

Outline

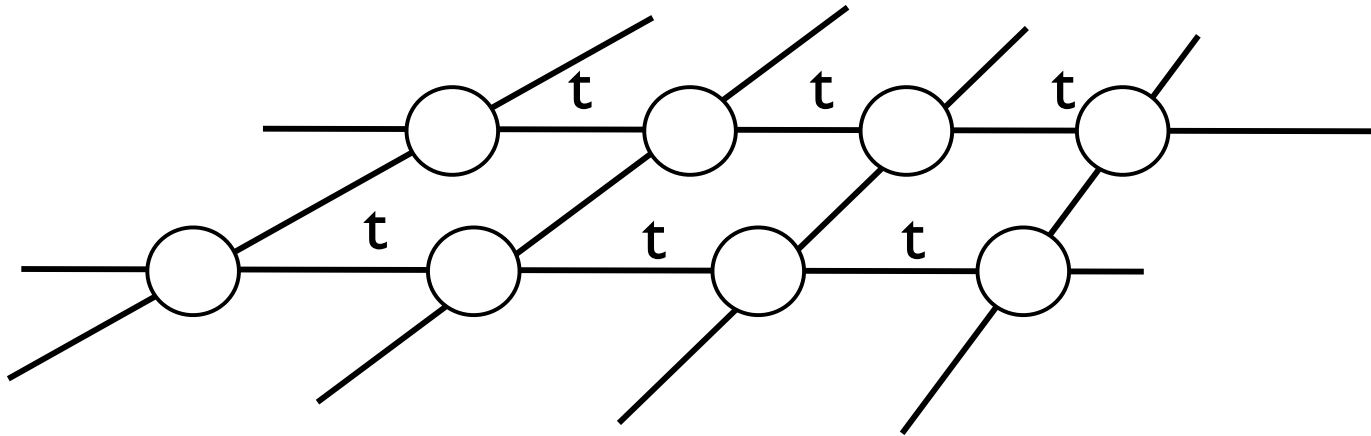
Overview of the the Kondo lattice model
Nonlinear canonical transformation on
even valence fermi systems
Transformations on odd valence insulators

Nonlinear canonical transformations applied to metal insulator transition

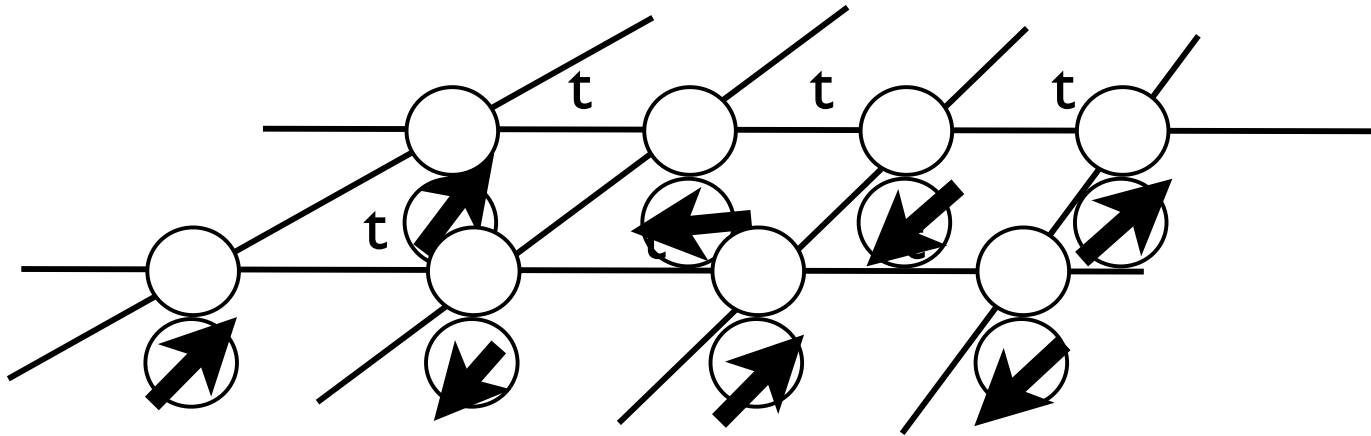
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Göteborgs Universitet

Nordita 2014-09-03

Kondo lattice model: begin with an ordinary tight binding lattice.



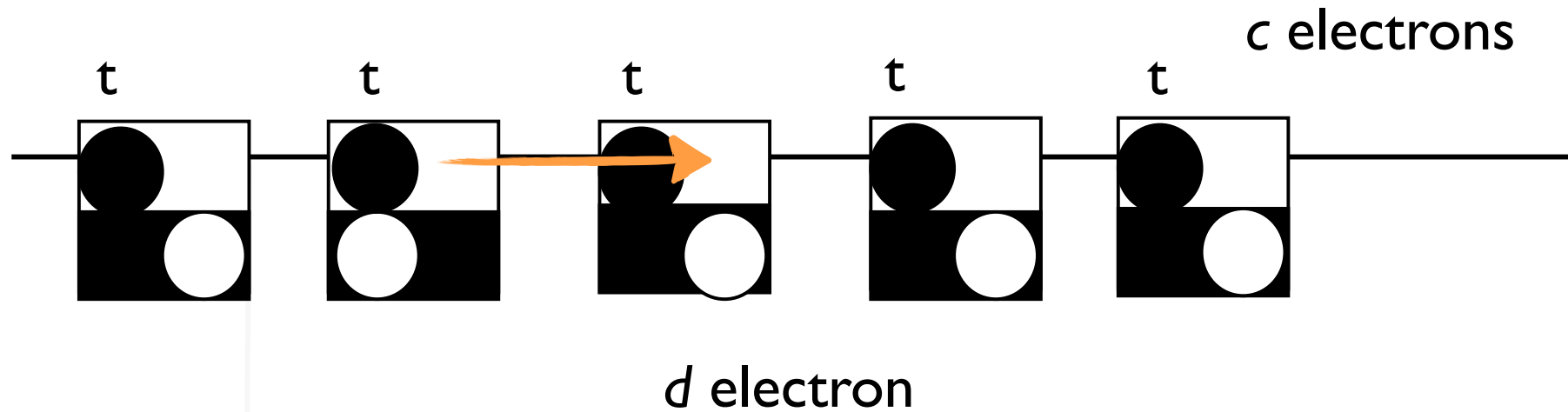
$$-t \sum_{rr'} (c_{\sigma r}^{\dagger} c_{\sigma r'} + C C)$$



and layer of deeper localized spins (electrons)

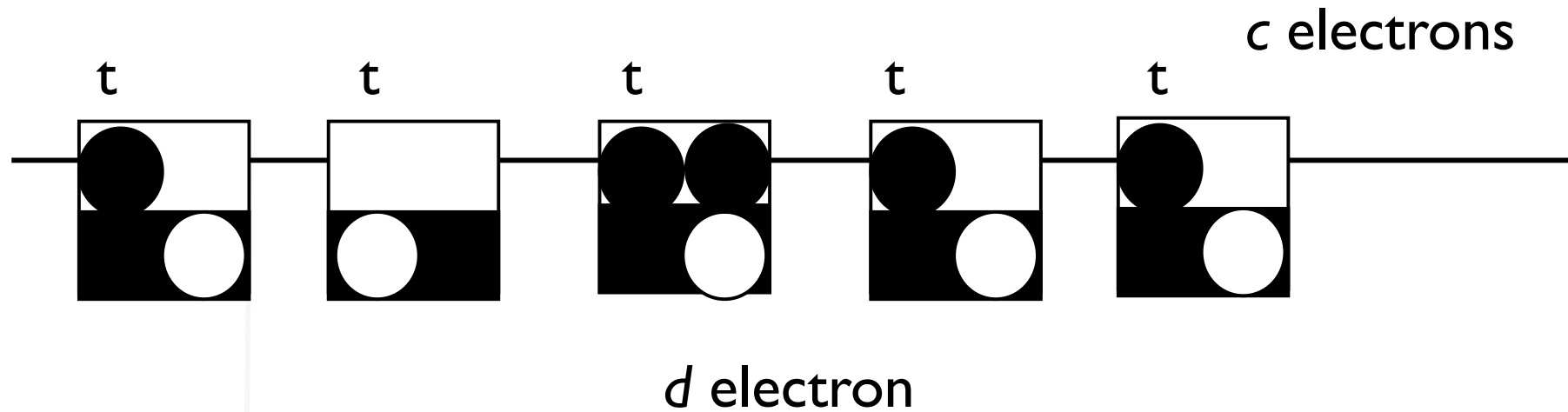
$$H_{Kondo} = \sum_r (J S_r \cdot S_{d,r} + U_d (n_d - 1)^2)$$

Kondo Lattice model:



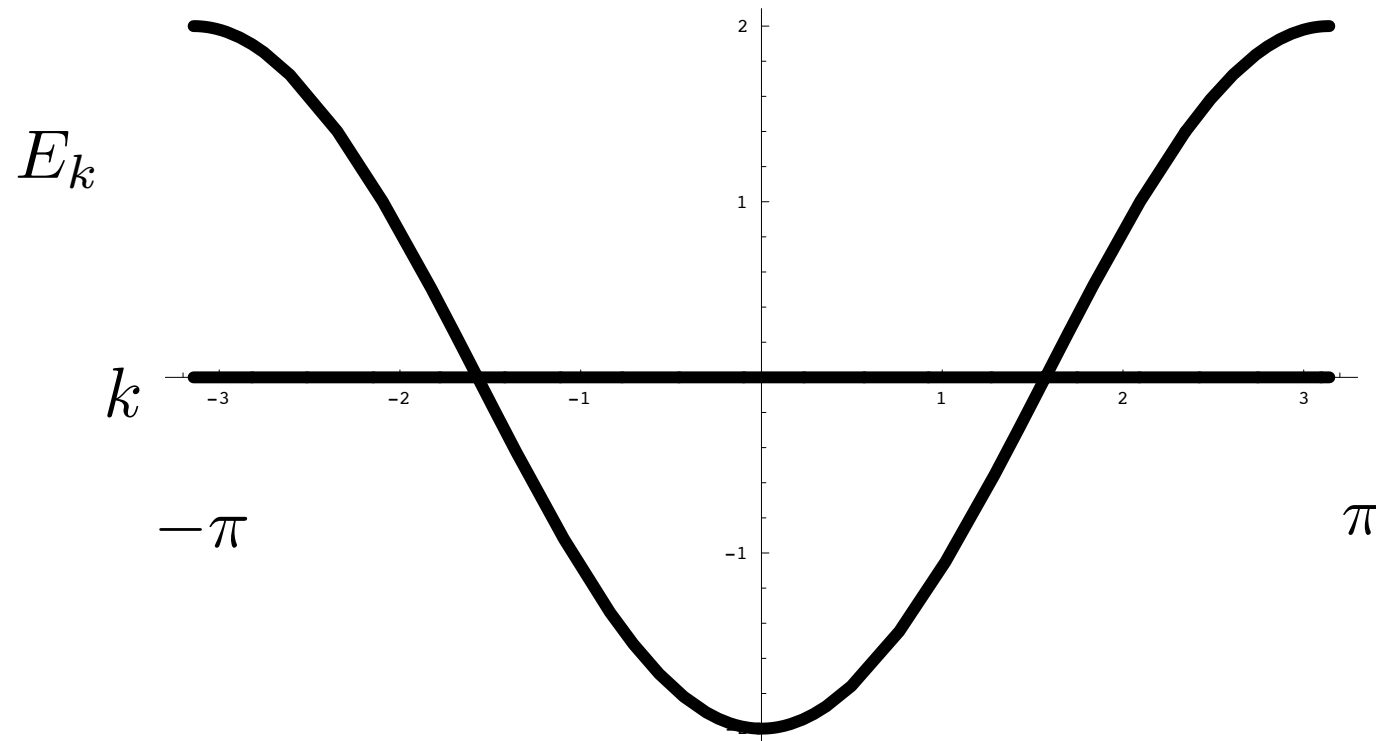
$$H_{KLM} = -t \sum_{rr'} (c_{\sigma r}^\dagger c_{\sigma r'} + CC) + \sum_r (JS_r \cdot S_{d,r} + U_d(n_d - 1)^2)$$

Kondo Lattice model:



$$H_{KLM} = -t \sum_{rr'} (c_{\sigma r}^\dagger c_{\sigma r'} + CC) + \sum_r (JS_r \cdot S_{d,r} + U_d(n_d - 1)^2)$$

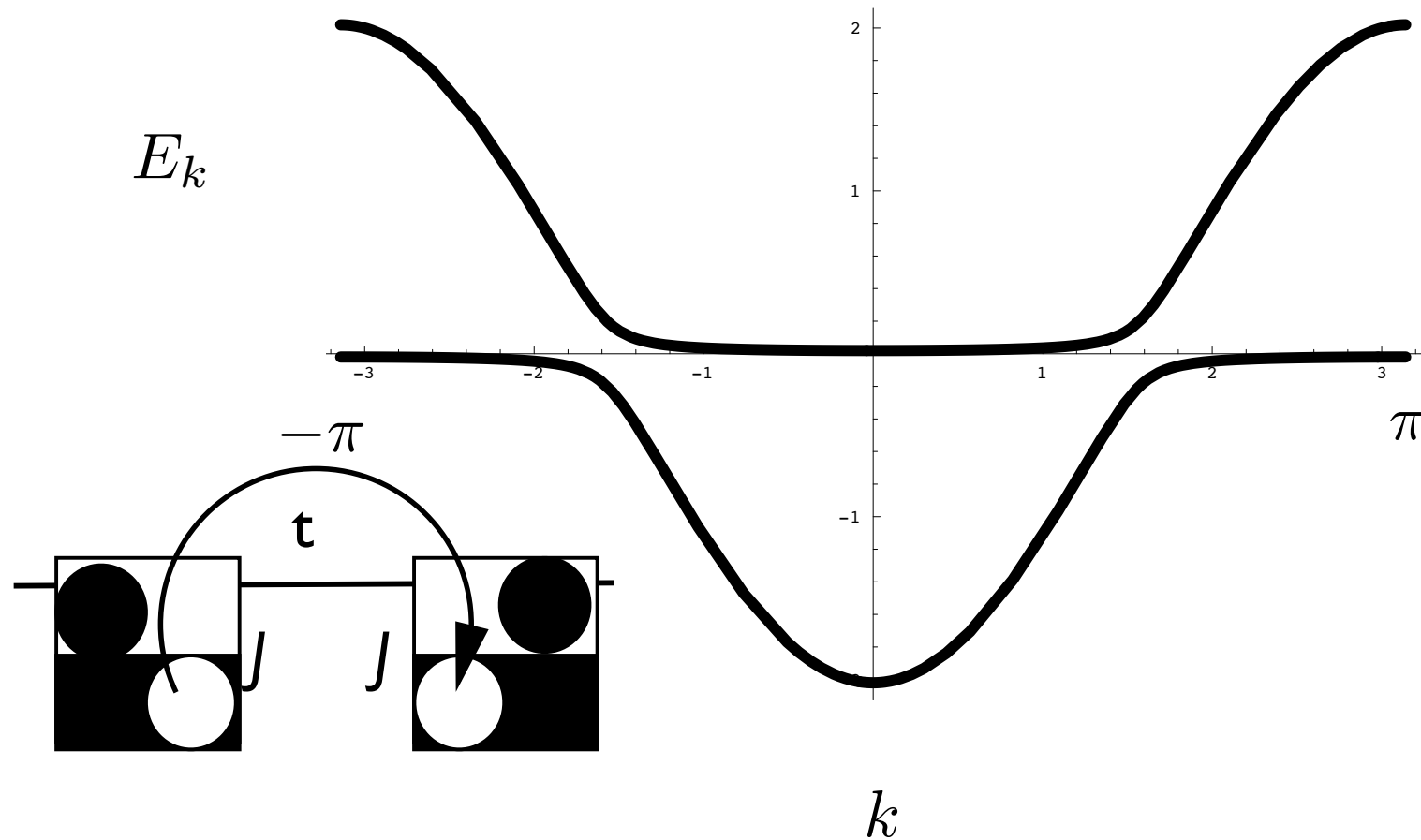
weak coupling $J=0$ “RKKY” regime



$$H_{KLM} =$$

$$-t \sum_{rr'} (c_{\sigma r}^\dagger c_{\sigma r'} + c_{\sigma r'} c_{\sigma r}) + \sum_r (J S_r \cdot S_{d,r} + U_d (n_d - 1)^2)$$

Kondo Lattice model -
weak coupling $J \sim 0$ “RKKY” regime

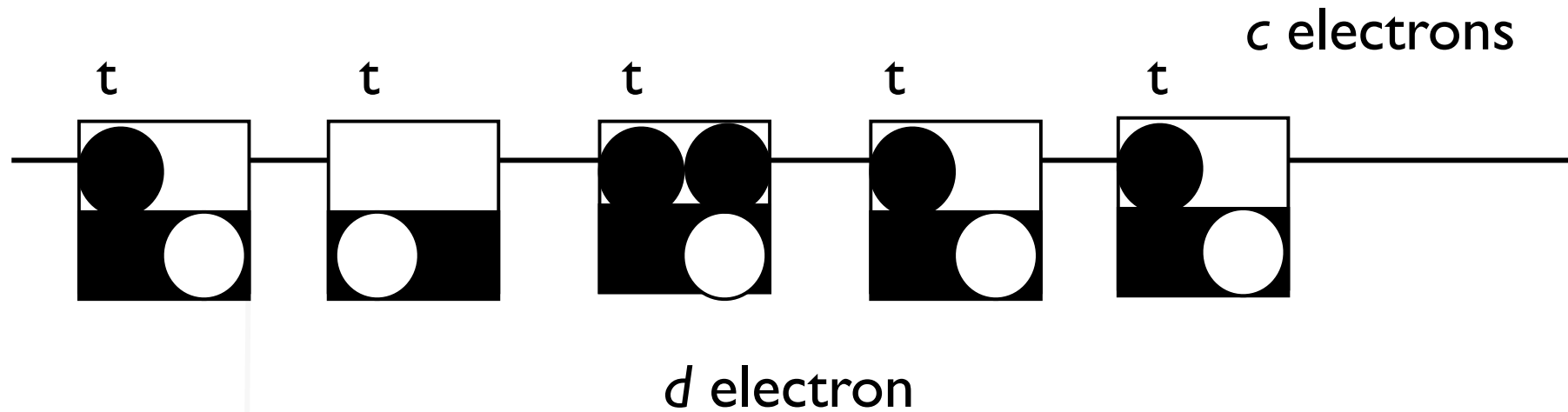


$$H_{int} = \sum_r (J S_r \cdot S_{d,r} + U_d (n_d - 1)^2)$$

Strong coupling limit is very different. $J \gg t$ gives a singlet ground state, with the energy gap of the singly occupied and single vacancy states determining the quasiparticle gap and is an insulator.

