Kraus operators for a pair of interacting qubits: a case study

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Abstract The Kraus form of the completely positive dynamical maps is appealing from the mathematical and the point of the diverse applications of the open quantum systems theory. Unfortunately, the Kraus operators are poorly known for the two-qubit processes. In this paper, we derive the Kraus operators for a pair of interacting qubit, while the strength of the interaction is arbitrary. One of the qubits is subjected to the x-projection spin measurement. The obtained results are applied to calculate the dynamics of the initial entanglement in the qubits system. We obtain the loss of the correlations in the finite time interval; the stronger the inter-qubit interaction, the longer lasting entanglement in the system.

1. Introduction

The "integral", i.e. so-called, Kraus form [1] of a completely positive dynamical map for an open quantum system [2, 3] is appealing for the mathematical reasons. Mathematical existence of the Kraus form for such processes is guaranteed by the Kraus theorem, universally [1-3]. On the other hand, a Kraus-form (KF) may be regarded as a solution to a differential master equation (ME) for the open system’s statistical operator (density matrix); a case when no ME exists for the process can be found e.g. in Refs. [4,5].

The Kraus operators are often constructed due to some physical assumptions or understanding of the underlying physical processes [6]. Nevertheless, such derivations may not provide the full physical (e.g. microscopic) details [7]. One way to obtain a proper KF for the open system’s dynamics is derivation from the related master equation for the process [7,8]–if such an ME exists [4,5]. To this end, it is important to note: phenomenological derivations of MEs may also be unreliable–often there appear certain subtleties of both mathematical and physical nature as well as unexpected pitfalls [9,10].

Having this in mind as well as the above-distinguished usefulness of KF, in this paper we derive the Kraus operators starting from a microscopically derived master equation for a pair of two-level systems (qubits). We are concerned with an ancilla qubit interacting with another qubit, which is subjected to a $S_x$ spin-projection quantum measurement. Usefulness of the
KF for the process is emphasized by application of our results in investigating 
the dynamics of entanglement in the qubits system. 

2. The master equation and the task

The total, isolated, system is described by the Hamiltonian:

\[ H = H_{10} + H_{20} + H_{E1\circ} + H_{12} + H_{1E1}, \]

where the symbol "\( \circ \)" stands for the subsystems self-Hamiltonians and the 
rest are the interaction terms. While the self-Hamiltonian terms are standard 
(see below), the qubits interaction is chosen \[6,11\]:

\[ H_{12} = \beta S_{1z} \otimes S_{2z}, \]

where the 1/2-spin operators \( S_{pz} = \sigma_{pz}/2 \) \( p = 1,2 \) and we take \( \hbar = 1 \), while 
the interaction with the environment:

\[ H_{1E1} = S_{1x} \otimes \int_0^{\nu_{\text{max}}} d\nu h(\nu)(a_{\nu}^\dagger + a_{\nu}) \equiv S_{1x} \otimes B_{E1}, \]

where appear the annihilation and creation operators satisfying the standard 
Bose-Einstein commutation \([a_{\nu}^\dagger, a_{\nu}] = -\delta(\nu - \nu')\).

For non-interacting qubits, i.e. for \( \beta = 0 \), the qubits can be described by 
mutually independent dynamics. However, for the interacting qubits \( \beta \neq 0 \), 
dynamics of the qubits cannot be mutually independent. Therefore we regard 
the pair of qubits, 1 + 2, as an open system subjected to the environment \( E_1 \) 
as described above.

