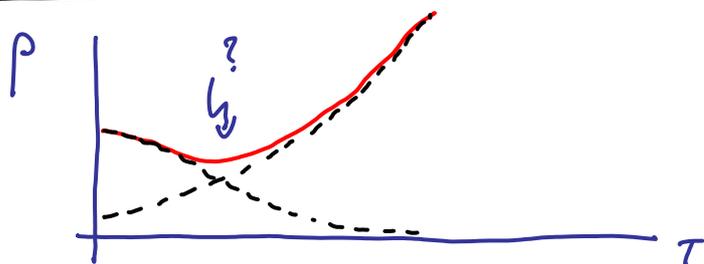


Equilibrium Review: cond-mat/9408101

Nonequilibrium: P. Metha, N. Andrei, PRL, 2005

Introduction:



Kondo, 1964: s-d-model for dilute random impurities

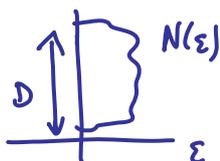
dimensionless

Description of magnetic impurities in metal:

$$H_0 = \sum_{ka} \epsilon_k c_{ka}^\dagger c_{ka}, \quad a = \pm 1/2$$

$$H_I = J \sum_{kk'} (c_{ka}^\dagger \bar{\sigma}_{ab} c_{k'b}) \cdot \bar{S}$$

$\bar{\sigma}_{el}(0) \leftarrow 1/2, 1, 3/2, \dots$



$D = \text{bandwidth} \approx 10,000 \text{ K} = \text{large}$

Kondo found  $\rho = \rho_0 + J^3 \ln T/D$ , logarithm was surprising! 2

Impurity Susceptibility:  $\chi^i = \frac{\mu^2}{T} [1 - J + J^2 \ln T/D - J^3 \ln^2 T/D]$

$$\chi = \left. \frac{\partial M_i}{\partial h} \right|_{h=0} (T)$$

Curie law, diverges for  $T \rightarrow 0$

Logarithms show up again: infrared instability.

Stimulated new theoretical approaches:

- resummation methods
- poor man's scaling (Anderson)
- numerical renormalization (Wilson)
- Bethe Ansatz (Andrei)
- CFT (Affleck Ludwig)
- bosonization
- Flow-equation Rg (Kohno)

resummation

$$\chi_i = \frac{\mu^2}{T} \left[ 1 - \frac{J}{1 + J \ln T/D} \right] = \frac{\mu^2}{T} \left[ 1 - \frac{1}{\ln T/T_K} \right]$$

$$T_K = D e^{-\frac{1}{J}} \quad : \text{low-energy scale that parametrizes theory.}$$

Summing subleading loop does not help!

For  $J > 0$  (AF coupling): IF-unstable

For  $J < 0$  (FM coupling): IF-stable

Anderson scaling: Start with  $H(D, J) \xrightarrow{\text{renormalize}} H(D', J')$  with  $D' < D$

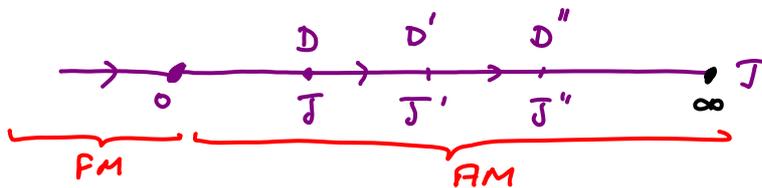
so  $T$  can become smaller before problems arise.

$$J \rightsquigarrow J' = J + J^2 \#$$

⊂ positive!

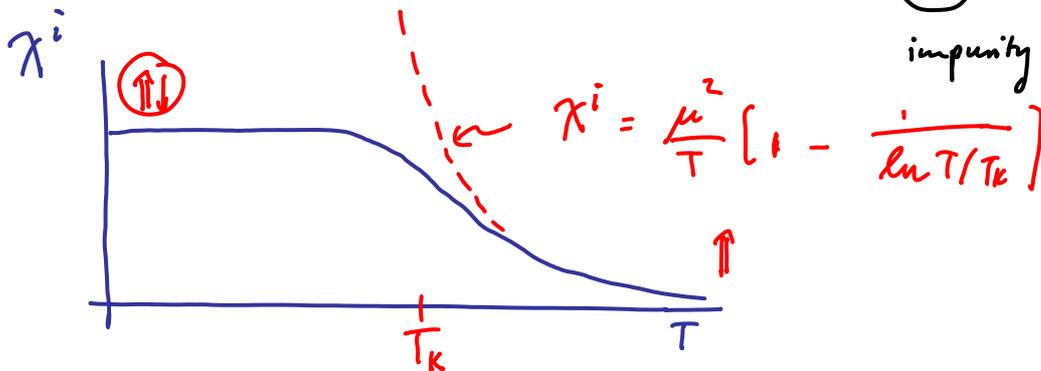
$J$  grows as  $D$  is reduced

Anderson, Wilson:



For  $S = 1/2$ :  $J'(D' \rightarrow 0) \rightarrow \infty$

$\uparrow \downarrow$  = singlet!  
impurity disappears!



