Transport-induced correlations in weakly interacting systems

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Received 19 June 2013 Accepted 29 July 2013 Published 30 August 2013

Online at stacks.iop.org/JSTAT/2013/P08015 doi:10.1088/1742-5468/2013/08/P08015

Abstract. We study spatial correlations in the transport of energy between two baths at different temperatures. To do this, we introduce a minimal model in which energy flows from one bath to another through two subsystems. We show that the transport-induced energy correlations between the two subsystems are of the same order as the energy fluctuations within each subsystem. The correlations can be either positive or negative and we give bounds on their values which are associated with a dynamic energy scale. The different signs originate as a competition between fluctuations generated near the baths and fluctuations of the current between the two subsystems. This interpretation sheds light on known results for spatially-dependent heat and particle conduction models.

Keywords: driven diffusive systems (theory), transport processes/heat transfer (theory), heat conduction

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1. Introduction

The physics of systems in and out of equilibrium can differ in dramatic ways. For example, in equilibrium, one-dimensional systems with short range interactions cannot show long-range correlations at positive temperatures. By contrast, in systems away from equilibrium long-range correlations are known to form when there is a steady-state current of a conserved quantity [1]–[4]. For diffusive systems this has been calculated for models of particle and heat transport [5]–[7], and measured in heat-transport experiments [4, 8].

In this paper, we study a minimal model for the formation of correlations during the transport of a conserved quantity. We focus primarily on energy, but also show an example in which the same model is used to describe the transport of particles. The model consists of two systems and two baths, arranged in a chain as shown in figure 1. The total energy of the system is $E_{\text{tot}} = E_1 + E_2 + E_{\text{int}}$, where the $E_{1,2}$ denote the energies of the subsystems and E_{int} is the interaction energy between the two subsystems, and also between the subsystems and the baths. We assume that the different components interact weakly, as is assured for example when the interactions are short range and the links between the components do not scale with system size. This means that the interactions allow energy to flow, but can be neglected in any energetic calculation. In equilibrium, i.e. when the temperatures of the two baths are equal, $T_0 = T_3$, the probability of subsystem 1 to be in state s_1 and subsystem 2 to be in state s_2 is given to leading order in system size by

$$P(s_1, s_2) = \frac{1}{Z} e^{-\beta E_{\text{tot}}} \approx \frac{1}{Z} e^{-\beta E_1(s_1)} e^{-\beta E_2(s_2)}$$

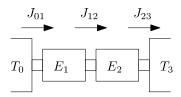


Figure 1. Heat conduction through a pair of systems. Subsystems 1 and 2 have energies E_1 and E_2 respectively. The baths have temperatures T_0 , T_3 . The currents of energy between the different parts are J_{01} , J_{12} , J_{23} .

where $\beta = 1/T_0$ and Z is the partition function. It follows immediately that the energy correlations vanish,

$$\langle E_1 E_2 \rangle_{\beta} - \langle E_1 \rangle_{\beta} \langle E_2 \rangle_{\beta} = 0, \tag{1}$$

where $\langle \cdots \rangle_{\beta}$ denotes a thermal average.

In this paper we use this minimal model to explain the mechanism by which longrange correlations develop in systems away from equilibrium, i.e. when $T_0 \neq T_3$, despite the negligible interaction energies. These correlations are related to an energy scale associated with the current traversing the system. We apply our formalism to several examples, both classical and quantum, and show that the correlations can be both positive and negative. We give a simple picture to explain the appearance and sign of the correlation, and which also sheds light on continuous diffusive systems.

The correlations discussed in this paper can in principle be measured in various systems. For example, correlations of the electric charge can be measured in double quantum dot experiments [9], and in small analog electronic circuits. The latter can also be used to measure heat fluctuations [10]. Such setups have been used in experimental studies of fluctuation relations [10, 11].

