Notes: Nevanlinna analytical Continuation Method

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Abstract

This is the abstract.

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I. THE ANALYTIC CONTINUATION PROBLEM

The analytic continuation problem seeks to extract real frequency dynamical information from imaginary-time correlation functions $C(\tau)$ data. Technically, this is a highly nontrivial task[1]. To see this, we use the relation between $C(\tau)$ and $A(\omega)$ [1, 2]:

$$C(\tau) = \int_{-\infty}^{\infty} d\omega \frac{e^{-\tau\omega}}{1 - \lambda e^{-\beta\omega}} A(\omega) = \int_{-\infty}^{\infty} d\omega K(\tau, \omega) A(\omega)$$
 (1)

where $K(\tau, \omega) = \frac{e^{-\tau \omega}}{1 - \lambda e^{-\beta \omega}}$ is the kernel, $\lambda = \pm 1$ for bosons/fermions respectively. One may consider to solve the problem by firstly discretize τ and ω and get:

$$C(\tau_i) = \sum_{j=1}^{N_\omega} K_{ij} A(\omega_j)$$
 (2)

Then do SVD decomposition of rectangular matrix K, write $K_i j = U_{il} \lambda_l V_{lj}$. Finally the spectral function reads

$$A(\omega_j) = \sum_{l=1}^{N_\tau} \frac{1}{\lambda_l} V_{ij} \sum_{i=1}^{N_\omega} C(\tau_i) U_{il}$$
(3)

It seems fine at the first glanse. However, if we consider the properties of $K(\tau, \omega)$, we would notice that it is highly sigular since it is exponentially small for large $|\omega|$, so small errors $C(\tau)$ would be amplified by exponentially small λ_l . This problem is well-known ill-posed[3, 4] and enormous efforts have been made[].

II. HOW TO SOLVE?

. . .

Here we introduce the recently developed Nevanlinna analytic continuation method[5].

III. NEVANLINNA ANALYTIC CONTINUATION METHOD

The Nevanlinna analytic continuation method[5] is an interpolation method. The key step is to build the conformal mappings from the open upper half of the complex plane C^+ to a closed unit disk $\bar{\mathcal{D}}$ in the complex plane and make use of the Schur algorithm [6–8] to do the interpolate.

A. Schur Algorithm

Schur Algorithm was introduced by I. Schur[9] in Section 1 of Ref.[6]. Here we list the main results we need while for a detailed introduction, see Ref.[8].

A Schur class(\mathcal{S}) consists of the Schur functions, which are the holomorphic functions from the open unit disk \mathcal{D} to the closed unit disk $\bar{\mathcal{D}}$. For a given Schur function $s_0(z)$, the Schur algorithm defines a set of $\{s_j(z) \in \mathcal{S}\}_{0 \leq j < \infty}$ starting from $s_0(z)$ by the recurrence relation:

$$zs_{j+1}(z) = \frac{s_j(z) - \gamma_j}{1 - \gamma_j^* s_j(z)}$$
(4)

where $s_j \in \mathcal{S}$ and $\gamma_j \equiv s_j(0)$ are called Schur parameters and $|\gamma_j| \leq 1$.

On the other hand, given an arbitrary strictly contractive sequence of Schur parameters $\{\gamma_0, \gamma_1, \ldots, \gamma_j, \ldots\} \subset \mathcal{D}$, one can construct a unique Schur function $s_0(z)$ by means of a continued fraction algorithm. In which we use the inverse relation of eq. (4)

$$s_j(z) = \frac{\gamma_j + z s_{j+1}(z)}{1 + \gamma_j^* z s_{j+1}(z)}$$
(5)

to construct the *n*-th Schur approximant, which we will denote by $s_0(z; \gamma_0, \gamma_1, \dots, \gamma_n)$. Namely, we write:

$$s_n(z;\gamma_n) = \gamma_n \tag{6}$$

$$s_j(z;\gamma_j,\gamma_{j+1},\ldots\gamma_n) = \frac{\gamma_j + zs_{j+1}(z;\gamma_{j+1},\ldots\gamma_n)}{1 + \gamma_j^* zs_{j+1}(z;\gamma_{k+1},\ldots\gamma_n)}$$
(7)

where $j = n - 1, n - 2, \dots, 1, 0$.