$$\frac{1}{\sqrt{2}} \left(\begin{array}{|c|} \hline \bullet \\ \hline \blacksquare \\ \hline \end{array} - \begin{array}{|c|} \hline \blacksquare \\ \hline \bullet \\ \hline \end{array} \right)$$

Generalize from Kondo to “Periodic Anderson” models



$$h_{atomic} = \epsilon_d n_d + U n_{d\uparrow} \cdot n_{d\downarrow} + V \sum_{\sigma} (c_{0,\sigma}^{\dagger} d_{\sigma} + CC)$$

We look more carefully at the ground state atomic singlet.

$$|\Psi_G\rangle = \frac{1}{\sqrt{2}} \left(\begin{array}{|c|} \hline \bullet \\ \hline \blacksquare \\ \hline \end{array} - \begin{array}{|c|} \hline \blacksquare \\ \hline \bullet \\ \hline \end{array} \right)$$

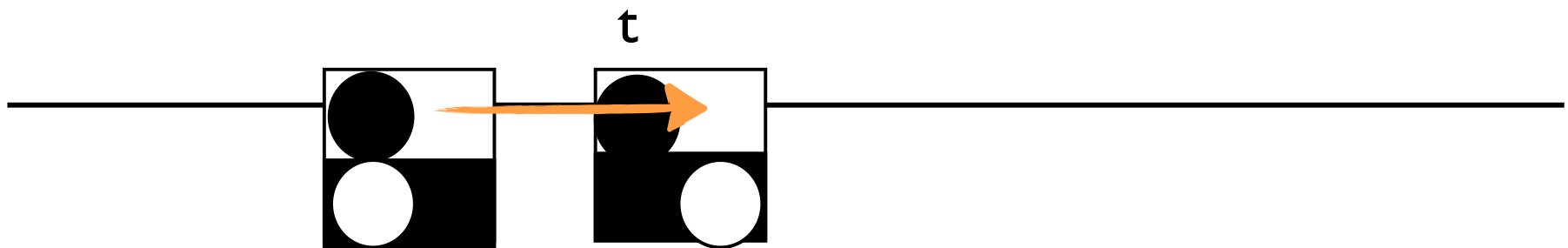
Can we create a local quasiparticle operator which creates an electron or hole on a singlet site?

Quasiparticles require:

A $\hat{c}_\alpha |\Psi_G\rangle = 0$

B $\hat{c}_\alpha \hat{c}_\alpha^\dagger |\Psi_G\rangle = |\Psi_G\rangle$

Do bare electron operators qualify?

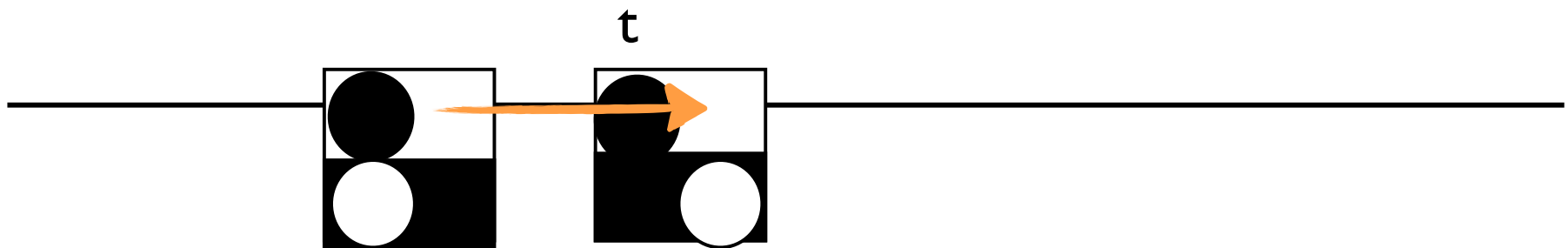


Quasiparticles require:

A $\hat{c}_\alpha |\Psi_G\rangle = 0$

$$c_{c,\uparrow} |\Psi_G\rangle = \frac{1}{\sqrt{2}} c_{f,\downarrow}^\dagger |0\rangle \neq 0$$

B $\hat{c}_\alpha \hat{c}_\alpha^\dagger |\Psi_G\rangle = |\Psi_G\rangle$

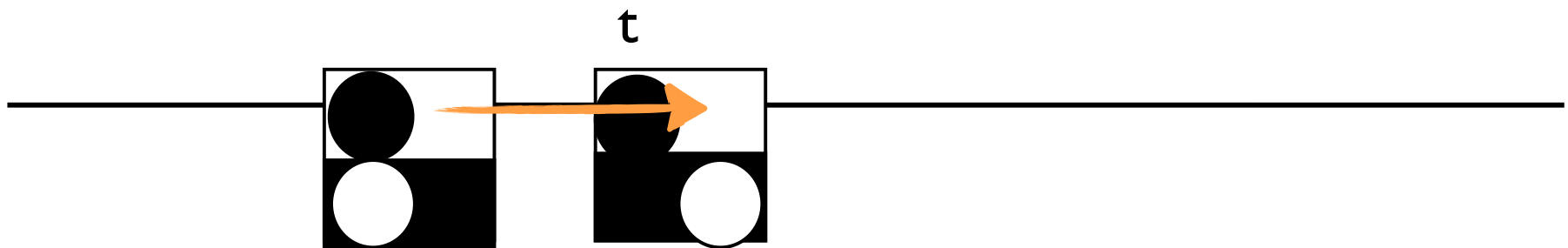


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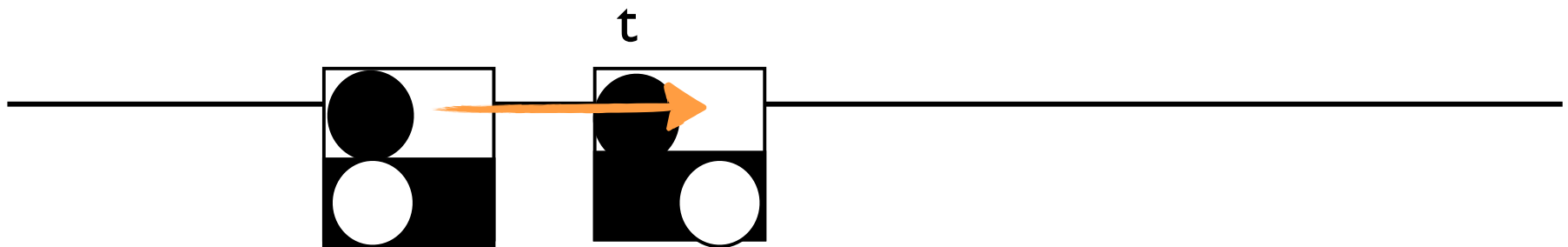


Quasiparticles require:

A $\hat{c}_\alpha |\Psi_G\rangle = 0$

~~$$c_{c,\uparrow} |\Psi_G\rangle = \frac{1}{\sqrt{2}} c_{f,\downarrow}^\dagger |0\rangle \neq 0$$~~

B $\hat{c}_\alpha \hat{c}_\alpha^\dagger |\Psi_G\rangle = |\Psi_G\rangle \quad c_{c,\uparrow} c_{c,\uparrow}^\dagger |\Psi_G\rangle = -\frac{1}{\sqrt{2}} c_{c,\downarrow}^\dagger c_{f,\uparrow}^\dagger |Vac\rangle \neq |\Psi_G\rangle$



Quasiparticles require:

A $\hat{c}_\alpha |\Psi_G\rangle = 0$

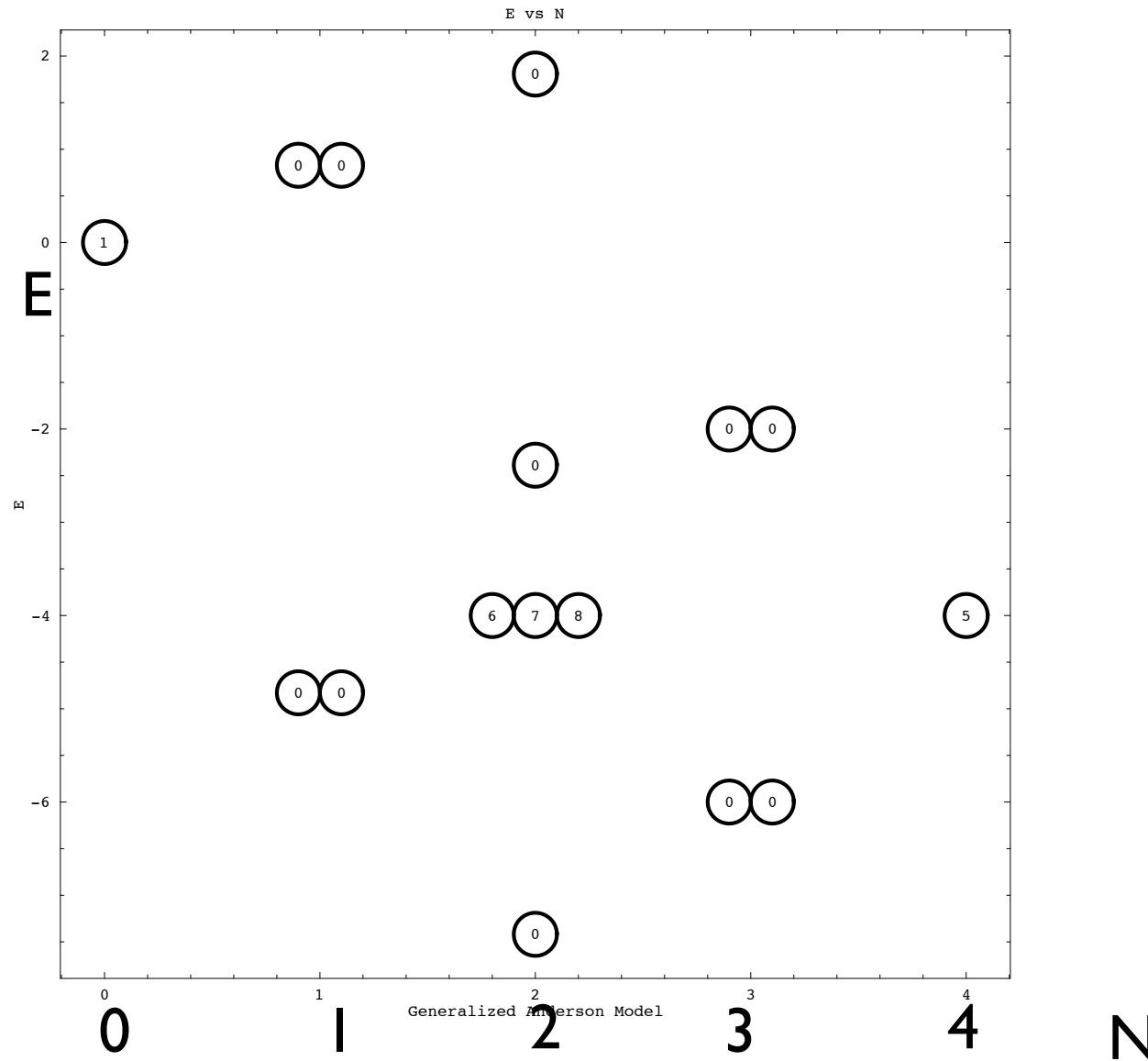
~~$$c_{c,\uparrow} |\Psi_G\rangle = \frac{1}{\sqrt{2}} c_{f,\downarrow}^\dagger |0\rangle \neq 0$$~~

B $\hat{c}_\alpha \hat{c}_\alpha^\dagger |\Psi_G\rangle = |\Psi_G\rangle$ ~~$c_{c,\uparrow} c_{c,\uparrow}^\dagger |\Psi_G\rangle = -\frac{1}{\sqrt{2}} c_{c,\downarrow}^\dagger c_{f,\uparrow}^\dagger |Vac\rangle \neq |\Psi_G\rangle$~~

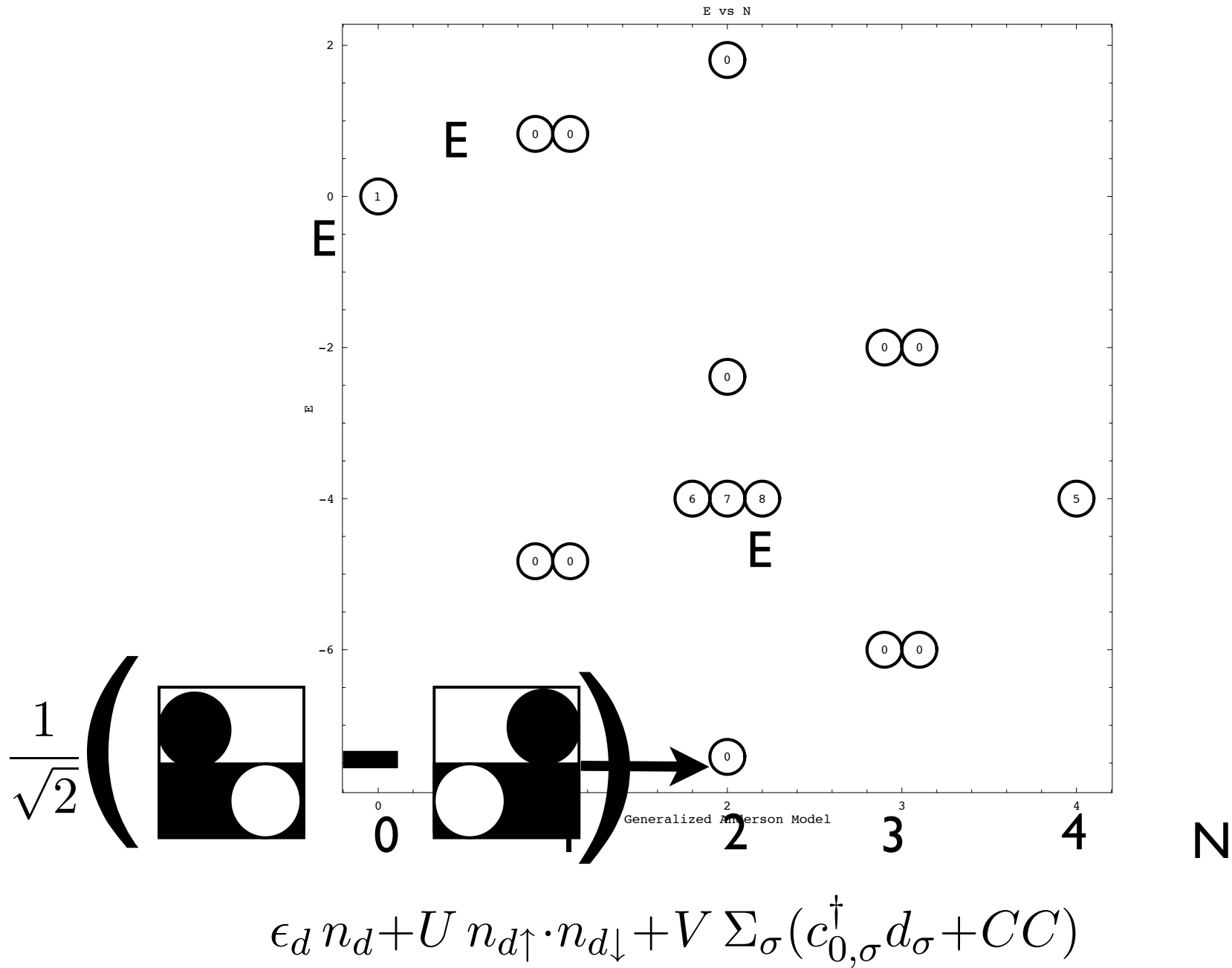
Bare electron operators are no good!

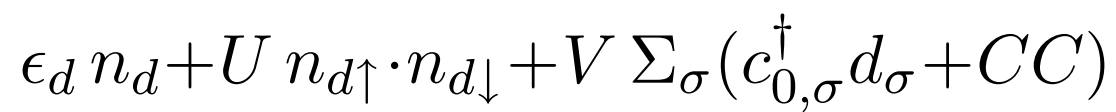
We now turn to dealing with this problem.

