar{\Psi}' = S^{3/2} \bar{\Psi}$ so

$$S_0 \rightarrow \int_{-\infty}^{\infty} \frac{d\tilde{\omega}}{2\pi} \int_{-\infty}^{\infty} \frac{d\tilde{k}}{2\pi} (1 - v\tilde{\omega} - \tilde{k}) \bar{\Psi}'(\tilde{\omega}, \tilde{k}) \Psi'(\tilde{\omega}, \tilde{k})$$

- NB - we chose rescaling factor for fields

in order to leave S_0 invariant

- this is appropriate when we are considering vicinity of free ($\lambda=0$) fixed point - which we saw was stable for $\lambda_0 < 0$

- in general we might choose some other field rescaling if we are studying vicinity of some other fixed point

- likewise, we don't always seek ω, k

by some power of S - more generally

$$\tilde{k} = k S, \quad \tilde{\omega} = \omega S^{\frac{1}{Z}}, \quad Z \equiv \text{"dynamical exponent"}$$

"

- fact that dispersion relation is (approximately)

linear, $S \propto (i\omega - V_F k)$ dictates
this choice in this case

- rescaling of $\vec{J}(\omega)$ simply determined
by Fourier transform:

$$\vec{J}(\omega) = \int_{-\infty}^{\infty} d\tau e^{i\omega\tau} \vec{J}(\tau)$$

$$\begin{aligned} \vec{J}(\tilde{\omega}) &= \int d\tau e^{i \frac{\tilde{\omega}}{S} \tau} \vec{J}(\tau) \\ &= \frac{1}{S} \int d\tilde{\tau} e^{i\tilde{\omega}\tilde{\tau}} \vec{J}(\tilde{\tau}) \end{aligned}$$

$$\vec{J}(\omega) \xrightarrow{A} S \cdot \vec{J}(\tilde{\omega})$$

- Kondo interaction is invariant under this
rescaling

$$\begin{aligned} &\int dk dk' d\omega d\omega' \vec{J}(\omega) \vec{J}(\omega') S^3 \vec{\varphi}' \cdot \vec{\varphi} S \vec{J}(\omega) \\ \rightarrow &\int d\tilde{k} d\tilde{k}' d\tilde{\omega} d\tilde{\omega}' \frac{S^4}{S^4} \vec{\varphi}' \cdot \vec{\varphi} S \vec{J}(\omega) \end{aligned}$$

- all factors of S cancel