2. The model

The currents between the different components of the system are denoted by J_{01} , J_{12} , J_{23} , see figure 1. Conservation of energy implies that

$$\frac{dE_1}{dt} = J_{01} - J_{12},
\frac{dE_2}{dt} = J_{12} - J_{23}.$$
(2)

We consider simple dynamics, where the current fluctuations are modeled by white noise. Separating fluctuations from average values, we write the currents as

$$J_{01} = A_{01}(E_1) + \sqrt{B_{01}(E_1)}\eta_{01}$$

$$J_{12} = A_{12}(E_1, E_2) + \sqrt{B_{12}(E_1, E_2)}\eta_{12}$$

$$J_{23} = A_{23}(E_2) + \sqrt{B_{23}(E_2)}\eta_{23}$$
(3)

where $A_{ij}(E_1, E_2) = \overline{J_{ij}}(E_1, E_2)$, $\overline{\eta_{ij}} = 0$, and $\overline{\eta_{ij}(t)\eta_{ij}(t')} = \delta(t - t')$, for indices ij = 01, 12, 23. Overbars denote averages at fixed E_1, E_2 , and B_{ij} are noise amplitudes.

The above model provides a good description of the dynamics when the energy flows are slow enough to allow subsystems 1 and 2 to constantly relax to their microcanonical equilibria at energies E_1 and E_2 . This assures that the state of each subsystem is well-described by its energy. The associated (microcanonical) inverse temperatures are $\beta_i(E_i) \equiv \partial S_i(E_i)/\partial E_i$ for i=1,2, where $S_i(E_i)$ is the entropy of the *i*th subsystem. This time scale separation also means that over the relaxation time of the entire system to its non-equilibrium steady-state, the current performs many independent fluctuations, and can therefore be modeled by white noise about the average.

The physical information about the systems and the energy flow is contained in the six functions A_{ij} and B_{ij} , which as we show below can be obtained for specific models. Under some conditions the functions A_{ij} and B_{ij} are related. Specifically, let w be the energy flow within the memory time of the noise (of the order of the subsystems' relaxation time, or shorter). If $(\Delta \beta)^2 \overline{w^3}_c \ll \overline{w}$, where $\overline{w^3}_c$ denotes the third cumulant and $\Delta \beta$ is the largest of $|\beta_1 - \beta_0|$, $|\beta_2 - \beta_1|$ and $|\beta_3 - \beta_2|$, then the following generalized fluctuation-dissipation relation holds:

$$2A_{ij} = (\beta_j - \beta_i) B_{ij}. \tag{4}$$

It follows by taking a cumulant expansion of the exchange fluctuation relation [12], and neglecting cumulants of order 3 and higher, according to the condition above, as discussed in detail in [13] (for a different approach, see [14]). It holds in general when the temperature differences are small (namely in the standard linear response regime) but can hold more generally in some cases even when the temperature differences are large (see [13] for examples). Therefore, equation (4) is a generalization of the Green–Kubo formula. In what follows we use this relation when applicable.

Finally, note that equations (2) and (3) describe [15], for a specific form of functions A_{ij} and B_{ij} , the dynamics of local energy transfers in a stochastic form [6, 16] of the discrete Kipnis, Marchioro and Presutti model [17]. We relate them to other models below.

3. Steady-state fluctuations

With the model defined, we turn to study the energy fluctuations at its steady-state. At the steady-state, $\langle dE_1/dt \rangle = \langle dE_2/dt \rangle = 0$, where angular brackets denote averages over the steady-state probability distribution. From equations (2) and (3), this implies $\langle A_{01} \rangle = \langle A_{12} \rangle = \langle A_{23} \rangle$.

As in equilibrium, fluctuations are expected to scale as $N^{1/2}$, where N is the system size. This will be shown self-consistently below. This motivates an expansion in small energy fluctuations $\delta E_1 \equiv E_1 - \langle E_1 \rangle$, $\delta E_2 \equiv E_2 - \langle E_2 \rangle$.

Using this we first find $\langle E_i \rangle$. To lowest order in δE_i , we have $\langle A_{ij}(E_1, E_2) \rangle = A_{ij}(\langle E_1 \rangle, \langle E_2 \rangle)$. Therefore, at $\delta E_1 = \delta E_2 = 0$,

$$A_{01}(\langle E_1 \rangle) - A_{12}(\langle E_1 \rangle, \langle E_2 \rangle) = 0,$$

$$A_{12}(\langle E_1 \rangle, \langle E_2 \rangle) - A_{23}(\langle E_2 \rangle) = 0,$$
(5)

which can be solved to obtain $\langle E_1 \rangle$ and $\langle E_2 \rangle$.

