Mode-coupling theory for multiple decay channels

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Abstract. We investigate the properties of a class of mode-coupling equations for the glass transition where the density mode decays into multiple relaxation channels. We prove the existence and uniqueness of the solutions for Newtonian as well as Brownian dynamics and demonstrate that they fulfil the requirements of correlation functions; in the latter case the solutions are purely relaxational. Furthermore, we construct an effective mode-coupling functional, which allows us to map the theory to the case of a single decay channel, such that the covariance principle found for the mode-coupling theory for simple liquids is properly generalized. This in turn allows establishment of the maximum theorem stating that long-time limits of mode-coupling solutions can be calculated as maximal solutions of a fixed-point equation without relying on the dynamic solutions.

Keywords: mode coupling theory, slow dynamics and ageing (theory), structural glasses (theory), stochastic processes (theory)
1. Introduction

The dynamic properties of a liquid are encoded in time-dependent correlation functions of suitable observables. The principal quantity characterizing the collective motion is the dynamic density–density correlation function or intermediate scattering function $S(q, t)$ [1], which either is experimentally accessible directly by scattering techniques on the molecular scale or can be calculated as a statistical average in particle-tracking methods for colloidal systems. Since these correlation functions are subject to some underlying stochastic process, say Newtonian dynamics, probability theory imposes constraints on the class of permissible functions [2], the most important one being that the corresponding power spectra are non-negative. While these properties are automatically fulfilled in an experiment or a computer simulation in the stationary state or thermal equilibrium, a theoretical description usually relies on a series of simplifying assumptions thereby necessarily involving approximations. It is hence a non-trivial question if the solutions of theoretical models respect general features of dynamic correlation functions within the framework of probability theory.

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In the case of colloidal particles which are model liquids on the micrometer scale, the microscopic dynamics is Brownian motion rather than Newtonian dynamics. The solvent friction overdamps the dynamics such that all colloidal correlation functions exhibit pure relaxations, i.e. they can be represented as positive superpositions of decaying exponentials only. The corresponding even more restricted class of permissible functions is well known in the mathematical literature [2] and the property is referred to as completely monotone functions. A manifestation of this property is that the power spectra are superpositions of Lorentzians centered at zero frequency.

For dense or supercooled liquids the structural relaxation slows down drastically and a microscopic theory for the plethora of phenomena associated with the glass transition is provided by the mode-coupling theory (MCT) of the glass transition for simple liquids developed by Götze and collaborators [3, 4]. The theory provides a set of integrodifferential equations for $S(q,t)$ and makes a series of non-trivial predictions in the vicinity of a dynamic bifurcation, for example the emergence of an extended plateau in the intermediate scattering functions, power-law relaxations at mesoscopic time windows, the time–temperature superposition principle, etc. Although the sharp dynamical transition predicted by MCT becomes smeared by ergodicity-restoring processes, various facets of the theory have been tested successfully [4, 5] in numerous experiments on supercooled liquids, colloids, and extensive computer simulations on hard spheres, Lennard-Jones mixtures, and Yukawa particles.

Many of the mathematical properties for the intermediate scattering function $S(q,t)$ in the case of simple one-component systems have been worked out rigorously; in particular, it has been demonstrated that solutions of the MCT equations uniquely exist for Newtonian [6] as well as for Brownian dynamics [7] compatible with the constraints provided by probability theory. In the Brownian case the solutions are completely monotone, thus reflecting the underlying relaxational dynamics [7] (see [8] for explicit numerical results for a distribution of relaxation rates for a hard-sphere system), and the existence of the long-time limits called glass form factors follows. Second, the MCT equations respect a covariance principle under certain transformations which entails a series of properties for the solutions [4]. For example, the glass form factor is the maximal solution of a fixed-point equation and it can be determined by employing a convergent monotone iteration scheme and can therefore be calculated without solving the dynamical equations. Glass-transition lines are identified with bifurcation singularities for the glass form factor and the class of possible bifurcations is completely characterized.

The simplest generalization of MCT is to consider multi-component mixtures, which requires us to employ matrix-valued correlation functions $S_{\mu\nu}(q,t)$, where the indices $\mu, \nu$ refer to the different species in the liquid. Then the equations of motion couple different species and the theory naturally has to deal with integrodifferential equations for matrices. The notion of positivity then has to be adapted to positive-definite matrices, and it has been shown that the covariance principle can be properly generalized [9] and all properties shown for the single-component system can be proven as well.

Most glass formers are molecular liquids. Rigid molecules display an orientational degree of freedom besides the translational one, and a mode-coupling theory for a molecular liquid [10, 11] as well as a single molecule in a solvent of spherical particles [12] has been worked out. Expanding the orientation in a complete set, i.e. spherical harmonics $Y_{\ell m}$ in the case of linear molecules or Wigner rotation matrix $D^{(\ell)}_{mn}$ for arbitrarily shaped...
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particles, again leads to matrix-valued correlation functions, yet a new phenomenon occurs: the density changes with both translation and orientation, and to describe the slow structural relaxation it is necessary to consider multiple relaxation channels. As a non-trivial prediction of these MCT equations, the glassy dynamics is independent of inertial parameters, i.e. the mass and moment of inertia of its molecular constituents [10]–[12]. However, the mathematical structure of these MCT equations differs explicitly from the case of multi-component systems; in particular, the covariance property appears to be lost.

Recently, liquids confined to a slit have been considered [13, 14], where translational invariance perpendicular to the walls is explicitly broken. Expansion into suitable Fourier modes for the perpendicular degree of freedom again requires us to deal with matrix-valued correlators \( S_{\mu\nu}(q, t) \), where the indices now refer to different modes. The fluctuating density decays by currents perpendicular or parallel to the walls, and the mathematical structure of the MCT is identical to that of molecules.

The goal of this work is to provide proofs that the MCT equations for multiple relaxation channels display unique solutions which reflect the constraints imposed on being correlation functions. This is achieved in the case of Newtonian as well as Brownian dynamics; in the latter case the theory ignores hydrodynamic interactions. We also formulate a covariance principle for the equations of motion in a wider sense than used for one- or multi-component systems, which is essential to connect to the techniques developed to prove the properties within the mode-coupling approach. We shall show that this generalized covariance principle is suited to introduce an effective mode-coupling functional such that the theory assumes the mathematical structure of a multi-component mixture. The properties of MCT equations for multiple relaxation channels follow from this mapping corroborating that the MCT approach is a robust strategy to arrive at physical correlation functions.

2. Matrix-valued correlation functions

The quantities of interest are the generalized intermediate scattering functions (ISFs)

\[
S_{\mu\nu}(q, t) = N^{-1} \langle \rho_{\mu}(\vec{q}, t)^* \rho_{\nu}(\vec{q}, 0) \rangle,
\]

as the time-correlation matrix of a set of density modes \( \rho_{\mu}(\vec{q}, t) \) labeled by a discrete mode index \( \mu \) and a continuous wavenumber \( \vec{q} \). Examples for such density modes will be given below. Furthermore, \( N \) is the number of particles and conventions are such that \( S_{\mu\nu}(q, t) \) becomes independent of the system size in the thermodynamic limit. The theory becomes more elegant by treating the generalized ISF as a matrix \( [S(q, t)]_{\mu\nu} = S_{\mu\nu}(q, t) \), and similarly for other correlation functions. To avoid technical complications we shall assume the matrices to be finite dimensional and the wavenumbers to be of a discrete finite set. We shall suppress the wavenumber \( q \) in the following if it merely serves as a label; the relations derived then hold for each wavenumber separately.

General properties of the ISF can be inferred from the fact that for any set of complex numbers \( y_{\mu} \) (and for each fixed wavenumber \( q \)) the quantity \( \sum_{\mu\nu} y_{\mu}^* S_{\mu\nu}(t) y_{\nu} \) is an autocorrelation function of a stationary stochastic process. By the spectral representation theorem [2] it can be written as the characteristic function \( \int e^{-i\Omega t} R(\Omega) \) of a finite spectral measure \( R(\Omega) \) corresponding to a non-decreasing, right-continuous function on the real
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line with \( R(-\infty) = 0, R(\infty) < \infty \). Within equilibrium statistical physics the time-reversed stochastic process obeys the same probabilistic law implying that the spectral measure is also symmetric.

The generalization to the matrix case is straightforward \([15]\),

\[ S(t) = \int e^{-i\Omega t} R(d\Omega), \]

such that \( R(\Omega) \) is a self-adjoint matrix-valued measure, i.e. \( R_{\mu\nu}(\Omega) \) is a complex finite Borel measure on the real line \( \Omega \in \mathbb{R} \) and a positive-semidefinite matrix for fixed \( \Omega \). Thus the matrix correlator fulfils \( S(-t) = S^\dagger(t) \) and by the equilibrium property it is also invariant under time-reversal \( S(-t) = S(t) \). Combining the two properties, one infers that the ISF is Hermitian, \( S(t) = S^\dagger(t) \).

We define the Fourier–Laplace transform by

\[ \hat{S}(z) = i \int_0^\infty e^{int} S(t) \, dt, \]

where the complex frequency \( z \) is confined to the upper complex half-plane \( \mathbb{C}_+ = \{ z \in \mathbb{C} | \text{Im}[z] > 0 \} \). The spectral representation theorem implies

\[ \hat{S}(z) = \int \frac{1}{\Omega - z} R(d\Omega), \quad z \in \mathbb{C}_+. \]

The following properties follow readily from the preceding relation for the Fourier–Laplace transforms of matrix correlation functions in equilibrium and complex frequencies \( z \in \mathbb{C}_+ \).

