

Markovian approximation for correlated initial states

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$$H_{3} = H_{5} + H_{R} + A I$$

Coupling parameter

activy on Hilbert space $\mathcal{H}_{SR} = \mathcal{H}_S \otimes \mathcal{H}_R$ Field of oscillators finite-dimensional Schwödinger dynamics $P_{SR}^{t} = e^{-itH_{A}} P_{SR} e^{itH_{A}}$ $r_{initial state}$ Reduced system state $P_{S}^{t} = H_{R} P_{SR}^{t}$ $r_{partial tace}$

Two kinds of Mihial states:

$$S_{SR} = S_S \otimes S_R$$
 mcorrelated
 S_{SR} not of moduct form correlated

Born approximation
$$S_{SR}^{t} \approx S_{SR}^{t} \approx S_{SR}^{t} \otimes S_{R}^{t}$$

(S_{R}^{t} stationary, $S \ll R \& A small$)

Markov approximation (R Coses memory quickly enough)

Master equation (Dynamical map for system)

$$P_{s}^{t} \approx e^{tz} P_{s}$$

 $P_{s} \mapsto V_{t} P_{s} = e^{tz} P_{s}$

 $\tau = \lambda^2 t$ fixed, $\lambda \rightarrow 0$, $t \rightarrow \infty$:

$$e^{\tau \mathcal{L}} \mathcal{P}_{S} = \lim_{A \to 0} \operatorname{tr}_{R} \begin{bmatrix} -i \overline{\lambda}_{2} H_{A} & i \overline{\lambda}_{2} H_{O} \\ e^{-i \overline{\lambda}_{2} H_{A}} & e^{i \overline{\lambda}_{2} H_{O}} & (\mathcal{P}_{S} \otimes \mathcal{P}_{R}) & e^{-i \overline{\lambda}_{2} H_{O}} & e^{i \overline{\lambda}_{2} H_{A}} \end{bmatrix}$$

interaction picture dynamics, rescaled

$$\mathcal{L} = \mathcal{L}(a) = -i \left[H_{S}, \cdot \right] + a^{2} \mathcal{K}$$

generates CPTP semigronp



SSR \$ IS OFR · Born approximation X

• Agnamical map X 55t



Carnot define dynamical map V_t but can ship analyze structure of reduced S dynamics

PHYSICAL REVIEW A 67, 062109 (2003)

Kraus representation in the presence of initial correlations

Hiroyuki Hayashi,* Gen Kimura,[†] and Yukihiro Ota[‡]

PHYSICAL REVIEW A **100**, 042120 (2019)

Dynamics of initially correlated open quantum systems: Theory and applications

Gerardo A. Paz-Silva,¹ Michael J. W. Hall,^{1,2} and Howard M. Wiseman¹

Works similar in spirit to ours





Annals of Physics 322 (2007) 631-656

ANNALS of PHYSICS www.elsevier.com/locate/aop





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Annals of Physics 322 (2007) 657-676

On the assumption of initial factorization in the master equation for weakly coupled systems 1: General framework

S. Tasaki ^a, K. Yuasa ^{b,*}, P. Facchi ^{c,d}, G. Kimura ^{b,1}, H. Nakazato ^b, I. Ohba ^b, S. Pascazio ^{e,d} On the assumption of initial factorization in the master equation for weakly coupled systems II: Solvable models

K. Yuasa ^{a,*}, S. Tasaki ^b, P. Facchi ^{c,d}, G. Kimura ^{a,1}, H. Nakazato ^a, I. Ohba ^a, S. Pascazio ^{e,d}

$$\frac{\mathrm{d}}{\mathrm{d}\tau}\rho_{\mathrm{I}}(\tau) = \mathcal{K}\rho_{\mathrm{I}}(\tau), \qquad \rho_{\mathrm{I}}(0) = \mathcal{P}\rho_{0} = \mathrm{tr}_{\mathrm{B}}\{\rho_{0}\} \otimes \Omega_{\mathrm{B}}.$$
(4.4)

That is, even if the initial state ρ_0 is not in a factorized form, but rather there is entanglement, or simply a classical correlation, between system S and reservoir B, all correlations disappear in van Hove's limit and system S behaves as if the total system started from the factorized initial state in (4.4) with a reservoir state Ω_B specified below.

