

MPPIKS Summer School Lecture

Boundary Conformal Field Theory Approach

to QTP's

- There is considerable interest in models involving a single localized quantum impurity interacting with a gasless ~~continuous~~ delocalized excitations
- eg Kondo model, multi-channel, multi-impurity [charge & spin varieties]
- magnetic impurity in spin system
- construction in quantum ~~wire~~ wire
- multi-terminal junction
- X-ray edge problems.
- magnetic monopoles (dyons) interacting with charged particles

- these models often exhibit interesting cross-over in behavior from high E to low E that can be described by RG
- getting an accurate description of full cross-over is difficult - sometimes done with Bethe Ansatz or NRG or ~~approx~~ with MFT's but all these methods have limitations
- high & low energy behavior is often relatively simple & can be accurately calculated
- very generally these high and low energy RG fixed points correspond to CIBCs in CFT's
- often they have equivalent but

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more familiar descriptions [eg. free fermions]

- but in some cases they correspond to NFL QCP's - eg. 2 channel (or multi-channel) Kondo, 2 impurity Kondo
- in these lectures I will review this general formalism and discuss the 2-channel Kondo example in some detail, including experiments by Goldhaber-Gordon group.
- the first step in mapping to a BCFT is to map into 1D
- this is possible, and convenient, due to fact that all interactions take place at $\vec{r}=0$ - short range

eg. - Kondo model

- for spherically symmetric case we can expand in spherical harmonics

- we ignore

$$H = \int d^3\vec{r} \psi_{\vec{r}}^\dagger \left(-\frac{\hbar^2 \nabla^2}{2m} \right) \psi_{\vec{r}} + J \psi_{\vec{r}=0}^\dagger \psi_{\vec{r}=0}$$

$$\epsilon = \frac{\hbar^2 k^2}{2m} - \mu$$

- we can expand in spherical harmonics

$$\psi_{\vec{r}} = \frac{1}{\sqrt{4\pi}} \psi_0(r) + \text{higher harmonics}$$

ℓ = 0 - wave

- only s-wave interacts with impurity
- NB - we ignore bulk interactions - justified?
- to study low energy physics (at $T \ll 1$)

where $V = \text{density of states} = mk_F / \pi^2$

we linearise $\epsilon(k)$ near k_F : $\epsilon = V_F(k - k_F)$

and define incoming and outgoing

waves ["left", "right" movers]

$$\psi_L(r) = \int_{-\infty}^{\infty} dk e^{\pm ikr} \psi_0(k_F + k)$$

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$$H = \frac{VF}{2\pi} i \int_0^\infty dr \left[\psi_L^\dagger \frac{d}{dr} \psi_L - \psi_R^\dagger \frac{d}{dr} \psi_R \right] \\ + VF \lambda \psi_L^\dagger(0) \vec{\sigma}_2 \psi_L(0) \cdot \vec{J} + \dots$$

where $\lambda = JV$, NB $\psi_L(0) = \psi_R(0)$

- a Lorentz-invariant ($VF \leftrightarrow c$)

Dirac fermion field theory in $(1+1)D$
with impurity interactions at $r=0$

- defined on half-line

at $\lambda=0$, this is an example of a CFT
with a boundary. [+ decoupled spin]

- fermions are massless $\Rightarrow$ no scale

imaginary time Lagrangian is

$$2\pi \mathcal{L} = \psi_L^\dagger \left(\frac{d}{dr} + iVF \frac{d}{dr} \right) \psi_L + \psi_R^\dagger \left(\frac{d}{dr} - iVF \frac{d}{dr} \right) \psi_R$$

let $Z = T + ir$ [and set $VF=1$]

- then, without boundary, $S = \int_{-\infty}^{\infty} dT \int_{-\infty}^{\infty} dr \mathcal{L}$

is invariant under conformal transformation

$$\psi_L \rightarrow \left(\frac{dw}{dz}\right)^{+1/2} \psi_L, \quad \psi_R \rightarrow \left(\frac{dw}{d\bar{z}}\right)^{+1/2} \psi_R$$

$$Z \rightarrow w(Z), \quad \psi_L \rightarrow \sqrt{w}'^{1/2} \psi_L, \quad \psi_R \rightarrow \sqrt{w}'^{1/2} \psi_R$$

where $\bar{z} = \bar{r} - i\tau$

$$\text{Since } S = \int dZ d\bar{Z} \left[\psi_L^\dagger \frac{d}{dZ} \psi_R + \psi_R^\dagger \frac{d}{d\bar{Z}} \psi_L \right]$$

- ∞ -dimensional set of transformations

$$w(z) = \sum_{n=0}^{\infty} a_n z^n, \quad a_n \in \mathbb{C}$$

$a_1 \in \mathbb{R}$ - scale transformation etc.