1. \( \hat{S}(z) \) is analytic.
2. \( \hat{S}(-z^*) = -\hat{S}^\dagger(z) \).
3. \( \lim_{\eta \to \infty} \eta \text{Im}[\hat{S}(z = i\eta)] \) is finite.
4. \( \text{Im}[\hat{S}(z)] \succeq 0 \).

Here, \( \text{Im}[\hat{S}(z)] = [\hat{S}(z) - \hat{S}^\dagger(z)]/2i \) is the proper generalization of the imaginary part for the matrix \( \hat{S}(z) \), and the symbol \( \succeq 0 \) indicates that the matrix is positive-semidefinite. Conversely, the Riesz–Herglotz representation theorem \([15]\) reveals that these four conditions are sufficient for \( \hat{S}(z) \) being the Laplace transform of an equilibrium matrix correlation function.

An important subclass of equilibrium correlation functions arises if the dynamics is purely relaxational, for example in the case of Brownian dynamics. Then all time-correlations consist of superpositions of relaxing exponentials in the following sense:

\[ S(t) = \int_0^\infty e^{-\gamma t} a(d\gamma), \quad t \geq 0, \]

where \( a(\gamma) \) is again a self-adjoint matrix-valued measure. Such functions \( S(t) \) are referred to as completely monotone and display the property that time-derivatives \( [-\partial_t]^\ell S(t) \succeq 0, \ell \in \mathbb{N} \) are positive-semidefinite for \( t > 0 \). By Bernstein’s theorem \([2]\) this condition is equivalent to the integral representation, equation (5).

The Fourier–Laplace transform of a completely monotone matrix-valued function can be extended to complex frequencies not located on the negative imaginary axis and is
represented as
\[ \hat{S}(z) = \int_0^\infty \frac{-1}{z + i\gamma} a(\text{d}\gamma), \quad z \in \mathbb{C} \setminus i\mathbb{R}^- . \] (6)

The following properties can be verified easily for \( z \in \mathbb{C} \setminus i\mathbb{R}^- \).

1. \( \hat{S}(z) \) is analytic.
2. \( \hat{S}(-z^*) = -\hat{S}^\dagger(z) \).
3. \( \lim_{\eta \to \infty} \hat{S}(z = i\eta) = 0 \).
4. \( \text{Re}\{\hat{S}(z)\} \geq 0 \) for \( \text{Re}\{z\} < 0 \).

Here, \( \text{Re}\{\hat{S}(z)\} = [\hat{S}(z) + \hat{S}^\dagger(z)]/2 \) now denotes the Hermitian part of the matrix \( \hat{S}(z) \). Conversely, these four properties guarantee a representation of the form of equation (5) (see [16], section 5, theorem 2.6).

3. Mode-coupling equations with multiple decay channels

We consider the dynamics of a generalized fluctuating density \( \rho_\mu(\vec{q}, t) \) where \( \vec{q} \) is a wavenumber along a direction in which translational invariance in the statistical sense holds. The additional mode index \( \mu \) selects further properties of the density mode. For example, in the case of multi-component mixtures it refers to the different species, whereas for linear molecules it represents a multi-index \( \mu = (\ell, m) \) characterizing the orientation of the molecule in terms of spherical harmonics \( Y_{\ell m}(\vartheta, \phi) \). For molecules of arbitrary shape Wigner rotation matrices \( D_{\ell mn}(\vartheta, \varphi, \chi) \) have to be employed, and correspondingly \( \mu = (\ell, m, n) \) encodes the dependence on the three Euler angles \( (\vartheta, \varphi, \chi) \) of the orientation of the molecule. In liquids confined to a slab of effective width \( L \) the wavenumber \( \vec{q} \) is only for directions parallel to the walls and \( \mu \in \mathbb{Z} \) indicates the discrete Fourier modes of wavenumbers \( Q_\mu = 2\pi\mu/L \) for the modulation perpendicular to the walls.

The density modes fulfil a continuity equation
\[ \dot{\rho}_\mu(\vec{q}, t) = i\sum_{\alpha=1}^r q^\alpha_\mu j^\alpha_\mu(\vec{q}, t), \] (7)

where we allow for \( r \in \mathbb{N} \) different decay channels. For one- or multi-component bulk liquids comprised of structureless particles only a single current is needed to be considered, \( r = 1 \), and the corresponding currents \( j^\alpha_\mu(\vec{q}, t) \) are identified with the species-dependent longitudinal mass transport, whereas the real-valued channel coupling \( q^\alpha_\mu \) corresponds to the magnitude of the wavenumber \( q = |\vec{q}| \). For linear molecules already two decay channels become relevant [10, 12], a translational one \( \alpha = T \) with \( q^T_\ell m = q \) and an orientational one \( \alpha = R \) where \( q^R_{\ell m} = \sqrt{\ell(\ell + 1)} \). General molecules require as many as six decay channels to be considered [17], three translational ones corresponding to the Cartesian directions and three orientational ones associated with the Cartesian components of the body-fixed frame. In the case of liquids in slit geometry [13, 14] the decay of the density mode occurs parallel and perpendicular to the walls, and characteristic couplings are \( q^0_\mu = q \) and \( q^\perp_\mu = Q_\mu \), respectively.

Correlating the density modes \( S_{\mu \nu}(q, t) = N^{-1}\langle \rho_\mu(\vec{q}, t)^*\rho_\nu(\vec{q}, 0) \rangle \) yields matrix-valued intermediate scattering functions, which should reflect the properties elaborated in
section 2. For Newtonian dynamics, a formally exact equation of motion can be derived relying on the Zwanzig–Mori projection operator formalism [1, 4], choosing as distinguished variables first the set of fluctuating densities \{\rho_\mu(\vec{q}, t)\} and then the currents \{j_\mu(\vec{q}, t)\}. The first equation of motion is found to be

$$\dot{S}(t) + \int_0^t K(t-t')S(t') \, dt' = 0,$$

where \(S = S(t = 0)\) is the initial value of the ISF and is identified as a generalized static structure factor. We again suppress the dependence on the wavenumber \(q\) if no confusion can occur, and all quantities appearing in the equations are to be read as wavenumber dependent. The explicit expression for the memory kernel \([K(q, t)]_{\mu\nu} = K_{\mu\nu}(q, t)\) reveals that it corresponds to a time-dependent correlation function of the variables \(\dot{\rho}_\mu(\vec{q}), \dot{\rho}_\nu(\vec{q})\) such that the dynamics is driven by a reduced Liouville operator [1, 4]. By Fourier–Laplace transform one can formally solve equation (8)

$$\hat{S}(z) = -\left[zS^{-1} + S^{-1} \hat{K}(z)S^{-1}\right]^{-1}.$$  

Employing the continuity equation, equation (7), the memory kernel naturally splits,

$$K_{\mu\nu}(q, t) = \sum_{\alpha=1}^{r} \sum_{\beta=1}^{r} q_\alpha^\beta K_{\mu\nu}^{\alpha\beta}(q, t)q_\nu^\beta.$$  

We indicate quantities associated with mode indices and channel indices by calligraphic symbols and again use matrix notation, e.g. \([K(q, t)]_{\alpha\beta} = K_{\alpha\beta}(q, t)\).

Employing the second Zwanzig–Mori projection step for the case of Newtonian dynamics, one finds the equation of motion for the currents

$$\hat{\mathcal{K}}(t) + JL^{-1}\mathcal{K}(t) + \int_0^t \mathcal{M}(t-t')\mathcal{K}(t') \, dt' = 0,$$

where \(\mathcal{J}^{\alpha\beta}(q) = N^{-1}\langle j_\mu^{\alpha}(q)j_\nu^{\beta}(q)\rangle\) is the static current correlator matrix, which also serves as initial condition \(\mathcal{K}(q, t = 0) = \mathcal{J}(q)\), and \([\mathcal{M}(q, t)]_{\alpha\beta} = M_{\mu\nu}^{\alpha\beta}(q, t)\) is a force kernel. Furthermore, we allow for an instantaneous damping term with positive-semidefinite matrices \(L^{-1}(q) \succeq 0\).

A formal solution is again obtained by Fourier–Laplace transform,

$$\hat{\mathcal{K}}(z) = -\left[z\mathcal{J}^{-1} + iL^{-1} + \hat{\mathcal{M}}(z)\right]^{-1}.$$  

Provided \(\hat{\mathcal{M}}(z) = \mathcal{O}(z^{-1})\) for \(z \to \infty\), the leading terms in the high-frequency expansion for the current correlator and the ISF do not involve the force kernel

$$\hat{K}(z) = -z^{-1}\mathcal{J} + z^{-2}\mathcal{I} + \mathcal{O}(z^{-3}),$$

$$\hat{S}(z) = -z^{-1}S - z^{-2}J + iz^{-4}\mathcal{I} + \mathcal{O}(z^{-5}), \quad z \to \infty,$$

where the contracted static current correlation matrix is denoted by \(J_{\mu\nu}(q) = \sum_{\alpha\beta} q_\mu^\alpha J_{\mu\nu}^{\alpha\beta}(q)q_\nu^\beta\) and similarly \([\mathcal{I}(q)]_{\mu\nu} = \sum_{\alpha\beta} q_\mu^\alpha [\mathcal{J}(q)L^{-1}(q)\mathcal{J}(q)]_{\alpha\beta}q_\nu^\beta\). The corresponding short-time expansion is readily inferred to be
Here the leading correction of order $O(t^4)$ depends on the force kernel $M(t=0)$.