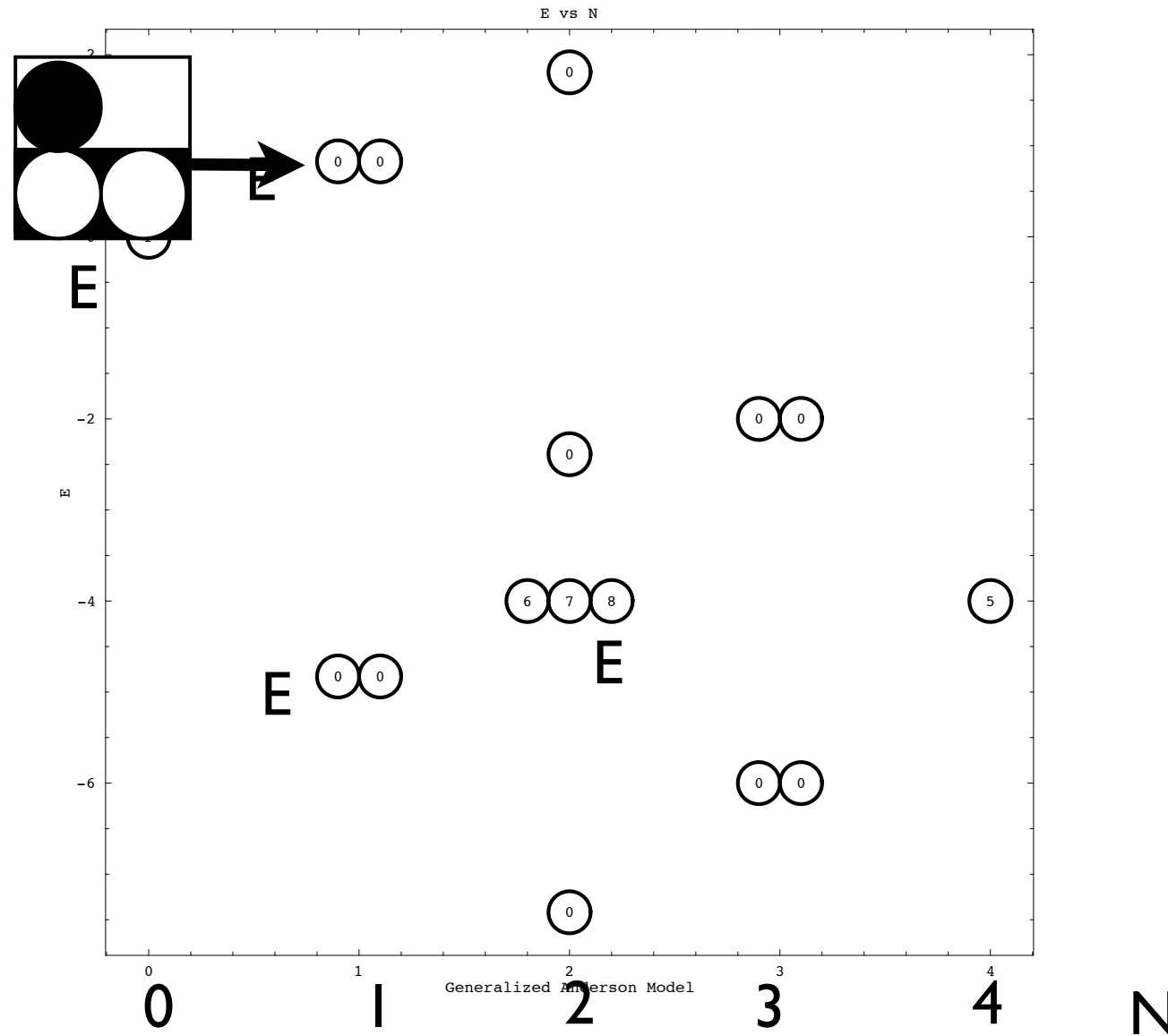
We begin by taking a “generic” value of the Anderson impurity parameters and plotting the energy values for different values of the impurity site occupation number.



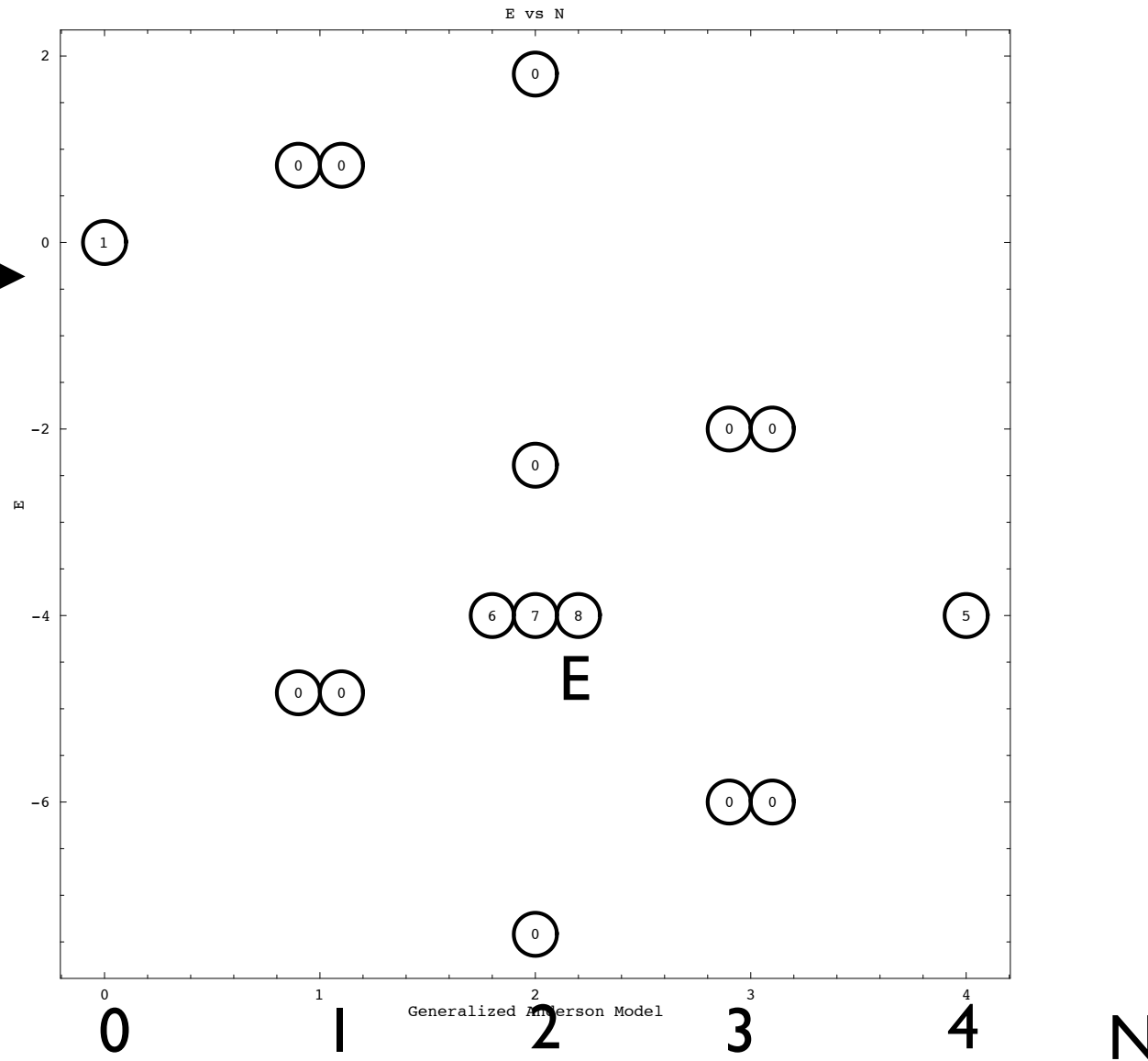
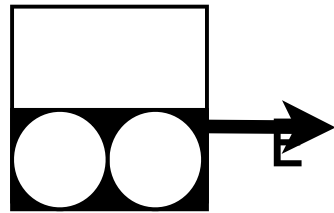
$$\epsilon_d n_d + U n_{d\uparrow} \cdot n_{d\downarrow} + V \sum_{\sigma} (c_{0,\sigma}^{\dagger} d_{\sigma} + C C)$$



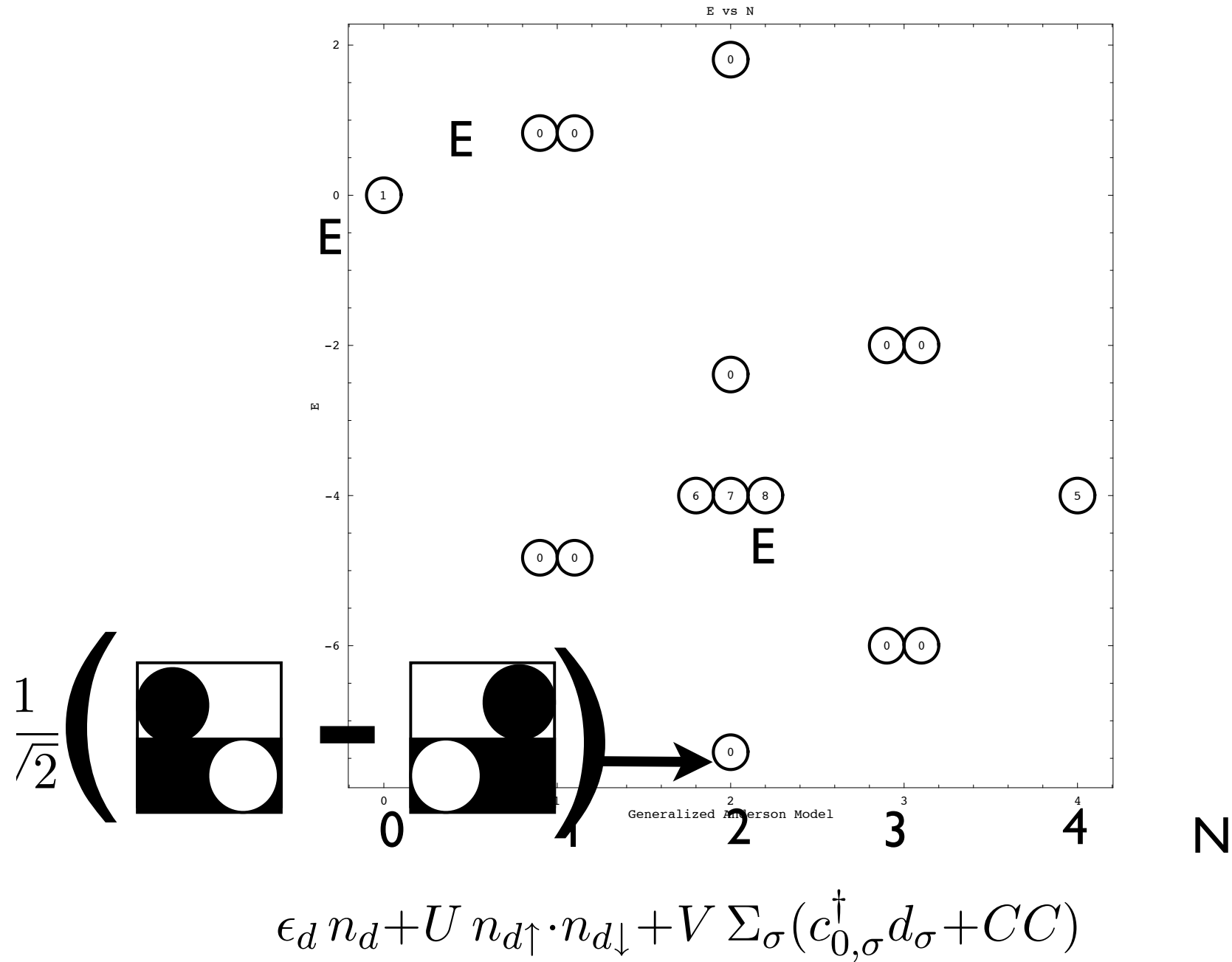


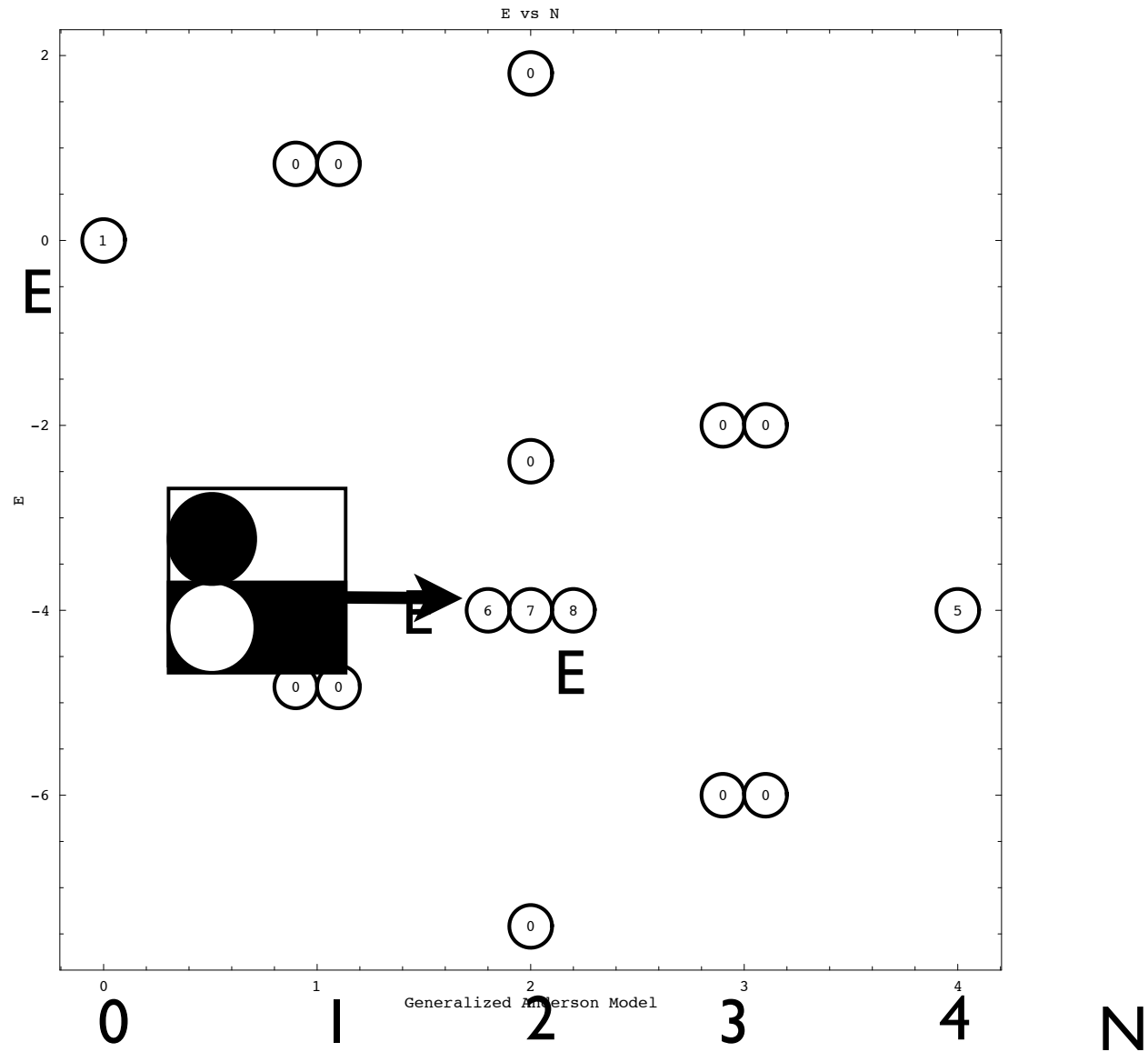


$$\epsilon_d n_d + U n_{d\uparrow} \cdot n_{d\downarrow} + V \sum_{\sigma} (c_{0,\sigma}^{\dagger} d_{\sigma} + C C)$$

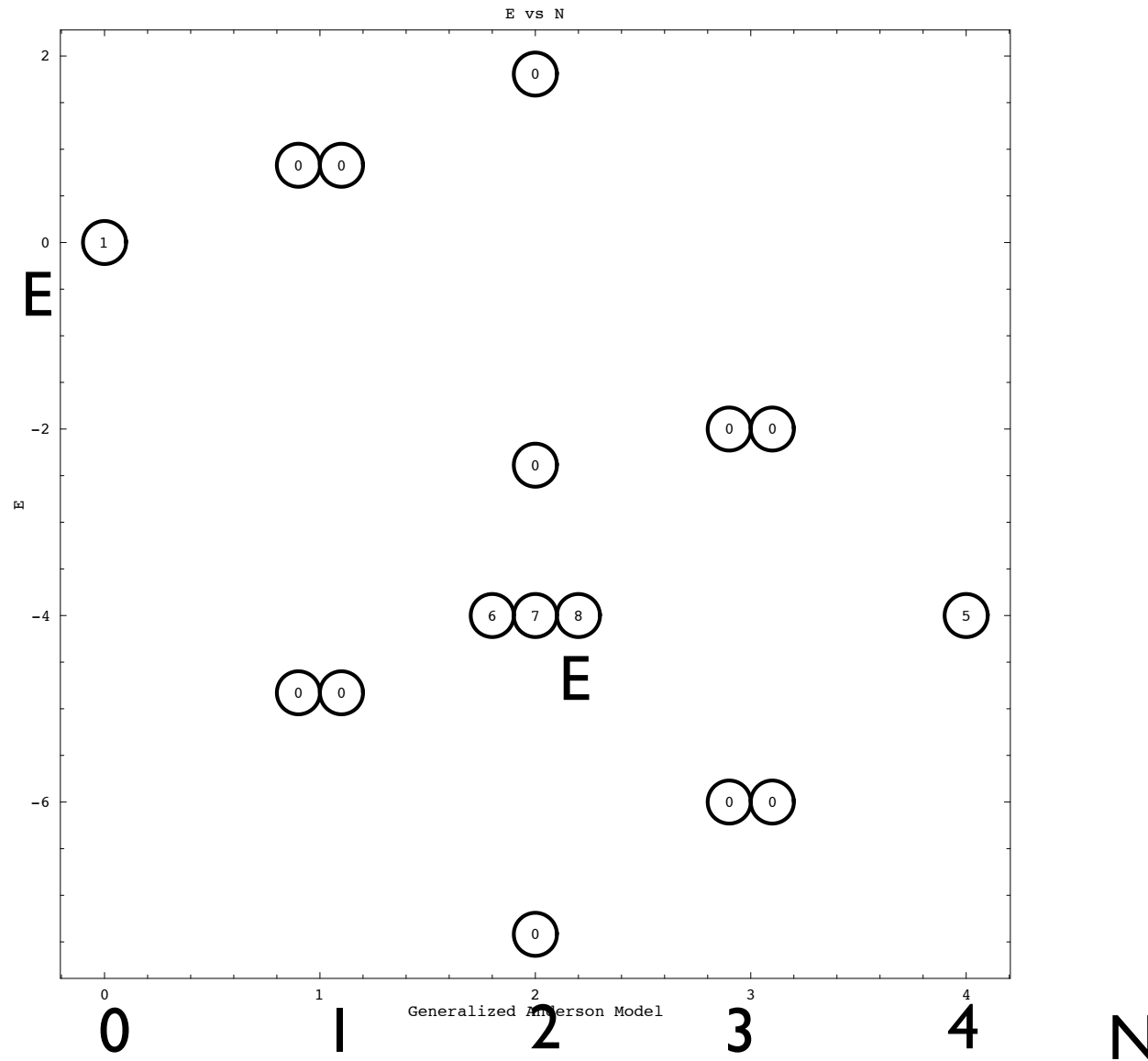


$$\epsilon_d n_d + U n_{d\uparrow} \cdot n_{d\downarrow} + V \sum_{\sigma} (c_{0,\sigma}^{\dagger} d_{\sigma} + C C)$$