$\Rightarrow$ symmetry has powerful consequences

- if system is defined on $1/2$ -disk $r > 0$,

$\text{Im } Z > 0$, full conformal symmetry

is broken but CIBC's preserve

∞ subset that maps boundary, $\text{Im } Z = 0$

into itself $a_n \in \mathbb{R} \Rightarrow W^\dagger(\tau) = w(\tau)$

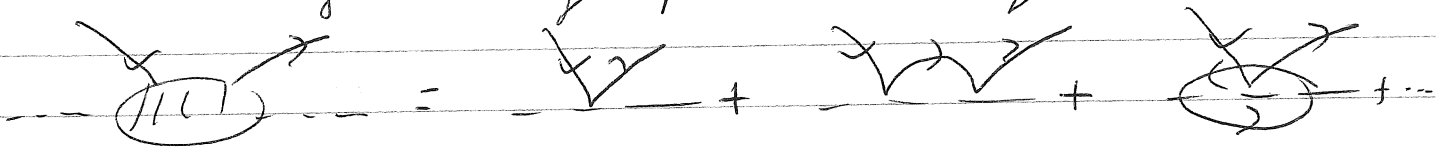
- in above example $\psi_L(r=0, \tau) = \psi_R(r=0, \tau)$

is a CIBC.

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of course Kondo interaction breaks BCI
- in fact it introduces a characteristic
energy scale - Kondo temperature

- Renormalization group (RG) eqns:



$$\frac{d\lambda}{d\ln D} = -\lambda^2 + \frac{\lambda^3}{2} + \dots$$

$$\Rightarrow \lambda(D) \text{ is } 0(D) \text{ at } T_K \pm D_0 e^{-\frac{1}{\lambda_0} \sqrt{\lambda_0}}$$

where D_0 is band-width-energy scale

at which bare coupling is evaluated

- but low energy behavior is again

described by a CIB(- at $E \ll T_K$

- this is most easily seen by considering

a tight-binding model with a large J

$$H = -t \sum_{j=0}^{\infty} (c_j^\dagger c_{j+1} + \text{h.c.}) + J \sum_i c_i^\dagger \frac{\sigma_z}{2} c_i \quad (\text{spin indices suppressed})$$

$$\begin{array}{ccccccc} \uparrow & & & & & & \\ H \rightarrow & 0 & 0 & 0 & 0 & & \\ & 0 & 1 & 2 & & & \end{array}$$

if $T \gg t$, 1 electron is trapped at site 0,
forming a singlet with impurity spin
- electrons on other sites can do anything

but go to site 0

- at $1/2$ -filling p-h symmetry, this
corresponds to a $\pi/2$ phase shift

bc modified from $\psi_L(i) = \psi_R(i)$ to $\psi_L(i) = -\psi_R(i)$

- this follows from writing $\psi_j = e^{i\frac{\pi}{2}j} \psi_R(j) + e^{-i\frac{\pi}{2}j} \psi_L(j)$

$$\begin{array}{lcl} \text{open bc: } \psi_{-1} = 0 \Rightarrow & & \psi_L = \psi_R \\ T \rightarrow \infty: \psi_0 = 0 \Rightarrow & & \psi_L = -\psi_R \end{array}$$

- another CIBC occurs [e impurity
spin disappears] - only change from
strong to weak coupling is in BC.

