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The local approach to quantum transport may violate the second law of thermodynamics

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Abstract – Clausius statement of the second law of thermodynamics reads: Heat will flow spontaneously from a hot to cold reservoir. This statement should hold for transport of energy through a quantum network composed of small subsystems each coupled to a heat reservoir. When the coupling between nodes is small, it seems reasonable to construct a local master equation for each node in contact with the local reservoir. The energy transport through the network is evaluated by calculating the energy flux after the individual nodes are coupled. We show by analyzing the most simple network composed of two quantum nodes coupled to a hot and cold reservoir, that the local description can result in heat flowing from cold to hot reservoirs, even in the limit of vanishing coupling between the nodes. A global derivation of the master equation which prediagonalizes the total network Hamiltonian and within this framework derives the master equation, is always consistent with the second law of thermodynamics.

Introduction. – Transport of energy in and out of a quantum device is a key issue in emerging technologies. Examples include molecular electronics, photovoltaic devices, quantum refrigerators and quantum heat engines [1–3]. A quantum network composed of quantum nodes each coupled to local reservoir and to other nodes constitutes the network. The framework for describing such devices is the theory of open quantum systems. The dynamics is postulated employing completely positive quantum master equations [4,5]. Solving the dynamics allows to calculate the steady-state transport of energy through the network.

It is desirable to have the framework consistent with thermodynamics. The first law of thermodynamics is a conservation law of energy; the energy of an isolated system is constant and can be divided into heat and work [6]. The dynamical version of the second law of thermodynamics states that for an isolated system the rate of entropy production is non-negative [7]. For a typical quantum device the second law can be expressed as

$$\frac{d}{dt}\Delta S = \frac{dS_{\text{int}}}{dt} + \frac{dS_{\text{m}}}{dt} - \sum_{i} \frac{J_i}{T_i} \geq 0,$$

where \( dS_{\text{int}}/dt \) is the rate of entropy production due to internal processes, expressed by the von Neumann entropy, \( dS_{\text{m}}/dt \) is the entropy flow associated with matter entering the system, and the last term is the contribution of heat flux, \( J_i \), from the reservoir \( i \).

Microscopic derivation of a global Markovian master equation (MME) of Lindblad-Gorini-Kossakowski-Sudarshan (LGKS) form [4,5], for the network is usually intricate. The local approach simplifies this task [8–14]. It is commonly considered that if the different parts of the network are weakly coupled to each other, a local master equation is sufficient to describe all the properties of the network. We will show that the local approach is only valid for local observables such as the population of each node, and is not valid for non-local observables describing energy fluxes.

The network model. – The simplest network model composed of two nodes shown in fig. 1 and is sufficient to demonstrate the distinction between the local and global approach. Heat is transported between two subsystems \( A \) and \( B \), where each is coupled to a single heat bath with temperature \( T_h \) and \( T_c \). The two subsystems are weakly coupled to each other. The global Hamiltonian is of the form

$$\mathbf{H} = \mathbf{H}_A + \mathbf{H}_B + \mathbf{H}_{AB} + \mathbf{H}_h + \mathbf{H}_c + \mathbf{H}_{Ah} + \mathbf{H}_{Bc}. \quad (2)$$

The bare network Hamiltonian, is \( \mathbf{H}_0 = \mathbf{H}_A + \mathbf{H}_B \) where the node Hamiltonians are \( \mathbf{H}_A = \omega_h \hat{a}^\dagger \hat{a} \) and \( \mathbf{H}_B = \omega_c \hat{b}^\dagger \hat{b} \).
which are composed of either two harmonic oscillators (HO) or of two-level systems (TLS), depending on the commutation relation.

\[
\begin{align*}
\hat{aa}^\dagger + \delta \hat{a}^\dagger \hat{a} &= 1, & \hat{aa} + \delta \hat{aa}^\dagger &= 0, \\
\hat{bb}^\dagger + \delta \hat{b}^\dagger \hat{b} &= 1, & \hat{bb} + \delta \hat{bb}^\dagger &= 0
\end{align*}
\]  

with \(\delta = 1\) for the TLS and \(\delta = -1\) for oscillators. The interaction between the system \(A\) and \(B\) is described by the swap Hamiltonian, \(\hat{H}_{AB} = \epsilon(\hat{a}^\dagger \hat{b} + \hat{b}^\dagger \hat{a})\), with \(\epsilon > 0\). The hot (cold) baths Hamiltonians are denoted \(\hat{H}_{h(c)}\), where \(T_h > T_c\). The system-bath interaction is given by, \(\hat{H}_{Ah} = \gamma_h \hat{a} \otimes \hat{R}_h\) and \(\hat{H}_{Bc} = \gamma_c \hat{b} \otimes \hat{R}_c\), with \(\hat{R}_{h(c)}\) operators belonging to the hot (cold) bath Hilbert space, and \(\gamma_{h(c)}\) are the system-baths coupling parameters.