We next expand equation (2) to leading order in the fluctuations to obtain

$$\frac{\mathrm{d}}{\mathrm{d}t} \begin{pmatrix} \delta E_1 \\ \delta E_2 \end{pmatrix} = \mathbf{R} \begin{pmatrix} \delta E_1 \\ \delta E_2 \end{pmatrix} + \mathbf{\Sigma} \begin{pmatrix} \eta_{01} \\ \eta_{12} \\ \eta_{23} \end{pmatrix}. \tag{6}$$

Here, **R** is a 2 × 2 matrix with elements $R_{ij} = \partial_{E_j} X_i|_{E_{1,2} = \langle E_{1,2} \rangle}$, where $X_1 = A_{01} - A_{12}$ and $X_2 = A_{12} - A_{23}$. The matrix **R** controls the relaxation of fluctuations, as $d(\overline{\delta E_i})/dt = \sum_j R_{ij} \overline{\delta E_j}$. **\(\Sigma\)** is a 3 × 2 matrix which specifies the strength of the fluctuations. Since the numbers B_{ij} do not generally vanish at $E_1 = \langle E_1 \rangle$, $E_2 = \langle E_2 \rangle$, we have to lowest order in the fluctuations

$$\Sigma \equiv \begin{pmatrix} \sqrt{B_{01}^s} & -\sqrt{B_{12}^s} & 0\\ 0 & \sqrt{B_{12}^s} & -\sqrt{B_{23}^s} \end{pmatrix},$$

with $B_{ij}^s = B_{ij} (\langle E_1 \rangle, \langle E_2 \rangle).$

Finally, the energy correlation matrix $C_{ij} = \langle E_i E_j \rangle - \langle E_i \rangle \langle E_j \rangle$ can be obtained by solving the Lyapunov equation [19]

$$\mathbf{RC} + \mathbf{CR}^T = -\mathbf{Q},\tag{7}$$

where $\mathbf{Q} \equiv \Sigma \Sigma^T$. In general, away from equilibrium the off-diagonal C_{12} does not vanish. In what follows, we will demonstrate this explicitly in two simple models.

It follows from equation (4) that A_{ij} and B_{ij} have the same scaling with system size. **R** is composed of derivatives of the type $\partial A_{ij}/\partial E_k$, and therefore scales as N^{-1} , where N is the system size. From equation (7) it follows that $\mathbf{C} \sim N$, as in equilibrium. The fluctuations in the energy then scale as $N^{1/2}$ and justify self-consistently the expansion in small $\delta E_1, \delta E_2$.

4. Linear A model

Consider a simple model, where subsystems 1 and 2 are identical, with microcanonical temperatures T(E). $E\left(T\right)$ will denote the inverse of this function. In addition, for the dynamics we take

$$A_{01} = \gamma \left[\varepsilon \left(T_0 \right) - \varepsilon_1 \right],$$

$$A_{12} = \gamma \left[\varepsilon_1 - \varepsilon_2 \right],$$

$$A_{23} = \gamma \left[\varepsilon_2 - \varepsilon \left(T_3 \right) \right].$$
(8)

Here ε are the energy densities in the baths ($\varepsilon(T_0)$ and $\varepsilon(T_3)$) and in the subsystems ($\varepsilon_{1,2} = E_{1,2}/N$), and γ is a rate constant, setting a time scale in the model. From the definition of **R**

$$\mathbf{R} = \frac{\gamma}{N} \begin{pmatrix} -2 & 1\\ 1 & -2 \end{pmatrix}.$$

The solution to equation (7) then reads

$$\mathbf{C} = \frac{N}{24\gamma} \begin{pmatrix} 7B_{01} + 4B_{12} + B_{23} & 2B_{01} - 4B_{12} + 2B_{23} \\ 2B_{01} - 4B_{12} + 2B_{23} & B_{01} + 4B_{12} + 7B_{23} \end{pmatrix}. \tag{9}$$

Note that at equilibrium $B_{01} = B_{12} = B_{23}$, so that **C** is diagonal, as expected from equation (1).