Within the MCT approach the memory kernel $M(q,t)$ is approximated as a functional that is local in time of the ISF as positive superpositions of products of the ISF [14]:

$$M_{\mu\nu}^{\alpha\beta}(q,t) = F_{\mu\nu}^{\alpha\beta}[S(t), S(t); q] + (M_{\text{reg}})^{\alpha\beta}_{\mu\nu}(q,t).$$

Explicitly, the MCT functional is represented as

$$F_{\mu\nu}^{\alpha\beta}[S(t), S(t); q] = \frac{1}{2N} \sum_{q_1, q_2} \sum_{\mu_1 \mu_2} \mathcal{Y}_{\mu_1 \mu_2}^{\alpha\beta}(q_1, q_2)$$

$$\times \mathcal{S}_{\mu_1 \nu_1}(q_1, t) \mathcal{S}_{\mu_2 \nu_2}(q_2, t) \mathcal{Y}_{\mu_1 \mu_2}^{\beta\alpha}(q_1, q_2),$$

and the vertices $\mathcal{Y}_{\mu_1 \mu_2}^{\alpha\beta}(q_1, q_2)$ are determined by static properties of the liquid only and assumed to be known smooth functions of control parameters; see [14]. The force correlator may contain an additional regular damping kernel $(M_{\text{reg}})(q,t)$ encoding faster dynamical processes, which are not captured by the MCT functional. This kernel is assumed to be known a priori and satisfies the constraints formulated in section 2 of an equilibrium matrix-valued correlation function.

Equations (8), (10), (11), and (16), (17) together with the initial conditions $S(t=0) = S, K(t=0) = J$ constitute a closed set of equations for the Newtonian case, and we shall prove the existence and uniqueness of solutions that are correlation functions in section 4.

For Brownian dynamics, equation (12) is replaced by

$$\dot{K}(z) = -\left[iD^{-1} + M(z)\right]^{-1},$$

such that $D$ is a positive-definite matrix. Here again, the memory kernel may also contain an a priori known regular damping, which satisfies the additional constraints of a purely relaxational equilibrium matrix-valued correlation function; see section 2. By the same arguments as above, the high-frequency expansion for the ISF is found to be

$$\dot{S}(z) = -z^{-1}S + z^{-2}iD + O(z^{-3}), \quad z \to \infty,$$

with the contracted relaxation rate matrix $D_{\mu\nu}(q) = \sum_{\alpha\beta} q_{\alpha\beta}^{\mu} D^{\alpha\beta}_{\mu\nu}(q) q_{\alpha\beta}^{\nu}$. In the temporal domain this implies

$$S(t) = S - Dt + O(t^2), \quad t \to 0,$$

and the positive-definite matrix $D$ determines the initial decay of the ISF.

From equation (18), one infers $\dot{K}(z) \to iD$ as $z \to \infty$, which suggests that $K(t)$ contains a $\delta$-function at the time origin. To arrive at integrodifferential equations with smooth functions, we split off the high-frequency limit, $\delta K(z) = K(z) - iD$, consistent with its contraction $\delta^* K(z) = K(z) - iD$. Then equation (9) translates back in the temporal domain to

$$\dot{S}(t) + DS^{-1}S(t) + \int_0^t \delta K(t-t')S(t') dt' = 0.$$
Second, equation (18) implies
\[
\left[ iD^{-1} + \hat{M}(z) \right] \delta \hat{K}(z) = -\hat{M}(z)iD, \tag{22}
\]
which yields an integral equation in the temporal domain
\[
\delta \hat{K}(t) + \int_0^t D\hat{M}(t-t')\delta \hat{K}(t') \, dt' = -D\hat{M}(t)D. \tag{23}
\]
Equations (21), (10), (23) and (16), (17) together with the initial conditions \(S(t = 0) = S\) constitute a closed set of equations for the Brownian case. The properties of their solutions are one of the central questions of this work.

4. Construction of MCT solutions

The existence of MCT solutions with the required properties is intimately connected to the representation of ISF in terms of continued fractions as prescribed by the Zwanzig–Mori approach, as well as with the closure via a microscopically derived MCT functional. We shall show that the force kernel obtained via an MCT functional of correlation functions and completely monotone functions has the same properties as correlation functions and completely monotone functions, respectively, and, given a force kernel with these properties, the ISF as obtained via the continued fraction has again the same properties. This observation will allow us to construct an iteration scheme not leaving the class of correlation functions or completely monotone functions. Since both classes are closed, the limiting functions, provided they exist, still display the desired properties. The scheme is then translated to the time domain and shown to be a constructive proof of a unique solution.

4.1. Properties of the MCT functional

In the first step we collect some properties of the MCT functional, some of which have already been discussed in [14]. To avoid technical problems we consider all matrices to be finite dimensional and the set of wavenumbers as a discrete finite set. Then the MCT functional maps a set of matrices in the mode space labeled by a wavenumber to matrices with a mode and channel indices with the same set of labels for the wavenumbers. Since the functional is local in time, time merely enters as a parameter.

It is convenient to introduce the pair-mode indices \(a := (\mu_1, \mu_2)\) and \(b := (\nu_1, \nu_2)\) and the double indices \(\gamma := (\alpha, \mu), \delta := (\beta, \nu)\). The MCT functional, equation (17), can then be represented as
\[
\mathcal{F}[F, E; q]^{\gamma\delta} = \frac{1}{4N} \sum_{ab} \sum_{\bar{q}_1, \bar{q}_2} Y_a^\gamma(\bar{q}, \bar{q}_1 \bar{q}_2) \times \left[ F(q_1) \otimes E(q_2) + E(q_1) \otimes F(q_2) \right]_{ab} Y_b^\delta(\bar{q}, \bar{q}_1 \bar{q}_2)^* \tag{24}
\]
where \(\otimes\) denotes the Kronecker product in the space of mode indices, \(E_{\mu_1\nu_1}(q_1)F_{\mu_2\nu_2}(q_2) = [E(q_1) \otimes F(q_2)]_{\alpha=\mu_1, \beta=\mu_2, \lambda=\nu_1, \gamma=\nu_2}\).

Positive definiteness of a Hermitian matrix \(E\) will be denoted \(E \succ 0\), and positive semidefiniteness by \(E \succeq 0\). Similarly, for two Hermitian matrices we write \(F \succ E\) if
\(F - E \succ 0\), etc. In the following, the label \(q\) for the wavenumber is suppressed and all operations are understood component-wise for each \(q\). Then one easily shows \cite{14} that \(E \succeq 0, F \succeq 0\) implies \(\mathcal{F}[F, E] \succeq 0\). Furthermore, the functional preserves ordering: \(F \succeq E\), \(\mathcal{F}[F, F] \succeq \mathcal{F}[E, E]\).

A matrix-valued correlation function \(S_{\mu\nu}(q, t)\) fulfills \(\sum_{ij} \sum_{\mu\nu} \xi_{ij}\, S_{\mu\nu}(q, t_i - t_j) \xi_{j\nu} \geq 0\) for any finite set of times \(t_i \in \mathbb{R}\) and complex numbers \(\xi_{ij} \in \mathbb{C}\). By Bochner’s theorem \cite{2}, this is an equivalent characterization of a (matrix-valued) correlation function. Then one concludes that if \(S_{\mu\nu}(q, t)\) and \(T_{\mu\nu}(q, t)\) are matrix-valued correlation functions so is the matrix \([S(q_1, t) \otimes T(q_2, t)]_{ab}\). This follows literally as in the case for scalars by decorating the complex numbers by indices. For complex numbers \(s_i^q\) then

\[
\sum_{ij} \sum_{\gamma\delta} s_i^q \mathcal{F}[S(t_i - t_j), T(t_i - t_j); q] s_j = \frac{1}{4N} \sum_{q_1, q_2} \left\{ \sum_{ij} \sum_{ab} \left( \sum_{\gamma} s_i^q \mathcal{Y}_{\mu\nu}(q q_1 q_2) \right) \times [S(q_1, t_i - t_j) \otimes T(q_2, t_i - t_j) + T(q_1, t_i - t_j) \otimes S(q_2, t_i - t_j)]_{ab} \times \left( \sum_{\delta} s_j^q \mathcal{Y}_{\mu\nu}^*(q q_1 q_2) \right) \right\} \geq 0,
\]

(25)

since the sums in the curly bracket are already non-negative by the characterization of correlation functions. Therefore, the MCT functional maps correlation functions to correlation functions.

In the case of Brownian dynamics the correlation functions \(S_{\mu\nu}(q, t)\) are completely monotone, and by Bernstein’s theorem \cite{2} can be uniformly approximated for \(t \geq 0\) by sums

\[
\sum_i e^{-\gamma_i t} \Delta a_{\mu\nu}^{(0)}(\gamma_i) \to S_{\mu\nu}(q, t),
\]

(26)

with \(\Delta a^{(0)}(\gamma_i) \geq 0\) and \(\gamma_i \geq 0\). Using such approximants for \(S(q, t)\) and \(T(q, t)\) a calculation similar to equation (25) shows that the approximants to \(\mathcal{F}[S(t), T(t); q] s^q\) are of the same type. Passing to the limits shows that the MCT functional maps completely monotone correlation functions onto completely monotone correlation functions.

4.2. Zwanzig–Mori equation of motion

Next we show that the representation of the ISF via memory kernels implies that correlation functions map to correlation functions, and in the Brownian case the property of complete monotonicity is also inherited. The proof is performed most easily in the Fourier–Laplace domain and follows essentially the argument for mixtures \cite{9}. Since the wavenumber merely serves as a label, we shall omit the dependence on \(q\) in the remainder of this section when convenient.