Response to magnetic field becomes finite as spins get bound into singlet.

Analogous to confinement in QCD: quarks look free when probed at high momenta, but are confined at small momenta.

$$H = \sum_{ka} \epsilon_k c_{ka}^\dagger c_{ka} + J \sum_{kk'} (c_{ka}^\dagger \bar{\sigma}_{aa'} c_{k'a'}) \cdot \bar{\sigma}_0$$



Bandwidth large, only low-energy properties are of interest:

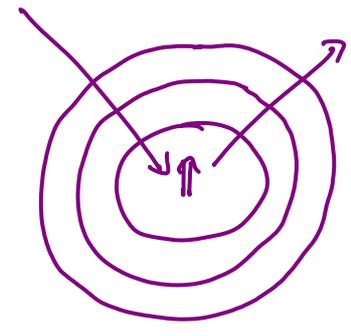
Assume  $T, \hbar \ll D \rightarrow \infty$ , universal results, independent of details of bandwidth.

Field-theoretic description:

$c_{\vec{k},a}$  good for translational invariance

↓ change to spherical basis

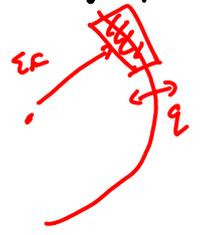
$c_{k\ell m, a}$  only  $\ell = \ell_0$  couples to impurity  
 ↑ angular momentum of local level



Typically one takes  $\ell_0 = 0, \ell = 0$

For  $\ell_0 > 0, |\ell| \leq \ell_0$  is possible, "multi-channel model"

Linearize free spectrum:  $\epsilon_k = \epsilon_F + q v_F$ , for  $|q| < D$



$c_{\vec{k},a} \rightarrow c_{k,a} \xrightarrow{\text{Fourier transform}} c_{q,a} \rightarrow \psi_a(x)$  with  $x \in \mathbb{R}$   
 $|q| < 0$   $x < 0$ : incoming  
 $x > 0$ : outgoing

Then:

$$\hat{H} = -i \int_{-\infty}^{\infty} dx \psi_a^\dagger(x) \partial_x \psi_a(x) + J \psi_a^\dagger(0) \bar{\sigma}_{ab} \psi_b(0) \cdot \bar{\sigma}_0$$

(only right-movers, chiral fields)

units:  $v_F = 1$

Dos:  $\nu = \frac{1}{\pi}$ ,  $J$  dimensionless.

Strictly speaking  $|x| < 1/D \rightarrow \infty$  as  $D \rightarrow 0$

Cutoffs will be introduced while diagonalizing  $H$ , in a way that preserves symmetries.

We are interested in finding eigenstates in Fock space, with

$$\hat{H} |F\rangle = E |F\rangle, \quad |F\rangle \text{ is many-body state}$$

Build up  $|F\rangle$  one electron at a time.

Electron number:  $\hat{N}_e = \int dx \psi^\dagger(x) \psi(x)$ ,  $[\hat{H}, \hat{N}_e] = 0$  7

We are interested in thermodynamic limit }  $N_e \rightarrow \infty$   
 $L \rightarrow \infty$ , with  $D = \frac{N_e}{L} = \text{fixed}$ .

② Scaling limit:  $D \rightarrow \infty$  will restore universality.

$N_e = 1$ :

general form of  $|F\rangle = \int dx F_{aa_0}(x) \psi_a^\dagger(x) |0\rangle \otimes |a_0\rangle$

where  $|0\rangle$  is "drained Fermi sea":  $\psi_a(x) |0\rangle = 0$   
 $|a_0\rangle$  is impurity state.