Interestingly, we see that the cross-correlation C_{12} can be either positive or negative, depending on the sign of $B_{01} - 2B_{12} + B_{23}$. We suggest the following simple interpretation. Current fluctuations transfer energy along the different bonds. The fluctuations relax by the average dynamics $d(\overline{\delta E_i})/dt = \sum_j R_{ij} \overline{\delta E_j}$. Consider a current fluctuation near a bath, say in J_{01} , which changes the energies by $(\delta E_1, \delta E_2) \propto (1,0)$. This fluctuation relaxes according to the sum of the modes of \mathbf{R} , $(\delta E_1, \delta E_2) \propto \mathrm{e}^{-\gamma t/N} (1,1) - \mathrm{e}^{-3\gamma t/N} (-1,1)$, for which δE_1 and δE_2 have the same sign. On the other hand, a fluctuation in the central bond J_{12} changes energies by $(\delta E_1, \delta E_2) \propto (-1,1)$, which decays as $(\delta E_1, \delta E_2) \propto -\mathrm{e}^{-3\gamma t/N} (-1,1)$. Therefore, current fluctuations near the baths promote positive correlations between the two subsystems, while current fluctuations in the center of the system contribute to negative correlation between the subsystems. The expression $C_{12} \propto B_{01} - 2B_{12} + B_{23}$ reflects the positive effect of the boundary noise B_{01} , B_{23} and the negative effect of the bulk noise B_{12} .

To demonstrate the different possible behaviors in different models consider first a specific example of the linear A model in which the subsystems are ideal gases with $E = C_v T$, with C_v a constant specific heat, and that equation (4) holds. This model of ideal gases satisfying Fourier's law is perhaps the simplest phenomenological model for energy transfer. To obtain B_{01}^s , B_{12}^s , B_{23}^s we solve equation (5) and use equation (4). Substituting into equation (9) we find

$$C_{12} = \frac{C_v}{27} (T_0 - T_3)^2.$$

Thus C_{12} is positive for any $T_0 \neq T_3$, and proportional to the system size.

This model can be considered as a 'boxed' version of standard models for heat or particle conduction with spatial dependence, known to have positive correlations in the continuum limit [5]. Indeed, a similar line of argumentation applies to continuum diffusive systems, which conduct particles or heat. If a fluctuation in the energy density E(x,t) decays according to a simple diffusion $\partial_t E = D\nabla^2 E$, with D a constant diffusion coefficient, a current fluctuation near the baths will contribute to positive correlations, while current fluctuations at the bulk will contribute to negative correlations. Therefore models with stronger noise near the boundary will have positive correlations, while models with stronger bulk noise will have negative correlations, as illustrated in figure 2. This gives a simple picture of the positive correlations found in standard heat conduction models [5, 6] as opposed to the negative correlation found, for example, in the simple symmetric exclusion process (SSEP) [1, 18], a standard particle conduction model.

Indeed, a boxed version of the SSEP can be constructed as follows. Consider a boxed model in which ε_i describes the density of particles in box i. A particle then hops from subsystem i to j with rate $\gamma \varepsilon_i (1 - \varepsilon_j)$ where 1 is the maximal density of particles that each subsystem can contain. This leads [19] to a linear $A = \gamma(\varepsilon_i - \varepsilon_j)$ model with a variance in the particle transfer between boxes $B_{ij} = \gamma(\varepsilon_i (1 - \varepsilon_j) + \varepsilon_j (1 - \varepsilon_i))$. Using the procedure outlined above we find here

$$C_{12} = \frac{-N}{27} (\varepsilon_0 - \varepsilon_3)^2. \tag{10}$$

Note that now the correlations are negative as opposed to positive in the previous example.

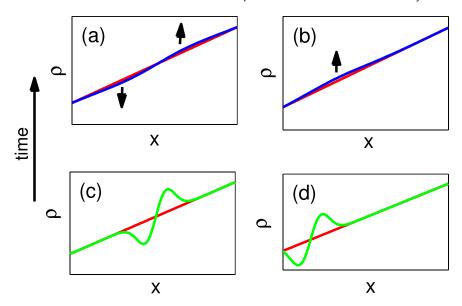


Figure 2. In continuum models with constant diffusivity, a current fluctuation in the center of the system (c) relaxes to a density fluctuation (a) which gives a negative contribution to spatial correlations. In contrast, a current fluctuation near the boundaries (d) adds a positive contribution to spatial correlations (b).