In the Newtonian case we shall assume that \(\mathcal{M}(t)\) is a correlation function which is equivalent to \(\hat{\mathcal{M}}(z)\) fulfilling properties (1)–(4) of section 2. The spectral representation shows that \(\hat{\mathcal{M}}(z = i\eta) = \mathcal{O}(\eta^{-1})\) as \(\eta \to \infty\). From the second equation of motion, equation (12), one infers readily that \(\hat{\mathcal{K}}(z)\) inherits property (1), since the denominator has positive imaginary part (in the matrix sense) for \(z \in \mathbb{C}_+\) and hence no zeros can occur. Second, property (2) is readily checked, and by \(\hat{\mathcal{K}}(z = i\eta) = i\eta^{-1} \mathcal{F} + \mathcal{O}(\eta^{-2})\) for \(\eta \to \infty\) property (3) follows. Decomposing the current correlator into real and imaginary parts.
with \( \text{Im}[z] > 0 \) yields
\[
\text{Im}[\hat{K}(z)] = \hat{K}^\dagger(z) \text{Im}[\hat{K}(z)^{-1}] \hat{K}(z) = \hat{K}^\dagger(z) \left\{ \text{Im}[z] J^{-1} + D^{-1} + \text{Im}[\hat{M}(z)] \right\} \hat{K}(z) \geq 0, \tag{27}
\]
which shows that property (4) is also fulfilled. Hence \( \hat{K}(z) \) is again the Fourier–Laplace transform of a correlation function and so is its contraction \( \hat{K}(z) \); in particular, \( \text{Im}[\hat{K}(z)] \geq 0 \) for \( \text{Im}[z] > 0 \). The first equation of motion, equation (9), shows by the same arguments as above that \( \hat{S}(z) \) again fulfills properties (1), (2), and (3). The same calculation as above,
\[
\text{Im}[\hat{S}(z)] = \hat{S}^\dagger(z) \left\{ \text{Im}[z] S^{-1} + S^{-1} \text{Im}[\hat{K}(z)] S^{-1} \right\} \hat{S}(z), \tag{28}
\]
demonstrates property (4), \( \text{Im}[\hat{S}(z)] \geq 0 \) for \( \text{Im}[z] > 0 \). Hence the Zwanzig–Mori equations for the Newtonian case yield correlation functions provided the force kernel is itself a correlation function. For a converse representation theorem see appendix A.

In the case of Brownian dynamics, \( \hat{M}(z) \) is assumed to fulfill \((1^*)–(4^*)\). Then equation (18) shows that \( -\delta \hat{K}(z) = iD - \hat{K}(z) \) also satisfies \((1^*), (2^*), \) and \((3^*)\). With the same calculation as in equation (27) one can show
\[
\text{Re}[-\delta \hat{K}(z)] = \hat{K}^\dagger(z) \text{Re}[\hat{M}(z)] \hat{K}(z). \tag{29}
\]
Hence for \( \text{Re}[z] < 0 \), \( \text{Re}[-\delta \hat{K}(z)] \geq 0 \), completing the argument that \( -\delta \hat{K}(z) \) is completely monotone and so is its contraction \( -\delta \hat{K}(z) \). The first Zwanzig–Mori equation, equation (9), shows that \( \hat{S}(z) \) also inherits properties \((1^*), (2^*), \) and \((3^*)\). Using \( \text{Re}[\hat{K}(z)] = \text{Re}[\delta \hat{K}(z)] \), the relation
\[
\text{Re}[\hat{S}(z)] = \hat{S}^\dagger(z) \left\{ -\text{Re}[z] S^{-1} + S^{-1} \text{Re}[-\delta \hat{K}(z)] S^{-1} \right\} \hat{S}(z), \tag{30}
\]
then proves also property \((4^*)\). Thus the Zwanzig–Mori equations for Brownian dynamics yield completely monotone correlation functions provided the force kernel is of the same type.

### 4.3. Iteration scheme

The fact that the mapping from (completely monotone) correlation functions yields a memory kernel by a mode-coupling functional that has the same properties, as well as the mapping from the memory kernel to the ISF again preserving the properties, suggests construction of an iteration scheme within the closed space of such functions. Such a scheme was first used by Götze and Lücke [18] for the dynamic structure factor of liquid helium. All operations are again to be understood for each wavenumber \( q \).

In the Newtonian case the sequence is initialized by \( (S^{(0)}(t), K^{(0)}(t)) = (S, J) \), which are trivially correlation functions. Equivalently, in the Fourier–Laplace domain \( (\hat{S}^{(0)}(z), \hat{K}^{(0)}(z)) = (-S/z, -J/z) \). Then the following mapping is employed:
\[
\hat{S}^{(n+1)}(z) = -\left[ zS^{-1} + S^{-1} \hat{K}^{(n)}(z) S^{-1} \right]^{-1}, \tag{31}
\]
Here $\hat{K}^{(n)}(z)$ is of course the contraction of $\hat{K}^{(n)}(z)$ and $\hat{M}^{(n)}(z)$ is the Fourier–Laplace transform of

$$\hat{M}^{(n)}(q, t) = \mathcal{F}[S^{(n)}(t), S^{(n)}(t); q] + (\mathcal{M}_{\text{reg}})(q, t).$$

Hence the coupling of different wavenumbers emerges via the MCT functional only.

In the case of Brownian dynamics the sequence is started by $(\hat{S}^{(0)}(z), \hat{K}^{(0)}(z)) = (-S/z, i\mathcal{D})$ and equation (32) is replaced by

$$\hat{K}^{(n+1)}(z) = -\left[i\mathcal{D} - \hat{M}^{(n)}(z)\right]^{-1}.$$ 

In section 4.2 it was shown that this corresponds to $-\delta \hat{K}^{(n+1)}(z) = i\mathcal{D} - \hat{K}^{(n+1)}(z)$, which is the Fourier–Laplace transform of a completely monotone correlation function.

Provided the sequence in the temporal domain $(S^{(n)}(t), K^{(n)}(t))$ converges uniformly on each finite time interval, the limits $S(t) = \lim_{n \to \infty} S^{(n)}(t)$ are matrix-valued correlation functions which are also completely monotone in the case of Brownian dynamics [2]. Furthermore, the limit represents a solution of the MCT equations.

4.4. Proof of convergence

The sequence constructed in section 4.3 will be translated to a sequence of functions in the temporal domain. The convergence shall then be demonstrated by providing estimates with respect to a suitable norm that we introduce below, following the strategy of Haussmann [6]. The main difference from the Picard iteration is that existence and uniqueness can be proven globally with uniform convergence on finite time intervals. The required ingredient is precisely the property that all approximants are correlation functions and therefore bounded by their respective initial values.

First, for Newtonian dynamics, equations (31) and (32) can be recast to integrodifferential equations, similar to equations (8) and (11). Integrating once,

$$S^{(n+1)}(t) = S - \int_0^t dt' \int_0^{t'} dt'' K^{(n)}(t' - t'') S^{-1} S^{(n+1)}(t''),$$

$$K^{(n+1)}(t) = \mathcal{J} - \int_0^t dt' \mathcal{J} D^{-1} K^{(n+1)}(t') - \int_0^t dt' \int_0^{t'} dt'' \mathcal{J} M^{(n)}(t' - t'') K^{(n+1)}(t''),$$

we obtain equations which can be analyzed by the techniques familiar from the Picard iteration scheme for ordinary differential equations. In contrast to this technique, at each iteration step a self-consistent integral equation requires to be solved to evaluate the subsequent approximant. This procedure ensures that each approximant satisfies the constraints of being a correlation function, as formulated in section 2.

Following [9], we denote the space of matrices with mode indices by $\mathcal{A}_0$ equipped with the standard operator norm $\| \cdot \|$. Then $\mathcal{A}_0$ is a $C^*$ algebra with respect to matrix multiplication and Hermitian conjugation as $*$-operation. For a finite set of $M$ wavenumbers, the matrices naturally form vectors $E = [E(q)]_{q=1,...,M} \in \mathcal{A}_0^M$. Equipping $\mathcal{A}_0^M$ with the maximum norm $\| E \| = \max_q \| E_q \|$ and taking all operations component-
wise, one easily shows that $\mathcal{A}_{0}^{M}$ is again a $C^*$ algebra [19]. Elements are called positive-semidefinite, $E \succeq 0$, if all components $E(q) \succeq 0$ are positive-semidefinite for each $q$. In particular, the norm preserves ordering; $E \succeq F$ implies $\|E\| \geq \|F\|$. The space of $M$-tuples of all matrices with both mode and channel indices denoted by $\mathcal{A}_{0}^{M}$ will be equipped by another $C^*$ algebra by the same procedure, and for simplicity we use the same symbol to indicate norms. Then contraction with the selectors induces a Lipschitz-continuous mapping of $\mathcal{A}_{0}^{M} \to \mathcal{A}_{0}^{M}$.

Since all approximants $S^{(n)}(t), K^{(n)}(t), K^{(n)}(t)$ are correlation functions, the spectral representation, equation (2), reveals that they cannot exceed their initial value, i.e. $S \succeq S^{(n)}(t), J \succeq K^{(n)}(t), J \succeq K^{(n)}(t)$, which implies $\|S^{(n)}(t)\| \leq \|S\|, \|K^{(n)}(t)\| \leq \|J\|, \|K^{(n)}(t)\| \leq \|J\|$.