The dynamics of the reduced system \(A + B\) is governed by the Master equation,

\[
\frac{d}{dt}\hat{\rho}_s = -i[\hat{H}_0 + \hat{H}_{AB}, \hat{\rho}_s] + L_h \hat{\rho}_s + L_c \hat{\rho}_s. 
\]  

With the LGKS dissipative terms, \(L_{h(c)}\), which differ for the local and global approaches. At steady state the heat flow from the hot (cold) bath is given by

\[
\mathcal{J}_{h(c)} = \text{Tr}[\{L_{h(c)} \hat{\rho}_s, \hat{H}_{h(c)} + \hat{H}_{AB}\}],
\]  

where \(\hat{\rho}_s\) is the steady-state density operator.

**Local approach.** – In the local approach it is assumed that the inter-system coupling does not affect the system bath coupling. Therefore in the derivation of the MME the Hamiltonian \(\hat{H}_{AB}\) is ignored and the dissipative terms take the form,

\[
\begin{align*}
L_h \hat{\rho}_s &= \gamma_h \left( \hat{a} \hat{\rho}_s \hat{a}^\dagger - \frac{1}{2} [\hat{a}^\dagger \hat{a}, \hat{\rho}_s] \right) \\
&\quad + e^{-\beta_h \omega_h} \left( \hat{a}^\dagger \hat{\rho}_s \hat{a} - \frac{1}{2} [\hat{aa}^\dagger, \hat{\rho}_s] \right), \quad \text{(6)} \\
L_c \hat{\rho}_s &= \gamma_c \left( \hat{b} \hat{\rho}_s \hat{b}^\dagger - \frac{1}{2} [\hat{b}^\dagger \hat{b}, \hat{\rho}_s] \right) \\
&\quad + e^{-\beta_c \omega_c} \left( \hat{b}^\dagger \hat{\rho}_s \hat{b} - \frac{1}{2} [\hat{bb}^\dagger, \hat{\rho}_s] \right). \quad \text{(7)}
\end{align*}
\]

when the node-to-node coupling is zero, \(\hat{H}_{AB} = 0\), each of the local master equations, eq. (6) and eq. (7), drives the local node to thermal equilibrium. The dynamics of the network is completely characterized by the expectation values of four operators: Two local observables \((\hat{a}^\dagger \hat{a}), (\hat{b}^\dagger \hat{b})\), and two \(AB\) correlations \(\langle \hat{X} \rangle \equiv \langle \hat{a}^\dagger \hat{b} + \hat{b}^\dagger \hat{a} \rangle\) and \(\langle \hat{Y} \rangle \equiv i(\hat{a}^\dagger \hat{b} - \hat{b}^\dagger \hat{a})\). For the dynamics we obtain

\[
\begin{align*}
\frac{d}{dt} \langle \hat{a}^\dagger \hat{a} \rangle &= -\gamma_h (1 + \delta e^{-\beta_h \omega_h}) \langle \hat{a}^\dagger \hat{a} \rangle + \gamma_h e^{-\beta_h \omega_h} - e \langle \hat{Y} \rangle, \\
\frac{d}{dt} \langle \hat{b}^\dagger \hat{b} \rangle &= -\gamma_c (1 + \delta e^{-\beta_c \omega_c}) \langle \hat{b}^\dagger \hat{b} \rangle + \gamma_c e^{-\beta_c \omega_c} + e \langle \hat{Y} \rangle, \\
\frac{d}{dt} \langle \hat{X} \rangle &= -\frac{1}{2} (\gamma_h (1 + \delta e^{-\beta_h \omega_h}) + \gamma_c (1 + \delta e^{-\beta_c \omega_c})) \langle \hat{X} \rangle \langle \hat{Y} \rangle - \langle \omega_h - \omega_c \rangle \langle \hat{Y} \rangle, \\
\frac{d}{dt} \langle \hat{Y} \rangle &= -\frac{1}{2} (\gamma_h (1 + \delta e^{-\beta_h \omega_h}) + \gamma_c (1 + \delta e^{-\beta_c \omega_c})) \langle \hat{X} \rangle + 2e \langle \langle \hat{a}^\dagger \hat{a} \langle \hat{b}^\dagger \hat{b} \rangle\rangle. 
\end{align*}
\]