Physically, eq.(3) describes a quantum measurement of the \( S_{1x} \) observable. We assume the initial tensor product state \( \rho_{12} \otimes \rho_{E1} \) and the weak 
coupling limit for eq.(3), while the environment \( E_1 \) being in the thermal state 
\( \rho_{E1} = \rho_{th} = \exp(-H_{E1\circ}/k_B T)/Z \) on temperature \( T \); \( k_B \) is the Boltzmann 
constant and \( Z \) is the normalization "statistical sum". That is, we consider 
the time-homogeneous, completely positive and trace preserving process for 
the 1 + 2 system and below we derive the proper master equation in the 
weak-coupling limit. Such physical situation is generally described by the 
following, Lindblad-form master equation (in the interaction picture) for the 
pair of qubits \[3\]:

\[ \frac{d\rho_{12}}{dt} = -i[H_{LS}, \rho_{12}] + \sum_{\nu,i,j} \left( \gamma_{ij}(\nu) \left[ A_j(\nu)\rho_{12}A_i(\nu) - \frac{1}{2}(A_i^\dagger(\nu)A_j(\nu), \rho_{12}) \right] \right) \]

(4)
Our task in this section is, starting from equations (1)-(3), to derive the explicit forms for the damping functions $\gamma_{ij}(\nu)$ and the Lindblad operators $A_i(\nu)$; for simplicity, but without loss of generality, we ignore the Lamb shift term $H_{LS}$.

The interaction picture is defined by the self-Hamiltonian

$$H_0 = H_{10} + H_{20} + H_{12} + H_{E_{10}} = \frac{\omega}{2}\sigma_{1z} + \frac{\omega}{2}\sigma_{2z} + \frac{\beta}{4}\sigma_{1z} \otimes \sigma_{2z} + H_{E_{10}},$$

(5)

where the environmental self-Hamiltonian: $H_{E_{10}} = \int_{0}^{\nu_{\text{max}}} d\nu a_\nu^\dagger a_\nu$ with the maximal frequency $\nu_{\text{max}}$. The alternative choice [3,13] of the interaction picture without the $H_{12}$ term in eq.(5) reduces the considerations to the weak qubits interaction, $\beta \ll 1$, which is a special case of our considerations.

From eq.(5) it readily follow the energy eigenvalues and eigenprojectors for the qubits system:

$$E_1 = \omega + \frac{\beta}{4}, \quad P_1 = |++\rangle\langle++|,$$

$$E_2 = -\frac{\beta}{4}, \quad P_2 = |+-\rangle\langle+--|,$$

$$E_3 = -\omega + \frac{\beta}{4}, \quad P_3 = |--\rangle\langle-+|.$$

(6)

where $|mn\rangle \equiv |m\rangle|n\rangle, m, n \in \{+,-\}$ and $\sigma_\pm|\pm\rangle = \pm|\pm\rangle$. From eq.(6) follows the set of the values for the parameter $\nu$ in eq.(4): $\{0, \nu_1 = E_1 - E_2 = \omega + \beta/2, \nu_2 = E_1 - E_3 = 2\omega, \nu_3 = E_2 - E_3 = \omega - \beta/2\}$.

The general expressions for the Lindblad operators [3]:

$$A_i(\nu) = P_n A_i P_m, \quad \nu = E_m - E_n$$

(7)

where $A_i = S_{1x}$, cf. eq.(3), while:

$$\gamma_{kl}(\nu) = 2\pi tr \left[ B_k(\nu) B_l \rho_{E_{10}} \right].$$

(8)

In eq.(8) [3]: $B_k(\nu) = \int_{-\nu}^{\nu} dt \exp(-\nu t)B_k(t)$ for the $B_k(t)$ representing the interaction-picture form of $B_k$. In eq.(3) there is only one such operator, $B_{E_{10}}$.

