10. Zwanzig-Mori Formalism

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Zwanzig-Mori formalism

Beginnings:
Two phenomenological approaches for the dynamics of systems close to or at thermal equilibrium:

▷ Phenomenological equations of motion for probability distributions (e.g. master equation, Fokker-Planck equation).
▷ Phenomenological equations of motion for dynamical variables (e.g. Langevin equation)

In these approaches, the focus is on selected degrees of freedom. All other degrees of freedom are taken into account summarily in the form of ad-hoc randomness.

Completions:
Microscopic foundations for these phenomenological approaches.

▷ Zwanzig (1960): Rigorous derivation of a generalized master equation from first principles, i.e. from the Liouville equation.
▷ Mori (1965): Rigorous derivation of a generalized Langevin equation from first principles, i.e. from the (quantum) Heisenberg equation or the (classical) canonical equations.

The focus is again on selected degrees of freedom but here the effects of the other degrees of freedom are taken into account on a basis that is exact and amenable to systematic approximation.

Variants:

▷ Zwanzig’s approach leads to a kinetic equation of a particular kind. There exist alternative ways to derive kinetic equations from the Liouville equations via systematic approximations (e.g. via BBGKY hierarchy).
▷ Mori’s approach has been formulated in more than one rendition. The version named projection operator formalism is most illuminating regarding the physical meaning of systematic approximations. The version named recursion method is most readily amenable to computational applications.
Time Dependence of Expectation Values

Quantum Hamiltonian system:

\[
\langle A \rangle_t = \text{Tr} \{ \rho(0) A(t) \} \quad \text{(Heisenberg representation)}
\]
\[
= \text{Tr} \{ \rho(0) e^{iHt/\hbar} A(0) e^{-iHt/\hbar} \} = \text{Tr} \{ e^{-iHt/\hbar} \rho(0) e^{iHt/\hbar} A(0) \}
\]
\[
= \text{Tr} \{ \rho(t) A(0) \} \quad \text{(Schrödinger representation)}
\]

- \[
\frac{dA}{dt} = i \hat{L} A = \frac{i}{\hbar} [\hat{H}, A] \quad \text{(Heisenberg equation)}
\]
- \[
\frac{d\rho}{dt} = -i \hat{L} \rho = \frac{1}{i\hbar} [\hat{H}, \rho] \quad \text{(quantum Liouville equation)}
\]
- \[
\hat{L} \equiv \frac{1}{\hbar} [\hat{H}, \ ] \quad \text{(quantum Liouville operator)}
\]
- \[
\{,\} \quad \text{(commutator)}
\]

Classical Hamiltonian system:

\[
\langle A \rangle_t = \int d^n q \int d^n p \rho(q_1, \ldots, q_n; p_1, \ldots, p_n; 0) A(q_1, \ldots, q_n; p_1, \ldots, p_n; t)
\]
\[
= \int d^n q \int d^n p \rho(q_1, \ldots, q_n; p_1, \ldots, p_n; t) A(q_1, \ldots, q_n; p_1, \ldots, p_n; 0)
\]

- \[
\frac{dA}{dt} = i \hat{L} A = -\{\hat{H}, A\} \quad \text{(Hamilton equation)}
\]
- \[
\frac{d\rho}{dt} = -i \hat{L} \rho = \{\hat{H}, A\} \quad \text{(classical Liouville equation)}
\]
- \[
\hat{L} \equiv \{\hat{H},\} = i \sum_{j=1}^n \left( \frac{\partial \hat{H}}{\partial q_j} \frac{\partial}{\partial p_j} - \frac{\partial \hat{H}}{\partial p_j} \frac{\partial}{\partial q_j} \right) \quad \text{(classical Liouville operator)}
\]
- \[
\{,\} \quad \text{(Poisson bracket)}
\]
Zwanzig’s method [nl/29]

Consider classical phase-space density \( \rho(t) = \rho(q_1, p_1; q_2, p_2; \ldots; q_n, p_n; t) \).