Go to "1st quantization":

$$H_0 |F\rangle = -i \int dy \psi_b^\dagger(y) \partial_y \psi_b(y) \int dx F_{aa_0}(x) \psi_a^\dagger(x) |0\rangle \otimes |a_0\rangle$$

$\delta_{ab} \delta(x-y)$

$$= -i \int dx \partial_x F_{aa_0}(x) \psi_a^\dagger(x) |0\rangle \otimes |a_0\rangle$$

8

$\Rightarrow N_e = 1$

$$h(x) = -i \partial_x + J \delta(x) \vec{\sigma} \cdot \vec{\sigma}_0$$

$N_e :$

$$h(x) = -i \sum_j \partial_{x_j} + J \delta(x_j) \vec{\sigma} \cdot \vec{\sigma}_0$$

Schrodinger Eq:

$$\left[ -i \partial_x + J \delta(x) \vec{\sigma}_{ab} \cdot (\vec{\sigma}_0)_{a_0 b_0} \right] F_{bb_0}(x) = E F_{aa_0}(x)$$

$E = k :$



only  $e^{ikx}$ , due to  $\partial_x$  (linearized spectrum)

$$F = e^{ikx} [A_{aa_0} \Theta(-x) + B_{aa_0} \Theta(x)]$$

Condition between A, B:

$$B_{aa_0} = \sum_{bb_0} S_{aa_0}^{bb_0} A_{bb_0}$$

S-matrix.

$$\begin{aligned} \hbar F = kF + -i [ -A_{aa_0} \delta(x) + B_{aa_0} \delta(x) ] & \quad \delta(x) \Theta(\pm x) = \frac{1}{2} \delta(x) \\ + \int \sigma_{ab}(\vec{\sigma}_0)_{a_0 b_0} \delta(x) e^{ikx} [ \underbrace{A_{bb_0} \Theta(-x) + B_{bb_0} \Theta(x)}_{\frac{1}{2}[A+B]} ] \end{aligned}$$

⇒ F is eigenstate if  $i(A - B) + \frac{1}{2} \int \sigma \cdot \vec{\sigma}_0 (A + B) = 0$ .

$$\Rightarrow \left( i + \frac{1}{2} \int \sigma_{ab}(\vec{\sigma}_0)_{a_0 b_0} \right) A_{bb_0} = \left( i - \frac{1}{2} \int \sigma_{ab}(\vec{\sigma}_0)_{a_0 b_0} \right) B_{bb_0}$$

$$\Rightarrow S_{aa_0}^{bb_0} = \frac{1}{i - \frac{1}{2} \int \vec{\sigma} \cdot \vec{\sigma}_0} \left[ i + \frac{1}{2} \int \vec{\sigma} \cdot \vec{\sigma}_0 \right]$$

Introduce:  $P_{aa_0}^{bb_0} = \frac{1}{2} (1 + \sigma \cdot \vec{\sigma}_0)_{aa_0}^{bb_0}$  claim  $\delta_a^{b_0} \delta_{a_0}^b$  (10)  
= exchange operator 

$$\Rightarrow (PA)_{ab} = A_{ba}$$

Proof:

$$\begin{aligned} \frac{1}{2} ( 1 + \sigma_{(1)}^z \cdot \sigma_{(2)}^z + \sigma_{(1)}^+ \sigma_{(2)}^- + \sigma_{(1)}^- \sigma_{(2)}^+ ) A_{\uparrow\uparrow} &= A_{\uparrow\uparrow} \\ A_{\downarrow\downarrow} &= A_{\downarrow\downarrow} \\ A_{\uparrow\downarrow} &= A_{\downarrow\uparrow} \\ A_{\downarrow\uparrow} &= A_{\uparrow\downarrow} \end{aligned}$$

⇒  $P^2 = 1$

$$\begin{aligned} S_{aa_0}^{bb_0} &= \frac{1}{i - \frac{1}{2} \int (\vec{\sigma} \cdot \vec{\sigma}_0 + 1) + \frac{1}{2} \int} \left[ i + \frac{1}{2} \int \vec{\sigma} \cdot \vec{\sigma}_0 \right] \\ &= \frac{1}{i + \frac{1}{2} \int - \int P} \frac{i + \frac{1}{2} \int + \int P}{(i + \frac{1}{2} \int) + \int P} \quad [ \quad ] \end{aligned}$$

$$= \frac{1}{(i + \frac{1}{2}J)^2 - J^2} \left[ (i + \frac{1}{2}J + JP) (i - \frac{1}{2}J + JP) \right]$$

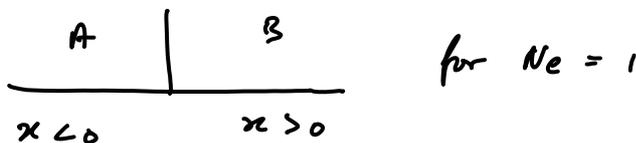
$$= a + ib P$$

$$\Rightarrow S^{10} = \frac{1 - icP}{1 - ic} \quad \text{acts on } A, \text{ gives } B,$$

$$c = \frac{J}{1 - \frac{3}{4}J^2} \sim J.$$

No,  $N_e = 1$  problem has been solved!