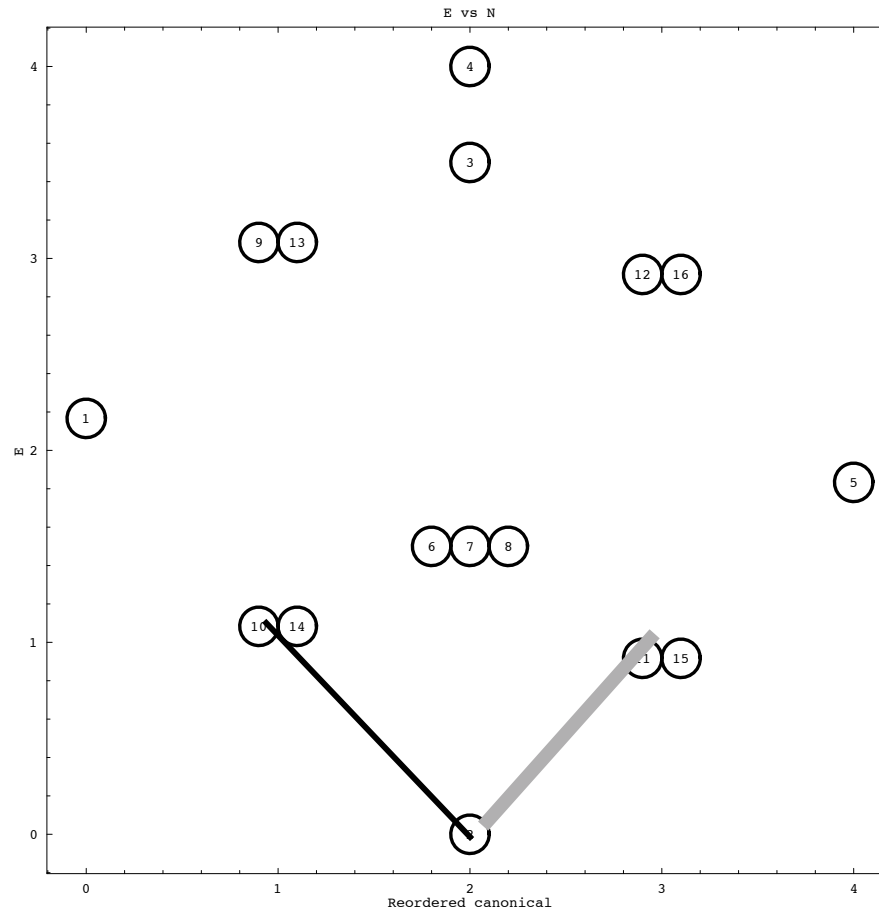


$$\epsilon_d n_d + U n_{d\uparrow} \cdot n_{d\downarrow} + V \sum_{\sigma} (c_{0,\sigma}^{\dagger} d_{\sigma} + C C)$$



$$\epsilon_d n_d + U n_{d\uparrow} \cdot n_{d\downarrow} + V \sum_{\sigma} (c_{0,\sigma}^{\dagger} d_{\sigma} + CC)$$

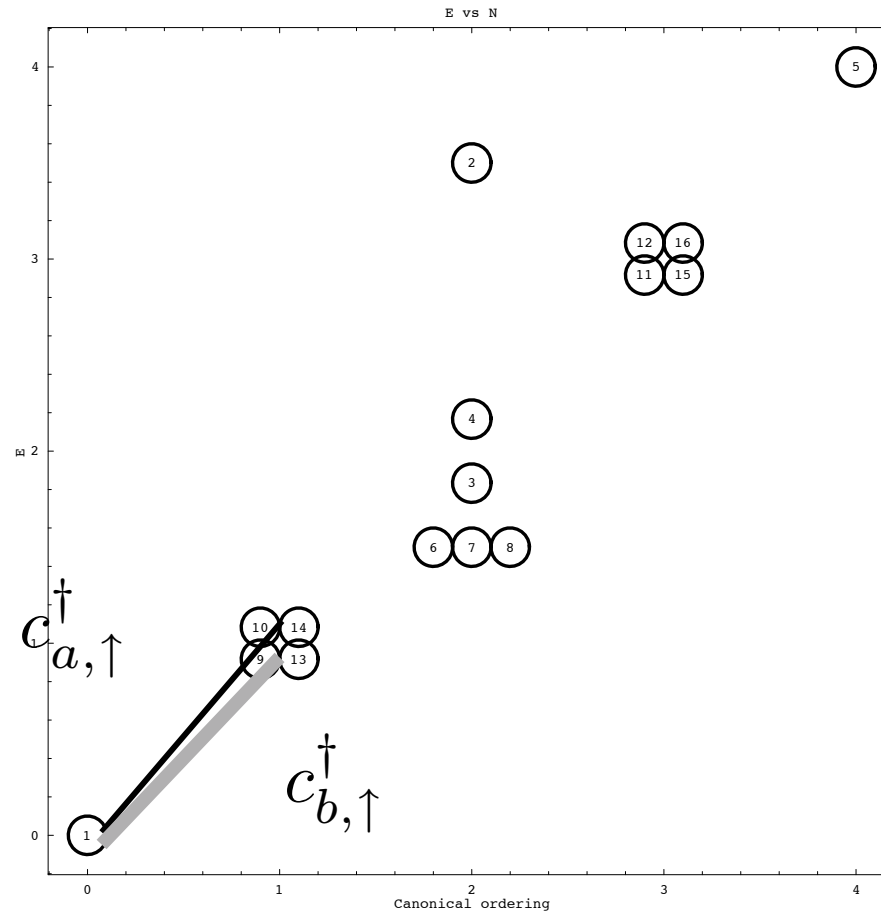
Specific Hamiltonian



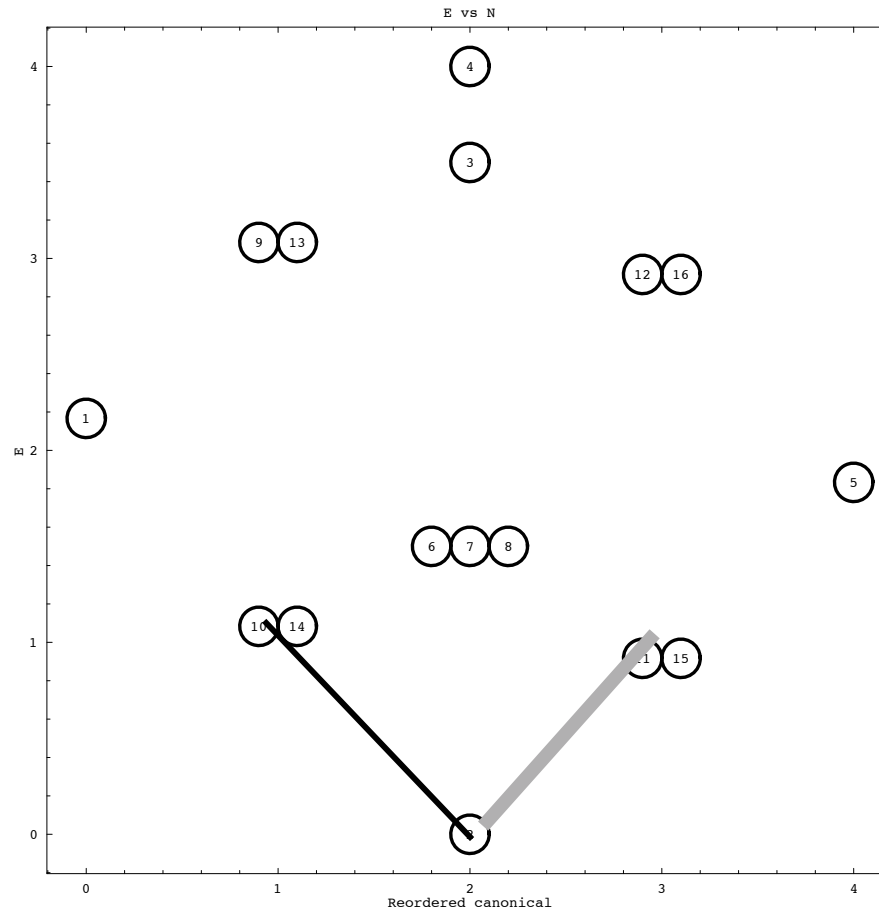
$$\epsilon_d = -\frac{1}{2}U \text{ and } V = 2, U = 8 .$$

$$\epsilon_d n_d + U n_{d\uparrow} \cdot n_{d\downarrow} + V \sum_{\sigma} (c_{0,\sigma}^{\dagger} d_{\sigma} + CC)$$

Another simple Fermi system

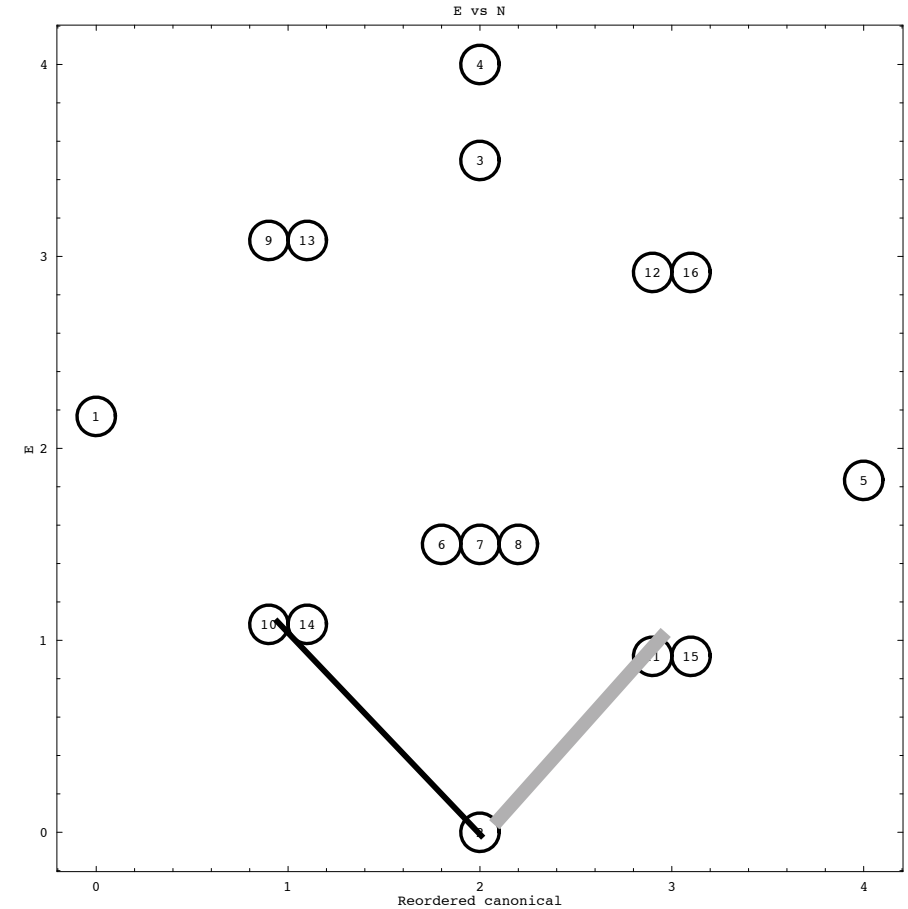
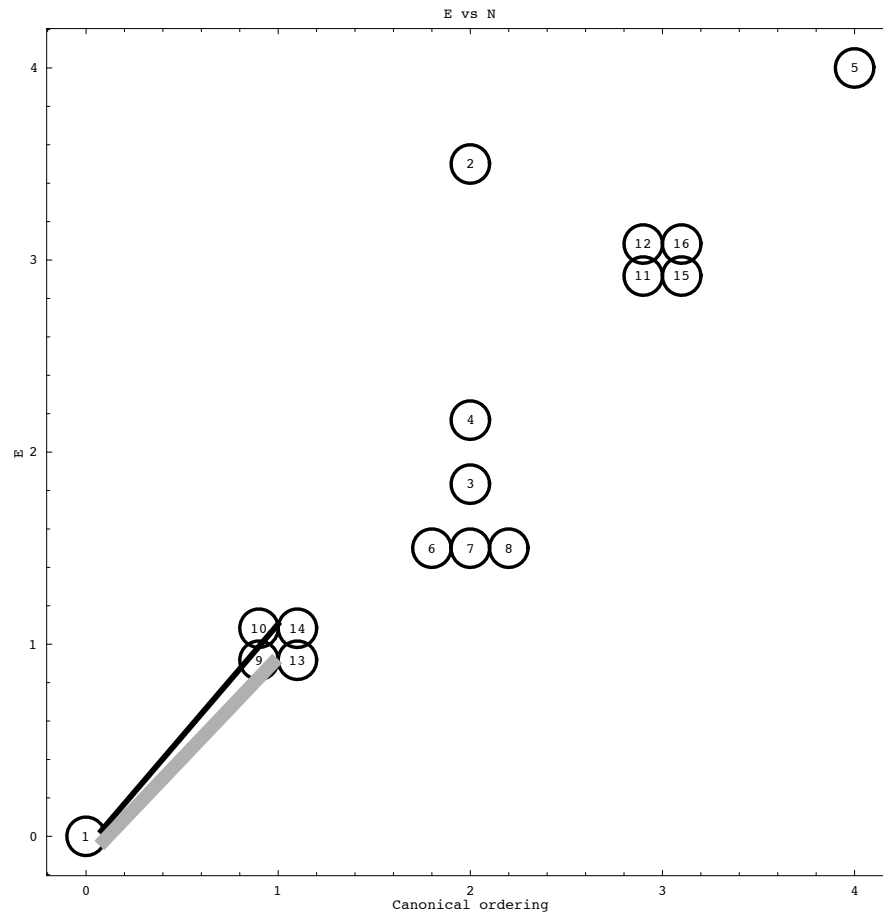


$$n_{tot} + \frac{1}{12}(n_a - n_b) - \frac{1}{2}S_a \cdot S_b$$



$$\epsilon_d = -\frac{1}{2}U \text{ and } V = 2, U = 8 .$$

$$\epsilon_d n_d + U n_{d\uparrow} \cdot n_{d\downarrow} + V \sum_{\sigma} (c_{0,\sigma}^{\dagger} d_{\sigma} + CC)$$