Distinguish system \((q_1, p_1)\) and heat bath \((q_2, p_2; \ldots; q_n, p_n)\).

Probability density of system: \( \rho_1(t) = \hat{P}\rho(t) \) via projection.

Probability density of heat bath: \( \rho_2(t) = \hat{Q}\rho(t) \), where \( \hat{Q} = 1 - \hat{P} \).

Implementation of projection:
\[
\rho_1(t) = \hat{P}\rho(q_1, p_1; q_2, p_2; \ldots; q_n, p_n; t) = \rho_{eq}(q_2, p_2; \ldots; q_n, p_n)\sigma(q_1, p_1; t),
\]
where \( \partial\rho_{eq}/\partial t = -\hat{L}\rho_{eq} = 0 \) and
\[
\sigma(q_1, p_1; t) = \int dq_2 dp_2 \cdots dq_n dp_n \rho(q_1, p_1; q_2, p_2; \ldots; q_n, p_n; t).
\]

Idempotency, \( \hat{P}\rho_1(t) = \rho_1(t) \), satisfied by construction.

**Liouville equation**, \( \partial \rho(t)/\partial t = -\hat{L}\rho(t) \), split into two coupled equations:
\[
\hat{P}\frac{\partial}{\partial t}\rho(t) = \frac{\partial}{\partial t}\rho_1(t) = -i\hat{P}\hat{L}[\rho_1(t) + \rho_2(t)], \quad (1)
\]
\[
\hat{Q}\frac{\partial}{\partial t}\rho(t) = \frac{\partial}{\partial t}\rho_2(t) = -i\hat{Q}\hat{L}[\rho_1(t) + \rho_2(t)]. \quad (2)
\]

Formal solution of (2) [nex68] to be substituted into (1):
\[
\rho_2(t) = e^{-i\hat{Q}\hat{L}t}\rho_2(0) - i \int_0^t d\tau e^{-i\hat{Q}\hat{L}\tau}\hat{L}\rho_1(t - \tau).
\]

**Zwanzig’s kinetic equation** (generalized master equation):
\[
\frac{\partial}{\partial t}\rho_1(t) = -i\hat{P}\hat{L}\rho_1(t) - i\hat{P}\hat{L}e^{-i\hat{Q}\hat{L}t}\rho_2(0) - \int_0^t d\tau \hat{P}\hat{L}e^{-i\hat{Q}\hat{L}\tau}\hat{L}\rho_1(t - \tau).
\]

#1 Autonomous part of system’s time evolution.
#2 Autonomous part of heat bath’s time evolution. Instantaneous effect of heat bath on system at time \( t \).
#3 Effect caused by system on heat bath at time \( t - \tau \) propagates in heat bath and feeds back into system at time \( t \).

First term often vanishes and second term (inhomogeneity) can be made zero by judicious choice of initial conditions.

Zwanzig’s kinetic equation can be used as the starting point for the derivation (via approximations) of a master equation or a Fokker-Planck equation.
Zwanzig’s kinetic equation.

In the derivation of Zwanzig’s kinetic equation,

\[ \frac{\partial}{\partial t} \rho_1(t) = -i \hat{P} \hat{L} \rho_1(t) - i \hat{P} \hat{L} e^{-iQLt} \rho_2(0) - \int_0^t d\tau \hat{P} \hat{L} e^{-iQ\tau} \hat{Q} \hat{L} \rho_1(t - \tau), \]  

(1)

from two projections of the Liouville equation,

\[ \hat{P} \frac{\partial \rho}{\partial t} = \frac{\partial \rho_1}{\partial t} = -i \hat{P} \hat{L} [\rho_1 + \rho_2], \quad \hat{Q} \frac{\partial \rho}{\partial t} = \frac{\partial \rho_2}{\partial t} = -i \hat{Q} \hat{L} [\rho_1 + \rho_2], \]  

(2)

we postulate the formal solution

\[ \rho_2(t) = e^{-iQLt} \rho_2(0) - i \int_0^t d\tau e^{-iQ\tau} \hat{Q} \hat{L} \rho_1(t - \tau). \]  

(3)

Verify that (3) is a solution of (2). The derivation involves one integration by parts.