$N_e = 2$ :

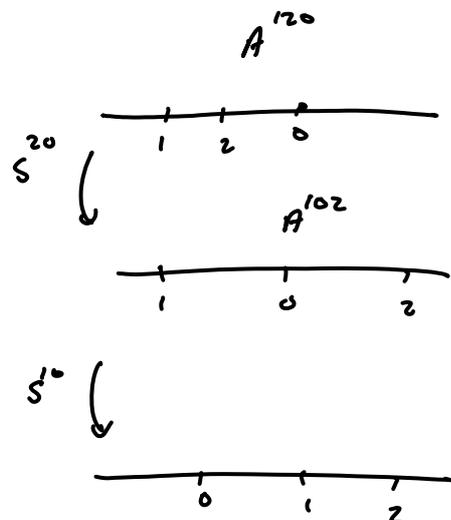
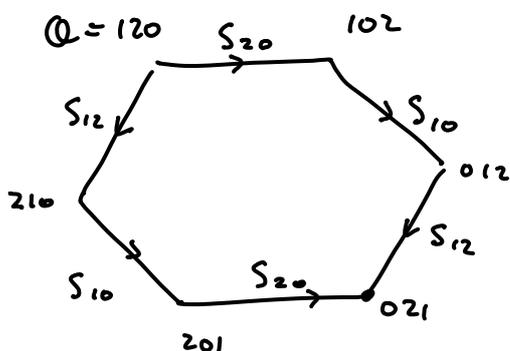


For  $N_e = 2$ ;  $3! = 6$  regions; labeled by  $\mathcal{Q}$ , for 6 ordering of  $x_1, x_2, 0$ .

$$F(x_1, x_2) = \sum_{\mathcal{Q}} A_{a_1 a_2 a_0}^{\mathcal{Q}} \Theta(x_{\mathcal{Q}})$$

$$S^{20} A^{120} = A^{102}$$

$$S^{10} A^{102} = A^{012}$$



two wrap of reading  $021$  from  $120$ :

$$S^{12} S^{10} S^{20} = S^{20} S^{10} S^{12}$$

Yang-Baxter equation.

If  $S$  satisfy YB, 2-electron wave-function can be constructed.

Deep assumption is made, (which usually fails):

$$F = e^{ik_1 x_1 + ik_2 x_2} \sum_Q A_{a_1 a_2 a_0} \Theta(x_Q) \quad , \quad E = k_1 + k_2$$

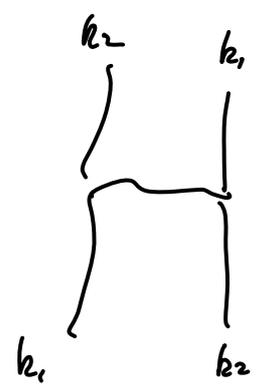
Assumption was made that the same  $k$ 's work in all  $\Theta$ 's.

more generally, we should use:  $k_1 + p_Q, k_2 - p_Q,$

where  $p_Q$  is not the same in each  $Q$ .

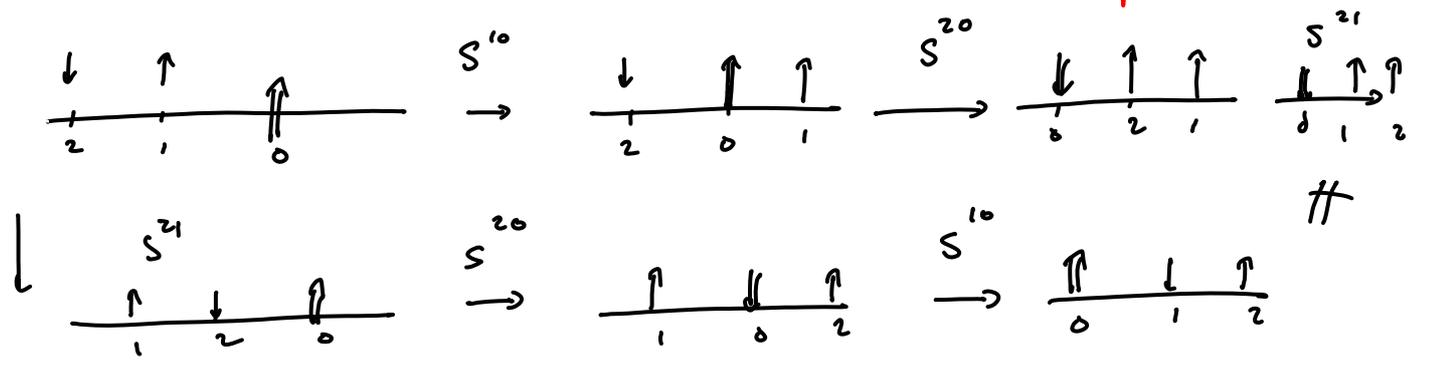
Bethe Ansatz is:  $p_Q = 0$  for all  $Q$