[+ elimination of impurity]

- starting with a small bare T , this
low energy description only becomes

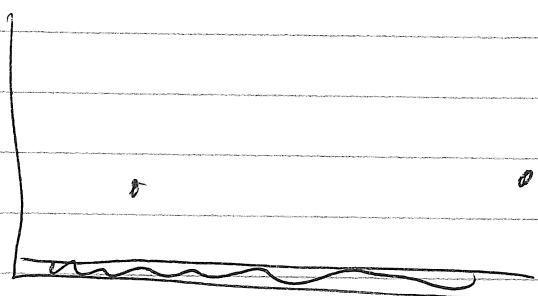
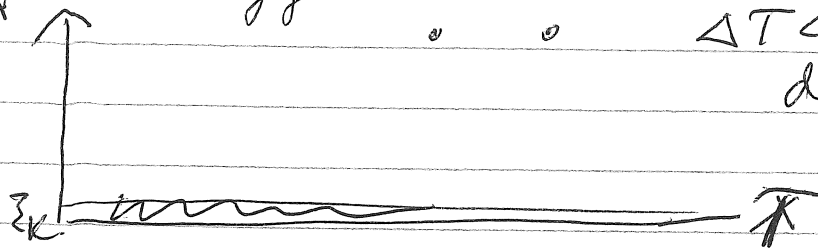
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valid at $E \ll T_K$

- ~~any~~^{low} "nearly" electron which forms singlet is at distance of $O\left(\frac{v_F}{T_K} \equiv \xi_K\right)$
- this is a simple example of a very general phenomenon.
- low energy effective Hamiltonian is a CFT with a CIBC for a wide range of problems. - tested above
- why? - no rigorous proof but -
- far from boundary compared to all microscopic length scales (such as ξ_K) and at long times ($\gg v_{TK}$) we might expect universal behavior characteristic of a critical point

2 different cases

$\Delta T \ll r$ - boundary doesn't matter



$\Delta T \gg r$
- boundary influences Green's functions

but only in a universal way consistent with system being critical - i.e. only ratios of lengths enter $\Delta T/r$, etc. + powers of length

$$e.g. \quad G(r_1, T_1; r_2, T_2) = \frac{1}{|T_1 - T_2|^n} f\left(\frac{r_1}{T_1 - T_2}, \frac{r_2}{T_1 - T_2}\right)$$

$$f \rightarrow \text{const} \quad \text{at} \quad \frac{r_1}{|T_1 - T_2|} \gg 1$$

- this type of scale-invariant long-distance behavior is described by CFT with a CIBC

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- NB- bulk critical theory doesn't change
due to ~~spacely~~ ^{short-range} interaction only at boundary
- all that happens is an RG flow from 1 CIBC to another
 - CIBC is a "termination" of bulk CI behavior as boundary is taken into account (but not approached too closely) that preserves a large~~st~~ subset of conformal symmetry
 - Some low energy, long distance properties
are ^{afternoon} ~~determined by the CIBC~~ completely by the conformally invariant Hamiltonian + CIBC - these include all long time & distance (e low T) Green's functions, the $T=0$ impurity entropy, $\ln g$, the $T=0$ ~~resistivity~~ resistivity (or conductance) for

dilute random array of $Kndr$ impurities
(or single quantum dot acting as a

$Kndr$ impurity) ...
[start here]

- other low energy long-distance properties

require consideration of leading irrelevant
operator (i.e. interactions) at fixed point

i.e. CFT & CIBC is exact description
of zero energy limit only

- it is a fixed point Hamiltonian ~~that~~
that system flows to

- at any finite energy scale H_{eff}
contains additional interacting terms which
break the conformal symmetry

- these:

1) don't involve the impurity degrees of freedom

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itself [eg. impurity spin] because it is screened at much higher energies [$\sim T_K$]

2) are ~~not~~ added to the model with

specified CIBC - the set of possible interactions is therefore determined by CIBC [+ other symmetries like $SO(2)$]

- of this ∞ set of interactions, all of which are irrelevant, the "least irrelevant" lowest dimension, LI^0 controls asymptotic behavior of various quantities: Kondo

- impurity susceptibility
- T -dependence of impurity entropy ($\Rightarrow$ specific heat)
- T -dependence of resistivity / susceptibility
- leading corrections to all Green's functions
- to calculate these effects we need

to know coupling constant of LIO in Hett

$$H = H_{\text{CIBC}}^{\text{CFT}} + g O_{\text{LI}}$$

- this cannot usually be determined exactly
[unless we can use Bethe ansatz or
a numerical method]

- but we can estimate it in order of
magnitude up to a factor of $\mathcal{O}(1)$

- it drops out in various ratios of
physical quantities [eg Wilson ratio
of impurity susceptibility & specific heat
or "logarithmic ratio" of resistivity to square
of susceptibility]
(to here)

- how are CIBCs specified?