We next turn to calculate the correlations in two quantum examples. In the first we consider coupled Fermi gases. In the second, treated in appendix A, we study systems which exchange energy through blackbody radiation. In both cases the correlations turn out to be positive.

5. Ideal Fermi gas

We now derive the correlation function starting from a quantum mechanical model of weakly interacting Fermi gases. To derive the average energy flow, we first consider two Fermi gases in adjacent boxes, with Hamiltonian $H = H_c + H_d + H'$, where $H_c = \sum_k \epsilon_k^c c_k^\dagger c_k$, $H_d = \sum_k \epsilon_k^d d_k^\dagger d_k$, and $H' = \lambda c_0^\dagger c_0 d_0^\dagger d_0$. In box c (d) we label fermion operators by $c, c^\dagger (d, d^\dagger)$, with dispersion ϵ_k^c (ϵ_k^d). In addition, there is a density–density interaction between the c and d fermions, localized at the contact point ('0') between the two boxes. This interaction allows for the exchange of energy between the boxes, but does not permit particle exchange between them.

For small λ we can treat H' as a weak perturbation. Its effect is to scatter a pair of electrons $\{c_{k_2}, d_{k_4}\}$ to a new pair $\{c_{k_1}, d_{k_3}\}$. Each such process leads to a transfer of energy $\epsilon_{k_3} - \epsilon_{k_4}$. The rate for this process is computed using the Fermi golden rule, leading to an average energy current $A_{cd} = f_{cd}^{(1)}$ and fluctuations $B_{cd} = f_{cd}^{(2)}$, where

$$f_{cd}^{(m)} = \frac{2\pi|\lambda|^2}{\hbar} \sum_{k_1...k_4} (\epsilon_{k_3} - \epsilon_{k_4})^m n_{k_2}^c n_{k_4}^d (1 - n_{k_1}^c) (1 - n_{k_3}^d) \delta\left(\epsilon_{k_2} + \epsilon_{k_4} - \epsilon_{k_1} - \epsilon_{k_3}\right)$$
(11)

where $n_k^c = 1/\exp((\epsilon_k - \mu_c)/T_c) + 1$ and $n_k^d = 1/\exp((\epsilon_k - \mu_d)/T_d) + 1$ are Fermi occupation functions and $\mu_{c,d}$ are chemical potentials. Note that we allow for different densities and temperatures in the two boxes.

We can replace the sums over k by integrals over energy, $\sum_k \to \int d\epsilon \, \nu(\epsilon)$. Furthermore, provided that $T_c, T_d \ll T_F$, the Fermi temperature, we can ignore the energy dependence of the density of states, and replace $\nu(\epsilon) \to \nu_{c,d}$ for each of the boxes. We then obtain

$$A_{cd} = \kappa_{cd} (T_c^4 - T_d^4) \tag{12}$$

where $\kappa_{cd} = 2\pi^5/15\hbar|\lambda|^2\nu_c^2\nu_d^2$. B_{cd} is found to be well-described by

$$B_{cd} \approx 4\kappa_{cd}(T_c^5 + T_d^5). \tag{13}$$

This expression is exact at $T_c = T_d$, and can be shown numerically to be within 1.5% of the exact result for all values of T_c and T_d . Equation (4) is satisfied when $T_c \approx T_d$. In what follows we will consider the case where the subsystems and the baths are all described by Fermi gases, and we will take $\kappa_{cd} = \kappa$ to be constant across all junctions.

Using equations (12) and (5) we solve for the average energies to find $T_1 = ((2T_0^4 + T_3^4)/3)^{1/4}$ and $T_2 = ((T_0^4 + 2T_3^4)/3)^{1/4}$. From this, following the procedure described above, and using the expression for E to lowest order in T/T_F , $E_i = \frac{3}{5}NT_F\left(1 + 5\pi^2/12\left(T_i/T_F\right)^2\right)$, we find

$$C_{12} = \frac{\pi^2 N \left(B_{01} - 2B_{12} + B_{23} \right)}{48\kappa T_{\rm F} \left(T_1^2 + T_2^2 \right)} \tag{14}$$

whose sign depends on the convexity of B_{cd} . Using equation (13) we find that correlations are always positive and increase with the temperature difference between the two baths.