Solution:
**Projection operator method** [nl31]

**Goal:** Determination of symmetrized time-correlation function (fluctuation function) for a dynamical variable $A(t)$ of a quantum or classical many-body Hamiltonian system $\mathcal{H}$ in thermal equilibrium.

Fluctuation function (real, symmetric, normalized):

$$C_0(t) \equiv \frac{\langle A(t) | A \rangle}{\langle A | A \rangle} = \frac{\langle A | A(-t) \rangle}{\langle A | A \rangle} = \frac{\langle A | e^{-\imath L t} | A \rangle}{\langle A | A \rangle}.$$

Dirac notation symbolizes inner product of choice as explained in [nl32]. Some properties of dynamic quantities depend on choice of inner product.

Relaxation function (via Laplace transform):\(^1\)

$$c_0(z) = \int_0^\infty dt \ e^{-zt} \frac{\langle A | e^{-\imath L t} | A \rangle}{\langle A | A \rangle} = \frac{1}{\langle A | A \rangle} \left\langle A \left| \frac{1}{z + \imath \epsilon} \right| A \right\rangle.$$

Projection operator method determines relaxation function via systematic approximation.

Inverse Laplace transform,

$$C_0(t) = \frac{1}{2\pi \imath} \int_{\cal c} dz \ e^{zt} c_0(z),$$

involves integral along straight path from $\epsilon - \imath \infty$ to $\epsilon + \imath \infty$ for $\epsilon > 0$.

In practical applications, the (real, symmetric) spectral density is inferred from the relaxation function as limit process,

$$\Phi_0(\omega) = 2 \lim_{\epsilon \to 0} \Re \{ c_0(\epsilon - \imath \omega) \},$$

and the fluctuation function via inverse Fourier transform,

$$C_0(t) = \int_{-\infty}^{+\infty} \frac{d\omega}{2\pi} e^{-\imath \omega t} \Phi_0(\omega).$$

\(^1\)The last bracket is also known as a Green’s function.
Kubo inner product

General properties of inner products:
\[ \langle A|B \rangle = \langle B|A \rangle^*, \quad \langle A|\lambda B \rangle = \lambda \langle A|B \rangle, \]
\[ \langle A|A \rangle = ||A||^2 \geq 0, \quad \langle A|B+C \rangle = \langle A|B \rangle + \langle A|C \rangle. \]

Kubo inner product for quantum system:
\[ \langle A|B \rangle \overset{1}{=} \frac{1}{\beta} \int_0^\beta d\lambda \langle e^{\lambda H}A^\dagger e^{-\lambda H}B \rangle, \]
where
\[ \langle A \rangle = \frac{1}{Z} \text{Tr} \{ e^{-\beta H}A \}, \quad Z = \text{Tr} \{ e^{-\beta H} \}, \quad \beta = \frac{1}{k_B T}. \]

Alternative inner product for quantum systems:
\[ \langle A|B \rangle \overset{2}{=} \frac{1}{2} \langle A^\dagger B + BA^\dagger \rangle. \]

Both inner products have the same classical limit:
\[ \langle A|B \rangle \overset{3}{=} \frac{1}{Z} \int d^nq dp e^{-\beta H(q,p)} A(q, p) B(q, p). \]

Inner products of [nln31] employ...