- surprisingly, in NFL cases we
don't usually explicitly specify a BC

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- instead a "boundary state" $|A\rangle$ is specified
- this allows all of above quantities to be calculated [e also finite size spectrum with arbitrary pair of CIBCs at 2 ends of finite system] $\Rightarrow$ it is equivalent to specifying BC explicitly
- FSS is very useful for comparing

CFT predictions to NRG calculations

- consider CI partition function, at inverse temperature β , for finite system of length l , with CIBCs A, B at 2 ends:

H_{AB}^l is corresponding Hamiltonian

- all eigenvalues are of form $\frac{\pi}{l} X$ for ~~some~~ dimensionless numbers X_i

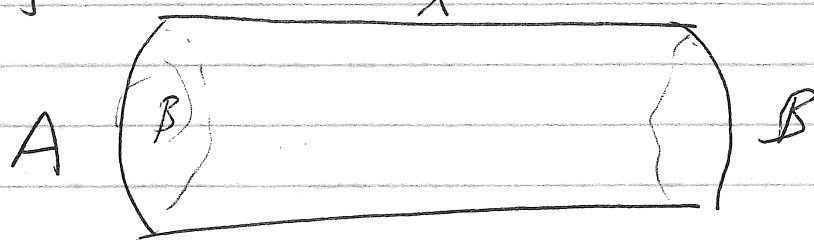
$$Z_{AB} = \text{tr } e^{-\beta H_{AB}^l}$$

- now we use a basic property of

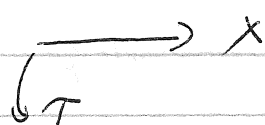
($d+1$) dimensional Lorentz invariant QFT's:

- we may interchange space and ^{imaginary} time
[modular transformation]

- now Z_A

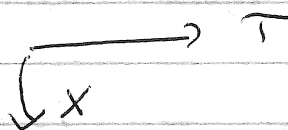


originally



Cylinder - periodic
in imaginary time

now consider



- now periodic in space (no boundary)

with system length β

- not periodic in T [not a trace]

- instead system propagates for time l

between initial & final states $|A\rangle, |B\rangle$

$$Z_{AB} = \langle A | e^{-l H_P^\beta} | B \rangle$$

H_P^β is Hamiltonian with PBC, length β

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- the boundary states $|A\rangle, |B\rangle$ determine Z_{AB} & in fact all other needed properties
- these states must obey certain properties to respect boundary conformal invariance = determined by J Cardy
- in some cases boundary conformal invariance + other symmetries allow only a finite set of possible boundary states
- then it may be relatively straightforward to determine boundary state corresponding to R/L fixed point of a given condensed matter QIP - eg. multi-channel Kndo (to p 20)

~~Simple Example - Quantum Ising Chain~~

Quantities Determined Completely by fixed point Hamiltonian (CIBC)

① Impurity Entropy

- how is it defined & measured experimentally
- for dilute Kondo impurities in a metal,
specific heat C_p is $C_p(T) = C_0(T) + M_{\text{imp}} C_{\text{imp}}(T) + \dots$
heat capacity

$C_{\text{imp}}(T)$ is extra entropy due to a single
impurity - it is $O(1)$ in volume limit