PHYSICAL REVIEW X 10, 041024 (2020)

Correlation-Picture Approach to Open-Quantum-System Dynamics S. Alipour[®],^{1,*} A. T. Rezakhani[®],^{2,†} A. P. Babu[®],¹ K. Mølmer,³ M. Möttönen[®],⁴ and T. Ala-Nissila^{1,5,‡} RindBlad-type NE mith jump operators depending on imitial correlation

Our main results : Perfurbation theory small 2, all times t >0.

- Inifial conclutions decay and then Markov approximation becomes valid for SR dynamics.
- Markovian approximation for reduced S dynamics is valid for all times.



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Quantum mocen tomography



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Witness for initial system-environment correlations in open-system dynamics

E.-M. LAINE^{1(a)}, J. PIILO^{1(b)} and H.-P. BREUER^{2(c)}

Initial SQ correlation can increase distripuishability of system stades.

$$\frac{4 \text{Incompled Hamiltonians}}{H_{S} = \sum_{j=1}^{N} E_{j} | \phi_{j} \rangle \langle \phi_{j} | , \qquad H_{R} = \int_{R^{3}} \omega(R) a^{*}(R) a(R) d^{3}R.$$

$$H_{S} = \int_{j=1}^{N} E_{j} | \phi_{j} \rangle \langle \phi_{j} | , \qquad H_{R} = \int_{R^{3}} \omega(R) a^{*}(R) a(R) d^{3}R.$$

$$|R|$$

$$\frac{3 \text{Incomplex for operator}}{I = G \otimes \varphi(g)}$$

$$any \text{ hermitian matrix}}$$

$$\varphi(g) = \int_{R^{3}} \left[a^{*}(g) + a(g) \right]$$

$$a^{*}(g) = \int_{R^{3}} g(R) \in L^{2}(R^{3}, d^{3}R)$$

$$\int_{T} g(R) \in L^{2}(R^{3}, d^{3}R)$$

Equilibrium states

S: Gibbs density matrix $P_{s,p} = \frac{e^{-pH_s}}{tr e^{-pH_s}}$

R: Infinitely extended (confinuous modes) KMS state WR, B Two-point function

$$\omega_{R_{i}\beta}\left(a^{*}(k)a(\ell)\right) = \frac{\delta(k-\ell)}{e^{\beta\omega(k)}-1}$$

Initial states

$$\int_{SA} (\cdot) = \Lambda \left(\int_{S}^{\infty} \otimes W_{R_{1}\beta} \right) = \sum_{\alpha} \int_{S}^{\infty} \otimes W_{R_{1}\beta} \left(K_{\alpha}^{*} \cdot K_{\alpha} \right)$$

$$\sum_{\alpha} K_{\alpha}^{*} K_{\alpha} = 1 \qquad \text{Krawo ops}$$

$$K_{\alpha} \in C \quad \text{`correlator algebra'}$$

Observable and correlator algebras

Test from chien spaces

 $L^{2}_{obs} : f \in L^{2}(\mathbb{R}^{3}, d^{3}\mathbb{R}) \quad 3 \text{ times differentiable w.r.t. } |\mathbb{R}| > 0 \quad \&$ $\cdot |\mathbb{R}| \sim 0 : \quad \mathbb{F}(\mathbb{R}) \sim |\mathbb{R}|^{P} \qquad p = -\frac{1}{2}, \frac{1}{2} \quad \text{or} \quad p > 1$ $\cdot |\mathbb{R}| \rightarrow \infty : \quad \mathbb{F}(\mathbb{R}) \sim |\mathbb{R}|^{-q} \qquad q > 3$