From equation (35) one derives the standard estimates for $n \in \mathbb{N}$,

$$\|S^{(n+1)}(t) - S^{(n)}(t)\| \leq \int_{0}^{t} dt' \int_{0}^{t'} dt'' \|J\| \|S^{-1}\| \|S^{(n+1)}(t'') - S^{(n)}(t'')\|$$

$$+ \int_{0}^{t} dt' \int_{0}^{t'} dt'' \|S\| \|S^{-1}\| \|K^{(n)}(t'') - K^{(n-1)}(t'')\|,$$

(37)

where the symmetry of the convolution has been exploited to shift the time argument in the last line. Similarly, equation (36) yields for $n \in \mathbb{N}$

$$\|K^{(n+1)}(t) - K^{(n)}(t)\| \leq \int_{0}^{t} dt' \|J\| \|D^{-1}\| \|K^{(n+1)}(t') - K^{(n)}(t')\|$$

$$+ \int_{0}^{t} dt' \int_{0}^{t'} dt'' \|J\| \|M\| \|K^{(n+1)}(t'') - K^{(n)}(t'')\|$$

$$+ \int_{0}^{t} dt' \int_{0}^{t'} dt'' \|J\|^2 \|M^{(n)}(t'') - M^{(n-1)}(t'')\|,$$

(38)

where $M = M(t = 0)$. Note that $\|M^{(n)}(t)\| \leq \|M\|$ has been used as well as $\|S^{(n)}(t)\| \leq \|S\|$.

By Lipschitz-continuity of the contraction $\|K^{(n+1)}(t) - K^{(n)}(t)\| \leq L_{1}\|K^{(n+1)}(t) - K^{(n)}(t)\|$ and Lipschitz-continuity of the mode-coupling functional $\|M^{(n+1)}(t) - M^{(n)}(t)\| \leq L_{2}\|S^{(n+1)}(t) - S^{(n)}(t)\|$ with suitable Lipschitz constants $L_{1} > 0, L_{2} > 0$. Then the growth of the scalar $X^{(n)}(t) = \|S^{(n+1)}(t) - S^{(n)}(t)\| + c\|K^{(n+1)}(t) - K^{(n)}(t)\|$, where $c > 0$ accounts for the different dimensional units, in the finite time interval $0 \leq t \leq T < \infty$ is controlled by

$$X^{(n)}(t) \leq L \int_{0}^{t} dt' [X^{(n)}(t') + X^{(n-1)}(t')], \quad n \in \mathbb{N}$$

(39)

with $L = T\|J\| \|S^{-1}\| + c\|J\| \|D^{-1}\| + cT\|J\| \|M\| + L_{1}T \|S\| \|S^{-1}\| + cL_{2}T\|J\|^2$. The preceding equation with the additional bound $X^{(n)}(t) \leq K := 2c\|J\| + 2\|S\|$ is essentially the same as studied in [4, 6, 7], and for completeness we repeat the argument. The sequence $X^{(n)}(t)$ is bounded by the $a^{(n)}(t)$ recursively defined by $a^{(n)}(t) = 2L \int_{0}^{t} dt' a^{(n-1)}(t'), a^{(0)}(t) = K$. First, $X^{(0)}(t) \leq a^{(0)}(t) = K, X^{(1)}(t) \leq a^{(0)}(t)$, and $X^{(1)}(t) \leq a^{(1)}(t) = 2LKt$ are obvious. Then $X^{(2)}(t) \leq 2L \int_{0}^{t} dt' a^{(0)}(t') = a^{(1)}(t)$, which in turn implies $X^{(2)}(t) \leq 2L \int_{0}^{t} dt' a^{(1)}(t') = a^{(2)}(t)$. The procedure can be continued to show
that $X^{(n)}(t) \leq a^{(m)}(t)$ for $m \leq n$. Explicit calculation shows $a^{(n)}(t) = K(2Lt)^n/n!$ such that $\|S^{(n)}(t) - S^{(m)}(t)\| \leq \sum_{k=m}^{n-1} \|S^{(k+1)}(t) - S^{(k)}(t)\| \leq \sum_{k=m}^{n} X^{(k)}(t) \leq \sum_{k=m}^{\infty} a^{(k)}(t) \to 0$ as $n > m \to \infty$, implying $S^{(n)}(t)$ to be Cauchy sequences uniformly on every finite time interval. Since $A_0^{(n)}$ is complete, the limit $S(t) = \lim_{n \to \infty} S^{(n)}(t)$ exists and represents again a correlation function [2], and similarly for the currents $\mathcal{K}(t) = \lim_{n \to \infty} \mathcal{K}^{(n)}(t)$.

Uniqueness is demonstrated similarly. Assuming two sets of solutions $S(t), \mathcal{K}(t)$ and $\tilde{S}(t), \tilde{\mathcal{K}}(t)$ with the same initial conditions, going through the same steps as above reveals that $Y(t) := \|S(t) - \tilde{S}(t)\| + c\|\mathcal{K}(t) - \tilde{\mathcal{K}}(t)\|$ satisfies $Y(t) \leq 2L \int_0^t dt' Y(t')$ and $Y(0) = 0$. Then, by the same argument as in the preceding paragraph, induction shows $0 \leq Y(t) \leq a^{(n)}(t) \to 0$ as $n \to \infty$ for all times $t$, hence the two solutions are equal.

In the case of Brownian dynamics, equations (35) and (36) will be replaced by (cf equations (21) and (23))

\[
S^{(n+1)}(t) = S - \int_0^t D S^{-1} S^{(n+1)}(t') dt' - \int_0^t dt' \int_0^t dt'' \delta \mathcal{K}^{(n)}(t') - t'' S^{-1} S^{(n+1)}(t''),
\]

\[
\delta \mathcal{K}^{(n+1)}(t) = -D \mathcal{M}^{(n)}(t) D - \int_0^t dt' D \mathcal{M}^{(n)}(t - t') \delta \mathcal{K}^{(n+1)}(t'),
\]

and it is clear that they correspond to the iteration scheme of section 4.3. Then $S^{(n)}(t), -\delta \mathcal{K}^{(n)}(t), -\delta \mathcal{K}^{(n)}(t)$ are completely monotone. By construction $S^{(n)}(t = 0) = S$, $-\delta \mathcal{K}^{(n)}(t = 0) = D \mathcal{M}(t = 0) D$ etc and one infers $\|S^{(n)}(t)\| \leq \|S\|, \|\delta \mathcal{K}^{(n)}(t)\| \leq \|D\|\|\mathcal{M}\| := K_1, \|\delta \mathcal{K}^{(n)}(t)\| \leq K_1 L_1$. Then the following bounds follow for $0 \leq t \leq T < \infty$:

\[
\|S^{(n+1)}(t) - S^{(n)}(t)\| \leq \|D\| \|S^{-1}\| \int_0^t dt' \|S^{(n+1)}(t') - S^{(n)}(t')\| \\
+ T \|S^{-1}\| \|S\| \int_0^t dt' \|\delta \mathcal{K}^{(n)}(t') - \delta \mathcal{K}^{(n-1)}(t')\| \\
+ TK_1 L_1 \|S^{-1}\| \int_0^t dt' \|S^{(n+1)}(t') - S^{(n)}(t')\|,
\]

\[
\|\delta \mathcal{K}^{(n+1)}(t) - \delta \mathcal{K}^{(n)}(t)\| \leq \|D\|\|\mathcal{M}\| \int_0^t dt' \|\delta \mathcal{K}^{(n+1)}(t') - \delta \mathcal{K}^{(n)}(t')\| \\
+ \|D\|\|\mathcal{M}\| \int_0^t dt' \|\delta \mathcal{K}^{(n+1)}(t') - \delta \mathcal{K}^{(n)}(t')\|,
\]

The presence of the first term on the rhs of equation (43) requires us to adapt the strategy to find suitable bounds. Upon substituting the bounds from equation (42) the combination of both equations yields estimates for $X^{(n)}(t) = \|S^{(n+1)}(t) - S^{(n)}(t)\| + c\|\mathcal{K}^{(n+1)}(t) - \mathcal{K}^{(n)}(t)\|$, which now satisfies the inequalities

\[
X^{(n)}(t) \leq L \int_0^t dt' [X^{(n)}(t') + X^{(n-1)}(t') + X^{(n-2)}(t')],
\]

for $n \geq 2$. Combining the two equations yields estimates as in equation (39), with the constant now replaced by $L = (\|D\|\|S^{-1}\| + TL_1\|S\|\|S^{-1}\| + TK_1 L_1 \|S^{-1}\|)(1 + c\|D\| L_2) + c\|D\|\|\mathcal{M}\| + c\|D\| K_1 L_2 < \infty$. The sequence $X^{(n)}(t)$ is bounded by the

\[
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\]

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constants $K = 2\|S\| + 2cK_1$ and can be controlled by a suitable functional series $a^{(n)}(t)$ satisfying the inequalities

$$a^{(n)}(t) \geq L \int_0^t [2a^{(n-1)}(t') + a^{(n-2)}(t')] dt', \quad (45)$$

and $a^{(0)}(t) = a^{(1)}(t) = K$. First, $X^{(0)}(t) \leq a^{(0)}(t)$, $X^{(1)}(t) \leq a^{(1)}(t)$, $X^{(2)}(t) \leq a^{(2)}(t)$ and $X^{(1)}(t) \leq a^{(1)}(t)$, $X^{(2)}(t) \leq a^{(1)}(t)$ are obvious. Then $X^{(2)}(t) \leq L \int_0^t dt' [2a^{(1)}(t') + a^{(0)}(t')] \leq a^{(2)}(t)$. The procedure can be continued to show that $X^{(n)}(t) \leq a^{(m)}(t)$ for $m \leq n$. It remains to find a suitable bound $a^{(n)}(t)$ satisfying equation (45) such that $\sum a^{(n)}(t) < \infty$.