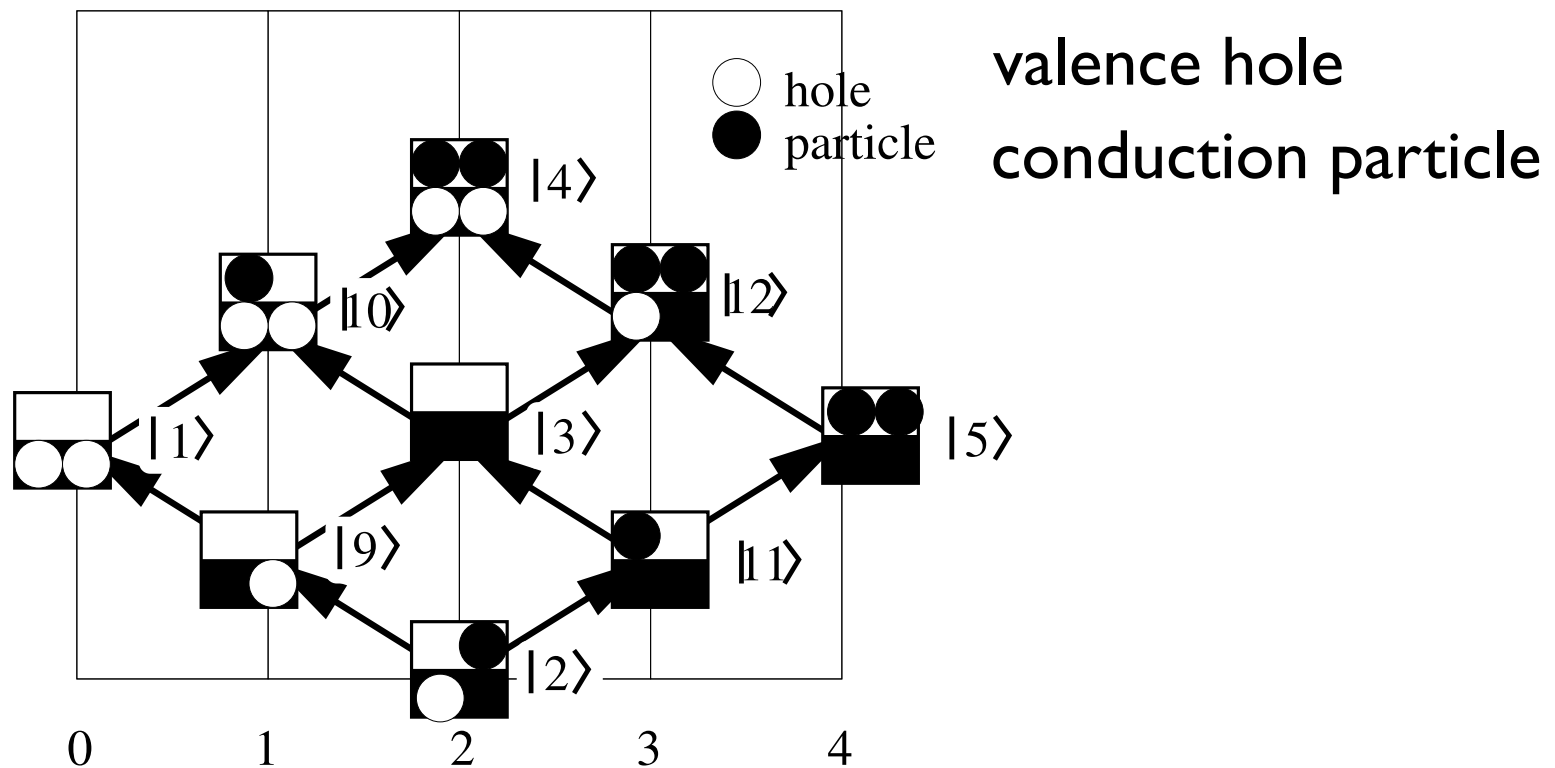


$$n_{tot} + \frac{1}{12}(n_a - n_b) - \frac{1}{2}S_a \cdot S_b$$

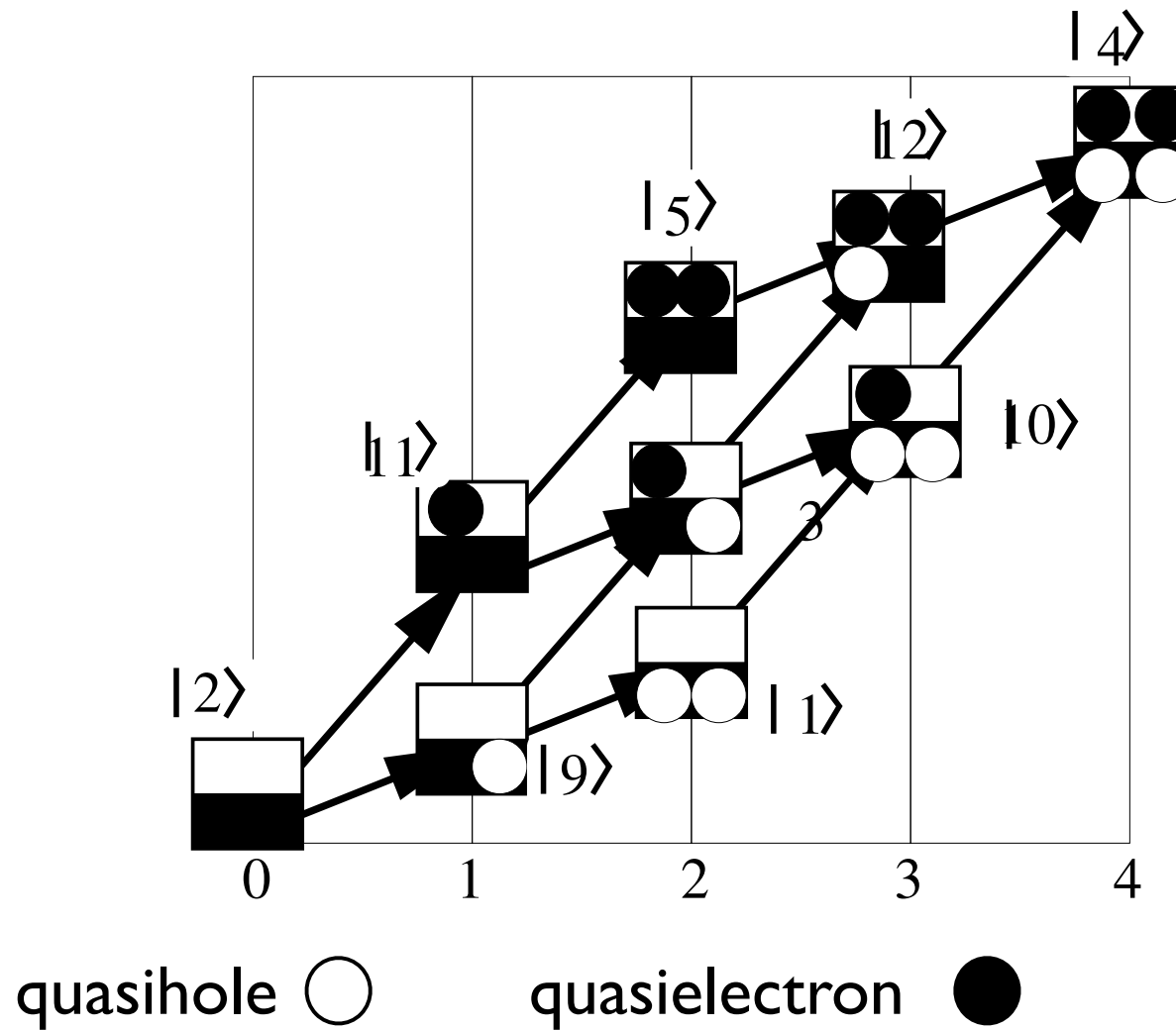
$$\epsilon_d = -\frac{1}{2}U \text{ and } V = 2, U = 8.$$

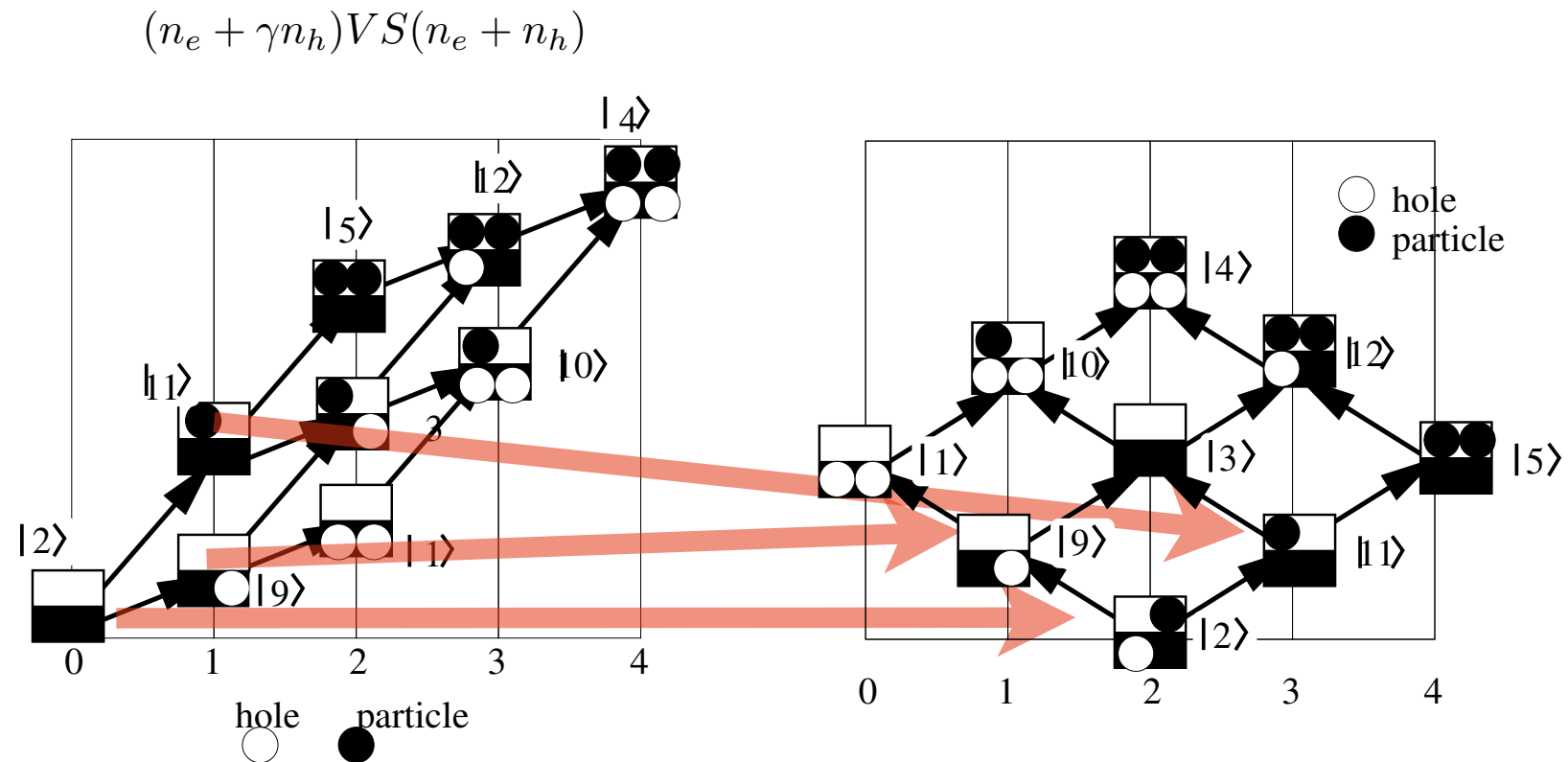
$$\epsilon_d n_d + U n_{d\uparrow} \cdot n_{d\downarrow} + V \sum_{\sigma} (c_{0,\sigma}^{\dagger} d_{\sigma} + CC)$$

Pictogram of the interacting state



$$(n_e + \gamma n_h)VS(n_e + n_h)$$





A permutation is a canonical transformation ** and can be represented by:

$$U = P(\hat{c}_e, \hat{c}_h, \hat{c}_e^\dagger, \hat{c}_h^\dagger)$$

Invert this

$$c_c^\dagger = U^\dagger \hat{c}_e^\dagger U$$

$$c_f = U^\dagger \hat{c}_h^\dagger U$$

This transformation is then done **exactly** with computer algebra

This reordering generates a canonical transformation that by construction maps states properly

$$\begin{aligned}
 c_{c,\uparrow}^\dagger = & \hat{c}_{h,\downarrow} \left(\tau_2 n_{e,\downarrow} + \frac{1}{2} \hat{c}_{h,\uparrow}^\dagger \hat{c}_{e,\downarrow}^\dagger - \tau_1 n_{e,\uparrow} - \frac{1}{\sqrt{2}} n_{h,\uparrow} \right) \\
 & + \hat{c}_{e,\uparrow}^\dagger \left(\tau_1 n_{h,\downarrow} + \frac{-1}{2} \hat{c}_{h,\uparrow} \hat{c}_{e,\downarrow} + \frac{1}{\sqrt{2}} n_{e,\downarrow} - \tau_2 n_{h,\uparrow} \right) \\
 & + \left(\frac{1}{\sqrt{2}} \hat{c}_{h,\downarrow} - \frac{1}{\sqrt{2}} \hat{c}_{e,\uparrow}^\dagger \right) + \dots,
 \end{aligned}$$

$$\tau_2 = (1 - 1/\sqrt{2}) \quad \text{and} \quad \tau_1 = \frac{1}{2}(\sqrt{2} - 1).$$

This reordering generates a canonical transformation that by construction maps states properly

$$\begin{aligned}
 c_{c,\uparrow}^\dagger = & \hat{c}_{h,\downarrow} \left(\tau_2 n_{e,\downarrow} + \frac{1}{2} \hat{c}_{h,\uparrow}^\dagger \hat{c}_{e,\downarrow}^\dagger - \tau_1 n_{e,\uparrow} - \frac{1}{\sqrt{2}} n_{h,\uparrow} \right) \\
 & + \hat{c}_{e,\uparrow}^\dagger \left(\tau_1 n_{h,\downarrow} + \frac{-1}{2} \hat{c}_{h,\uparrow} \hat{c}_{e,\downarrow} + \frac{1}{\sqrt{2}} n_{e,\downarrow} - \tau_2 n_{h,\uparrow} \right) \\
 & + \left(\frac{1}{\sqrt{2}} \hat{c}_{h,\downarrow} - \frac{1}{\sqrt{2}} \hat{c}_{e,\uparrow}^\dagger \right) + \dots,
 \end{aligned}$$

“Bogoliubov”

$$\tau_2 = (1 - 1/\sqrt{2}) \quad \text{and} \quad \tau_1 = \frac{1}{2}(\sqrt{2} - 1).$$

$$\begin{aligned} c_{f,\uparrow}^\dagger = & \hat{c}_{e,\uparrow}^\dagger \left(\frac{1}{2} n_{h,\downarrow} + n_{e,\downarrow} + \frac{-1}{2} \hat{c}_{h,\uparrow} \hat{c}_{e,\downarrow} + \frac{1}{\sqrt{2}} \hat{c}_{h,\uparrow}^\dagger \hat{c}_{e,\downarrow}^\dagger \right) \\ & + \hat{c}_{h,\downarrow} \left(\frac{-1}{2} n_{e,\uparrow} + \frac{1}{2} \hat{c}_{h,\uparrow}^\dagger \hat{c}_{e,\downarrow}^\dagger - \frac{1}{\sqrt{2}} \hat{c}_{h,\uparrow} \hat{c}_{e,\downarrow} - n_{h,\uparrow} \right) + \cdots . \end{aligned}$$

Note no Bogoliubov-like terms

We get an exact mapping to electrons and holes

$$n_{tot} = 2 + \hat{c}_{e,\uparrow}^\dagger \hat{c}_{e,\uparrow} + \hat{c}_{e,\downarrow}^\dagger \hat{c}_{e,\downarrow} - \hat{c}_{h,\downarrow}^\dagger \hat{c}_{h,\downarrow} - \hat{c}_{h,\uparrow}^\dagger \hat{c}_{h,\uparrow}$$

The local atomic Kondo interaction becomes

$$\begin{aligned} H_{KLM} = & 3Jn_e + \frac{1}{4}n_e(n_e - 1)(-6J + U_f) & + \\ & \frac{1}{8}n_e(n_e - 1)n_h(6J - U_f) + \frac{1}{8}n_en_h(-18J + U_f) & + \\ & \frac{1}{8}n_en_h(n_h - 1)(6J - U_f) + 3Jn_h & + \\ & \frac{1}{4}n_h(n_h - 1)(-6J + U_f) + \frac{1}{8}(S_c \cdot S_f)(2J - U_f) \end{aligned}$$

The original Kondo Lattice model

$$H_{KLM} = -t \sum_{rr'} (c_{c,\sigma}^\dagger(r) c_{c,\sigma}(r') + CC) + \sum_r (J S_c(r) \cdot S_f(r) + U_f(n_f(r) - 1)^2)$$

becomes a **dilute** fermi gas which is then truncated to a free fermion model in electron and hole operators

$$\begin{aligned} H_{KLM}^{eff} = & J(n_e(r) + n_h(r)) + \\ & \sum_s (-1)^s t ((\hat{c}_{e,s}^\dagger(r) \hat{c}_{h,-s}^\dagger(r') - \hat{c}_{e,s}(r') \hat{c}_{h,-s}(r) + CC)) + \\ & \sum_s (t(\hat{c}_{e,s}^\dagger(r) \hat{c}_{e,s}(r') - \hat{c}_{h,s}^\dagger(r) \hat{c}_{h,s}(r')) + CC) \end{aligned}$$

(Note the the Hamiltonian is approximate but the transformation is exact. This is important when computing expectation values of functions of bare electron operators!)

For the record, the exact transformation can be written down and used but the details are not interesting

$$c_{c,\uparrow}^\dagger = \hat{c}_{h,\downarrow} (\tau_1 n_{h,\uparrow} n_{e,\uparrow} + \tau_2 n_{e,\downarrow} + \tau_1 n_{h,\uparrow} n_{e,\downarrow} + \tau_1 n_{e,\uparrow} n_{e,\downarrow} + \frac{1}{2} \hat{c}_{h,\uparrow}^\dagger \hat{c}_{e,\downarrow}^\dagger - \tau_1 n_{e,\uparrow} - \frac{1}{\sqrt{2}} n_{h,\uparrow}) + \\ \hat{c}_{e,\uparrow}^\dagger (\tau_1 n_{h,\downarrow} + \frac{-1}{2} \hat{c}_{h,\uparrow} \hat{c}_{e,\downarrow} + \frac{1}{\sqrt{2}} n_{e,\downarrow} - \tau_1 n_{h,\downarrow} n_{e,\downarrow} - \tau_1 n_{h,\uparrow} n_{e,\downarrow} - \tau_1 n_{h,\uparrow} n_{h,\downarrow} - \tau_2 n_{h,\uparrow}) + \\ (\frac{1}{\sqrt{2}} \hat{c}_{h,\downarrow} - \frac{1}{\sqrt{2}} \hat{c}_{e,\uparrow}^\dagger)$$

$$c_{f,\uparrow}^\dagger = \hat{c}_{e,\uparrow}^\dagger (\frac{1}{2} n_{h,\downarrow} + \frac{-1}{2} n_{h,\uparrow} n_{h,\downarrow} + n_{e,\downarrow} + \frac{-1}{2} n_{h,\uparrow} n_{e,\downarrow} + \frac{-1}{2} n_{h,\downarrow} n_{e,\downarrow} + \frac{-1}{2} \hat{c}_{h,\uparrow} \hat{c}_{e,\downarrow} + \frac{1}{\sqrt{2}} \hat{c}_{h,\uparrow}^\dagger \hat{c}_{e,\downarrow}^\dagger) + . \\ \hat{c}_{h,\downarrow} (\frac{-1}{2} n_{e,\uparrow} + \frac{1}{2} n_{h,\uparrow} n_{e,\uparrow} + \frac{1}{2} n_{h,\uparrow} n_{e,\downarrow} + \frac{1}{2} n_{e,\uparrow} n_{e,\downarrow} + \frac{1}{2} \hat{c}_{h,\uparrow}^\dagger \hat{c}_{e,\downarrow}^\dagger - \frac{1}{\sqrt{2}} \hat{c}_{h,\uparrow} \hat{c}_{e,\downarrow} - n_{h,\uparrow})$$

$$\tau_2 = (1 - 1/\sqrt{2}) \quad \tau_1 = \frac{1}{2}(\sqrt{2} - 1)$$

Ground state energies accurate to $(t/J)^4$ *

$$E_k = \frac{3J}{4}\Delta_k \pm te_k \pm \mu \text{ where } \Delta_k = \sqrt{1 + \left(\frac{4te_k}{3J}\right)^2}.$$

$$\alpha = \frac{1}{2} \iiint_{T^d} (1 - \Delta_k^{-1}) \left(\frac{d\theta}{2\pi}\right)^d$$

$$\beta = \iiint_{T^d} e_k^2 / \Delta_k \left(\frac{d\theta}{2\pi}\right)^d \approx \frac{d}{2} - \frac{d}{3}\lambda^2 + O(t^4)$$