 L_{cor}^2 : $f \in L_{obs}^2$ and also $e^{\beta |\mathbf{k}|} f(\mathbf{k}) \in L^2(\mathbf{R}^3, d^3\mathbf{k})$

$$\int_{SR} = \frac{1}{2} e^{\mathcal{E}} \left(\int_{S}^{\infty} \omega_{R,\beta} \right) e^{\mathcal{E}^{*}}$$

$$\mathcal{E} = \sum_{j} B_{j} \otimes a^{*}(f_{j}) + \sum_{\ell} D_{\ell} \otimes a(h_{\ell})$$

Assumption on effectiveness of SR coupling.

Fermi Golden Anle Condition

spectral density. of reservoir $\langle \phi_{m}, G, \phi_{n} \rangle \Im (|E_{m}-E_{n}|) \neq 0 \quad \forall E_{m} \neq E_{n}$

Second order fauntion processes are not expressed

Main result

Theorem 3.4 (SR dynamics for correlated initial states) There is a constant $\lambda_0 > 0$ such that if $|\lambda| < \lambda_0$, then the following holds true. Let $\rho_{SR} = \Lambda(\rho_S \otimes \omega_{R,\beta})$ and let ρ_S be its reduction to S. Then for all $t \ge 0$, $O \in \mathcal{O}$,

$$\rho_{\mathrm{SR}}\left(e^{\mathrm{i}tH_{\lambda}}Oe^{-\mathrm{i}tH_{\lambda}}\right) = \left(e^{t\mathcal{L}_{\mathrm{S}}(\lambda)}\rho_{\mathrm{S}}\otimes\omega_{\mathrm{R},\beta}\right)(O) + \chi(\lambda,t,O) + R(\lambda,t,O),$$

where $\mathcal{L}_{S}(\lambda)$ is the Davies generator and

$$|R(\lambda, t, O)| \le C(O) |\lambda|^{1/4}.$$

The correlation term χ satisfies

$$\begin{split} \chi(\lambda, t, O) &= 0 \quad if \quad \rho_{\mathrm{SR}} = \rho_{\mathrm{S}} \otimes \omega_{\mathrm{R},\beta} \\ \chi(\lambda, t, O_{\mathrm{S}} \otimes \mathbb{1}_{\mathrm{R}}) &= 0 \\ |\chi(\lambda, t, O)| &\leq C(O) \, (1 + t^2)^{-3/2}. \end{split}$$

The constants C(O) depend on O as well as the initial state ρ_{SR} (but are independent of t, λ). For $O = O_S \otimes \mathbb{1}_R$ we have $C(O_S \otimes \mathbb{1}_R) \leq c \|O_S\|$ with c independent of O_S .

Consequences



2

The system dynamics is Markorian at all times:

$$t_{R}(p_{SR}^{t}) = e^{tZ}(t_{SR}^{o}) + O(|a|'_{4})$$

Ontline of proof

$$\begin{split} & \omega_{0} \equiv & \int_{S_{1}\beta} \otimes \omega_{R_{1}\beta} \\ & \alpha_{\chi}^{t}(\cdot) \equiv & e^{it\#_{\lambda}} \cdot e^{-it\#_{\lambda}} \end{split}$$

mcoupled equilibrium state

Heitenberg, dynamics

KMS condition $\omega(A \propto_{o}^{i\beta}(B)) = \omega_{o}(BA)$

ω

$$\omega_{0} \left(\begin{array}{c} \mathsf{K}^{*} & \alpha_{\lambda}^{t} \left(0 \right) \\ \mathsf{K}^{*} & \alpha_{\lambda}^{t} \left(0 \right) \\ \end{array} \right) \\ = & \omega_{0} \left(\begin{array}{c} \alpha_{0}^{-i\beta}(\mathsf{K}) \\ \alpha_{0}^{*} \\ \end{array} \right) \\ = & \omega_{0} \left(\begin{array}{c} \mathcal{C} \\ \alpha_{\lambda}^{t} \\ \end{array} \right) \\ \end{array} \right)$$