The desired master equation can be shortly presented as:

$$\frac{d\rho_{12}}{dt} = -\alpha^2 \left( \sum_{i=1}^{2} \left( \gamma_i A_i + \gamma_i^- A_i^\dagger \right) \right) [\rho_{12}],$$

(9)
where $\alpha$ presents the weak coupling constant of the system-environment interaction; if $\gamma \equiv \gamma(\nu)$, then $\gamma^- \equiv \gamma(-\nu)$, with all superoperators satisfying:

$$A[\rho_{12}] = A\rho_{12}A^\dagger - \frac{1}{2}\{A^\dagger A, \rho_{12}\}, \quad A^\dagger[\rho] = A^\dagger\rho_{12}A - \frac{1}{2}\{AA^\dagger, \rho_{12}\}. \quad (10)$$

From (6) and (7) straightforwardly follow the non-zero Lindblad operators in eq.(10):

$$A_1 = \frac{1}{8}\sigma_1^- \otimes (I_2 - \sigma_{2z}),$$
$$A_2 = \frac{1}{8}\sigma_1^- \otimes (I_2 + \sigma_{2z}). \quad (11)$$

Calculation of the damping functions $\gamma_{kl}(\nu)$ is straightforward; technical details are presented in [7]. For completeness, we provide a few main steps.

Due to the only one term in eq.(3), the general expression eq.(8) reduces to:

$$\gamma_{xx}(\nu) = \frac{2\pi h(\nu)}{\nu \max} \int_0^{\nu \max} d\nu' h(\nu') tr(a_{\nu}(a_{\nu'} + a_{\nu'}^\dagger)\rho_{th}). \quad (12)$$

With the use of expressions for the thermal averages, i.e. when the environment is in thermal equilibrium [14]:

$$\overline{a_{\nu}a_{\nu}^\dagger} = 0 = \overline{a_{\nu}^\dagger a_{\nu}}, \quad (13)$$
$$\overline{a_{\nu}^\dagger a_{\nu}^\dagger} = \delta(\nu' - \nu)(1 + \bar{n}(\nu')) \quad (14)$$

and

$$\overline{a_{\nu}^\dagger a_{\nu}} = -\delta(\nu' - \nu)\bar{n}(\nu'), \quad (15)$$

follow the expressions for $\gamma_{xx}(\nu)$ and $\gamma_{xx}(-\nu)$:

$$\gamma_{xx}(\nu) = 2\pi J(\nu)(1 + \bar{n}(\nu)), \quad \gamma_{xx}(-\nu) = 2\pi J(\nu)\bar{n}(\nu), \quad (16)$$

where $\nu$ takes the above distinguished values and the average number of bosons in thermal state $\bar{n}(\nu) = (e^{-\nu/T} - 1)^{-1}$. We choose the standard Ohmic spectral density $J(\nu) = \alpha\nu e^{-\nu/\nu_c}$ with the cutoff $\nu_c$.

In the high temperature limit, which we are concerned with, $\bar{n}(\nu) \gg 1$ and therefore $\gamma(\nu) \approx \gamma(-\nu) = 2\pi J(\nu)\bar{n}(\nu)$, which reduces the list of the damping functions to only two of them (for fixed $\nu$): $\gamma_1 \equiv \gamma(\nu_1)$ and $\gamma_2 \equiv \gamma(\nu_3)$. 


Substitution of eqs. (10) and (11) in eq.(9), after simple calculation, gives
the following relation of the damping functions to the Lindblad operators
appearing in eq.(9):

\[
\begin{align*}
\gamma_1 &\equiv 4\pi J(\omega + \beta/2)\bar{n}(\omega + \beta/2), \quad A_2, A_2^\dagger, \\
\gamma_2 &\equiv 4\pi J(\omega - \beta/2)\bar{n}(\omega - \beta/2), \quad A_1, A_1^\dagger,
\end{align*}
\]

that completes the master equation (9).

3. Derivation of the Kraus operators

In this section we derive the interaction-picture Kraus operators for the
qubits dynamics described by the master equation (9). We use a method
recently developed in Ref.[8].

3.1 A brief overview of the method

In Ref. [8], the authors developed a general procedure for deriving a Kraus
decomposition from a known master equation and vice versa, regarding the
finite-dimensional quantum systems. The only assumption is that the master
equation is local in time.