- we can calculate it from single impurity
model [eg 1D version] as

$$C_{\text{imp}}(T) = \lim_{h \rightarrow 0} \left[C_e(T, J) - C_e^{(0)}(T_h) \right]_{T=0}$$

where $C_e^{(0)}$ is ~~spec~~ heat capacity

~~without impurity~~

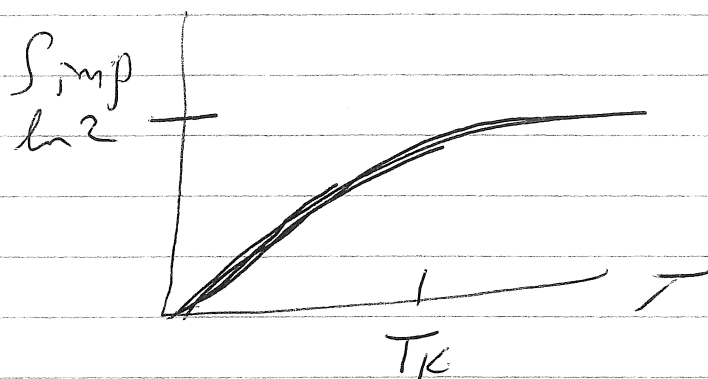
$$\text{NB } C_e(T, J=0) = C_e^{(0)}(T) + \ln(2S+1)$$

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impurity entropies is $S_{\text{imp}}(T) = \int_0^T \frac{dT}{T} C_{\text{imp}}(T)$

- for usual $\beta = \frac{1}{2}$, single channel

Kondo model



in scaling limit where $\log T \rightarrow 0$

$S_{\text{imp}} = \ln 2$ at $T \gg T_K$ since

the $T \rightarrow 0$ $T \rightarrow 0$ we get entropies
of free spin $= \ln 2$ [2 states]

at $T \rightarrow 0$, impurity is screened

e doesn't contribute $\Rightarrow S_{\text{imp}}(0) = 0$

KB - can also define $S_{\text{imp}}(T) = \lim_{T \rightarrow \infty} [S(T) - S^{(0)}(T)]$

- a general formula can be found for

S_{imp} at $T \rightarrow \infty$, $T \rightarrow 0$ - property of

RG fixed points in ultra-violet
e infrared for any CIBC.

- it is a pure number [setting $k_B = 1$]
on dimensional grounds [~~for $T \rightarrow 0$ or ∞~~
~~there is no~~]

- it can be expressed in terms of a simple
matrix element of B. state $|A\rangle$

② S-matrix in single particle sector

[resistivity & inductance in Kondo model]

- in 1D Dirac fermion theory we

calculate $\langle \psi_L^\dagger(\tau + i\epsilon) \psi_R(\tau - i\epsilon') \rangle$

or more simply $\langle \psi_L^\dagger(r) \psi_R(r) \rangle$

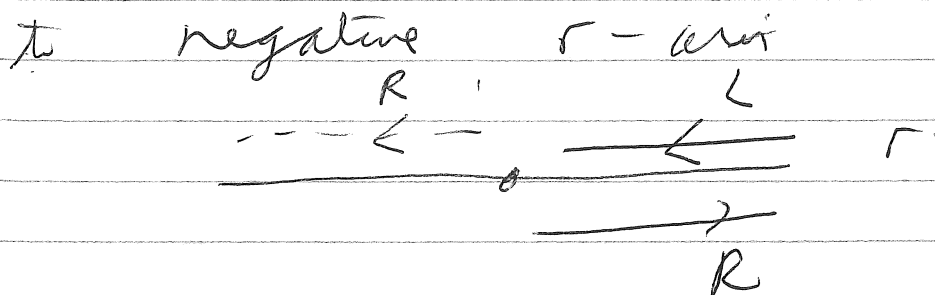
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- for $\lambda = 0$, bc $\psi_R(0) = \psi_L(0)$

implies that we may consider

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$\psi_R(r) = \psi_L(-r)$ - analytic continuation



$$R_1 \quad \langle \psi_L^+(r) \psi_R(r) \rangle = \langle \psi_L^+(r) \psi_L(-r) \rangle$$

$$= \frac{1}{2i\pi}$$

- in single-channel 1D model with

simple bc

$$\psi_R(r) = -\psi_L(r)$$

$$\psi_R(r) = -\psi_L(-r)$$

$$R_2 \quad \langle \psi_L^+(r) \psi_R(r) \rangle = -\frac{1}{2i\pi}$$

- more generally, when we break p-h

symmetry, we could have an arbitrary

phase shift, δ , at the Fermi surface.