6. Bound on the energy correlations

It is natural to ask how the magnitude of the correlations is related to the different energy scales in the problem: k_BT_i for i=0–3, and a dynamic energy scale $\langle J \rangle \tau$, composed of the average current $\langle J \rangle$ multiplied by the relaxation time of fluctuation to the steady-state, τ . The last energy scale is directly related to the non-equilibrium steady-state.

In appendix B we consider systems in which equation (4) holds, and show that, if the average currents grow with the difference between E_i and E_j (i.e. $\partial_{E_i} A_{ij} > 0$ and $\partial_{E_j} A_{ij} < 0$), then

$$|C_{12}| \le 18 \langle J \rangle \tau \max_{i=0,1,2} \left(|\beta_i - \beta_{i+1}|^{-1} \right).$$
 (15)

This gives an upper bound on the size of the correlations. The bound is proportional to the dynamic energy scale $\langle J \rangle \tau$, demonstrating how correlations disappear upon approach to equilibrium.

The bound above can be understood as follows. Correlations have units of energy squared, and vanish at equilibrium. One therefore expects them to scale as $\langle J \rangle \tau$, multiplied by an energy scale derived from the energies $k_B T_i$. This is expressed by the Lyapunov equation (7). The matrix \mathbf{R} sets the time scale τ for relaxation, with $\tau^{-1} = \lambda_{\text{max}}$, the larger of the eigenvalues of \mathbf{R} . The elements of the matrix \mathbf{Q} are sums of B_{ij} terms. According to equation (4), $B_{ij} = 2A_{ij}/(\beta_i - \beta_j)$, these are proportional to the averaged currents A_{ij} multiplied by a combination of the energy scales $k_B T_i$. The solution to the Lyapunov equation therefore scales as $\langle J \rangle \tau$ multiplied by a combination of the energies $k_B T_i$.

Interestingly, the energy scale $\langle J \rangle \tau$ also appears in an exact relation for the non-equilibrium fluctuations in a simpler model, where a single system is connected to two heat baths. In [13] it was shown that when equation (4) holds, energy fluctuations in the system are given by $\langle J \rangle \tau((\beta_2 - \beta_1)/(\beta_0 - \beta_1)(\beta_2 - \beta_0))$.

Finally, we comment that the large deviation functional (LDF) of the energy or density profile in non-equilibrium systems has recently been the subject of close attention [3, 15, 16], [20]–[25]. In particular, it has been shown that the non-locality of the LDF is directly related to the non-equilibrium long-range correlations [3, 26]. The model studied here, being low dimensional, might serve as a good template for understanding their general properties, in particular, in light of the insights gained into the sign of the correlations in driven diffusive systems. To date, the LDF has been determined for a few models in the macroscopic limit [16, 27] and obeys a variational principle which remains to be understood for generic systems.

Acknowledgments

We are very grateful to Luca D'Alessio for many useful discussions and important comments. We would like to acknowledge the support of BSF and ISF grants, the EU under grant agreement no. 276923—MC–MOTIPROX, NSF DMR-0907039 and the SCHePS Paris 7 interdisciplinary project.

Appendix A. Coupled blackbodies

Let us consider two coupled cavities at different temperatures. Each cavity has a thermal gas of photons, and the photons can leak from one cavity to the other at the interface between the two. The Hamiltonian is

$$H = \sum_{\mathbf{k},\alpha} \hbar c k a_{\mathbf{k},\alpha}^{\dagger} a_{\mathbf{k},\alpha} + \sum_{\mathbf{k},\alpha} \hbar c k b_{\mathbf{k},\alpha}^{\dagger} b_{\mathbf{k},\alpha} + H_{\text{int}}.$$
 (A.1)