One can show by induction that $a^{(n)}(t) = K(\bar{L}/T)^{(n-1)/2}I((n+1)/2)$, $n \in \mathbb{N}$ will do, provided the constant $\bar{L} \geq L + 1/T$ is chosen large enough. The remainder of the proof follows the Newtonian case.

5. Generalized covariance principle and the maximum theorem

In this section we demonstrate that the MCT equations satisfy a covariance principle under a linear transformation of $S(q,t)$. This covariance allows us to prove the maximum theorem, which states that the long-time limit $S(q,t \to \infty)$, provided it exists, can be calculated as the maximal solution of a fixed-point equation. Solutions where these limits are identically zero are called ergodic and identified with fluid behavior, while non-ergodic ones correspond to (idealized) glass states. Then the maximum theorem allows calculation of the phase diagram and in particular the glass-transition lines without solving the dynamic equations explicitly. Furthermore, since the fixed-point equation is purely algebraic, the glass-transition lines originate as bifurcations associated with singular behavior in its vicinity.

For the case of a simple liquid these properties have been proven rigorously for Brownian dynamics [7], and the proofs can also be adapted for the Newtonian case, assuming the existence of a long-time limit of the ISF. For matrix-valued ISF and Brownian dynamics all properties can be properly generalized [9], relying on the notion of $C^*$ algebras. Here we shall show that for multiple decay channels the covariance principle can be extended such that the maximum theorem still holds, yet the transform for the memory kernels becomes nonlinear and non-local in time.

5.1. Fixed-point equation

In section 5, we have shown that the solutions $S(t)$ of the MCT equations are correlation functions; in particular, they are positive-semidefinite for all times. Then the long-time limit, also referred to as the glass form factor,

$$F(q) = \lim_{t \to \infty} S(q, t) \succeq 0, \quad (46)$$

inherits this property, provided the limit exists. While this is guaranteed for Brownian dynamics (the solutions are non-negative, monotonically decreasing), in the Newtonian case all numerical solutions suggest that the limit also exists, yet no rigorous proof appears to be available. To proceed, we shall hence assume that the limit exists also for Newtonian dynamics.
By the MCT functional, the memory kernel also displays a long-time limit
\[ \mathcal{N}(q) := \lim_{t \to \infty} \mathcal{M}(q,t) = \mathcal{F}[F,F;q] \geq 0, \] (47)
which is again positive-semidefinite, as it should be since microscopically the memory kernel is also a correlation function. In the Laplace domain a non-vanishing long-time limit corresponds to a simple pole at zero frequency \( \mathcal{M}(z) = -z^{-1} \mathcal{N}[1 + o(z^0)] \) for \( z \to 0 \), where we again, here and in the following, suppress the dependence on the wavenumber. Then the representation of the current correlator for Newtonian, equation (12), as well as for Brownian dynamics, equation (18), leads to the low-frequency behavior \( \hat{K}(z) = z \mathcal{N}^{-1} + o(z) \). Thus the contraction reads to leading order \( \hat{K}(z) = z \mathcal{N}^{-1} + o(z) \), with
\[ \mathcal{M}^{-1} = \sum_{\alpha \beta} q_{\alpha} \mathcal{M}^{-1}_{\alpha \beta} q_{\beta} \geq 0. \] (48)

Considered as a functional of the long-time limits, \( \mathcal{N}[F] \) displays the properties of an effective static mode-coupling functional in the space of matrices with mode indices \( \mu, \nu \), as shown in [14]. In particular, it preserves ordering \( \mathcal{N}[E] \geq \mathcal{N}[F] \) if \( E \geq F \).

The first equation of motion, equation (9), shows that \( \hat{S}(z) = -z^{-1} \mathcal{F}[1 + o(z^0)] \) for \( z \to 0 \), where the glass form factor now satisfies
\[ \mathcal{F} = \mathcal{S}^{-1} + \mathcal{N}[F]^{-1} \mathcal{S}^{-1} \] (49)

The equations (47), (48) and (49) constitute a closed set of equations called fixed-point equations and the long-time limit has to represent one of the solutions of these equations. The coupling of different wavenumbers originates now from the effective static mode-coupling functional. In general, the fixed-point equations allow for many solutions; in particular, zero is always a solution. Hence, a criterion for which of the solutions represents the long-time limit of the dynamic MCT equations is needed. For the case of single decay channels, the answer is provided by the maximum theorem, stating that the long-time limit is represented by the maximal solution [4, 5, 9]. More precisely, out of all solutions \( \mathcal{F} \geq 0 \) that are positive-semidefinite there is a unique solution \( \mathcal{F} \) that is larger than any of the other \( \mathcal{F}, \mathcal{F} \geq \mathcal{F} \), and this is the one corresponding to the glass form factor. The maximal solution can be found by a monotone iteration scheme.

For multiple decay channels, it has been shown [14] that the fixed-point equations admit a maximal solution, which again can be determined by an iteration scheme. Here we prove that it also corresponds to the long-time limit of the dynamic MCT equations.

5.2. Effective mode-coupling theory functional and generalized covariance principle

The strategy consists of properly generalizing the covariance principle demonstrated for one-component systems [7] and mixtures [9]. Therefore, we shall construct an effective dynamic mode-coupling functional \( \mathcal{M}(t) \) acting only in the space of matrices with mode indices. The effective dynamic functional will be shown to be a proper extension of the effective static functional used in section 5.1 to preserve the properties of correlation functions or completely monotone functions. This effective dynamic functional is then the suitable starting point to apply the covariance idea.
First, we consider Newtonian dynamics and define the kernel $\tilde{M}(z)$ in the Laplace domain via

$$zJ^{-1} + \tilde{M}(z)\allowbreak^{-1} := -\tilde{K}(z).$$

(50)

In order to keep the argument simple, the regular damping has been discarded, $D^{-1} \equiv 0$. In appendix A it is shown that $\tilde{M}(z)$ is well defined and again corresponds to a correlation function. The current kernel $\tilde{K}(z)$ itself is the contraction of $\tilde{K}(z)$, which is obtained from the memory kernels $M(z)$ by equation (12). Then $\tilde{M}(z)$ can be viewed itself as a functional of the intermediate scattering functions, and we refer to it as the effective dynamic MCT functional. Specializing equation (50) to the limit $z \to 0$ and going through the arguments of section 5.1 again shows that the effective dynamic functional reduces to the effective static functional $N$. Casting equation (50) in the temporal domain,

$$\int_0^t JM(t')K(t - t')\, dt' = -\partial_t K(t),$$

(51)

shows that given $K(t')$ for times $t' \leq t$ the effective MCT functional can be determined up to time $t$ by solving a standard (matrix-valued) Volterra integral equation of the first kind. Thus the effective dynamic MCT functional is a causal smooth functional of the intermediate scattering functions. Whereas the original memory kernel is local in time, this is in general no longer the case for the effective MCT functional originating from multiple relaxation kernels, as can be checked by a high-frequency expansion. To emphasize the dependence on the time-dependent past of the intermediate scattering function, we indicate the effective mode-coupling functional by $\tilde{J}$.

The next step is to reformulate the iteration scheme of section 5.1 in terms of the effective memory kernel. A simple calculation shows that the map

$$\mathcal{H} : (\tilde{S}(z), \tilde{M}(z)) \mapsto (-S/z + [zS^{-1} - z^2J^{-1} - z^2\tilde{M}(z)]^{-1}, \tilde{M}[\{S\}](z)),$$

(52)

generates the same sequence of approximants as above, provided it is initialized with $(\tilde{S}(0)(z) = -S/z, \tilde{M}(0)(z) = 0)$. In particular, the sequence was shown never to leave the space of correlation functions and to converge to the fixed-point solution of

$$\mathcal{H}(\tilde{S}(z), \tilde{M}(z)) = (\tilde{S}(z), \tilde{M}(z)),$$

which is the unique solution of the MCT equations.

Assume now that $\bar{F}$ with $S \geq \bar{F} \geq 0$ is a solution of the fixed-point equation, equations (47), (48) and (49), then we define a mapping

$$\unrhd : (S(t), M[\{S\}](t)) \mapsto (\bar{S}(t) = S(t) - \bar{F}, \bar{M}[\{\bar{S}\}](t) = M[\{S\}](t) - N[\bar{F}]).$$

(53)

To make progress, we make the assumption that long-time limits, $F = \lim_{t \to -\infty} S(t), \lim_{t \to -\infty} M(t)$, of the correlation functions $S(t), M(t)$ emerging as fixed points of the map, equation (52), exist. Analyzing the low-frequency behavior of equation (50) and employing the definition in equation (48) reveals that $\lim_{t \to -\infty} M(t) = N[F]$. The spectral representation, equation (2), further shows that they correspond to a jump in the spectral measure at zero frequency, $\lim_{t \to -\infty} S(t) = \mathcal{R}\{0\} \geq 0$, and is positive-semidefinite, and correspondingly for $M(t)$. Considering $\tilde{M}[\{\bar{S}\}]$ as a functional of $\bar{S}(t)$, we have to show that it maps correlation functions $\bar{S}(t)$ to correlation functions $\tilde{M}(t).$
Since the spectral measures $R_M, R_{\tilde{M}}$ associated with $M(t)$ and $\tilde{M}(t)$ differ only by a jump at zero frequency $R_{\tilde{M}}(\{0\}) = R_M(\{0\}) - \mathcal{N}[\vec{F}]$, what remains to be shown is that $\lim_{t \to \infty} \tilde{M}(t) = \lim_{t \to \infty} M(t) - \mathcal{N}[\vec{F}] = \mathcal{N}[F] - \mathcal{N}[\vec{F}] \succeq 0$. However, by assumption, $\tilde{S}(t)$ is also a correlation function, thus its long-time limit fulfills $\lim_{t \to \infty} \tilde{S}(t) = F - \vec{F} \succeq 0$. Then the desired property follows since the functional $\mathcal{N}[F]$ preserves ordering.

