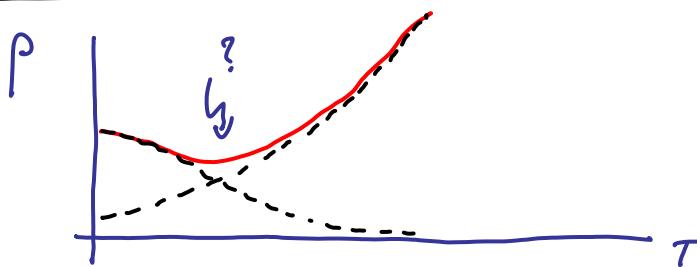


Equilibrium Review: cond-mat/9408101

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

### Introduction:

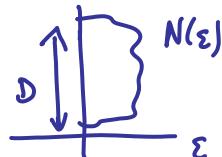


Kondo, 1964: s-d-model for dilute random impurities

Description of magnetic impurities in metal:

$$H_0 = \sum_{ka} \varepsilon_k c_{ka}^+ c_{ka}, \quad H_I = J \sum_{k k' b} (c_{ka}^+ \bar{\sigma}_{ab} c_{k'b}) \cdot \vec{S}$$

$a = \pm \gamma_2$



dimensionless

$$\bar{\sigma}_{el}(0)$$

$\gamma_2, 1, 3/2, \dots$

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

Kondo found  $\rho = \rho_0 + J^3 \ln T/D$ , logarithm was surprising! L2

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

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

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 (Kehrein)

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resummation  $\rightarrow$

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

$T_K = D e^{-\frac{1}{J}}$  : 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

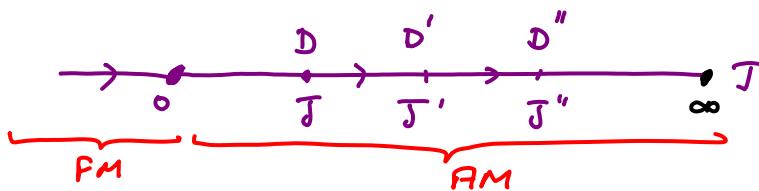
Anderson scaling: Start with  $H(D, J) \xrightarrow{\text{renormalize}} H(D', J')$  with  $D' < D$   
so  $T$  can become smaller before problems arise.

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

$\approx$  positive!

J grows as D is reduced

Anderson, Wilson:

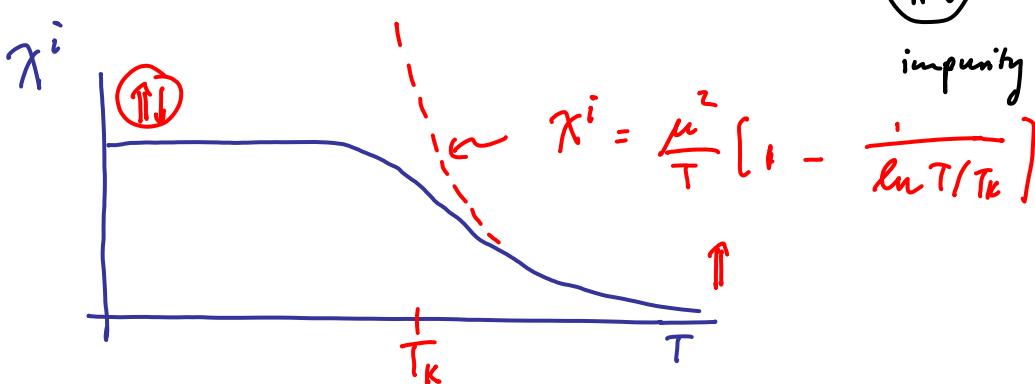


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For  $S = \gamma_2$  :  $J'(D' \rightarrow 0) \rightarrow \infty$ .

$\uparrow\downarrow$  = 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.

$$H = \sum_{ka} \varepsilon_k c_{ka}^\dagger c_{ka} + J \sum_{kk'} (\bar{c}_{ka}^\dagger \bar{\sigma}_{aa'} c_{k'a'}). \bar{\sigma}$$



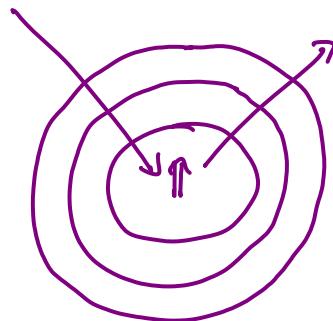
Bandwidth large, only low-energy properties are of interest:

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

Field-theoretic description:

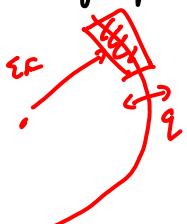
$\tilde{c}_{b,a}$  good for translational invariance  
 { change to spherical basis

$c_{klm,a}$  only  $l = l_0$  couples to impurity  
 angular momentum of local level



Typically one takes  $l_0 = 0$ ,  $l=0$   
 For  $l_0 > 0$ ,  $|m| \leq l_0$  is possible, "multi-channel model"

Linearize free spectrum:  $\varepsilon_k = \varepsilon_F + q v_F$ , for  $|q| < D$



$\tilde{c}_{k,a} \rightarrow c_{k,a} \rightarrow c_{q,a} \xrightarrow{|q| < 0}$  Fourier transform with  $x \in \mathbb{R}$   
 $x \leftarrow 0$ : incoming  
 $x \rightarrow \infty$ : 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  $|x| < /D \rightarrow \infty$  as  $D \rightarrow 0$

Cutoff 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.