The rates \(\gamma_{h(c)} > 0\) depend on the specific properties of the bath and its interaction with the system. Equations (8) fulfill the dynamical version of the first law of thermodynamics: The sum of all energy (heat) currents at steady state is zero, \(\mathcal{J}_h + \mathcal{J}_c = 0\). The heat flow from the hot heat bath can be cast in the form (see [15] for details)

\[
\mathcal{J}_h = (e^{\delta \omega_h} - e^{-\delta \omega_h}) \mathcal{F},
\]

where \(\mathcal{F}\) is a function of all the parameters of the system, which is always positive, and is different for the HO and TLS medium. The Clausius statement for the second law of thermodynamics implies that heat can not flow from a cold body to a hot body without external work being performed on the system. It is apparent from eq. (9), that the direction of heat flow depends on the choice of parameters. For \(\frac{\omega_c}{\omega_h} < \frac{\omega_h}{\omega_c}\) heat will flow from the cold bath to the hot bath, thus the second law is violated even at vanishing small \(AB\) coupling, cf. fig. 2. The breakdown of the second law has been examined in several models, see [16] and references therein. In [16] a Fermionic transport model between two heat baths at the same temperature was studied in the weak system-bath coupling limit MME and was compared to a solution within the formalism of nonequilibrium Green functions. At steady state, the current between the baths according to the weak coupling MME is nonzero, which implies a violation of the second law in the sense that heat flows constantly between two heat baths at the same temperature. This sort of violation can also be observed in eq. (9) when taking \(T_h = T_c\). It was claimed in [16] that the violation of the second law is a consequence of neglecting higher-order coherent processes between the system and the baths due to the weak coupling limit. In fact, the treatment introduced in [16] corresponds to the local approach described above. Next, we introduce a proper weak coupling MME, which always obeys the second law of thermodynamics.

**Global approach.** – The global approach is based on the holistic perception where the MME is derived in the
The red zones corresponds to negative entropy production rate, a clear violation of thermodynamics. The global MME, by construction, therefore we restrict the analysis to the harmonic nodes. Following the standard weak coupling limit, in the regime therefore we restrict the analysis to the harmonic nodes. Following the standard weak coupling limit, in the regime 

\[
\mathcal{H}_h = \frac{1}{\sin^2(\theta) e^{\beta\omega_-} - e^{\gamma_c \omega_-}} \gamma_c \gamma_h \omega_- \\
+ \frac{e^{\beta\omega_-} (-1 + e^{\gamma_c \omega_-}) \gamma_c + e^{\gamma_c \omega_-} (-1 + e^{\beta\omega_-}) \cos^2(\theta) \gamma_h}{\sin^2(\theta) e^{\beta\omega_-} - e^{\gamma_c \omega_-}} \gamma_c \gamma_h \omega_+ \\
+ e^{\beta\omega_-} (-1 + e^{\gamma_c \omega_-}) \cos^2(\theta) \gamma_c + \sin^2(\theta) e^{\beta\omega_-} (-1 + e^{\beta\omega_-}) \gamma_h.
\]

(13)

where \( \omega_- > 0 \) the dissipative terms of the MME reads,

\[
\mathcal{L}_h \rho_s = \gamma_h^+ \cos^2(\theta) \left( \hat{d}_+ \rho_s \hat{d}_+^\dagger - \frac{1}{2} \{ \hat{d}_+^\dagger \hat{d}_+, \rho_s \} \right) \\
+ e^{-\beta \omega_-} \left( \hat{d}_- \rho_s \hat{d}_-^\dagger - \frac{1}{2} \{ \hat{d}_-^\dagger \hat{d}_-, \rho_s \} \right) \\
+ \gamma_h^- \sin^2(\theta) \left( \hat{d}_- \rho_s \hat{d}_-^\dagger - \frac{1}{2} \{ \hat{d}_-^\dagger \hat{d}_-, \rho_s \} \right) \\
+ e^{-\beta \omega_-} \left( \hat{d}_+ \rho_s \hat{d}_+^\dagger - \frac{1}{2} \{ \hat{d}_+^\dagger \hat{d}_+, \rho_s \} \right)
\]

(11)

and

\[
\mathcal{L}_r \rho_s = \gamma_c^+ \sin^2(\theta) \left( \hat{d}_+ \rho_s \hat{d}_+^\dagger - \frac{1}{2} \{ \hat{d}_+^\dagger \hat{d}_+, \rho_s \} \right) \\
+ e^{-\beta \omega_-} \left( \hat{d}_- \rho_s \hat{d}_-^\dagger - \frac{1}{2} \{ \hat{d}_-^\dagger \hat{d}_-, \rho_s \} \right) \\
+ \gamma_c^- \cos^2(\theta) \left( \hat{d}_- \rho_s \hat{d}_-^\dagger - \frac{1}{2} \{ \hat{d}_-^\dagger \hat{d}_-, \rho_s \} \right) \\
+ e^{-\beta \omega_-} \left( \hat{d}_+ \rho_s \hat{d}_+^\dagger - \frac{1}{2} \{ \hat{d}_+^\dagger \hat{d}_+, \rho_s \} \right)
\]