The next step is to construct a mapping $\mathcal{H}$ such that the diagram

$$
(\hat{S}(z), \hat{M}(z)) \xrightarrow{\mathcal{H}} \mathcal{H}\{\hat{S}(z), \hat{M}(z)\}
$$

commutes. Explicit calculation shows that this is fulfilled with

$$
\mathcal{H} : (\hat{S}(z), \hat{M}(z)) \mapsto (-\hat{S}/z + [z\hat{S}^{-1} - z^3J^{-1} - z^2\hat{M}(z)]^{-1}, \hat{M}\{\hat{S}\}(z)),
$$

and $\hat{S} = S - \vec{F}$, as follows from equation (53). The preceding equations show that the MCT equations are covariant with respect to the linear transform $\sim$, i.e. all equations assume the same form and acquire merely $\sim$. Note that the current matrix $J$ remains the same, which becomes obvious upon inspection of the short-time expansion, equation (15). Comparison with [14] shows that this mapping is the proper generalization of the renormalized mode-coupling functional to the case of finite frequencies $z$.

For Brownian motion the argument runs essentially along the same path. We define the effective memory kernel $\tilde{M}(z)$ in the Laplace domain for the Brownian case via

$$
\left[iD^{-1} + \tilde{M}(z)\right]^{-1} := -\hat{K}(z) = -iD - \delta \hat{K}(z).
$$

It is shown in appendix B that $\tilde{M}(z)$ is well defined and corresponds to a completely monotone function. Again, the effective memory kernel should be viewed as a non-local causal smooth functional of the intermediate scattering functions, $M\{\{S\}\}(t)$. However, its initial value fulfills that $D^{-1}M(t = 0)D^{-1}$ is merely the contraction of $D^{-1}M(t = 0)D^{-1}$. We also note that the definitions of the effective kernel for the Newtonian and Brownian cases differ from each other, nevertheless, considered as functionals of $S(t)$, they are expected to display the same long-time behavior. By construction they both reduce to the effective static functional in the long-time limit.

The mapping $\mathcal{H}$ differs only by notation from the Newtonian case

$$
\mathcal{H} : (\hat{S}(z), \hat{M}(z)) \mapsto (-\hat{S}/z + [z\hat{S}^{-1} - z^3iD^{-1} - z^2\hat{M}(z)]^{-1}, \hat{M}\{\hat{S}\}(z)).
$$

A fixed-point solution $\vec{F}$ induces the mapping $\sim$, as in equation (53). Provided $\tilde{S}(t)$ is again completely monotone, then $\tilde{M}(t)$ is also completely monotone, as can be seen in the following. By the representation theorem, $M(t) - \lim_{t \to \infty} M(t)$ is completely monotone (note that the long-time limit is guaranteed to exist), thus $\tilde{M}(t)$ is completely monotone if $\lim_{t \to \infty} M(t) - \mathcal{N}[\vec{F}] \succeq 0$. The remainder of the argument is as in the Newtonian case.

The covariance of the MCT equation expressed in terms of a commuting diagram, equation (54), is achieved by
\( \tilde{\mathcal{H}} : (\tilde{S}(z), \tilde{M}(z)) \mapsto (-\tilde{S}/z + [z\tilde{S}^{-1} - z^2 D^{-1} - z^2 \tilde{M}(z)]^{-1}, \tilde{M}[[\tilde{S}]](z)), \) (58)

and again \( \tilde{S} = S - \bar{F}. \)

5.3. Maximum principle

The covariance property shows that the problem of finding a solution \( S(t) \) of the MCT equations is the same as finding a solution \( \tilde{S}(t) = S(t) - \bar{F} \) for the mapped problem, provided \( \bar{F} \geq 0 \) is a solution of the fixed-point equations. Remember that for the Newtonian case we had to use the additional assumption that the long-time limit exists. However, the existence and uniqueness of such a solution has been shown by the iteration scheme in the temporal domain. Hence we conclude that \( \tilde{S}(t) \) is uniquely determined and corresponds to a correlation function (where the long-time limit exists by assumption) in the case of Newtonian dynamics and to a completely monotone function for Brownian dynamics.

For the long-time limits \( F = \lim_{t \to \infty} S(t) \geq 0, \tilde{F} = \lim_{t \to \infty} \tilde{S}(t) \geq 0 \) this implies the conclusion \( F \geq \tilde{F}. \) Hence the long-time limits are not smaller than any solution of the fixed-point equations. Yet the long-time limits are a solution of the fixed-point equations themselves, and therefore correspond to the maximal solution \( F \equiv \tilde{F}. \) The observation is referred to as the maximum property, and allows us to determine the long-time limit of the MCT equations by solving algebraic equations rather than coupled integrodifferential equations.

By using the effective MCT functional we have reduced the problem of discussing the mathematical properties of the solutions of the MCT to the case of a single decay channel. The corresponding results can be taken over directly. In particular, consider the linearization \( \Psi[\delta F] \) as obtained from

\[
[S^{-1} + N[F]]^{-1} - [S^{-1} + N[F + \delta F]]^{-1} = \Psi[\delta F] + O(\delta F)^2,
\]

(59)

where \( F \geq 0 \) is the maximal solution of the fixed-point equation. One readily shows that \( \Psi : \mathcal{A}_0^M \to \mathcal{A}_0^M \) is a linear positive map on a \( C^* \) algebra, i.e. it fulfils \( \Psi[\delta F] \geq 0 \) if \( \delta F \geq 0. \) In [9] it was shown by a proper extension of the Frobenius–Perron theorem that \( \Psi \) has a maximum non-degenerate eigenvalue not exceeding unity in the generic case where \( \Psi \) is irreducible. The manifold in the phase diagram where this eigenvalue becomes unity is identified with the glass-transition lines, and the associated bifurcation behavior has to be of the \( A_2 \) type according to the classification of Arnol’d [20]. The generic case corresponds to the fold bifurcation scenario \( A_2 \), and explicit expressions for the exponent parameter have been worked out [21, 22]. For the case of mixtures, asymptotic solutions with scaling properties hold in the vicinity of the glass-transition lines, similarly to the one-component case [21, 23].

6. Summary and conclusions

In general, a density mode can decay into more than one channel such that the associated particle current density naturally splits into different parts. The representation in terms of memory kernels then suggests we consider parallel relaxation, which implies a
mathematically different structure of the equations of motion. Within the mode-coupling approach, the force kernels are again approximated by positive superpositions of products of density correlation functions at the same instant of time. Our work provides proofs that the solutions of this generalized MCT still respect the constraints of probability theory and purely relaxational dynamics, extending the ideas introduced in the one-component case with a single decay channel [4, 6, 7] and their generalization to matrix-valued correlation functions [9].

The proof of existence and uniqueness of solutions relies on an iteration scheme similar to ordinary differential equations, and the generalization to multiple decay channels requires only mild adaptations. Similarly, the ideas to show that the iteration does not leave the space of correlation functions or completely monotone functions are directly transferable from the matrix case. The shown properties underline that the MCT approach is robust and encodes a series of natural requirements on any theory of dynamical phenomena without fine tuning.

Much stronger conclusions follow from a property observed by Götze known as the covariance principle [4, 24]. First a fixed-point equation is derived whose solutions are candidates for the long-time limits, i.e. glass form factors, then it is shown that the equations of motion determining the dynamics ‘on top’ of this candidate are of the same form as the original problem. Since the solutions of the transformed problem are again correlation functions or completely monotone, one can define a semi-ordering for glass form factors such that the long-time limits of the dynamic equations coincide with the maximal solution of the fixed-point equations.

While the matrix-valued theory displays this covariance property in a straightforward generalization of the single-component theory, this is no longer the case for multiple decay channels. In fact, one can convince oneself that application of a linear transform to the density correlation function and the MCT kernels does not lead to form-invariant equations of motion. The key idea has been to introduce an effective (single decay channel) MCT functional to map the theory to a case where the covariance principle is known to apply. On the level of the glass form factors, this was implemented in [14], and it was shown that the effective MCT functional preserves ordering. The extension to the dynamics comes at the price that the functional is no longer local in time, yet it remains causal in the sense that to determine the effective kernel at a given time knowledge of the density correlation functions at earlier times is sufficient. With this prerequisite, we have shown a generalized covariance principle, such that essentially all of the consequences elaborated for the one-component case can be taken over. In particular, the maximum property follows, stating that the long-time limits of the dynamic theory can be calculated without solving the dynamic equations explicitly. Rather, it is sufficient to compute the maximal solution of the fixed-point equation. This in turn permits us to distinguish ergodic liquid-like solutions from non-ergodic idealized states and thus to construct a non-equilibrium state diagram. The glass-transition lines are identified with bifurcations of the fixed-point equation, and they have to be of $A_\ell$ type in the classification by Arnol’d [20], the simplest being the $A_2$ fold bifurcation.

We have not pursued reproduction of properties such as a non-zero radius of convergence for a short-time expansion, the existence of a power series for small frequencies for the control parameter off the glass-transition lines, or the emergence of scaling laws in the vicinity of the glass transition. Since the effective mode-coupling functional is non-local
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in time, the proofs presented for the one- and multi-component liquids [4, 7, 9, 12] do not readily apply; however, it appears promising to repeat the essential steps on the level of the multiple channel description.