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

We are interested in the thermodynamic limit  $\left. \begin{array}{l} \\ \end{array} \right\} \text{ : } \begin{array}{l} N_e \rightarrow \infty \\ L \rightarrow \infty \end{array}$ , with  $D = \frac{N_e}{L} = \text{fixed}$ .

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

$$\underline{N_e = 1}:$$

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

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

Go to "1st quantization":

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

$\uparrow \qquad \uparrow$

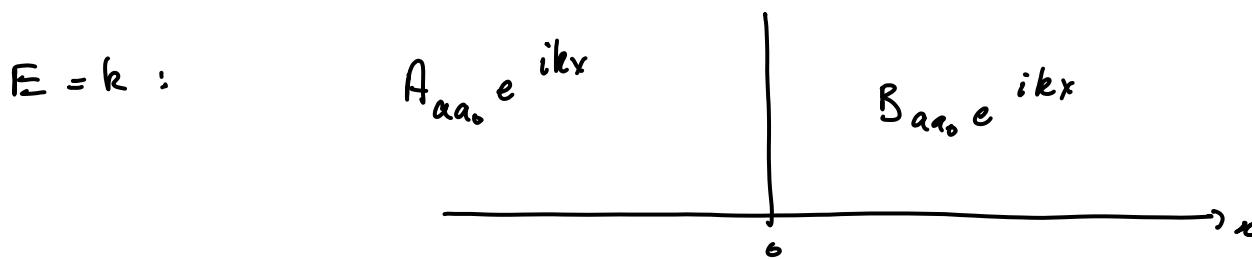
$\delta_{ab} \delta(x-y)$

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

$$\Rightarrow N_e = 1 \quad h(x) = -i \partial_x + \mathcal{T} \delta(x) \vec{\sigma} \cdot \vec{\sigma}_0$$

$$N_e : \quad h(x) = -i \sum_j \partial_{x_j} + \mathcal{T} \delta(x_j) \vec{\sigma} \cdot \vec{\sigma}_0$$

$$\text{Schrödinger Eq: } \left( -i \partial_x + \mathcal{T} \delta(x) \vec{\sigma}_{ab} \cdot (\vec{\sigma}_0)_{a_0 b_0} \right) F_{bb_0}(x) = E F_{aa_0}(x)$$



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:

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

$\hookrightarrow$  S-matrix.

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

$$\Rightarrow F \text{ is eigenstate if } i[A - B] + \frac{1}{2} J \sigma \cdot \vec{\sigma}_0 (A + B) = 0.$$

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

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

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

$\vec{\sigma} \cdot \vec{\sigma}_0$  = exchange operator



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

$$\text{Proof: } \frac{1}{2} \left( 1 + \vec{\sigma}_{(1)}^z \cdot \vec{\sigma}_{(2)}^z + \vec{\sigma}_{(1)}^+ \vec{\sigma}_{(2)}^- + \vec{\sigma}_{(1)}^- \vec{\sigma}_{(2)}^+ \right) A_{\uparrow\uparrow} = A_{\uparrow\uparrow}$$

$$A_{\downarrow\downarrow} = A_{\downarrow\downarrow}$$

$$A_{\uparrow\downarrow} = A_{\downarrow\uparrow}$$

$$\Rightarrow P^2 = 1 \qquad \qquad \qquad A_{\downarrow\uparrow} = A_{\uparrow\downarrow}$$

$$S_{aa_0}^{bb_0} = \frac{i}{i - \frac{1}{2} J(\vec{\sigma} \cdot \vec{\sigma}_0 + 1) + \frac{1}{2} J} [i + \frac{1}{2} J \vec{\sigma} \cdot \vec{\sigma}_0]$$

$$= \frac{i}{i + \frac{1}{2} J - \frac{1}{2} J} \frac{i + \frac{1}{2} J + \frac{1}{2} J}{(i + \frac{1}{2} J) + \frac{1}{2} J} [ \quad ]$$

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

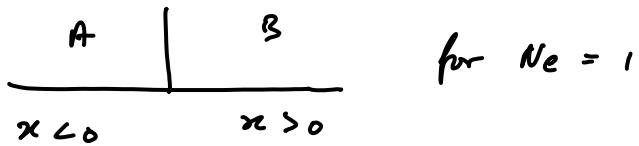
$$= a + ib P$$

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

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

So,  $N_e = 1$  problem has been solved!