GNS representation

$$= \langle \mathcal{L}_{\circ}, \pi(\mathcal{C}) \pi(a_{\mathcal{A}}^{\ell}(\circ)) \mathcal{L}_{\circ} \rangle$$

Liouville operator

=
$$\langle \Omega_0, \pi(\mathcal{C}) e^{itL_A} \pi(0) e^{-itL_A} \Omega_0 \rangle$$

Perturbation Heavy of KMS states: $\| \mathcal{D}_{o} - \mathcal{D}_{a} \| \leq C |a|$ Invariance $L_{a} - \mathcal{D}_{a} = 0$

=
$$(\Pi_0, \Pi(\mathcal{C}) e^{itL_{\mathcal{A}}} \Pi(0) \Pi_0 + O(1a1)$$

indep. of t
Use resonance expansion [KM 2017, M 2021]

$$e^{itL_{A}} = e^{itM(A)} \otimes P_{A} + P_{A}^{L} e^{itL_{A}^{L}} P_{A}^{L} + O(|A|^{1/4})$$
weakly on dense set of rectors
$$P_{A} = 1 \Omega_{0} \times \Omega_{0}|$$

$$p_{0} = 0 \times \Omega_{0}|$$

$$P_{0} = \Omega_{0} \otimes \Omega_{0}$$

$$P_{0} = \Omega_{0} \otimes \Omega_{0}$$

$$P_{0} = \Omega_{0} \otimes \Omega_{0}$$

$$P_{0} = (\Omega_{0}, \pi(C)) \times (e^{itM(A)} \otimes P_{A}) \pi(0) \Omega_{0} + \chi(A_{0}t_{0})$$

$$+ O(|A|^{1/4})$$

Term along P_2:

$$(e^{itM(a)} \otimes P_{\Omega}) \pi (O_{S} \otimes O_{R}) \Lambda_{o}$$

$$= \omega_{R_{1}\beta} (O_{R}) \cdot (e^{itM(a)} \otimes 4_{R}) \pi (O_{S} \otimes 4_{R}) \Lambda_{o}$$

$$= \omega_{R_{1}\beta} (O_{R}) \cdot \pi ((e^{t\mathcal{L}^{*}}O_{S}) \otimes 4_{R}) \Lambda_{o}$$

$$Definer \mathscr{L}^{*}$$

$$\begin{split} & \left(\begin{array}{c} \Omega_{0}, \pi \left(\varepsilon \right) \left(\varepsilon^{it} H(\varepsilon) \right) \left(\varepsilon^{i} \Omega_{0} \right) \pi \left(\varepsilon^{0} \Omega_{0} \right) \pi \left(\varepsilon^{0} \left(\varepsilon^{0} \right) \right) \right) \\ & = & \left(\left(\varepsilon^{0} \Omega_{0} \right) \left(\left(\varepsilon^{0} \Omega_{0} \right) \right) \pi \left(\varepsilon^{0} \left(\varepsilon^{0} \Omega_{0} \right) \right) \pi \left(\varepsilon^{0} \left(\varepsilon^{0} \Omega_{0} \right) \right) \\ & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \pi \left(\varepsilon^{0} \left(\varepsilon^{0} \Omega_{0} \right) \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \left(\varepsilon^{0} \Omega_{0} \right) \right) \\ & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \left(\varepsilon^{0} \Omega_{0} \right) \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \left(\varepsilon^{0} \Omega_{0} \right) \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \left(\varepsilon^{0} \Omega_{0} \right) \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \left(\varepsilon^{0} \Omega_{0} \right) \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \left(\varepsilon^{0} \Omega_{0} \right) \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \left(\varepsilon^{0} \Omega_{0} \right) \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \\ & = & \left(\varepsilon^{0} \Omega_{0} \right) \left(\varepsilon^{0} \Omega_{0} \right) \\ & = & \left(\varepsilon^{0$$

Thank you 🙏