Given the initial data consisting of N points $\{\mathcal{Y}_0, \mathcal{Y}_1, \dots, \mathcal{Y}_{N-1}\} \subset \mathcal{D}$ and target data $\{\mathcal{C}_0, \mathcal{C}_1, \dots, \mathcal{C}_{N-1}\} \subset \mathcal{D}$, we can find a holomorphic function $s(z) : \mathcal{D} \to \bar{\mathcal{D}}$ such that $s(\mathcal{Y}_j) = \mathcal{C}_j$ for all j by combining eq. (6), eq. (7) and the linear fractional transform:

$$\xi(z, \mathcal{Y}_j) = \frac{z - \mathcal{Y}_j}{1 - z\mathcal{Y}_j^*} \tag{8}$$

and s(z) has the following recursion relation:

$$\tilde{s}_{j}(\tilde{z}) = \frac{C_{j} + \xi(\tilde{z}, \mathcal{Y}_{j})\tilde{s}_{j+1}(\tilde{z})}{1 + C_{j}^{*}\xi(\tilde{z}, \mathcal{Y}_{j})\tilde{s}_{j+1}(\tilde{z})}$$

$$(9)$$

Proof. eq. (8) builds a one to one correspondence of a Schur function s(z) with another Schur function $\tilde{s}(z)$.

$$\tilde{s}(\xi^{-1}(z,\mathcal{Y})) = s(z), \text{ or } s(\xi(z,\mathcal{Y})) = \tilde{s}(z)$$
 (10)

From eq. (10), we have:

$$\gamma_j = s_j(0) = \tilde{s}_j(\xi^{-1}(0, \mathcal{Y})) = \tilde{s}_j(\mathcal{Y}_j) = \mathcal{C}_j$$
(11)

inserting eq. (10) and eq. (11) to eq. (5), we have:

$$\tilde{s}_{j}(\xi^{-1}(z, \mathcal{Y}_{j})) = \frac{C_{j} + z\tilde{s}_{j+1}(\xi^{-1}(z, \mathcal{Y}_{j}))}{1 + C_{i}^{*}z\tilde{s}_{j+1}(\xi^{-1}(z, \mathcal{Y}_{j}))}$$
(12)

replace $\xi^{-1}(z, \mathcal{Y}_j) \in \mathcal{D}$ by $\tilde{z} \in D$, we have:

$$\tilde{s}_{j}(\tilde{z}) = \frac{C_{j} + \xi(\tilde{z}, \mathcal{Y}_{j})\tilde{s}_{j+1}(\tilde{z})}{1 + C_{j}^{*}\xi(\tilde{z}, \mathcal{Y}_{j})\tilde{s}_{j+1}(\tilde{z})}$$

$$(13)$$

Therefore we have:

$$s_{N-1}(z; \mathcal{C}_{N-1}) = \frac{\mathcal{C}_{N-1} + \xi(z, \mathcal{Y}_{N-1}) s_N(z)}{1 + \mathcal{C}_{N-1}^* \xi(z, \mathcal{Y}_{N-1}) s_N(z)}$$
(14)

$$s_j(z; \mathcal{C}_j, \mathcal{C}_{j+1}, \dots \mathcal{C}_N) = \frac{\mathcal{C}_j + \xi(z, \mathcal{Y}_j) s_{j+1}(z; \mathcal{C}_{j+1}, \dots \mathcal{C}_N)}{1 + \mathcal{C}_i^* \xi(z, \mathcal{Y}_j) s_{j+1}(z; \mathcal{C}_{k+1}, \dots \mathcal{C}_N)}$$
(15)

where j = N - 2, N - 3, ..., 1, 0 and $s_0(z) \equiv s(z)$. In eq. (14), we notice that there is an degrees of freedom to choose an arbitrary $s_N(z) \in \mathcal{S}$, eq. (6) correponds the special case $s_{n+1}(z) = 0$.

G. Pick and R. Nevanlinna studied the interpolation problem independently in 1917[10] and 1919[11] respectively, showing that an interpolating function exists if and only if the Pick matrix

$$P_{jk} = \frac{1 - \mathcal{C}_k^* \mathcal{C}_j}{1 - \mathcal{Y}_j^* \mathcal{Y}_k} \tag{16}$$

is positive semi-definite. Furthermore, the function s(z) is unique if and only if the Pick matrix has zero determinant. It is called the Nevanlinna-Pick theorem.

B. Generalized Schur Algorithm

Schur algorithm can be modified to expand all contractive functions $(\in \mathcal{B})[12]$, which are holomorphic functions mapping from the upper half plane \mathcal{C}^+ to $\bar{\mathcal{D}}$.