If the dynamical map for the process eq.(9) is formally presented as:

\[
\rho_{12}(t) = \Phi_t[\rho_{12}(0)],
\]

and the master equation eq.(9) is shortly presented as:

\[
\frac{d\rho_{12}(t)}{dt} = \Lambda_t[\rho_{12}(0)],
\]

then (for the time-independent superoperator \(\Lambda\)), the following matrix rela-
tion is fulfilled:

\[
F = e^{Lt}.
\]

The matrices \(F = (F_{ij})\) and \(L = (L_{ij})\) are well defined for the finite-
dimensional systems, and representations for the map \(\Phi_t\) and for the super-
operator \(\Lambda\), respectively, in a chosen orthonormalized basis \(\{G_i\}\) of hermitian
operators acting on the system’s Hilbert state space.

Introducing the so-called Choi matrix [8]:

\[
S_{nm} = \sum_{r,s} F_{rs} tr (G_r G_n G_s G_m)
\]

and its non-negative (real) eigenvalues \(d_i\), follow the desired Kraus operators:
where the unitary matrix $U = (u_{ij})$ diagonalizes the Choi matrix $S = (S_{nm})$. That is, the procedure provides a Kraus form of the process:

$$\rho_{12}(t) = \sum_i K_i(t)\rho_{12}(0)K_i^\dagger(t).$$

(23)

The trace preservation implies:

$$\sum_i K_i^\dagger(t)K_i(t) = I, \forall t.$$

(24)

### 3.2 The Kraus operators for eq.(9)

Distilled from Section 3.1, the procedure for derivation of the Kraus operator is as follows: First, from the master equation eq.(9), the $L$ matrix is derived. Then due to eq.(20), the $F$ matrix follows that, in accordance with eq.(21), provides the Choi matrix $S$. Finally, diagonalization of the Choi matrix gives rise to the Kraus operators, eq.(22).

We proceed by first obtaining the $\Lambda$ operator in the standard representation of the $\sigma_i \otimes \sigma_j/2, i, j = 0, 1, 2, 3$, operators; $\sigma_i = I/\sqrt{2}$, while for $i > 0$, the $\sigma_i/\sqrt{2}$'s represent the standard (normalized) Pauli operators. The nonzero matrix elements are as follows: $-8(\gamma_1 + \gamma_2) = L_{2,2} = L_{3,3} = L_{5,5} = L_{6,6} = L_{8,8} = L_{9,9} = L_{10,10} = L_{11,11} = L_{12,12} = L_{13,13} = L_{14,14} = L_{15,15}$, $-b(\gamma_1 - \gamma_2) = L_{2,14} = L_{3,15} = L_{14,2} = L_{15,3}$, $-16(\gamma_1 + \gamma_2) = L_{4,4} = L_{16,16}$ and $-16(\gamma_1 - \gamma_2) = L_{4,16} = L_{16,4}$.

From the $L$ matrix easily follows the $F$ matrix with the following nonzero entries: $1 = F_{1,1} = F_{7,7}$, $(\exp(-16t\gamma_1) + \exp(-16t\gamma_2))/2 = F_{2,2} = F_{3,3} = F_{14,14} = F_{15,15}$, $(\exp(-16t\gamma_1) - \exp(-16t\gamma_2))/2 = F_{2,14} = F_{3,15} = F_{14,2} = F_{15,3}$, $(\exp(-8t(\gamma_1 + \gamma_2)) = F_{5,5}$, $F_{6,6} = F_{8,8} = F_{9,9} = F_{10,10} = F_{11,11} = F_{12,12} = F_{13,13}$, $(\exp(-32t\gamma_1) + \exp(-32t\gamma_2))/2 = F_{4,4} = F_{16,16}$ and $(\exp(-32t\gamma_1) - \exp(-32t\gamma_2))/2 = F_{4,16} = F_{16,4}$.