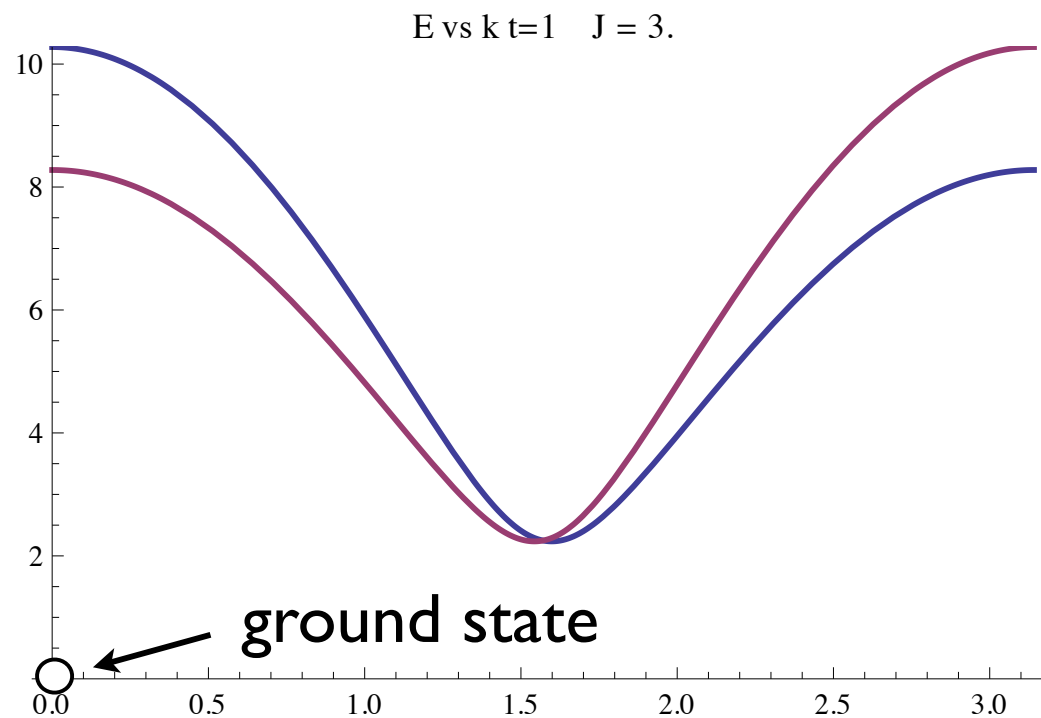
$$\begin{aligned} \langle H_{KLM} \rangle_{free} = & \frac{-3}{4} + 3\alpha + \frac{-15}{4}\alpha^2 + \frac{3}{2}\alpha^3 + \frac{-8}{3}\beta\lambda^2 + \\ & \alpha^3\beta\lambda^2(72 + -48\sqrt{2}) + \alpha\beta^3d^{-2}\lambda^4\left(\frac{64}{3} + \frac{-128}{9}\sqrt{2}\right) + \\ & \alpha\beta\lambda^2(16 + -8\sqrt{2}) + \beta^5d^{-4}\lambda^6\left(\frac{-64}{27} + \frac{128}{81}\sqrt{2}\right) + \\ & \beta^3d^{-2}\lambda^4\left(\frac{-64}{9} + \frac{32}{9}\sqrt{2}\right) + \alpha^2\beta^3d^{-2}\lambda^4\left(\frac{-64}{3} + \frac{128}{9}\sqrt{2}\right) + \\ & \alpha^4\beta\lambda^2(-36 + 24\sqrt{2}) + \alpha^2\beta\lambda^2(-52 + 32\sqrt{2}) \end{aligned}$$

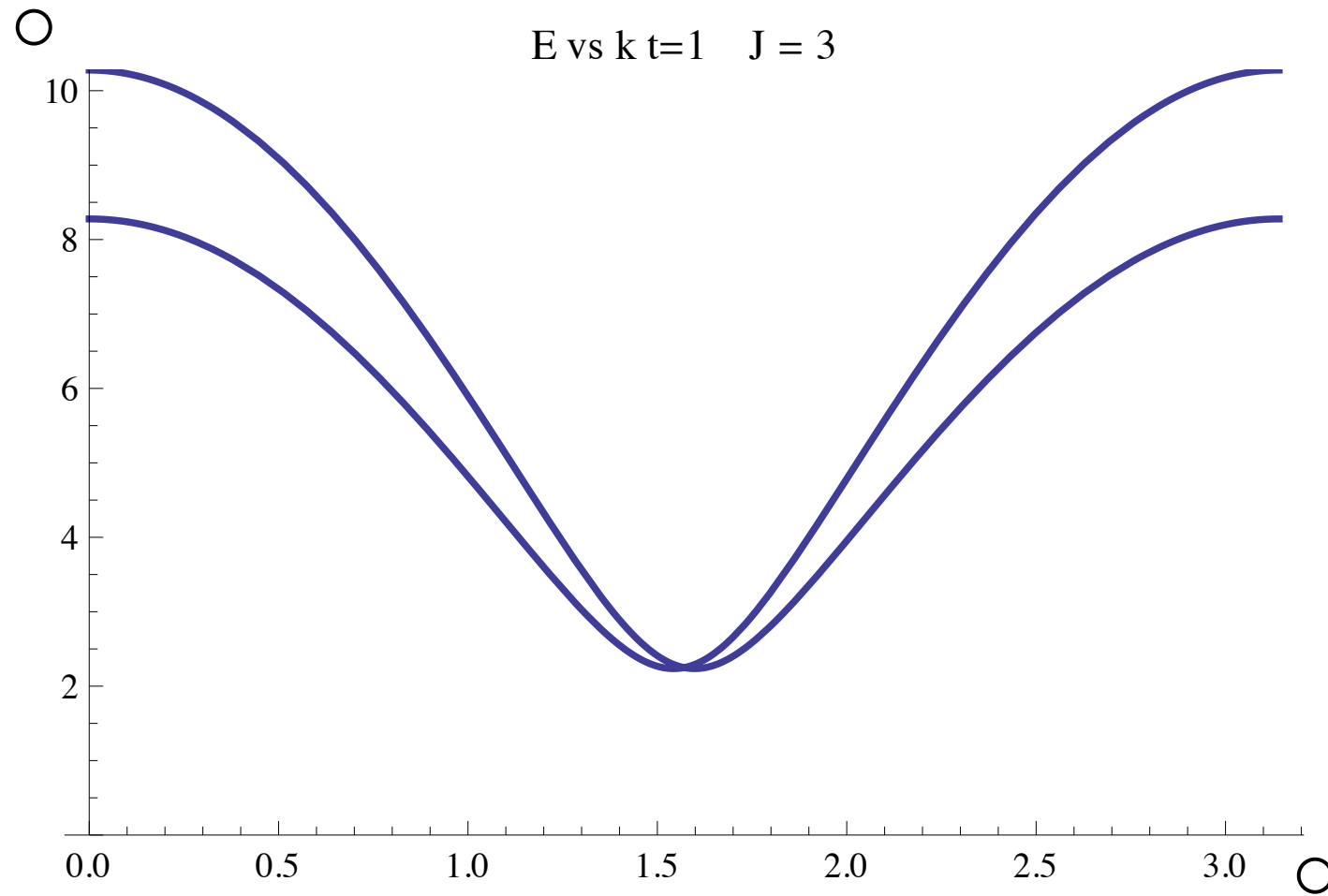
* Zheng and Oitmaa, 2003

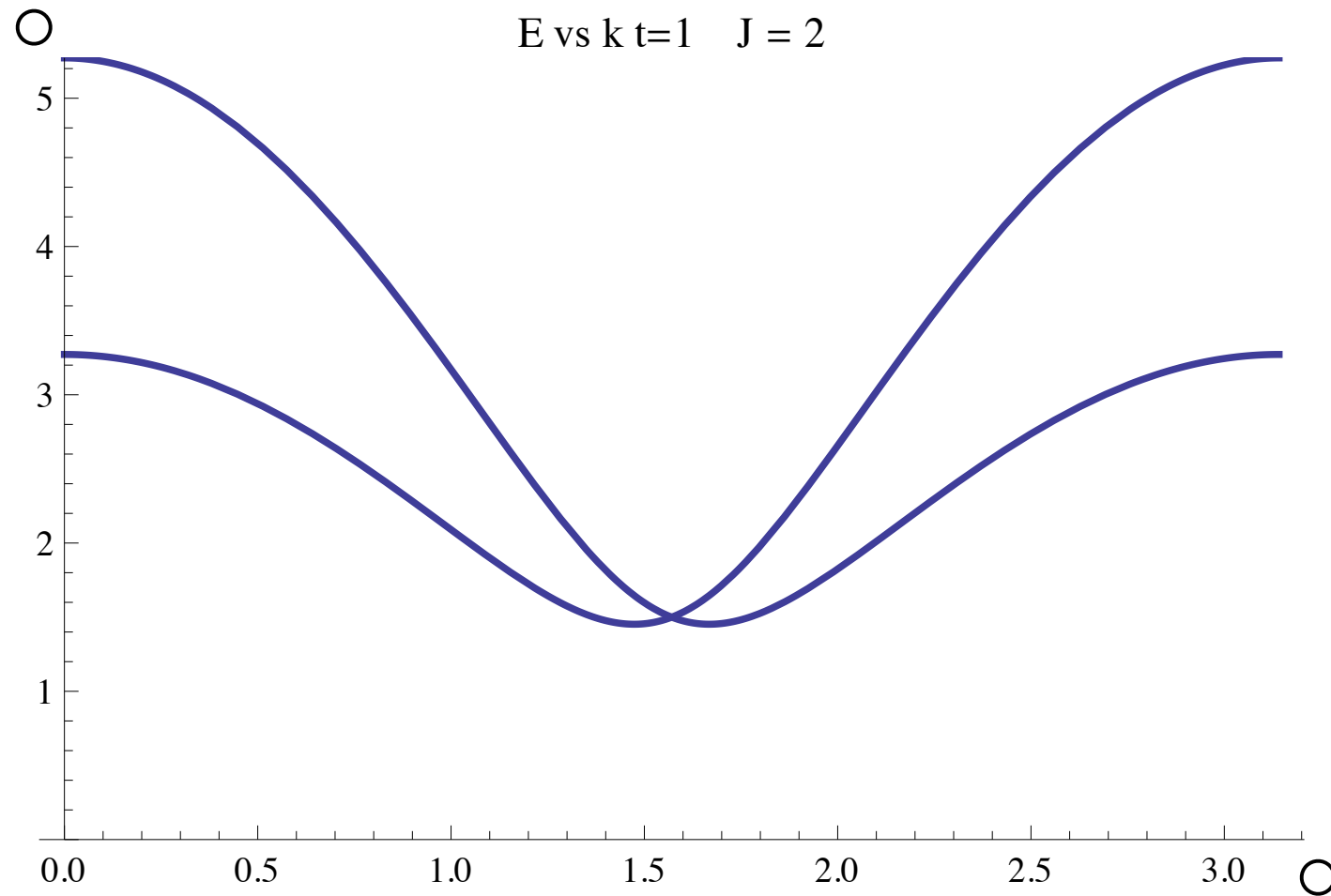
Spectrum and spectral weights can be computed with the exact canonical transformation but with the approximate Hamiltonian.

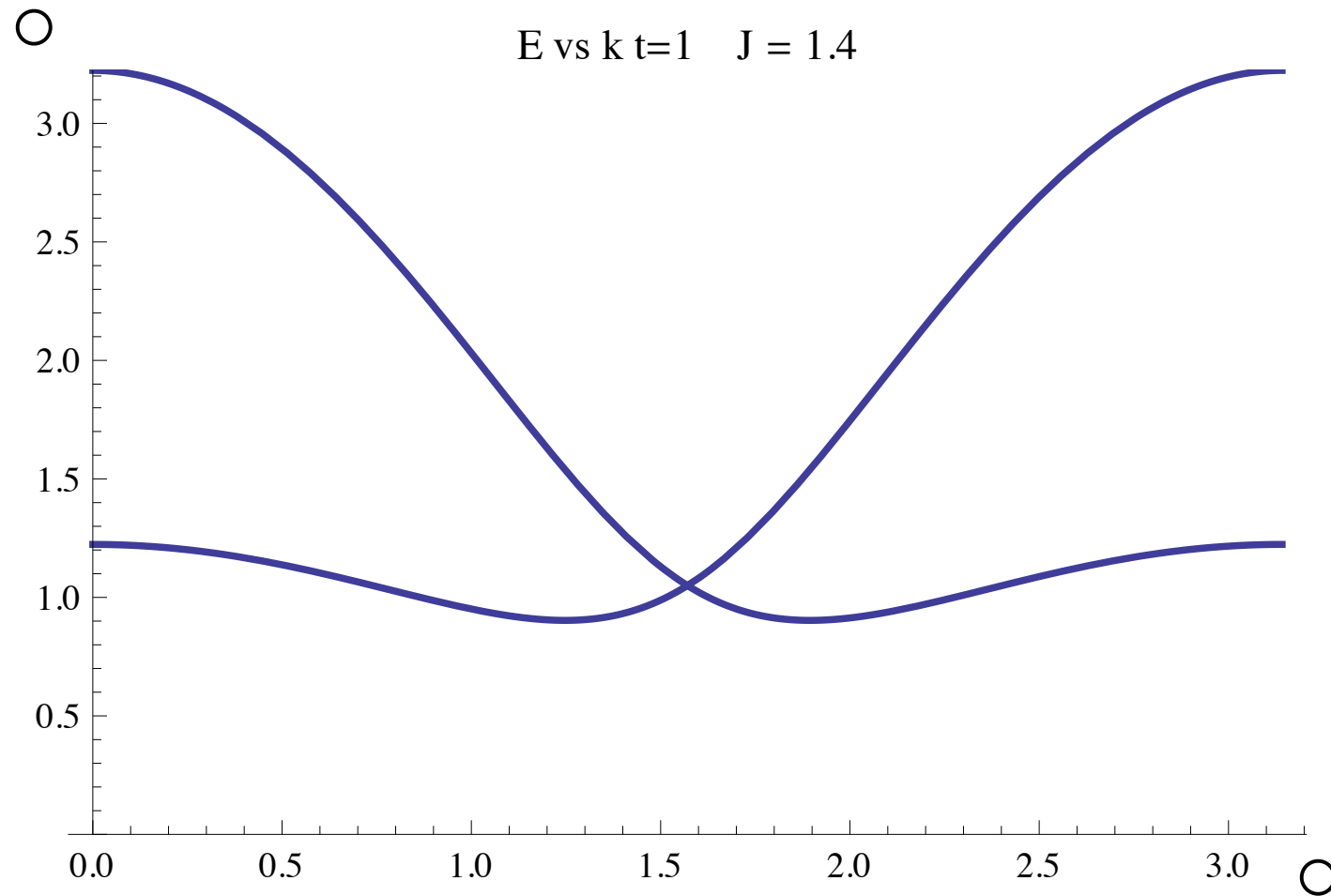
$$\Delta_k = \sqrt{1 + \left(\frac{4te_k}{3J}\right)^2}.$$

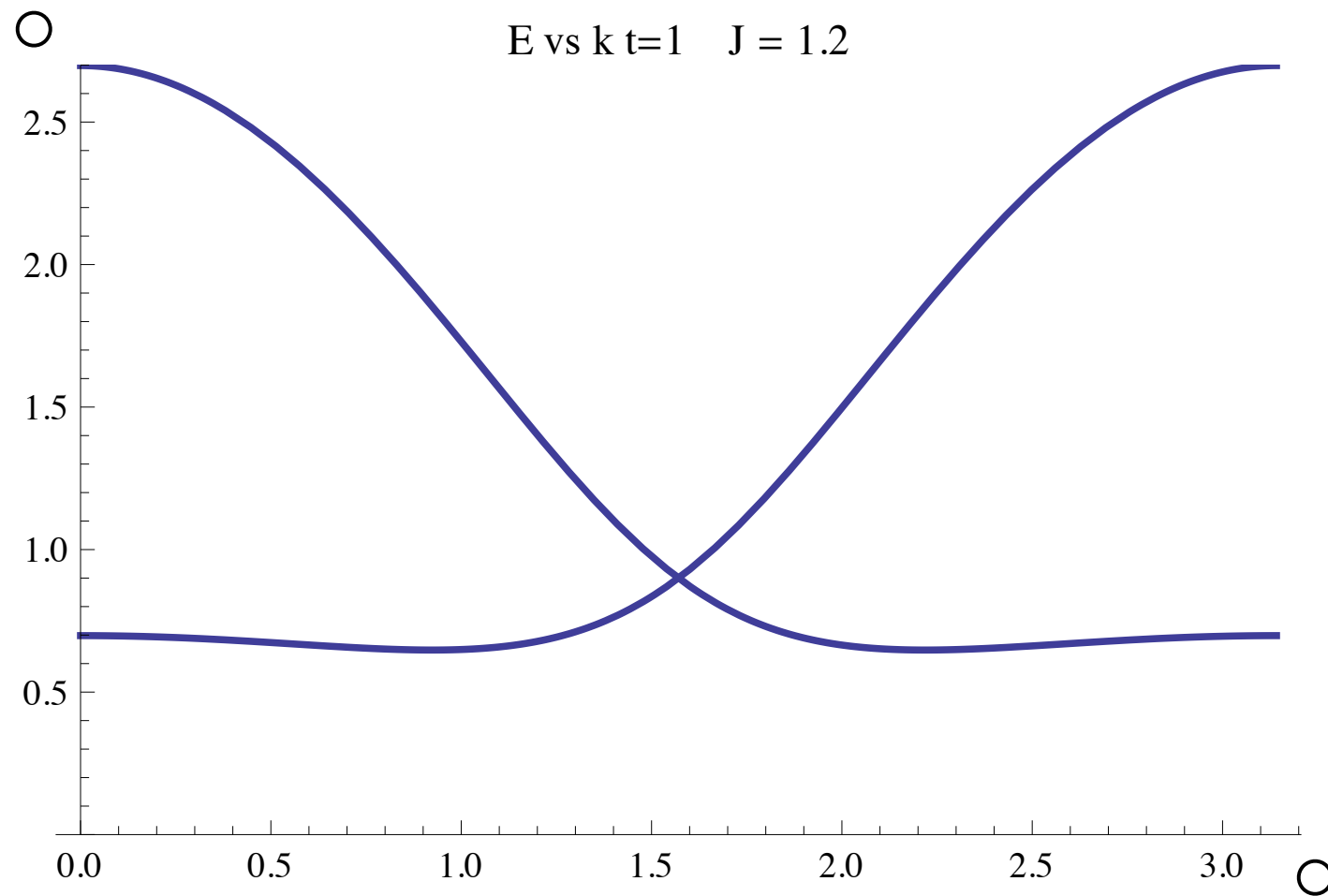
$$E_k = \frac{3J}{4}\Delta_k \pm te_k \pm \mu$$