Given the initial data consisting of N points $\{\mathcal{Y}_0, \mathcal{Y}_1, \dots, \mathcal{Y}_{N-1}\} \subset \mathcal{C}^+$ and target data $\{\mathcal{C}_0, \mathcal{C}_1, \dots, \mathcal{C}_{N-1}\} \subset \bar{\mathcal{D}}$, in order to find a holomorphic function $\theta(z) \in \mathcal{B}$ such that $\theta(\mathcal{Y}_j) = z_j$

for all j, we should make use of the Mobius transform $h(z, \mathcal{Y}) = \frac{z-\mathcal{Y}}{z-\mathcal{Y}^*}$ which maps $\mathcal{C}^+/\bar{\mathcal{C}}^+$ to $\mathcal{D}/\bar{\mathcal{D}}$, which means it establishes a one-to-one correspondence of $\theta(z)$ to a schur function s(z) with:

$$\theta(h^{-1}(z,\mathcal{Y})) = s(z), \text{ or } s(h(z,\mathcal{Y})) = \theta(z)$$
 (17)

We denote $h(z, \mathcal{Y}_j)$ as $h_j(z)$ form now on.

The recursion relation between $\theta_j(z)$ and the next contractive function $\theta_{j+1}(z)$ can be easily build as follows. From eq. (17), we have:

$$s_j(0) = \theta_j(h_j^{-1}(0)) = \theta_j(\mathcal{Y}_j) = \gamma_j$$
(18)

Let $\theta_{j+1}(z) = s_{j+1}(h_j(z))$, then use the recursion relation eq. (4), we have:

$$z\theta_{j+1}(h_j^{-1}(z)) = \frac{\theta_j(h_j^{-1}(z)) - \gamma_j}{1 - \gamma_i^* \theta_j(h_j^{-1}(z))} \stackrel{\text{def}}{=} \phi_j(h_j^{-1}(z))$$
(19)

Form the first and the third terms of eq. (19) we have:

$$\phi_j(h_j^{-1}(z)) = z\theta_{j+1}(h_j^{-1}(z)) = h_j(h_j^{-1}(z))\theta_{j+1}(h_j^{-1}(z))$$
(20)

replace $h_j^{-1}(z)$ with $z \in \mathcal{D}$ by $z \in \mathcal{C}^+$, we have

$$\phi_i(z) = h_i(z)\theta_{i+1}(z) \tag{21}$$

We can read from eq. (21) that $\phi_j(z) \in \mathcal{B}$ and $\phi_j(\mathcal{Y}_j) = 0$.

Form the second and the third terms of eq. (19) we can read:

$$\phi_j(z) = \frac{\theta_j(z) - \gamma_j}{1 - \gamma_j^* \theta_j(z)} \tag{22}$$

Together with eq. (21) we get the recursion relation between $\theta_j(z)$ and $\theta_{j+1}(z)$:

$$\theta_j(z) = \frac{\phi_j(z) + \gamma_j}{\gamma_j^* \phi_j(z) + 1} = \frac{h_j(z)\theta_{j+1}(z) + \gamma_j}{\gamma_j^* h_j(z)\theta_{j+1}(z) + 1}$$
(23)

The recursive final $\theta(z)$ can conveniently be written in a matrix form:

$$\theta(z)[z;\theta_N(z)] = \frac{a(z)\theta_N(z) + b(z)}{c(z)\theta_N(z) + d(z)}$$
(24)

where

$$\begin{pmatrix} a(z) & b(z) \\ c(z) & d(z) \end{pmatrix} = \prod_{j=1}^{N-1} \begin{pmatrix} h_j(z) & \gamma_j \\ \gamma_j^* h_j(z) & 1 \end{pmatrix}$$
 (25)

with j increasing from left to right. Like in eq. (6), there is a alse freedom to choose $\theta_N(z)$.

C. Interpolation of Green's functions

The retared Green's function $G^R(\omega + i\eta)$ and the Masubara Green's function $G(i\omega_n)$ can be expressed consistently by replacing the variables $i\omega_n$ and $\omega + i\eta$ with a single complex variable z. G(z) is analytic in the upper half plane \mathcal{C}^+ . Our problem is that once we have Masubara frequencies $\{i\omega_n\} \subset \mathcal{C}^+$ and target data $\{G(i\omega_n)\} \subset \mathcal{C}$, where \mathcal{C} is the complex plane, how can we get interpolate them and get the holomorphic function $G(z): \mathcal{C}^+ \to C$?