The rest of the calculation is also straightforward but rather involved. Hence we just give the final expressions for the non-zero Kraus operators:

$$K_1 = \frac{1 - e^{-32t\gamma_2}}{2} \begin{pmatrix} 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & t \\ 0 & 0 & 0 & 0 \\ 0 & -t & 0 & 0 \end{pmatrix}, K_2 = \frac{1 - e^{-32t\gamma_2}}{2} \begin{pmatrix} 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & -1 \\ 0 & 0 & 0 & 0 \\ 0 & -1 & 0 & 0 \end{pmatrix},$$

(25)
The unital character of the map \( \Phi_t \) for \( t = 0 \) is described by eq. (9), that is, \( \sum_i K_i I K_i = I \); equivalently \( \Phi_t[I] = I \), see eq. (18), i.e. \( \Lambda_t[I] = 0 \), see eq. (19).

From equations (25)-(30) follows the completeness relation eq. (24) for every instant of time \( t \). Hermiticity of the Kraus operators, \( K_i \), \( \forall i \), implies the unital character of the map \( \Phi_t \) described by eq. (9), that is, \( \sum_i K_i I K_i = I \); equivalently \( \Phi_t[I] = I \), see eq. (18), i.e. \( \Lambda_t[I] = 0 \), see eq. (19).
In the initial instant of time \( t = 0 (\tau = 0) \), \( B = 8 \), and therefore \( A = 0 \) while \( A' = 1/16 \). Therefore \( K_i(0) = 0, i = 1, 2, ..., 7 \), while \( K_8(0) = -I \), thus satisfying the initial condition, \( \rho_{12}(0) = K_8(0)\rho_{12}(0)K_8(0) \). From eqs.(25)-(27) it follows the time independence of the Kraus operators \( K_i, i = 1, 2, ... 6 \) in the asymptotic limit. That is

\[
\lim_{t \to \infty} \sum_{i=1}^{6} K_i(t)K_i(t) = I/4.
\]

Then, due to the completeness relation eq.(24), it follows the dominant contribution from the \( K_7 \) and \( K_8 \) operators:

\[
\lim_{t \to \infty} \sum_{i=7}^{8} K_i(t)K_i(t) = 3I/4.
\]

Of course, this does not imply existence of a stationary state, which is defined in the Schrödinger picture, in which the Kraus operators are defined as:

\[
U_{(12)}^{(12)}(t)K_i(t), i = 1, 2, ..., 8,
\]

where \( U_{(12)}^{(12)}(t) = \exp(-it(H_o - H_{E_1})) \) and \( H_o \) is defined by eq.(5).

### 3.3 Entanglement sudden death for the pair of qubits

As an application of the results of Section 3.2, we investigate entanglement dynamics for the pair of qubits described by the master equation (9). We assume the initial maximally entangled state for the pair of qubits, \( (| + - \rangle + | - + \rangle)/\sqrt{2} \), and choose the following set of the values for the parameters appearing in eq.(17): \( \omega = 0.1, \alpha = 0.02, T = 100, \nu_c = 100 \). As a measure of quantum entanglement we use the standard and well-studied measure of \textit{concurrence} [15]

\[
C(\rho(t)) = \max\{0, \Lambda(t)\},
\]

where \( \Lambda(t) = \sqrt{\lambda_1(t)} - \sqrt{\lambda_2(t)} - \sqrt{\lambda_3(t)} - \sqrt{\lambda_4(t)} \) with the eigenvalues \( \lambda_1 > \lambda_2 > \lambda_3 > \lambda_4 \) of

\[
\rho(t)(\sigma_{1y} \otimes \sigma_{2y})\rho^*(t)(\sigma_{1y} \otimes \sigma_{2y})
\]