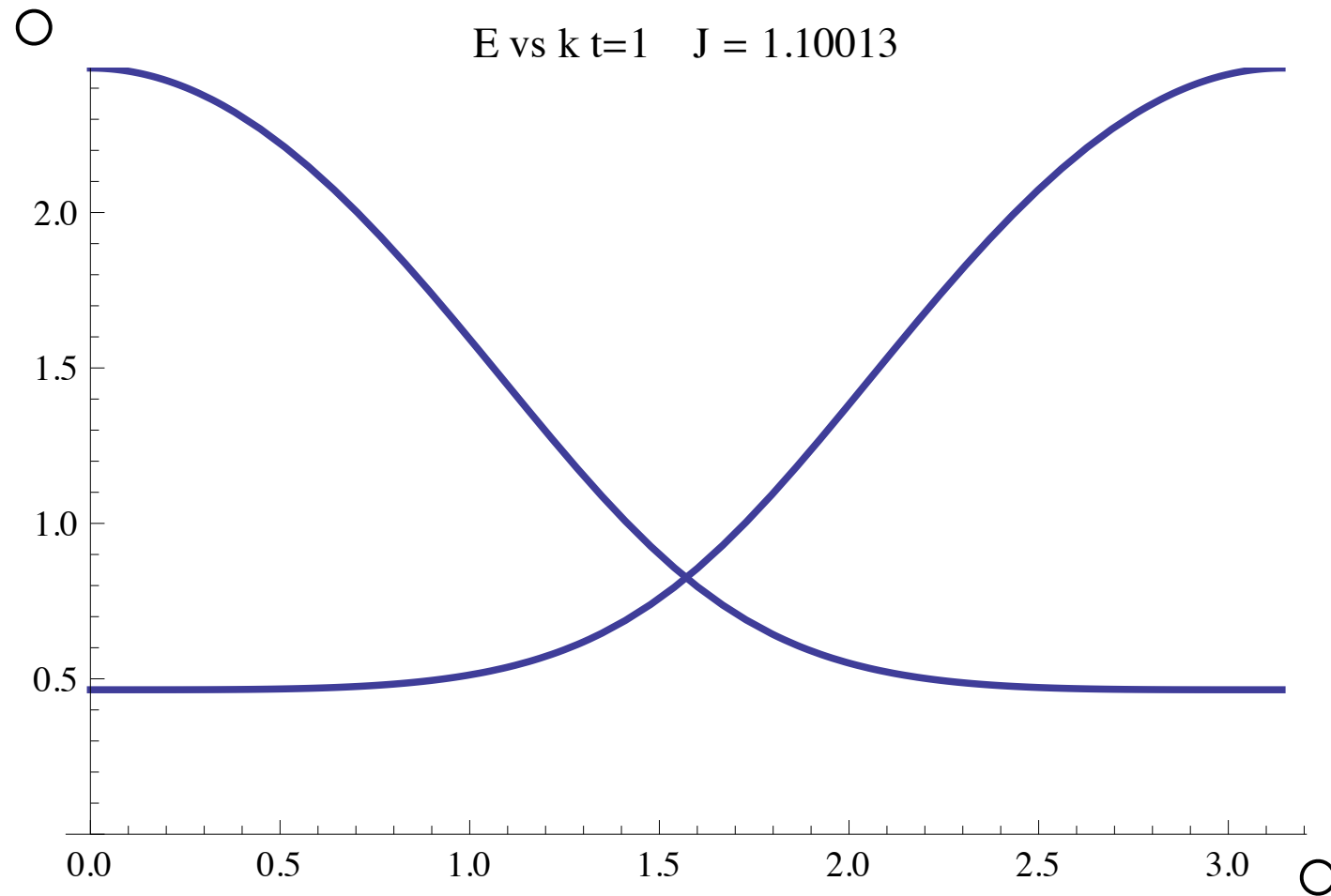


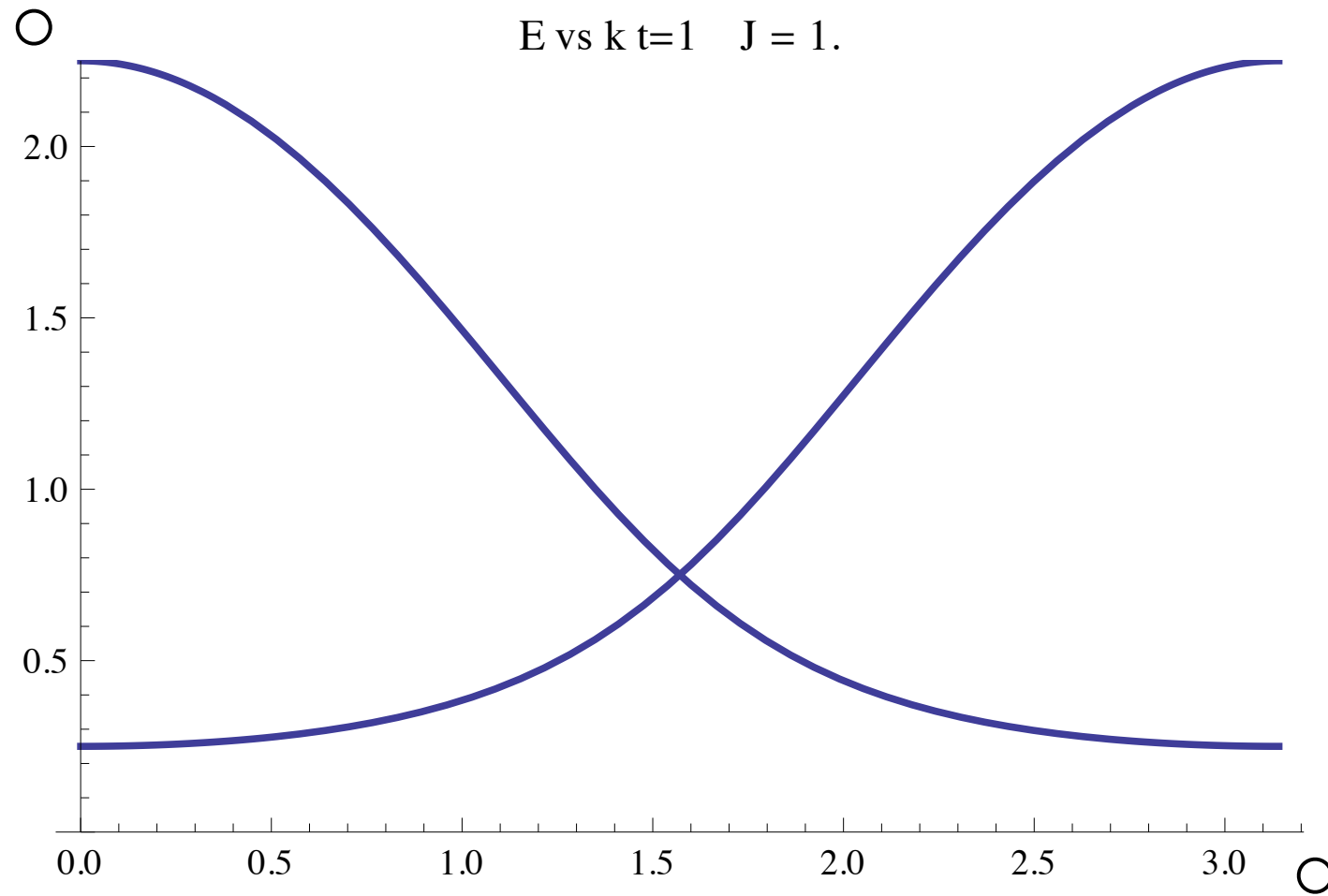


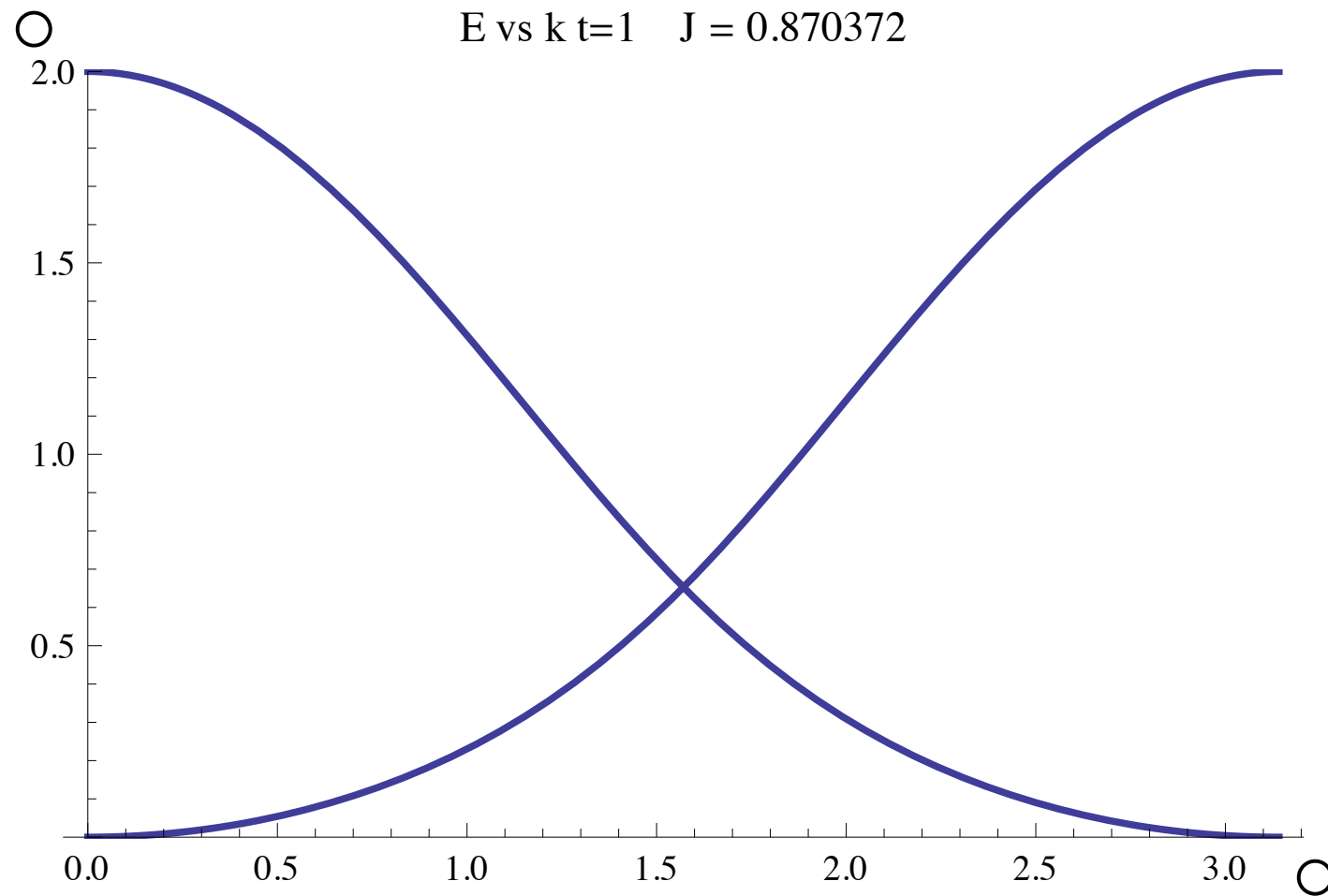


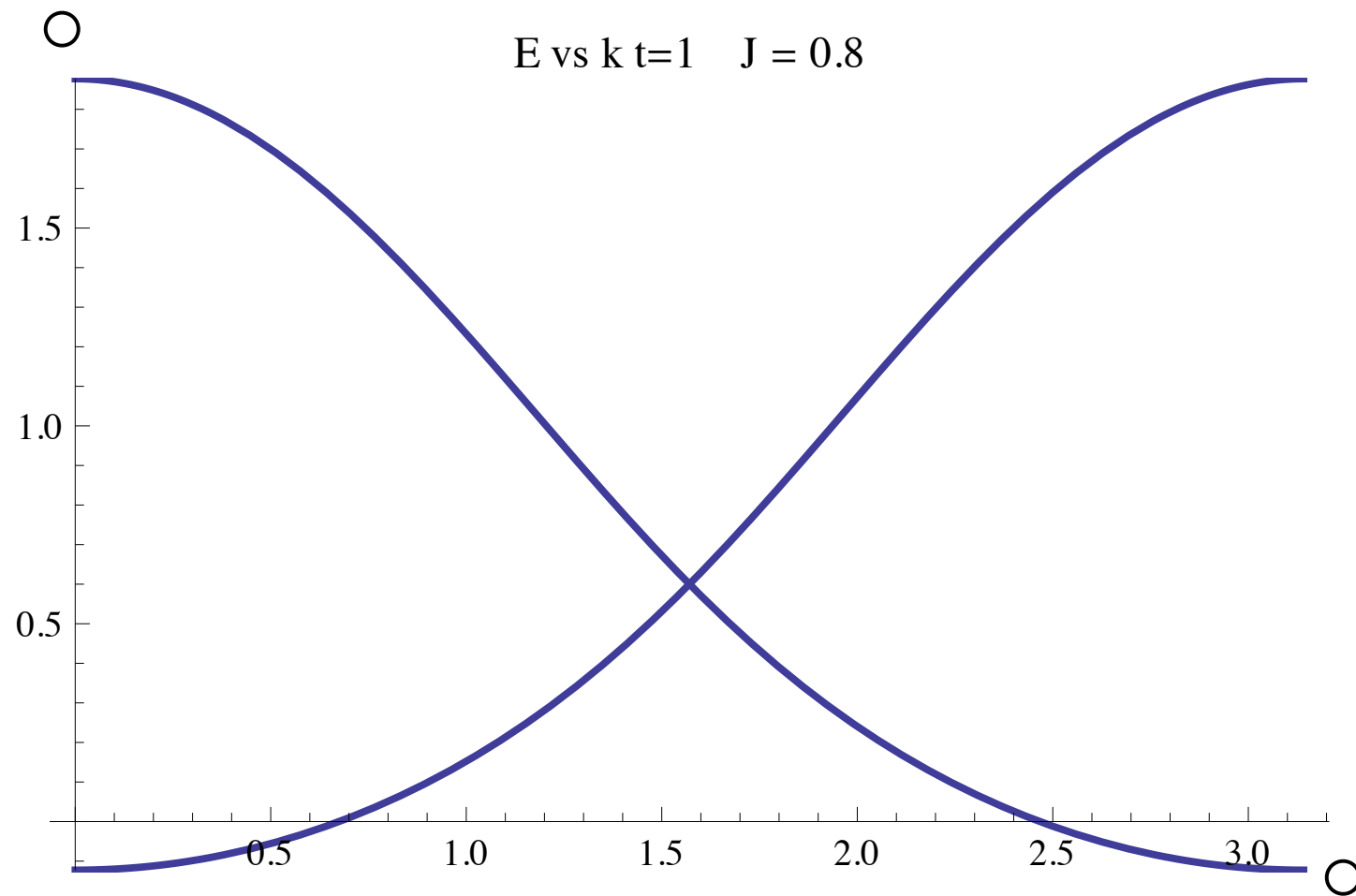


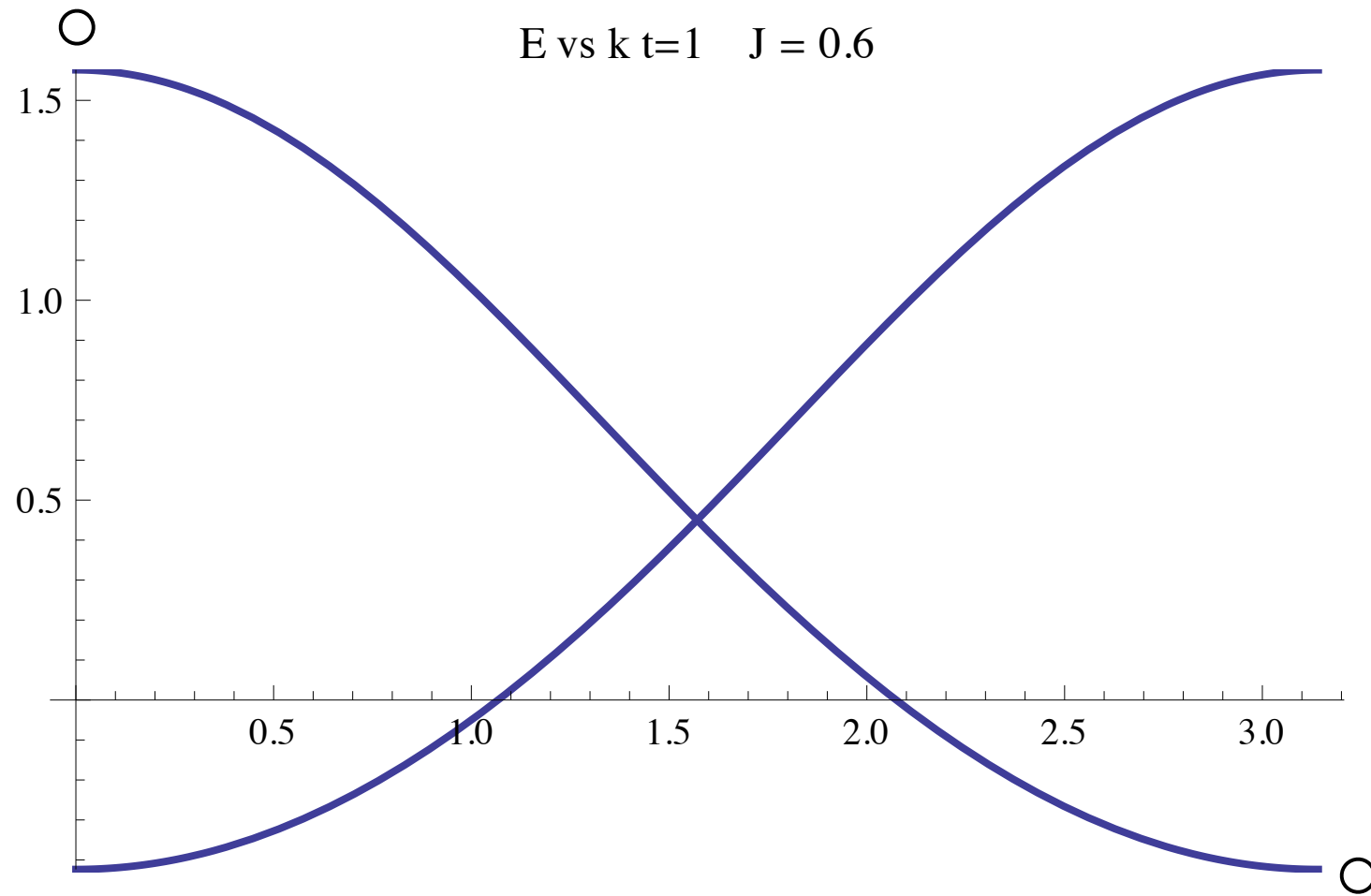


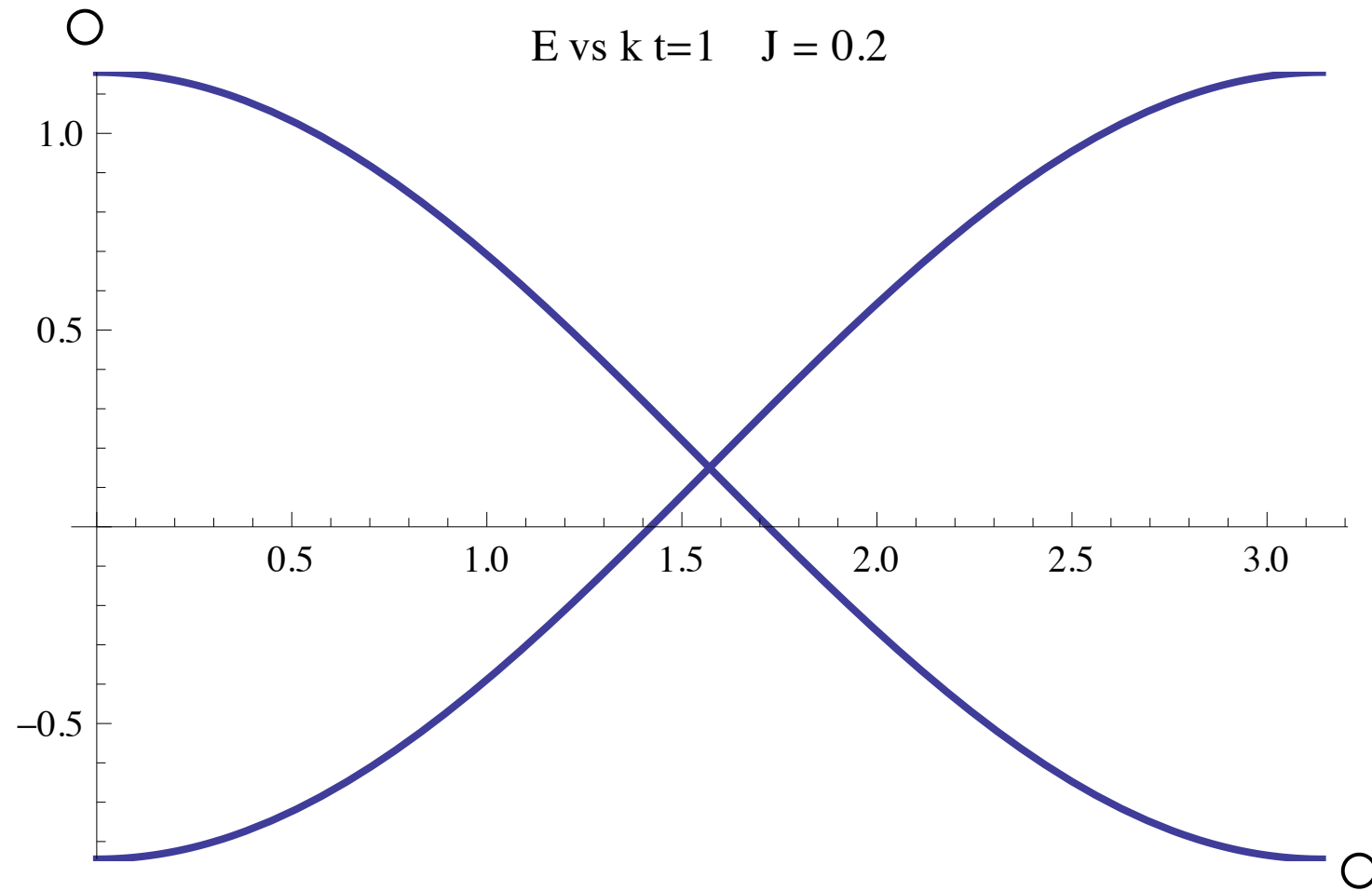


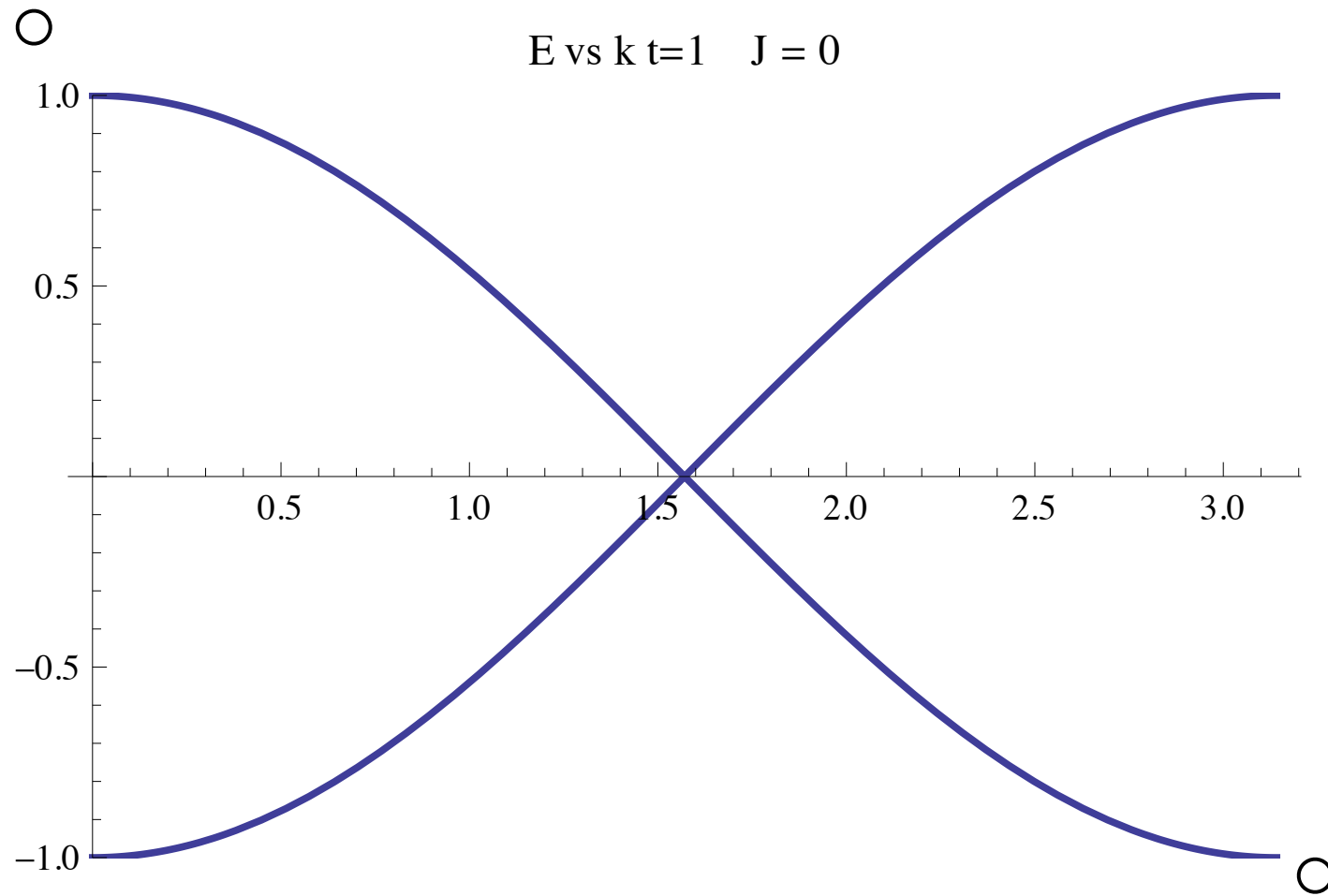




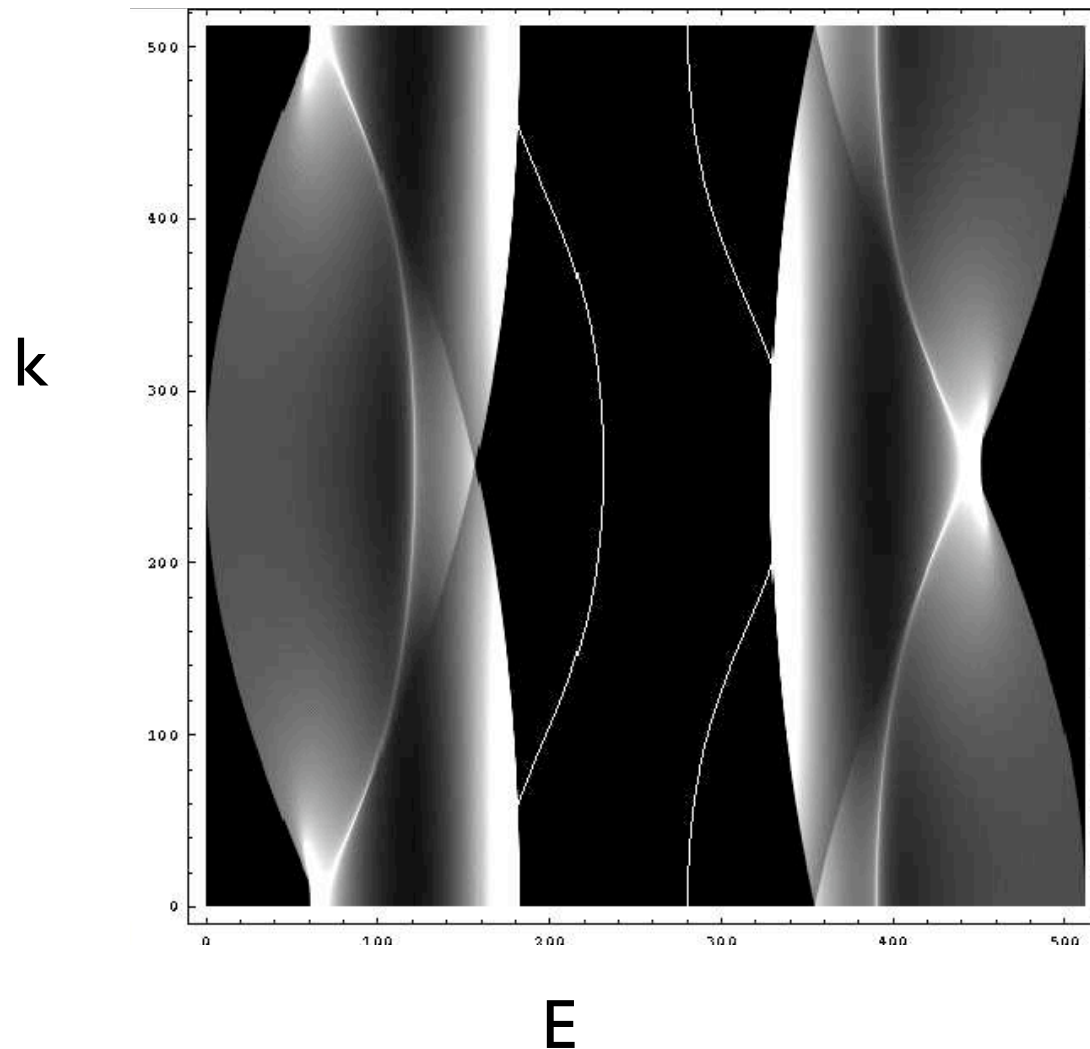








c-electron spectral weight



Note delta function component together with a continuous scattering part.

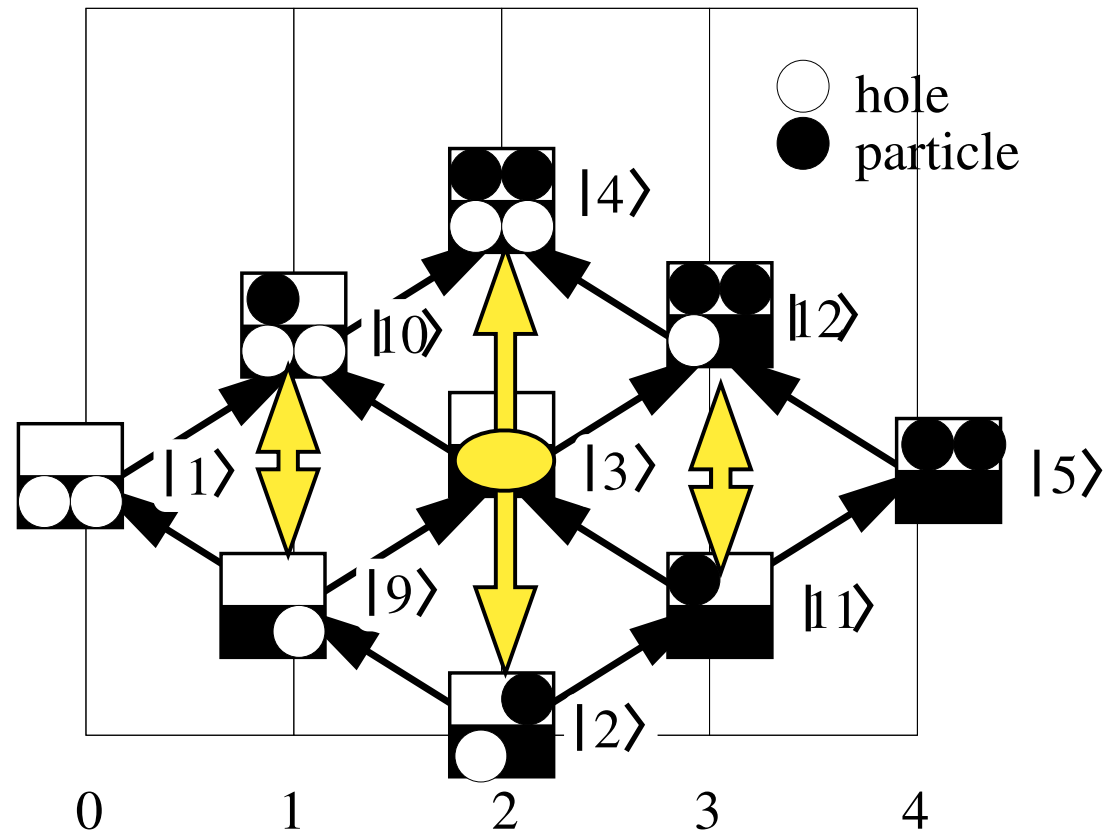
Generalizations:

Transformations can be generalized from fixed permutation to

$$SU(2) \otimes SU(3) \otimes SU(2)$$

that can continuously rotate from the identity transformation (weak coupling) to strong coupling permutation

Transformations can be generalized to $SU(2) \otimes SU(3) \otimes SU(2)$



$$SU(2) \otimes SU(3) \otimes SU(2)$$

An approximate free quasiparticle model can be constructed which can describe a type of metal-insulator transition for which spectrum and bare electron spectral weights can be computed exactly

Can we apply these ideas to Hubbard model?

Previous method does not work for odd valence insulator.

$$H = -t \sum_{\langle rr' \rangle} (c_{s,r}^\dagger c_{s,r'} + CC) + U \sum_r \frac{1}{2} (n_r - 1)^2$$

The ground state consists of one electron per site. Consider the Neel ground state. Can we make a unitary transformation that maps one of the two **one-particle** states to the “Fermi” vacuum consisting of zero particles.

Guided by previous studies of ‘nonlinear’ canonical transformation, it turns out that the following ***noncanonical*** transformation accomplishes this without restricting the Hilbert space:

$$\hat{c}_r = c_{\downarrow, r}^\dagger (1 - n_{\uparrow, r}) + (-1)^r c_{\downarrow, r} n_{\uparrow, r}$$

$$q_r^+ = (c_{\downarrow, r}^\dagger - (-1)^r c_{\downarrow, r}) c_{\uparrow, r}$$

$$q_r^- = (q_r^+)^{\dagger}$$

$$q_r^z = \frac{1}{2} - n_{\uparrow, r}$$

This exactly converts the algebra generated by two Fermionic spin degrees of freedom exactly to a ‘quasicharge’ Fermionic ‘ CP ’ and a ‘quasispin’ Bosonic $SU(2)$

$$\{\hat{c}_r, \hat{c}_{r'}^\dagger\} = \delta_{r,r'}$$

$$\{\hat{c}_r^\dagger, \hat{c}_{r'}^\dagger\} = 0$$

$$[\hat{c}_r^\dagger, q_{r'}^i] = 0 ;$$

$$[q_r^i, q_{r'}^j] = i\delta_{rr'} \sum_k \epsilon_{ijk} q_r^k$$

The hopping splits into three pieces

$$H = t (T_0 + T_1 + T_{-1}) + U h_U$$

$$T_0 = \frac{1}{2} \sum_{\langle r, r' \rangle} (1 + 4q_r q_{r'}) (\hat{c}_r^\dagger \hat{c}_{r'} + C C)$$

$$T_1 = \frac{1}{2} \sum_{\langle r, r' \rangle} -1^r (1 - 4q_r q_{r'}) (\hat{c}_r^\dagger \hat{c}_{r'}^\dagger)$$

$$T_{-1} = \frac{1}{2} \sum_{\langle r, r' \rangle} -1^r (1 - 4q_r q_{r'}) (\hat{c}_{r'} \hat{c}_r)$$

$$h_U = \frac{1}{2} \sum_r \hat{c}_r^\dagger \hat{c}_r$$

The Hubbard interaction has become a chemical potential for the quasicharge operators.

Conclusions

- These arguments suggest that the free electrons must be nonperturbatively rewritten as composite operators in a strongly interacting system.
- Nonlinear canonical transformations are natural tools to study doped even values Mott insulators at least in the strong coupling limit.
- A natural generalization of this leads to exact nonunitary transformations that give charge-like fermionic and bosonic spin-like degrees of freedom.