(12)

with \( \gamma_h^{\pm} = \gamma_{h(c)} (\omega_{\pm}) \). The calculated steady-state heat flow from the hot bath is given by

\[
\text{see eq. (13) above}
\]

which is positive for all physical choice of parameters. Rewriting eq. (11) and eq. (12) in the local basis, the effective coupling of subsystem \( A \) with the cold bath and of subsystem \( B \) with the hot bath is immediately apparent (see [15] for details). These equations converge to eq. (6) and eq. (7) for \( \epsilon = 0 \).

To further study the dynamics of \( A \) and \( B \), the explicit form of heat baths is specified, characterizing the rates \( \gamma \) [20]:

\[
\gamma_l \equiv \gamma_l (\Omega) = \pi \sum_k |g_l(k)|^2 \delta(\omega(k) - \Omega)[1 - e^{-\beta \omega(k)}]^{-1},
\]

(14)

where \( \omega(k) \) are the frequencies of the baths modes. For the case of a 3-dimensional phonon bath with a linear dispersion relation the relaxation rate can be expressed as

\[
\gamma_l (\Omega) = \kappa \Omega^2 [1 - e^{-\beta \Omega}]^{-1},
\]

(15)

where \( \kappa > 0 \) embodies all the constants and is proportional to the square of the system-bath coupling.

The steady-state observables of the local and global approached are compared in fig. 3 as a function of the node-to-node coupling strength \( \epsilon \). For local observables
such as the local population $\hat{n}_A \equiv \langle \hat{a}^\dagger \hat{a} \rangle$ the two approaches converge to the thermal population when $\epsilon \ll \{\omega_h, \omega_c, \sqrt{\omega_h - \omega_c}\}$. However, the non-local observables such as the current $J_h$ deviate qualitatively. In the local approach when $\frac{\omega_h}{T_c} < \frac{\omega_c}{T_h}$ the second law is violated: the heat flow becomes negative for all values of the coupling $\epsilon$ while for the global approach $J_h$ is always positive, cf. fig. 3.

The local approach is also not reliable even for parameters where the second law is obeyed: $\frac{\omega_c}{T_c} > \frac{\omega_h}{T_h}$. Deviations from the exact global approach appear in the favorable domain of small $\epsilon$, as seen in fig. 4 displaying $J_h$ for a wide range of $\omega_c$. It is noteworthy that the behavior of the heat flows observed in fig. 4 will be the same for all $\epsilon$, also when $\epsilon \ll \kappa$. The only domain where the global approach breaks down is on resonance, when $\omega_h = \omega_c$ and $\epsilon < \kappa$. At this point, the secular approximation is not justified since the two Bohr frequencies $\omega_h$ are not well separated, and on the time scale $1/\kappa \omega^2$, one can not neglect rotating terms such as $\epsilon^2 \omega^2$ [21].

Additional insight is obtained when examining the covariance matrix for the two-mode Gaussian state (see [15] for details). The correlations between subsystems $A$ and $B$ is fully determined by the set of correlation functions \{\text{cor}(x_A, x_B), \text{cor}(x_A, p_B), \text{cor}(p_A, x_B), \text{cor}(p_A, p_B)\}. \{x, p\}$ are the position and momentum coordinates of the subsystems. In both approaches \text{cor}(x_A, x_B) and \text{cor}(p_A, p_B) are equal for small $\epsilon$. The two additional correlations, \text{cor}(x_A, p_B) and \text{cor}(p_A, x_B), vanish at steady state in the global approach, where in the local approach they remain finite. Thus, in the local approach the nodes are over correlated compared to the global approach. It should be noted that in steady state none of the approaches generate entanglement. The two-mode Gaussian state is a separable state according to the separability criterion for continuous variable systems [22,23].

To summarize: As expected, the local dynamical approach is incorrect for strong coupling between the subsystems. In the weak coupling limit, local observables converge to their correct value. The non-local observables such as heat currents are qualitatively and quantitatively erroneous in the local MME. A strong indication is the violation of the second law of thermodynamics. The completely positive LGKS generator is a desired form for the master equation. However, for consistency with the physical world, a microscopic global derivation of the master equation is required. Such approaches are consistent with thermodynamics [24–28].

**REFERENCES**

Quantum transport and violation of the second law of thermodynamics