Based on the knowledge of Schur algorithm, if we can find a one-to-one correspondence of G(z) and a contractive function $\theta(z) \in \mathcal{B}$, then we can futher generalize the algorithm in section III B.

To do this, we firstly introduce the Nevanlinna functions $f(z) \in \mathcal{N}$. In complex analysis, a Nevanlinna function is a complex function that is analytic in the open upper half plane \mathcal{C}^+ and has non-negative imaginary part, i.e., maps into $\bar{\mathcal{C}}^+$ (the overline denotes inclusion of the boundary). The invertible Möbius transform $h(z) = \frac{z-i}{z+i}$ maps Nevanlinna functions one to one to contractive functions:

$$\theta(z) = h(f(z)), \text{ or } f(z) = h^{-1}(\theta(z))$$
 (26)

Given the initial data consisting of N points $\{\mathcal{Y}_0, \mathcal{Y}_1, \dots, \mathcal{Y}_{N-1}\} \subset \mathcal{C}^+$ and target data $\{C_0, C_1, \dots, C_{N-1}\} \subset \bar{\mathcal{C}}$, The only thing we needed is to let γ_j in eq. (23) be $\gamma_j \equiv h(C_j)$.

Moreover, the corresponding Pick matrix is generalized to:

$$P_{jk} = \frac{1 - h(C_k)^* h(C_j)}{1 - h(\mathcal{Y}_j)^* h(\mathcal{Y}_k)}$$
(27)

The aforementioned conformaling mappings are shown in fig. 1.

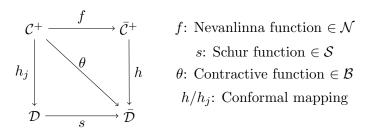


FIG. 1. Conformal mappings

Now we are ready to disscuss the Green's functions. The Lehmann representation of

Green's function G(z) is:

$$G(z) = \frac{1}{Z} \sum_{nm} |A_{nm}|^2 \frac{e^{-\beta E_n} \pm e^{-\beta E_m}}{z - E_m + E_n}$$
 (28)

where the "+" sign is for fermionic Green's functions and "-" sign is for bosonic Green's functions.

1. Fermionic case

In fermionic case, if we take z=x+iy with y>0, i.e. $z\in\mathcal{C}^+$, we can easily prove that $\mathrm{Im}G(z)\leq 0$. Therefore $-G(z)\in\mathcal{N}$ is a Nevanlinna function.

Proof.

$$ImG(x+iy) = -\frac{1}{Z} \sum_{nm} \frac{Q_{nm}^F y}{(x - E_m + E_n)^2 + y^2}$$
 (29)

where

$$Q_{nm}^{F} = |A_{nm}|^{2} (e^{-\beta E_{m}} + e^{-\beta E_{n}}) \ge 0$$
(30)

$$\therefore \operatorname{Im} G(x+iy) \leq 0 \text{ for } y > 0.$$

2. Bosonic case

The bosonic case is less trivial. The imaginary part of bosonic Green's function is:

$$ImG(x+iy) = -\frac{1}{Z} \sum_{nm} \frac{Q_{nm}^B y}{(x - E_m + E_n)^2 + y^2}$$
(31)

where

$$Q_{nm}^{B} = |A_{nm}|^{2} (e^{-\beta E_{m}} - e^{-\beta E_{n}}) = \operatorname{Sgn}(E_{m} - E_{n}) |Q_{nm}^{B}|$$
(32)

 $\operatorname{Im} G(x+iy)$ doesn't have a definite sign.

However, we can prove that $\tilde{G}(z) = -zG(z)$ is a Nevanlinna function.

Proof.

$$\operatorname{Im} \tilde{G}(x+iy) = -\operatorname{Im}[(x+iy)G(x+iy)]$$

$$= -x\operatorname{Im}G(x+iy) - y\operatorname{Re}G(x+iy)$$
(33)

where:

$$ReG(z) = \frac{1}{Z} \sum_{nm} \frac{Q_{nm}^B (x - E_m + E_n)}{(x - E_m + E_n)^2 + y^2}$$
(34)

For given (n, m)-term:

$$\tilde{\mathcal{Q}}_{nm}^{B} = xy\mathcal{Q}_{nm}^{B} - (x - E_m + E_n)y\mathcal{Q}_{nm}^{B}$$

$$= y(E_m - E_n)\mathcal{Q}_{nm}^{B}$$

$$= y|E_m - E_n||\mathcal{Q}_{nm}^{B}|$$
(35)

 $\therefore \tilde{\mathcal{Q}}_{nm}^B \ge 0 \text{ for } y > 0.$

$$\therefore \operatorname{Im} \tilde{G}(x+iy) = \frac{1}{Z} \sum_{nm} \frac{\tilde{\mathcal{Q}}_{nm}^B}{(x-E_m+E_n)^2 + y^2} \ge 0$$
 (36)

and $\tilde{G}(z)$ is a Nevanlinna function.