Essence of integrability: only momenta are interchanged, not shifted!!



$S^{12} = 1$  ?

does not work, since  $S^{10} S^{20} \neq S^{20} S^{10}$



Different order of swapping leaves us in different state, since scatterer has internal degrees of freedom  $\Rightarrow$  Noncommutativity!

So  $S^{12} = 1$  does not work. Reverse question: does  $S^{12}$  exist such that YB holds.

Claim:  $S^{12} = P^{12} \Rightarrow P^{12} S^{10} S^{20} = S^{20} S^{10} P^{12}$ , YB holds.

But is it reasonable to use  $S^{12} = P^{12}$ ? Does it introduce interactions among electrons? (We may be solving the wrong model!) Answer: No!

Consider 2 free electrons:  $\hbar = -i(\partial_1 + \partial_2)$

a solution:  $F_{a_1 a_2} = e^{i(k_1 x_1 + k_2 x_2)} A_{a_1 a_2}, \quad E = k_1 + k_2$

another solution:  $F_{a_1 a_2}(x_1, x_2) = \sum_q e^{i(k_1 + q)x_1 + i(k_2 - q)x_2} A_q$   
 $= e^{i(k_1 x_1 + k_2 x_2)} \sum_q e^{iq(x_1 - x_2)} A_q$   
 $= e^{i(k_1 x_1 + k_2 x_2)} f(x_1 - x_2) \leftarrow \text{infinite}$

$\Rightarrow$  the level  $E$  is infinitely degenerate.

antisymmetrize  $\rightarrow$   $F_{a_1 a_2} = A e^{i(k_1 x_1 + k_2 x_2)} \left( A_{a_1 a_2} \Theta(x_1 - x_2) + \frac{SA}{A_{a_2 a_1}} \Theta(x_2 - x_1) \right)$ .  
 any  $S$  can be used!  
 no continuity concerns, since  $\partial_x$  is linear!

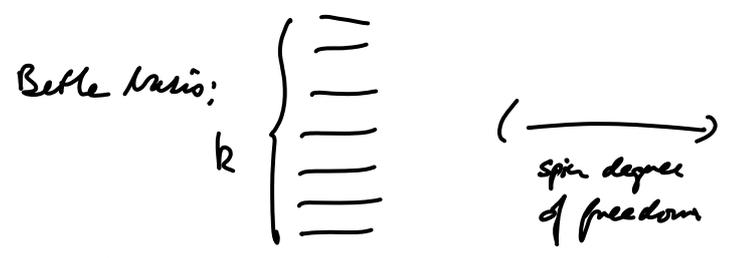
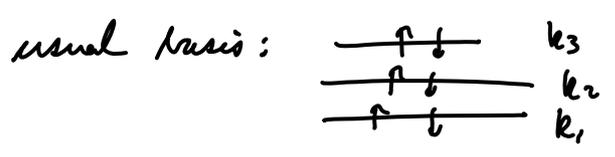
To perturb degenerate levels, choose special basis, such that

$\langle i | V | j \rangle = V \delta_{ij}$ , else  $\frac{\langle i | V | j \rangle}{E_i - E_j}$  blows up

So, the choice  $S^{12} = P^{12}$  corresponds to choice of basis in the degenerate space of  $k_1 + k_2 = E$  two-electron states.

$S^{12} = P^{12}$  corresponds to charge-spin decoupling:

$F_{a_1 a_2} = \begin{pmatrix} e^{i(k_1 x_1 + k_2 x_2)} & e^{i(k_1 x_1 + k_2 x_2)} \\ - & - \end{pmatrix} \left( A_{a_1 a_2} \Theta(x_1 - x_2) + A_{a_2 a_1} \Theta(x_2 - x_1) \right)$



but two electrons with same  $k$  would give 0.

charge degree of freedom