$N_e = 2$ :

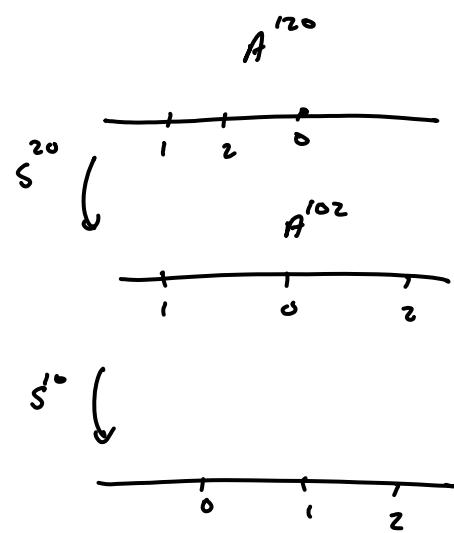
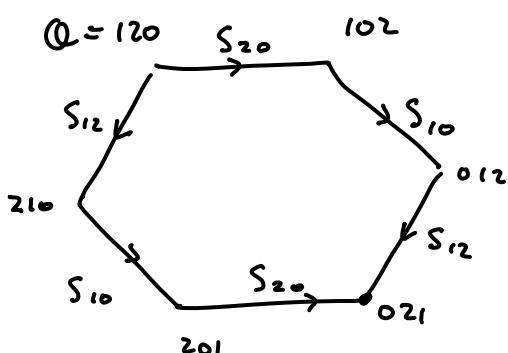


For  $N_e = 2$ ,  $3! = 6$  regions ; labeled by  $\Omega$  , for 6 ordering of  $x_1, x_2, 0$ .

$$F(x_1, x_2) = \sum_{\Omega} A_{a_1 a_2 a_0}^{\Omega} \theta(x_{\Omega})$$

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

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



two ways of reaching 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. 13

Deep assumption is made, (which usually fails):

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

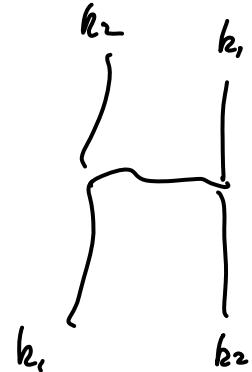
Assumption was made that the same  $k$ 's work in all Q's.

more generally, we should use:  $k_1 + p_0$ ,  $k_2 - p_0$ ,

where  $p_0$  is not the same in each Q.

Bethe Ansatz is:  $p_0 = 0$  for all Q

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

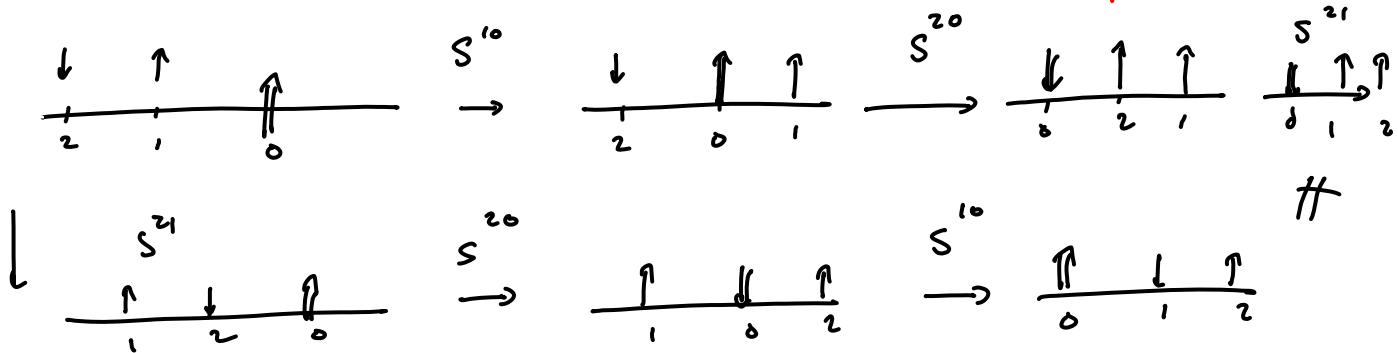


$$S^{12} = 1 ?$$

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

one possible outcome.



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:  $\hat{h} = -i(\partial_1 + \partial_2)$

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

$\Rightarrow$  The level  $E$  is infinitely degenerate.  
 antisymmetrize  $\rightarrow$

any  $S$  can be used!

$$F_{a_1 a_2} = A e^{i(k_1 x_1 + k_2 x_2)} \left( A_{a_1 a_2} \delta(x_1 - x_2) + \frac{SA}{A_{a_1 a_2}} \delta(x_2 - x_1) \right).$$

no continuity concerns, since  $\partial_x$  is linear!

To perturb degenerate levels, choose special basis, such that

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

Hence, 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:

$$F_{a_1 a_2} = \left( e^{i(k_1 x_1 + k_2 x_2)} - e^{i(k_1 x_2 + k_2 x_1)} \right) \left( A_{a_1 a_2} \delta(x_1 - x_2) + A_{a_2 a_1} \delta(x_2 - x_1) \right).$$

usual basis:

Bethe basis:

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

charge degree of freedom