\[ \triangleright \text{quantum Liouville operator: } L = \frac{1}{\hbar}[L, ], \]
Heisenberg equation of motion:
\[ \frac{dA}{dt} = \frac{i}{\hbar} \{ H, A \} = iLA. \]

\[ \triangleright \text{classical Liouville operator: } L = i\{ H, \} = i \sum_{j=1}^n \left( \frac{\partial H}{\partial q_j} \frac{\partial}{\partial p_j} - \frac{\partial H}{\partial p_j} \frac{\partial}{\partial q_j} \right), \]
Hamilton’s equation of motion:
\[ \frac{dA}{dt} = -\{ H, A \} = iLA. \]

\[ ^1\text{Designed to satisfy classical fluctuation-dissipation theorem in [nln39].} \]
\[ ^2\text{Designed to satisfy quantum fluctuation-dissipation theorem in [nln39].} \]
\[ ^3\text{Option for all inner products: subtract } \langle A^\dagger \rangle \langle B \rangle. \]
Projection operators

The relaxation function \( c_0(z) \) is determined recursively by a succession of subdivisions of the many-body dynamics into components that are treated rigorously and a remainder that is treated phenomenologically. It is expected that the remainder diminishes in importance as the number of rigorous components is increased systematically.

The time evolution of the dynamical variable \( A(t) \) can be conceived as a “pirouette” performed by the vector \( |A(t)\rangle \) through the Hilbert space.

The subdivisions are implemented by a sequence of projections onto one-dimensional Hilbert subspaces traversed by \( |A(t)\rangle \).

Initial condition: \( |f_0\rangle = |A(0)\rangle = |A\rangle \).
Projection operators \( P_n \) and \( Q_n = 1 - P_n \), \( n = 0, 1, \ldots \),

\[
P_0 = |f_0\rangle \langle f_0|, \quad P_0^2 = P_0, \quad (P_0)^\dagger = P_0, \quad P_0Q_0 = Q_0P_0 = 0.
\]
Orthogonal direction:\(^1\)

\[
|f_1\rangle = iL|f_0\rangle, \quad \langle f_0|f_1\rangle = 0, \quad P_0|f_1\rangle = 0, \quad Q_0|f_1\rangle = |f_1\rangle - P_0|f_1\rangle = |f_1\rangle,
\]

\[
P_1 = |f_1\rangle \langle f_1|, \quad Q_1 = 1 - P_1.
\]
The systematic generation of further orthogonal direction will be discussed in the context of the recursion method.

Successive projections filter out particular aspects of the many-body dynamics to be taken into account rigorously. The filters are applied in series. What passes through \( n \) filters is the remainder to be treated phenomenologically.

The physical content of this process can be gleaned from the first two projections carried out in detail:

- First projection [nlh33],
- Second projection [nlh34].

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\(^1\)Unitary transformation \( e^{iLt} \) makes \( iLA \) orthogonal to \( A \), implying \( \langle A|iLA\rangle = 0 \).
First projection \[\text{[nln34]}\]

Rewrite relaxation function from \[\text{[nln31]}\] with projection operators from \[\text{[nln33]}\] and apply Dyson identity \((X+Y)^{-1} = X^{-1} - X^{-1}Y(X+Y)^{-1}\).\(^1\)

\[
(f_0|f_0)c_0(z) = \frac{1}{z + iL} \left[ 1 + \left( \frac{-i}{z} \right) LQ_0 + \frac{(-i)^2}{z^2} LQ_0LQ_0 + \cdots \right] f_0 = \frac{1}{z} (f_0|f_0).
\]

\[
\left( f_0 \frac{1}{z + iLQ_0} iL \right) f_0 = \left( f_0 \left[ 1 - (z+iLQ_0) \frac{1}{z + iLQ_0} + \frac{z}{z + iLQ_0} \right] iL \right) f_0 = \left( f_0 \left[ (-i)Q_0 - iQ_0LQ_0Q_0L \right] \right) f_0 = \left( f_1 \frac{1}{z + iL_1} \right) f_1: \quad |f_1\rangle = Q_0|f_0\rangle = Q_0L|f_0\rangle, \quad L_1 = Q_0LQ_0.
\]

Relaxation function after first projection expressed via memory function:

\[
c_0(z) = \frac{1}{z + \Sigma_1(z)}, \quad \Sigma_1(z) = \frac{1}{(f_0|f_0)} \left( f_1 \left[ 1 - \frac{1}{z + iL_1} \right] f_1 \right).
\]

Memory function \(\Sigma_1(z)\) of original problem, \(\{L, |f_0\rangle\}\), can be reinterpreted as the (as yet non-normalized) relaxation function of a new dynamical problem, \(\{L_1, |f_1\rangle\}\).