Here, a, a^{\dagger} and b, b^{\dagger} denote creation and annihilation operators of photons with momentum \mathbf{k} and polarization α in each of the two cavities and $k = |\mathbf{k}|$. We impose reflecting boundary conditions at the surface z = 0 separating the two cavities, so that k_z in the sums is restricted to positive values. The interaction term

$$H_{\text{int}} = \lambda \hbar c \sum_{\alpha=1}^{2} \int dx \, dy \, a_{(x,y,0),\alpha}^{\dagger} b_{(x,y,0),\alpha} + \text{h.c.}$$
(A.2)

effectively converts outgoing a photons from the first cavity into b photons in the second cavity. The prefactor λ plays the role of the transmission through the barrier separating the two systems and we assume that the barrier does not radiate, i.e. it is effectively at zero temperature. Usual blackbody radiation corresponds to $\lambda = 1$.

To calculate the energy transfer between the two systems we will use the Fermi golden rule treating H_{int} as a perturbation. Then, the matrix element squared for a photon of a given polarization with momentum $\mathbf{k} = (k_x, k_y, k_z)$ in box a to transfer to box b with

momentum $\mathbf{k}' = (k_x, k_y, k_z')$ is

$$\frac{|\lambda|^2 \hbar^2 c^2}{L^2} n_{\mathbf{k}}^{(a)} (1 + n_{\mathbf{k}'}^{(b)}) \tag{A.3}$$

where $n_{\mathbf{k}}^{(a,b)} = 1/(\mathrm{e}^{-\beta_{a,b}\hbar ck} - 1)$ are the Bose occupation factors for the two cavities, and L is the linear size of the cavities along the contact surface. Each such process leads to an energy transfer $\hbar ck$. Then, using the Fermi golden rule and summing over initial and final states, we find the energy transfer rate from box a to b to be

$$W_{a\to b} = \frac{|\lambda|^2 \hbar c^2 L^2}{(2\pi)^3} \sum_{\alpha=1}^2 \int d^3 \mathbf{k} \int dk'_z \, \hbar c k \, n_{\mathbf{k}}^{(a)} (1 + n_{\mathbf{k}'}^{(b)}) \delta(\hbar c k - \hbar c k').$$

Taking into account that $k_z > 0$ and $k_z' > 0$, and using $\delta(\hbar ck - \hbar ck') = (k_z/\hbar ck)\delta(k_z - k_z')$, we obtain

$$W_{a\to b} = \frac{|\lambda|^2 \hbar c^2 L^2}{4\pi^2} \int_0^\infty dk \, k^3 n_k^{(a)} (1 + n_k^{(b)}). \tag{A.4}$$

This reduces to the usual result for blackbody radiation when $\beta_b = \infty$ and $\lambda = 1$.

The net energy transfer rate from a to b is then

$$A_{ab} = W_{a \to b} - W_{b \to a} = |\lambda|^2 \sigma (T_a^4 - T_b^4) \tag{A.5}$$

where $\sigma = \pi^2 L^2 / 60c^2 \hbar^3$ is the Stefan constant. Note that since the energy density of a blackbody is proportional to T^4 the blackbody radiation results in a linear A model. We can compute the energy fluctuations in a similar fashion,

$$B_{ab} = \frac{|\lambda|^2 \hbar c^2 L^2}{(2\pi)^3} \sum_{\alpha=1}^2 \int d^3 \mathbf{k} \int dk' (\hbar ck)^2 \, \delta(\hbar ck - \hbar ck') \left[n_{\mathbf{k}}^{(a)} (1 + n_{\mathbf{k}'}^{(b)}) + n_{\mathbf{k}}^{(b)} (1 + n_{\mathbf{k}'}^{(a)}) \right]$$
$$= \frac{|\lambda|^2 \hbar^2 c^3 L^2}{(2\pi)^2} \int dk \, k^4 \left[n_k^{(a)} (1 + n_k^{(b)}) + n_k^{(b)} (1 + n_k^{(a)}) \right].$$

Numerically we find that this expression is well approximated by

$$B_{ab} = 8|\lambda|^2 \sigma \left[(T_a T_b)^{5/2} + \frac{45}{\pi^4} \left(T_a^{5/2} - T_b^{5/2} \right)^2 \right]$$
 (A.6)

which satisfies equation (4) when $T_a \approx T_b$. Applying our formalism in a similar fashion to that described in the discussion of the Fermi gas, we find that the correlations are always positive.