and "*" denoting the complex-numbers conjugate. The density matrix \( \rho(t) \) in eq.(35) is the 1 + 2-system’s state eq.(23) for the Kraus operators given by eq.(33).
In Fig.1 we can see dependence of $C$ on both, time $t$ and the strength of the inter-qubits interaction $\beta$. Expectably, we observe dynamical decrease of $C$ and hence of the entanglement in the qubits 1 + 2 system for every chosen value of $\beta$. Fig.2 depicts slower decrease of $C$ for larger $\beta$, for every instant of time $t$. That is, the largest (negative) value of $\Lambda(t)$ increases with the increase of $\beta$. Physically, Figures 1 and 2 reveal detrimental influence of the environment on the initial entanglement for the pair of qubits; the stronger inter-qubits interaction (the larger $\beta$) the qubits more efficiently ”hold together”. Our results reveal the phenomenon of the so-called entanglement sudden death [16]: instead of the expected smooth, asymptotic approach to $C = 0$, we obtain dynamical change of the concurrence from the initial $C(0) = 1$ to the final $C(t) = 0$ value for the finite time interval $t$.

4. Discussion

Procedure presented in this paper is universal, in that it formally equally applies to arbitrary physical situation for a pair of qubits that is subjected
to a completely positive dynamical process. For both qubits monitored by their respective environments (or by a common environment), the technicalities are much more involved but without any conceptual or methodological obstacles or open issues. The use of the Kraus operators is a straightforward application of the matrix (linear algebra) calculus.

Tracing out a qubit from eq.(9) universally leads to a master equation for the other qubit. However, this procedure is trivial only for the initial tensor-product state $\rho_1 \otimes \rho_2$ for the pair of qubits. A qubit can be regarded as a part of the (extended) environment of the other qubit. This gives rise to alternative bipartitions of the total system.

Tracing out that includes the qubit 2 (the qubit 1) regards the $1 + E'_1 \equiv 1 + (E_1 + 2)$ (the $2 + E'_2 \equiv 2 + (1 + E_1)$) bipartition. Only if there is not correlations in the initial state of the qubits, the bipartition of the total system, $(1 + 2) + E_1$, which is assumed in Section 2 and leads to eq.(4), is interchangeable with the alternative bipartitions of the total system: $(1 + E_1) + 2$ and $1 + (E_1 + 2)$. The presence of the initial correlations in the qubits $1 + 2$ system implies initial correlations [17] in both $1 + E'_1$ and $2 + E'_2$ bipartitions, and hence [3] non-complete-positivity of the processes for the individual qubits; for this reason, in general, independent derivations of the master equations for the individual qubits are required [18,19]. That is, the initial correlations in the qubits system breaks the symmetry between the different bipartitions of the total system. For this reason we only regarded the bipartition $(1 + 2) + E_1$ and eq.(4) as the reliable basis for the analysis of the qubits dynamics—the absence of the initial correlations in this bipartition is supposed from the very start, $\rho_{12} \otimes \rho_{E_1}$, Section 2.

For a pair of qubits, every Kraus operator can be written $K_k = \sum_{i,j} c_{ij} A_{1i} \otimes B_{2j}$; an example is given by eq.(11). Then the state for the pair of qubits reads:

$$\rho_{12}(t) = \sum_{k,i,i',j,j'} c_{ij}^k c_{ij'}^k A_{1i} \otimes B_{2j} \rho_{12}(0) A_{1i'}^\dagger \otimes B_{2j'}^\dagger,$$  

(36)

which after the tracing out the qubit 2 (the qubit 1) gives a master equation for the qubit 1 (for the qubit 2). As emphasized above, only for the initial tensor-product state $\rho_{12} = \rho_1 \otimes \rho_2$, one can obtain a Kraus form for a single-qubit dynamics—that guarantees [1] complete positivity of the process the qubit is subjected to. Then eq.(36) gives, e.g.:

$$\rho_1 = tr_2 \rho_{12} = \sum_{i,i'} A_{1i} \rho_1 A_{1i'}^\dagger \left( tr_2 \sum_k (\sum_j c_{ij}^k B_{2j}) \rho_2 (\sum_j c_{ij}^k B_{2j})^\dagger \right).$$  