$$\langle \psi_L^+(r) \psi_R(r) \rangle = \frac{e^{2i\delta}}{2i\pi}$$

- above example was $\delta = \pi/2$

~~from~~ or $\langle \psi_k^\dagger(r) \psi_k(r) \rangle = \frac{S_{(1)}^{(1)}}{2i\pi}$

where $S_{(1)}^{(1)}$ is S-matrix element for single particle to single particle scattering

- in general [eg ~~in~~ multi-channel (Kndr)] we find $|S_{(1)}^{(1)}|$ can be < 1

- we can interpret this as a finite probability for 1 particle $\rightarrow$ 1 particle + n particle-hole pair scattering at

$E \rightarrow 0$ - this violates usual Fermi liquid behavior pointed out by Landau.

T-matrix is $T = -\frac{i}{2\pi V} [1 - S_{(1)}^{(1)}]$

- resistivity at $T=0$ for dilute (Kndr) scatterers [or conductance thru quantum dot in Kndr regime]

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is proportional to $T_2 (1 - t_1)$

- so maximum resistance $\left[\frac{2e^2}{h} \text{ conductance} \right]$
occurs for single channel case

- it is reduced at NFL critical point

Explicit Realization of 2 Channel Kondo QCP

$$H = \frac{VF}{2\pi} i \int_0^{\infty} dr \left[\psi_L^\dagger + \frac{\alpha_a}{2} \frac{d}{dr} \psi_L - \psi_R^\dagger + \frac{\alpha_a}{2} \frac{d}{dr} \psi_R \right] \\ + VF \lambda \left[\psi_L^\dagger \frac{\sigma_a}{2} \psi_L + \psi_R^\dagger \frac{\sigma_a}{2} \psi_R \right], \quad a=1, 2$$

λB - equal Kondo couplings to both channels.

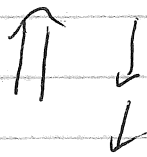
- why does this have a NFL infrared fixed point?

- To consider stability of FL fixed point we can again consider lattice model with a strong bare coupling, $T \gg t$

- but how 2 electrons get trapped at

site $j=0$ (1 from each channel) going into a $S=1$ state

- large AF coupling to impurity spin (of size $\frac{1}{2}$) leads to an $S=\frac{1}{2}$ ground state



- there is now a weak^{Kondo} coupling $\sim t^c/J$ of this $S=\frac{1}{2}$ spin to site $j=1$

- also turns out to be AF

- so, it is not consistent to assume $\frac{J}{t}$ renormalizes to ∞

- instead, we get an intermediate

- coupling fixed point (to p. 14)

- we found CIBC - checked against

Bethe Ansatz & NRG

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skip
[$g = \frac{1}{2}$ - fractional "ground state degeneracy"]

$S_{11} = 0$ - maximal inelastic scattering

in this sense

$\Rightarrow T = -\frac{i}{2\pi\nu} [1 - S_{11}]$ has

$\frac{1}{2}$ value for single channel case

$G = \frac{e^2}{h}$ not $2\frac{e^2}{h}$ at $T = V = 0$

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- Channel asymmetry is a relevant
perturbation of dimension $1/2$

- this was established (IA e A Ludwig)

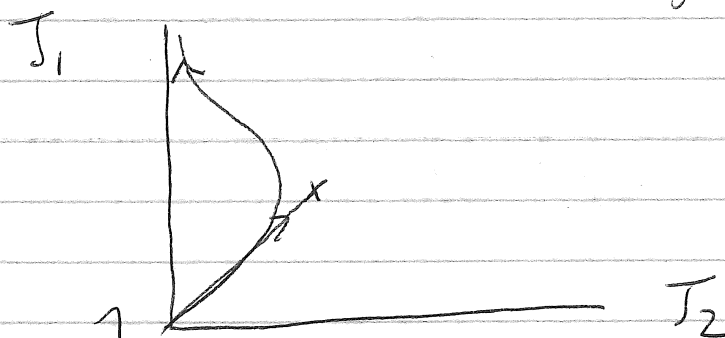
by finding most relevant operator allowed
by symmetry when $T_1 \neq T_2$