Moreover, $-i\omega_n G(i\omega_n)$ is the Fourier transform of $\frac{\mathrm{d}G(\tau)}{\mathrm{d}\tau}$.

We can perform analytical continuation $z = \omega + i\eta$ on $\tilde{G}(z)$:

$$2\operatorname{Im}\tilde{G}(\omega + i\eta) = \omega[-2\operatorname{Im}G(\omega + i\eta)] - 2\eta\operatorname{Re}G(\omega + i\eta)$$

$$\xrightarrow{\eta \to 0^{+}} \omega A(\omega)$$
(37)

where $A(\omega)$ is the spectral function.

IV. HARDY BASIS OPTIMIZATION

Appendix A: Conformal transforms

1. The linear fractional transform

The linear fractional transform is:

$$\xi(z, \mathcal{Y}) = \frac{z - \mathcal{Y}}{1 - z\mathcal{Y}^*} \tag{A1}$$

It is a one to one mapping of the open unit disk \mathcal{D} onto itself and a one to one mapping of the unit circle \mathcal{T} . It maps point \mathcal{Y} to the center of \mathcal{D} .

2. The Mobius transform

The mapping from $C^+/\bar{C^+}$ to $D^+/\bar{D^+}$ is called Mobius transform. It has the form:

$$h(z, \mathcal{Y}) = \frac{z - \mathcal{Y}}{z - \mathcal{Y}^*} \tag{A2}$$

where $\mathcal{Y} \in \bar{\mathcal{C}}^+$ and $\mathcal{Y} \neq 0$. We can easily prove that $|h(z,\mathcal{Y})| \leq 1$ for $z \in \bar{\mathcal{C}}^+$ and $|h(z,\mathcal{Y})| = 1$ if z is real. $h(z,\mathcal{Y})$ maps $\mathcal{Y} \in \bar{\mathcal{C}}^+$ to the center of the unit disk \mathcal{D} and the real axis as the edge of $\bar{\mathcal{D}}$, the rest part of upper half complex plane is wrapped inside the unit disk. If $\tilde{z} \in \mathcal{D}$, the inverse transform is:

$$h^{-1}(\tilde{z}, \mathcal{Y}) = \frac{\mathcal{Y} - \tilde{z}\mathcal{Y}^*}{1 - \tilde{z}}$$
(A3)

Angin one can prove $\operatorname{Im} h^{-1}(\tilde{z}, \mathcal{Y}) = (\operatorname{Im} \mathcal{Y})(1 - |\tilde{z}|^2) > 0.$

Proof of $|h(z,\mathcal{Y})| \leq 1$ for $z \in \bar{\mathcal{C}}^+$ and $|h(z,\mathcal{Y})| = 1$ if z is real. We already know that $\mathrm{Im} z \geq 0, \mathrm{Im} \mathcal{Y} > 0$.

$$|h(z,\mathcal{Y})|^{2} = \frac{z - \mathcal{Y}}{z - \mathcal{Y}^{*}} \frac{z^{*} - \mathcal{Y}^{*}}{z^{*} - \mathcal{Y}} = \frac{|z|^{2} + |\mathcal{Y}|^{2} - z\mathcal{Y}^{*} - z^{*}\mathcal{Y}}{|z|^{2} + |\mathcal{Y}|^{2} - z\mathcal{Y} - z^{*}\mathcal{Y}^{*}}$$

$$= \frac{|z|^{2} + |\mathcal{Y}|^{2} - 2(\operatorname{RezRe}\mathcal{Y} + \operatorname{Im}z\operatorname{Im}\mathcal{Y})}{|z|^{2} + |\mathcal{Y}|^{2} - 2(\operatorname{Re}z\operatorname{Re}\mathcal{Y} - \operatorname{Im}z\operatorname{Im}\mathcal{Y})}$$
(A4)

If Im z = 0, $|h(z, \mathcal{Y})|^2 = 1$. If Im z > 0, $|h(z, \mathcal{Y})|^2 < 1$. And we notice that if $\text{Im} \mathcal{Y} = 0$, we map all points in $\bar{\mathcal{C}}$ to point 1 except for point \mathcal{Y} itself.

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