Projection operator \(Q_0\) acts as filter on the Liouvillian \(L\), absorbing that part of dynamics dealt with explicitly in first projection. Explicit information contained in normalization constant of \(\Sigma_1(z)\).

\(^1\text{Direct consequence of operator identity } (X+Y)(X+Y)^{-1} = X(X+Y)^{-1}+Y(X+Y)^{-1} = 1, \text{ here with } X = z + iLQ_0, \quad Y = iLP_0.\)
Second projection \[nln35\]

Rewrite memory function from \[nln34\] with projection operators from \[nln33\] and apply Dyson identity:

\[
\langle f_0|f_0 \rangle \Sigma_1(z) = \left\langle f_1 \left| \frac{1}{z+iL_1} \right| f_1 \right\rangle = \left\langle f_1 \left| \frac{1}{z+iL_1P_1+iL_1Q_1} \right| f_1 \right\rangle \\
= \left\langle f_1 \left| \frac{1}{z+iL_1Q_1} \right| f_1 \right\rangle - \left\langle f_1 \left| \frac{1}{z+iL_1Q_1} \right| f_1 \right\rangle \frac{\langle f_0|f_0 \rangle}{\langle f_1|f_1 \rangle} \Sigma_1(z),
\]

where simplifications analogous to \[nln34\] are carried out.

\[
\Rightarrow \Sigma_1(z) = \frac{\langle f_1|f_1 \rangle / \langle f_0|f_0 \rangle}{z + \langle f_1|f_1 \rangle \left\langle f_1 \left| \frac{z}{z+iL_1Q_1} \right| f_1 \right\rangle},
\]

\[
\left\langle f_1 \left| \frac{z}{z+iL_1Q_1} \right| f_1 \right\rangle = \cdots = \left\langle f_1 \left| \frac{(-i)L_1Q_1}{z+iQ_1L_1Q_1} \right| f_1 \right\rangle \\
= \left\langle f_2 \left| \frac{1}{z+iL_2} \right| f_2 \right\rangle,
\]

with \(|f_2\rangle = Q_1iL_1|f_1\rangle\), \(L_2 = Q_1L_1Q_1\).

Memory function (first termination function) after second projection expressed via second termination function:

\[
\Sigma_1(z) = \frac{\Delta_1}{z + \Sigma_2(z)}, \quad \Sigma_2(z) = \frac{1}{\langle f_1|f_1 \rangle} \left\langle f_2 \left| \frac{1}{z+iL_2} \right| f_2 \right\rangle
\]

with continued-fraction coefficients \(\Delta_1 = \langle f_1|f_1 \rangle / \langle f_0|f_0 \rangle\).

The \(n^{th}\) projection yields

\[
\Sigma_{n-1}(z) = \frac{\Delta_{n-1}}{z + \Sigma_n(z)}, \quad \Sigma_n(z) = \frac{1}{\langle f_{n-1}|f_{n-1} \rangle} \left\langle f_n \left| \frac{1}{z+iL_n} \right| f_n \right\rangle
\]

with \(\Delta_{n-1} = \langle f_{n-1}|f_{n-1} \rangle / \langle f_{n-2}|f_{n-2} \rangle\)

and \(|f_n\rangle = Q_{n-1}iL_{n-1}|f_{n-1}\rangle\), \(L_n = Q_{n-1}L_{n-1}Q_{n-1}\).
Continued-fraction representation