(37)
Since the terms
\[ b_{ii'} = tr_2 \sum_k (\sum_j c_{kj}^i B_{2j}) \rho_2 (\sum_j c_{kj}^i B_{2j})^\dagger \equiv \sum_k tr_2 B_{2ik} \rho_2 B_{2i'k}^\dagger \]  
constitute a positive semi-definite matrix, diagonalization of the \((b_{ii'})\) matrix, 
\[ b_{ii'} = b_k w_{ik} w_{i'k}^*, b_k \geq 0, \forall k, \]
gives rise to the Kraus operators for the qubit 1:
\[ K_k^1 = \sum_i \sqrt{b_k} w_{ik} A_{1i}, \]  
where the unitary matrix \((w_{ik})\) diagonalizes the \((b_{ii'})\) matrix. As emphasized above, this procedure, and complete positivity of the qubit’s dynamics, breaks for any kind of the initial correlations in the 1 + 2 system [3, 13]. To this end, a more detailed analysis with an emphasis on the subtleties regarding the very concept of complete positivity [3, 13, 20] will be presented elsewhere.

An important extension of the standard procedure for Markovian dynamics is considered for a pair of weakly interacting damped harmonic oscillators [3, 13]; in our considerations, this is the \(\beta \ll 1\) case. Then the interaction term \(H_{12}\), eq.(2), may be regarded as a perturbation and an alternative interaction picture can be used by omitting \(H_{12}\) in eq.(5). Then \(H_{12}\) appears in the commutator term of the master equation with the Lindblad operators, which are obtained for the case \(\beta = 0\), cf. eq.(B.1) in Ref. [13]. Nevertheless, this is just a special case of our considerations. Due to the commutation \([H_{10} + H_{20}, H_{12}] = 0\), the Lindblad operators eq.(11) are the same for both \(\beta \neq 0\) and \(\beta = 0\). Of interest is the interaction picture state defined as \(\tilde{\rho}_o(t) = U_o^\dagger \rho(t) U_o\), where \(U_o = \exp(-it(H_{10} + H_{20} + H_{E_1})); \rho(t)\) is in the Schrödinger-picture. The exact interaction picture state \(\tilde{\rho}(t) = U^\dagger \rho(t) U\), where \(U(t) = \exp(-it(H_{10} + H_{20} + H_{12} + H_{E_1})))\). Keeping the terms of the first order in (small) \(\beta\), the approximation \(U \approx U_o (I - it H_{12})\) easily gives:
\[ \tilde{\rho}_o (t) \approx \tilde{\rho}(t) - it[H_{12}, \tilde{\rho}_o(t)]. \]  
(40)

Taking the time derivative of eq.(40) easily follows, in the new interaction picture:
\[ \frac{d\tilde{\rho}_o}{dt} \approx -i[H_{1E_1}, \tilde{\rho}(t)] - i[H_{12}, \tilde{\rho}_o(t)] \approx -i[H_{1E_1} + H_{12}, \tilde{\rho}_o(t)], \]  
(41)
which is eq.(B.1) in Ref. [13]; \(H_{12}\) is of the same form for both pictures. We also note that, while eq.(40) is universal (for sufficiently small \(\beta\)), in general, the Lindblad operators for the two interaction pictures are not identical; nevertheless, even in the more general cases, our approach (approximation of
the above unitary operator $U(t)$, in the zeroth order of the time-independent perturbation, leads to eq. (B.1) of Ref. [13].

5. Conclusion

Derivation of the Kraus operator-sum from the microscopic Hamiltonian model may be technically involved. This may be the reason behind the lack of the explicit forms of the Kraus operators for most of the non-single-qubit processes. Nevertheless, the use of the Kraus form of the dynamical map is often technically simple. Our results aim at reducing this gap and exhibiting the technical advantage of the use of a Kraus form for certain basic tasks in open systems and quantum information theory contexts.

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