An additional term linear in the intermediate scattering function may be added in the mode-coupling functional, as it occurs naturally for the interaction of liquid particles with a frozen disorder; see for example [25]–[28]. Provided, the full functional still satisfies the constraints formulated in section 4.1, all conclusions for the general properties of the MCT solutions for the multichannel MCT remain unaffected. More generally, all nonlinear functionals of the intermediate scattering function which map correlation functions in the case of Newtonian dynamics, or purely relaxational functions for Brownian dynamics, onto functions with the same properties can be included without modification. Similarly, the extension of the proofs to the incoherent dynamics coupled to the collective motion should not require new ideas.

The MCT equations may be equipped by an additional \(a \text{ priori}\) known regular damping term, see equation (16), which captures fast processes that are not included within the MCT functional and are smoothly embedded into our analysis. Generically, intriguing glass-transition scenarios may occur, if one of the channels dominates the caging mechanism. For example partial freezing of only the translational degrees of freedom for dipolar hard spheres [10] or for the collective rotational degrees of freedom for elongated ellipsoids [29] has been uncovered.

The MCT equations for multiple decay channels may challenge the replica theory for structural glasses [32], which is a purely static theory. Besides predicting a thermodynamic glass transition of Kauzmann type, it also allows us to deduce a singularity for a dynamical transition by deriving an equation for the glass form factors. For a hard-sphere liquid in high dimensions \(d \to \infty\) it has recently been shown that the critical volume fraction \(\varphi^{\text{replica}}(d)\) from replica theory [33] and \(\varphi^{\text{MCT}}(d)\) from MCT [30, 31] do not coincide. Using a different closure for the calculation of the static input into replica theory, Szamel [34] has demonstrated that replica theory yields exactly the same equation for the glass form factors as MCT does. However, since replica theory lacks a dynamical origin, it seems impossible to prove that the long-time limit of the intermediate scattering functions is equal to the maximum solution for the glass form factors. In addition, since replica theory only considers density and not current density, its form probably does not involve a channel index. Consequently, its fixed-point equation for the glass form factors will not depend on the number of decay channels, in strong contrast to the corresponding equation from MCT. How replica theory and MCT in the case of more than one decay channel could be reconciled is far from being obvious.

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Appendix A. Representation theorem for Newtonian dynamics

In this appendix we show that a matrix-valued correlation function can be represented in the Laplace domain via another correlation function.

**Theorem 1.** If $\hat{K}(z)$ is the Fourier–Laplace transform of a matrix-valued correlation function with asymptotic expansion

$$
\hat{K}(z) = -z^{-1}J + O(z^{-3}), \quad z \to \infty, (A.1)
$$

with $J \succ 0$, then there exists a representation

$$
\hat{K}(z) = -[zJ^{-1} + \hat{M}(z)]^{-1}, (A.2)
$$

such that $\hat{M}(z)$ corresponds again to a correlation function.

**Proof.** The proof is a generalization of the corresponding property in the scalar case (see Akhiezer [35], p 111, lemma 3.3.6).

By the spectral representation theorem,

$$
\hat{K}(z) = \int (\Omega - z)^{-1} R_K(d\Omega), (A.3)
$$

with $R_K(\Omega)$ the associated self-adjoint matrix-valued measure. From the asymptotic expansion, equation (A.1), one infers $J = \int R_K(d\Omega)$ and by assumption $J$ is positive-definite. Next, solve formally for $\hat{M}(z)$

$$
\hat{M}(z) = -zJ^{-1} - \hat{K}(z)^{-1}, \quad z \in \mathbb{C}_+. (A.4)
$$

In order for this to be well defined, $\hat{K}(z)$ has to be invertible in the complex upper half-plane. Assume for the moment the contrary. Then, there is a vector $y$ not identically to zero and for a $z_1 \in \mathbb{C}_+$ with $\hat{K}(z_1)y = 0$, implying that the complex analytic function $y^\dagger \hat{K}(z)y$ displays a zero in the upper half-plane. However, the representation

$$
\text{Im}[\hat{K}(z)] = \int \frac{\text{Im}[z] R_K(d\Omega)}{|\Omega - z|^2} \geq 0 (A.5)
$$

reveals that $\text{Im}[y^\dagger \hat{K}(z)y] \geq 0$ is non-negative and harmonic, and by the mean-value property a zero can occur only if $y^\dagger \hat{K}(z)y \equiv 0$ for all $z \in \mathbb{C}_+$. The latter case is excluded by the assumption on the asymptotic expansion, equation (A.1).

Thus $\hat{M}(z)$ is well defined in the upper half-plane. To demonstrate that it corresponds to a correlation function, we show again properties (1)–(4) of section 2. The explicit expression, equation (A.4), then shows that $\hat{M}(z)$ is analytic in $\mathbb{C}_+$, proving property (1). Property (2) follows immediately from equation (A.4) and $\hat{K}(-z^*) = -\hat{K}(z)^\dagger$, since $K(t)$ is a correlation function. From the asymptotic expansion, equation (A.1), one finds

$$
\hat{M}(z) = -zJ^{-1} - [z^{-1}J + O(z^{-3})]^{-1}
= -zJ^{-1} + zJ^{-1}[1 + O(z^{-2})]^{-1} = O(z^{-1}), (A.6)
$$

which implies property (3). From, equation (A.4), the imaginary part can be represented as

$$
\text{Im}[\hat{M}(z)] = -J^{-1} \text{Im}[z] + \hat{K}^\dagger(z)^{-1} \text{Im}[\hat{K}(z)] \hat{K}(z)^{-1}. (A.7)
$$

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The right-hand side is positive-semidefinite as can be seen in the following. For any vector \( y \) define the vector
\[
\textbf{\( x(\Omega) = (\Omega - z)^{-1} \hat{\textbf{K}}(z)^{-1} y J^{-1} y \).}
\]
(A.8)

Then \( \int x(\Omega)^\dagger R_K (d\Omega)x(\Omega) \geq 0 \). Upon expanding, one finds
\[
\textbf{\( y^\dagger \hat{\textbf{K}}^\dagger(z)^{-1} \int \frac{R_K (d\Omega)}{|\Omega - z|^2} \hat{\textbf{K}}(z)^{-1} y - y^\dagger J^{-1} y \geq 0 \).}
\]
(A.9)

Last, the representation of \( \text{Im}[\hat{\textbf{K}}(z)] \), equation (A.5), shows that the kernel \( \hat{\textbf{M}}(z) \) displays a positive-semidefinite imaginary part, and property (4) also holds. □

Appendix B. Representation theorem for Brownian dynamics

In this appendix we show that the effective memory kernel
\[
\hat{\textbf{M}}(z) = -\hat{\textbf{K}}(z)^{-1} - iD^{-1},
\]
(B.1)
is well defined for frequencies \( z \in \mathbb{C} \setminus i\mathbb{R}^- \) and is completely monotone provided \( -\delta \hat{\textbf{K}}(z) = iD - \hat{\textbf{K}}(z) \) shares the same property and \( \text{Im}[\hat{\textbf{K}}(z)] > 0 \) for \( z \in \mathbb{C}_+ \).

Proof. If \( \delta \hat{\textbf{K}}(z) \equiv 0 \) then \( \hat{\textbf{M}}(z) \equiv 0 \) and nothing needs to be shown.

First, by assumption \( \text{Im}[\hat{\textbf{K}}(z)] > 0 \) for \( z \in \mathbb{C}_+ \) and \( \hat{\textbf{K}}(z) \) is invertible also for \( z \in \mathbb{C} \setminus i\mathbb{R}^- \). Assume for the moment the contrary. Then there is a non-zero vector \( y \) with \( [iD + \delta \hat{\textbf{K}}(z)] y = 0 \). However, since \( -\delta \hat{\textbf{K}}(z) \) corresponds to a completely monotone function, there is a self-adjoint complex measure \( \mathbf{a}(\gamma) \) with
\[
\text{Re}\left[\int_0^\infty \frac{-1}{z + 1\gamma} \mathbf{a}(d\gamma)\right] = \int_0^\infty \frac{\text{Re}[z]}{|z + 1\gamma|^2} \mathbf{y}\dagger \mathbf{y} \mathbf{a}(d\gamma) \neq 0 \quad \text{for Re}[z] \neq 0
\]
(B.2)

which shows that \( \hat{\textbf{K}}(z)^{-1} \) is well defined now for all \( z \in \mathbb{C} \setminus i\mathbb{R}^- \).

The analytic properties (1*), (2*) are inherited from the corresponding ones for \( \delta \hat{\textbf{K}}(z) \). Since \( \lim_{z \to \infty} \hat{\textbf{K}}(z) = iD \), one finds \( \lim_{z \to \infty} \hat{\textbf{M}}(z) = 0 \), which implies (3*). Last, (4*) follows from
\[
\text{Re}[\hat{\textbf{M}}(z)] = \hat{\textbf{K}}_\dagger(z)^{-1} \text{Re}[-\delta \hat{\textbf{K}}(z)] \hat{\textbf{K}}(z)^{-1} \succeq 0,
\]
(B.4)
for \( \text{Re}[z] < 0 \), since \( -\delta \hat{\textbf{K}}(z) \) corresponds to a completely monotone function. By the representation theorem, properties (1*)–(4*) for \( \hat{\textbf{M}}(z) \) imply that it corresponds to the Fourier–Laplace transform of a completely monotone (matrix-valued) function. □

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