- stable fixed point corresponds to
most strongly coupled channel screening
impurity spin & other one decouples
at low E - i.e. $\delta_1 = \pi/2$, $\delta_2 = 0$

- relevant channel symmetry breaking perturbation at the QCP has dimension $1/2$

$$\Rightarrow (T_1, -T_2)_{\text{eff}}^{(\tau)} = (T_1, -T_2)_0 \left(\frac{D}{T_1} \right)^{1/2} \text{ for}$$

an appropriate UV cut off scale, D



(to part. 12)

- right at the QCP, LIO has dimension $3/2$

$$\Rightarrow \text{coupling constant } \left[\frac{1}{\sqrt{TK}} \right]$$

- various experimental realizations of 2CKE were proposed over the years:
- atoms with a ^{low-lying} quadrupolar state or other

orbital degrees of freedom playing the role of

"spin" with actual electron spin playing

the role of "channel"

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- atomic impurities in tunnel barrier
- hi-Tc superconductor,
- all were questioned
- D. Grahame-Gordon et al. finally produced an indisputable exptl. realization
- a double quantum dot device
- you might think that a single q.d. device corresponded to a 2CK system with left & right leads being 2 channels
- actually, this doesn't work



$$H_A \sim H_L^0 + H_R^0 + (t_L^\dagger \psi_L^\dagger(u) + t_R \psi_R(u) + h.c.) + U(n_d - 1)^2$$

$$H_K \sim \frac{1}{D} (\psi_L^\dagger + \psi_R^\dagger) + \sum_{\sigma} (\psi_L^\dagger + \psi_R^\dagger) \cdot \vec{J}$$

$$J \sim t^2/U$$

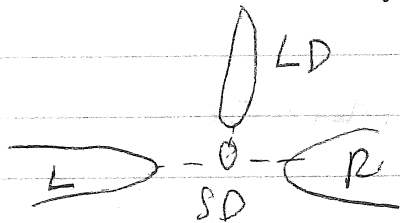
- a particular linear combination acts as a

single channel in Kondo model

[Glasman-Raikh], other channel

decouples

- in GG expt [Y Oreg e GG] a large 2^{M_1} adt
acts as a 2^{K_1} channel



- again a linear combination of 3 "channels"

$$\text{occurs: } \left(\sum_{i=1}^3 t_i \psi_i \right)^+ \rightarrow \sum_{i=1}^3 t_i' \psi_i \cdot \vec{S}$$

- however, due to its finite size, charging

energy of LD is important ($\epsilon \approx \psi_3$)

$$\delta H = \frac{U_d}{2} (\hat{n}_3 - n_0)^2$$

- at energy scales below U_d tunnelling

on or off LD is suppressed, leaving $2CKM$

$$: \psi_3, \frac{\epsilon_1 \psi_1 + \epsilon_2 \psi_2}{\sqrt{\epsilon_1^2 + \epsilon_2^2}}$$

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- size of LD must be small enough that $T_K \ll U$ but large enough that finite-size ~~low~~ single particle level spacing $\ll T_K$
- then must tune gates so $T_1 \approx T_2$
- CFT methods predict dependence of conductance on T e source-drain voltage of form:

$$G \equiv \frac{dI}{dV_{sd}} = \frac{e^2}{h} \left[1 - \left(\frac{T}{T_K} \right)^{1/2} F\left(\frac{eV_{sd}}{T}\right) \right]$$

(where F is not known, in general)

$\frac{e^2}{h}$ factor $\left[\frac{1}{2} \text{ of } \frac{2e^2}{h} \right]$ & exponent $\frac{1}{2}$ follow simply from CFT results

- cross-over energy scale is determined

from combining weak & strong coupling RG:

$$T_c \propto T_K \frac{(\lambda_1 - \lambda_2)^2}{(\lambda_1 + \lambda_2)^4} \quad (\text{Pustelnik, Borda, Blazman, von Delft})$$

- Conclusion: BCFT is a general approach to studying low energy RG fixed points of QIP's
- impurity degrees of freedom are integrated out, and we are left with a CIBC
- this method can reproduce in a systematic way FL f.p.'s found ~~earlier~~ as well as NFL QCP's that occur in some models