Relaxation function after \( n \) successive projections:
\[
c_0(z) = \frac{1}{z + \frac{\Delta_1}{z + \frac{\Delta_2}{z + \ldots + \frac{\Delta_{n-1}}{z + \Sigma_n(z)}}}}
\]

The explicit dynamical information extracted from the original many-body system, \( \{L, |f_0\rangle\} \), in the first \( n \) projections is contained in the continued-fraction coefficients:
\[\Delta_1, \ldots, \Delta_n.\]

Each projection adds a layer of projection operators around the original Liouvillian:
\[L_n = Q_{n-1} \cdots Q_0 L Q_0 \cdots Q_{n-1}.\]

Expectation (somewhat naively):

- If \( n \) is sufficiently large, all distinctive spectral features of \( L \) will have been filtered out and incorporated explicitly into the relaxation function via the \( \Delta_i \).
- Whatever features of \( L \) still shine through the \( n \) filters are adequately represented by a source of white noise.
- The memory function associated with white noise is a constant, commonly represented by a relaxation time:
  \[\Sigma_n(z) = \frac{1}{\tau_n} = \text{const.}\]
- This completion of the continued fraction is known under the name \( n \)-pole approximation. The relaxation function is characterized by \( n \) poles in the complex frequency plane.

Options used in practical applications:

- A number of continued-fraction coefficients are determined on phenomenological grounds along with a terminating relaxation time \( \tau_n \). Examples: classical relaxator (1 pole), classical oscillator (two poles).
- A number of continued-fraction coefficients are calculated from the original many-body system via the recursion method along with a termination function \( \Sigma_n(z) \) inferred from an extrapolation scheme.
**n-Pole Approximation** [nln87]

Relaxation function after \( n \) successive projections (Mori formalism [nln36]) or \( n \) iterations (recursion method [nln84]),

\[
c_0(z) = \frac{1}{z + \Delta_1},
\]

\[
c_0(z) = \frac{1}{z + \Delta_2},
\]

\[
c_0(z) = \frac{1}{z + \cdots}
\]

\[
c_0(z) = \frac{1}{z + \frac{\Delta_{n-1}}{1}},
\]

has termination function represented by a relaxation time \( \tau_n \):

\[
\Gamma_n(z) = \Delta_n \frac{c_n(z)}{c_{n-1}(z)} = \frac{1}{\tau_n}.
\]

Consequences:

- Singularity structure of \( c_0(z) \) reduced to \( n \) poles in complex \( z \)-plane.
- Frequency moments \( M_{2k} = \langle \omega^{2k} \rangle \) divergent for \( k \geq n \).

Limiting cases:

- \( \tau_n \to \infty \):
  All poles approach \( \omega \)-axis: \( z = \epsilon - i\omega \) with \( \epsilon \to 0 \).
  Spectral density \( \Phi_0(\omega) \) is a sum of \( \delta \)-functions.

- \( \tau_n \to 0 \):
  Transition to \( (n-1) \)-pole approximation: \( \frac{\Delta_{n-1}}{z + 1/\tau_n} \to \frac{1}{\tau_{n-1}} \).

Illustrations for \( n = 1, 2, 3 \):

- relaxation function: \( c_0(z) = \frac{1}{z + \Sigma(z)} \),
- memory function: \( \Sigma(z) \),
- spectral density: \( \Phi_0(\omega) = 2 \lim_{\epsilon \to 0} \text{Re}[c_0(\epsilon - i\omega)] \).
1-pole spectral approximation

Memory function representing white noise:

\[ \Sigma(z) = \frac{1}{\tau_1} = \text{const.} \]

Relaxation function:

\[ c_0(z) = \frac{1}{z + \frac{1}{\tau_1}} = \frac{\tau_1}{\tau_1 z + 1}. \]

Spectral density:

\[ \Phi_0(\omega) = \frac{2\tau_1}{1 + \tau_1^2\omega^2}. \]
2-pole spectral approximation