As a side remark we note that the same result for the blackbody radiation can be obtained using a different type of perturbation

$$H_{\text{int}} = \frac{\lambda \hbar c}{\mathrm{i}} \sum_{\alpha=1}^{2} \int \mathrm{d}x \,\mathrm{d}y \,a_{(x,y,0),\alpha}^{\dagger} \partial_{z} b_{(x,y,z),\alpha} \bigg|_{z=0} + \text{h.c.}$$
(A.7)

At $\lambda = 1$ this perturbation is nothing but the energy flux operator of photons. The easiest way to check that this perturbation gives the same result as equation (A.2) is to discretize the Hamiltonian along the z-direction,

$$H_{\text{int}} = \lambda \frac{\hbar c}{\mathrm{i}d} \sum_{\alpha=1}^{2} \int \mathrm{d}x \,\mathrm{d}y \, \left[a_{(x,y,0),\alpha}^{\dagger} b_{(x,y,d),\alpha} - \text{h.c.} \right], \tag{A.8}$$

where d is the lattice spacing, and then make the gauge transformation $b \to b \mathrm{e}^{-\mathrm{i}\pi/2}$. This gauge transformation obviously does not affect H_0 , while H_{int} reduces to equation (A.2) in the continuum limit, using $1/d \to \delta(z)$. Hence we recover the equivalence of the two choices for the perturbation.

Appendix B. Derivation of equation (15)

Here we derive the bound given in equation (15). We assume that $\partial_{E_i} A_{ij} > 0$ and $\partial_{E_j} A_{ij} < 0$, meaning that the average currents grow when the difference between E_i and E_j grows. The matrix **R** is then given by

$$\mathbf{R} = \begin{pmatrix} -\gamma_1 - \gamma_2 & \gamma_3 \\ \gamma_2 & -\gamma_3 - \gamma_4 \end{pmatrix},$$

where $\gamma_1 = -\partial_{E_1} A_{01}$, $\gamma_2 = \partial_{E_1} A_{12}$, $\gamma_3 = -\partial_{E_2} A_{12}$, and $\gamma_4 = \partial_{E_2} A_{23}$ are all positive numbers. As $|\operatorname{Tr}(\mathbf{R})| = |\gamma_1 + \gamma_2 + \gamma_3 + \gamma_4|$ then $\max \gamma \equiv \max \{\gamma_i\} \leq |\operatorname{Tr}(\mathbf{R})|$.

The correlation is given by

$$C_{12} = \frac{B_{23}(\gamma_1 + \gamma_2)\gamma_3 + B_{01}\gamma_2(\gamma_3 + \gamma_4)}{2\text{Tr}(\mathbf{R})\det(\mathbf{R})} - \frac{B_{12}(\gamma_2\gamma_4 + \gamma_1\gamma_3 + 2\gamma_1\gamma_4)}{2\text{Tr}(\mathbf{R})\det(\mathbf{R})},$$

so that

$$|C_{12}| \leq \frac{9 \max (B_{ij}) (\max \gamma)^{2}}{2 |\operatorname{Tr} (\mathbf{R})| \det (\mathbf{R})}$$

$$\leq 9 \langle J \rangle \frac{|\operatorname{Tr} (\mathbf{R})|}{\det (\mathbf{R})} \max_{i=0,1,2} \left(|\beta_{i} - \beta_{i+1}|^{-1} \right),$$

where in the second equality we used equation (4), that at the steady-state $\langle J \rangle = A_{01} = A_{12} = A_{23}$, and that $(\max \gamma)^2 \leq [\operatorname{Tr}(\mathbf{R})]^2$. As $|\operatorname{Tr}(\mathbf{R})| / \det(\mathbf{R}) = \lambda_1^{-1} + \lambda_2^{-1} \leq 2\tau$, one obtains equation (15)

$$|C_{12}| \le 18 \langle J \rangle \tau \max_{i=0,1,2} (|\beta_i - \beta_{i+1}|^{-1}).$$

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