Memory function representing classical relaxator:

\[ \Sigma(z) = \frac{\Delta_1}{z + \frac{1}{\tau_1}}. \]

Relaxation function:

\[ c_0(z) = \frac{\tau_2 z + 1}{\tau_2 z^2 + z + \Delta_1 \tau_2}. \]

Spectral density:

\[ \Phi_0(\omega) = \frac{2 \Delta_1 \tau_2}{\tau_2^2 (\Delta_1 - \omega^2)^2 + \omega^2}. \]
3-pole spectral approximation

Memory function representing damped classical oscillator:

$$\Sigma(z) = \frac{\Delta_1}{z} + \frac{\Delta_2}{z + \frac{1}{\tau_3}}.$$  

Relaxation function:

$$c_0(z) = \frac{\tau_3 z^2 + z + \Delta_2 \tau_3}{\tau_3 z^3 + z^2 + (\Delta_1 + \Delta_2) \tau_3 z + \Delta_1}.$$  

Spectral density:

$$\Phi_0(\omega) = \frac{2\Delta_1 \Delta_2 \tau_3}{[\tau_3 \omega (\omega^2 - \Delta_1 - \Delta_2)]^2 + (\omega^2 - \Delta)^2}.$$
Relaxation function with uniform continued-fraction coefficients.

Find closed-form expressions for the relaxation function $c_0(z)$, the spectral density $\Phi_0(\omega)$, and the fluctuation function $C_0(t)$, of some physical system if we know that the (infinite) sequence of continued-fraction coefficients is

(a) uniform: $\Delta_1 = \Delta_2 = \ldots = \frac{1}{3} \omega_0^2$,
(b) almost uniform: $\Delta_1 = \frac{1}{3} \omega_0^2$, $\Delta_2 = \Delta_3 \ldots = \frac{1}{7} \omega_0^2$.

Solution:
Link to Green’s Function Formalism

Retarded Green’s function (two varieties):

\[ \tilde{G}_\pm(t - t') \doteq \langle A(t); B(t') \rangle_\pm = -i\theta(t - t') \langle [A(t), B(t')]_\pm \rangle. \] (1)

Equations of motion are hierarchical in nature:

\[ i\frac{\partial}{\partial t} \langle A(t); B(t') \rangle_\pm = \delta(t - t') \langle [A, B]_\pm \rangle + \langle [A(t), \mathcal{H}_-; B(t')]_\pm \rangle. \] (2)

Frequency-dependent Green’s functions via Fourier transform:

\[ G_\pm(\omega) \doteq \langle A; B \rangle_\pm = \int_{-\infty}^{+\infty} dt \ e^{i\omega t} \tilde{G}_\pm(t), \] (3)

\[ \omega \langle A; B \rangle_\omega = \langle [A, B]_\pm \rangle + \langle [A(t), \mathcal{H}_-; B(t')]_\pm \rangle. \] (4)

Spectral representations relate the Green’s functions \( G_\pm(\zeta) \) to the functions \( S(\omega), \Phi(\omega), \) and \( \chi''(\omega) \) as defined in [nln39]:

\[ G_\pm(\zeta) = \int_{-\infty}^{+\infty} d\tilde{\omega} \ \frac{S(\tilde{\omega})}{2\pi \zeta - \tilde{\omega}} [1 \pm e^{-\beta \tilde{\omega}}], \quad \zeta = \omega + i\epsilon, \quad \epsilon > 0, \] (5a)

\[ G_+(\zeta) = 2 \int_{-\infty}^{+\infty} d\tilde{\omega} \ \frac{\Phi(\tilde{\omega})}{2\pi \zeta - \tilde{\omega}}, \quad G_-(\zeta) = 2 \int_{-\infty}^{+\infty} d\tilde{\omega} \ \frac{\chi''(\tilde{\omega})}{2\pi \zeta - \tilde{\omega}}. \] (5b)

Inverse relations:

1. \( \Phi(\omega) = -\lim_{\epsilon \to 0} \text{Im} \left[ G_+(\omega + i\epsilon) \right] \) (spectral density)
2. \( \chi''(\omega) = -\lim_{\epsilon \to 0} \text{Im} \left[ G_-(\omega + i\epsilon) \right] \) (dissipation function)
3. \( S(\omega) = \frac{\Phi(\omega)}{2(1 + e^{-\beta \omega})} = \frac{\chi''(\omega)}{2(1 - e^{-\beta \omega})} \) (structure function)

Relation to relaxation function for \( i\zeta = -z \) (see [nln31] and [nln84]):

\[ c_0(z) = \int_0^{+\infty} dt \ e^{\epsilon t} \frac{\Phi(t)}{\Phi(0)} = \frac{1}{2\Phi(0)} \int_{-\infty}^{+\infty} d\omega \ \frac{\Phi(\omega)}{2\pi \zeta - \omega} = \frac{1}{2\Phi(0)} G_+(\zeta). \]

1Here we set \( \hbar = 1 \) as is common practice.
2The bracket \( [\ , \ ]_- \) stands for commutators and the bracket \( [\ , \ ]_+ \) for anticommutators.
3One variety coincides with Kubo’s response function from [nln27]: \( \tilde{G}_-(t) = -\tilde{\chi}_{AB}(t) \).
[nex71] Structure function of harmonic oscillator I.

Consider the quantum harmonic oscillator (for $\hbar = 1$),

$$H = \frac{p^2}{2m} + \frac{1}{2}m\omega_0^2 q^2 = \omega_0 \left( a^\dagger a + \frac{1}{2} \right),$$

where $q = (a^\dagger + a)/\sqrt{2m\omega_0}$, $p = i\sqrt{m\omega_0/2}(a^\dagger - a)$ relate the position and momentum operators ($[q, p] = i$) to the boson creation and annihilation operators ($[a, a^\dagger] = 1$).

Use the recursion method with inner product $\langle A|B \rangle = \frac{1}{2}(\langle AB \rangle + \langle BA \rangle)$ to calculate the structure function $S_{qq}(\omega)$ for the position variable at temperature $T$, where $\langle a^\dagger a \rangle = n_B = (e^{\beta\omega} - 1)^{-1}$, $\beta = 1/k_B T$.

Solution:
Consider the classical harmonic oscillator,

\[ H = \frac{p^2}{2m} + \frac{1}{2} m\omega_0^2 q^2. \]

Use the recursion method with inner product \( \langle A|B \rangle = \langle AB \rangle \) to calculate the structure function \( S_{qq}(\omega) \) for the position variable at temperature \( T \), where \( \langle p^2 \rangle / 2m = \frac{1}{2} m\omega_0^2 \langle q^2 \rangle = \frac{1}{2} k_B T \), according to the equipartition theorem.

**Solution:**
Consider the quantum harmonic oscillator (for $\hbar = 1$),

$$H = \frac{p^2}{2m} + \frac{1}{2}m\omega_0^2 q^2 = \omega_0 \left(a^\dagger a + \frac{1}{2}\right),$$

where $q = (a^\dagger + a)/\sqrt{2m\omega_0}$, $p = i\sqrt{m\omega_0}/2(a^\dagger - a)$ relate the position and momentum operators ($[q, p] = i$) to the boson creation and annihilation operators ($[a, a^\dagger] = 1$).

Solve the equation of motion

$$\zeta \langle\langle A; B\rangle\rangle^\pm = \langle[A, B]_\pm\rangle + \langle\langle A, H\rangle; B\rangle^\pm$$

for (a) the Green’s function $\langle\langle q; q\rangle\rangle^\pm$ and (b) the Green’s function $\langle\langle q; q\rangle\rangle^\mp$. Infer from each Green’s function the structure function $S_{qq}(\omega)$ for the position variable at temperature $T$.

Solution: