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Quantum metrology with unitary parametrization processes

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Jing Liu¹, Xiao-Xing Jing¹ & Xiaoguang Wang^{1,2}

¹Zhejiang Institute of Modern Physics, Department of Physics, Zhejiang University, Hangzhou 310027, China, ²Synergetic Innovation Center of Quantum Information and Quantum Physics, University of Science and Technology of China, Hefei, Anhui 230026, China.

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Correspondence and
requests for materials
should be addressed to
X.G.W. (xgwang@
zimp.zju.edu.cn)

Quantum Fisher information is a central quantity in quantum metrology. We discuss an alternative representation of quantum Fisher information for unitary parametrization processes. In this representation, all information of parametrization transformation, i.e., the entire dynamical information, is totally involved in a Hermitian operator \mathcal{H} . Utilizing this representation, quantum Fisher information is only determined by \mathcal{H} and the initial state. Furthermore, \mathcal{H} can be expressed in an expanded form. The highlights of this form is that it can bring great convenience during the calculation for the Hamiltonians owning recursive commutations with their partial derivative. We apply this representation in a collective spin system and show the specific expression of \mathcal{H} . For a simple case, a spin-half system, the quantum Fisher information is given and the optimal states to access maximum quantum Fisher information are found. Moreover, for an exponential form initial state, an analytical expression of quantum Fisher information by \mathcal{H} operator is provided. The multiparameter quantum metrology is also considered and discussed utilizing this representation.

How to precisely measure the values of physical quantities, such as the phases of light in interferometers, magnetic strength, gravity and so on, is always an important topic in physics. Obtaining high-precision values of these quantities will not only bring an obvious advantage in applied sciences, including the atomic clocks, physical geography, civil navigation and even military industry, but also accelerate the development of fundamental theories. One vivid example is the search for gravitational waves. Quantum metrology is such a field attempting to find optimal methods to offer highest precision of a parameter that under estimation. In recently decades, many protocols and strategies have been proposed and realized to improve the precisions of various parameters^{1–22}. Some of them can even approach to the Heisenberg limit, a limit given by the quantum mechanics, showing the power of quantum metrology.

Quantum Fisher information is important in quantum metrology because it depicts the theoretical lowest bound of the parameter's variance according to Cramér-Rao inequality^{23,24}. The quantum Fisher information for parameter α is defined as $F = \text{Tr}(\rho L^2)$, where ρ is a density matrix dependent on α and L is the symmetric logarithmic derivative (SLD) operator and determined by the equation $\partial_\alpha \rho = (\rho L + L \rho)/2$. For a multiparameter system, the counterpart of quantum Fisher information is called quantum Fisher information matrix \mathcal{F} , of which the element is defined as $\mathcal{F}_{\alpha\beta} = \text{Tr}(\rho \{L_\alpha, L_\beta\})$, where L_α, L_β are the SLD operators for parameters α and β , respectively.

Recently, it has been found²⁷ that quantum Fisher information can be expressed in an alternative representation, that all information of parametrization process in quantum Fisher information is involved in a Hermitian operator \mathcal{H} . This operator characterizes the dynamical property of the parametrization process, and totally independent of the selection of initial states. Utilizing this representation, the quantum Fisher information is only determined by \mathcal{H} and the initial state.

In this report, we give a general expression of quantum Fisher information and quantum Fisher information matrix utilizing \mathcal{H} operator. For a unitary parametrization process, \mathcal{H} can be expressed in an expanded form. This form is particularly useful when the Hamiltonian owns a recursive commutation relation with its derivative on parameter estimation. We calculate the specific expression of \mathcal{H} in a collective spin system, and provide an analytical expression of quantum Fisher information in a spin-half system for any initial state. Based on this expression, all optimal states to access maximum quantum Fisher information are found in this system. Furthermore, considering this spin-half system as a multiparameter system, the quantum Fisher information matrix, can be easily obtained by the known form of \mathcal{H} in single parameter estimations. On the other hand,



inspired by a recent work²⁸, for an exponential form initial state, we provide an analytical expression of quantum Fisher information using \mathcal{H} operator. A demonstration with a spin thermal initial state is given in this scenario. The maximum quantum Fisher information and the optimal condition are also discussed.

Results

Quantum Fisher information with \mathcal{H} operator. For a general unitary parametrization transformation, the parametrized state $\rho(\alpha)$ is expressed by $\rho(\alpha) = U(\alpha)\rho_0U^\dagger(\alpha)$, where ρ_0 is a state independent of α . In this paper, since the parameter α is only brought by $U(\alpha)$, not the initial state ρ_0 , we use U instead of $U(\alpha)$ for short. Denote the spectral decomposition of ρ_0 as $\rho_0 = \sum_{i=1}^M p_i |\psi_i\rangle\langle\psi_i|$, where p_i and $|\psi_i\rangle$ are the i th eigenvalue and eigenstate of ρ_0 and M is the dimension of the support of ρ_0 . It is easy to see that p_i and $U|\psi_i\rangle$ are the corresponding eigenvalue and eigenstate of $\rho(\alpha)$, respectively. The quantum Fisher information of $\rho(\alpha)$ can then be expressed by^{29,30}

$$F = \sum_{i=1}^M 4p_i \langle \Delta^2 \mathcal{H} \rangle_i - \sum_{i \neq j} \frac{8p_i p_j}{p_i + p_j} |\langle \psi_i | \mathcal{H} | \psi_j \rangle|^2, \quad (1)$$

where^{25,26}

$$\mathcal{H} := i(\partial_\alpha U^\dagger)U \quad (2)$$

is a Hermitian operator since the equality $(\partial_\alpha U^\dagger)U = -U^\dagger(\partial_\alpha U)$. Meanwhile,

$$\langle \Delta^2 \mathcal{H} \rangle_i = \langle \psi_i | \mathcal{H}^2 | \psi_i \rangle - \langle \psi_i | \mathcal{H} | \psi_i \rangle^2 \quad (3)$$

is the variance of \mathcal{H} on the i th eigenstate of ρ_0 . When $\partial_\alpha U$ commutes with U , \mathcal{H} can be explained as the generator of the parametrization transformation²⁷. The expression (1) of quantum Fisher information is not just a formalized representation. The operator \mathcal{H} is only determined by the parametrization process, that is the dynamics of the system or the device. For a known dynamical process of a parameter, i.e., known system's Hamiltonian, \mathcal{H} is a settled operator and can be obtained in principle. In this representation, the calculation of quantum Fisher information is separated into two parts: the diagonalization of initial state and calculation of \mathcal{H} . For a general 2-dimensional state, the quantum Fisher information reduces to

$$F_{\text{qubit}} = 4(2\text{Tr}\rho^2 - 1) \langle \Delta^2 \mathcal{H} \rangle_{1(2)}. \quad (4)$$

The subscript of the variance can be chosen as 1 or 2 as any Hermitian operator's variances on two orthonormal states are equivalent in 2-dimensional Hilbert space. For a pure state, the quantum Fisher information can be easily obtain from Eq. (4) with taking the purity $\text{Tr}\rho^2 = 1$ and the variance on that pure state, i.e., ref. 27

$$F_{\text{pure}} = 4 \langle \Delta^2 \mathcal{H} \rangle_{\text{in}}. \quad (5)$$

Namely, the quantum Fisher information is proportional to the variance of \mathcal{H} on the initial state. In this scenario, denote the initial state $\rho_0 = |\psi_0\rangle\langle\psi_0|$, the quantum Fisher information can be rewritten into $F_{\text{pure}} = \langle \psi_0 | L_{\text{eff}}^2 | \psi_0 \rangle$, with the effective SLD operator

$$L_{\text{eff}} = i2[\mathcal{H}, |\psi_0\rangle\langle\psi_0|]. \quad (6)$$

For a well applied form of parametrization transformation $U = \exp(-itH_\alpha)$ ²⁷, where \hbar has been set as 1 in Planck unit, and being aware of the equation

$$\partial_\alpha e^A = \int_0^1 e^{sA} (\partial_\alpha A) e^{(1-s)A} ds, \quad (7)$$

\mathcal{H} can then be expressed by

$$\mathcal{H} = - \int_0^t e^{isH_\alpha} (\partial_\alpha H_\alpha) e^{-isH_\alpha} ds. \quad (8)$$

Defining a superoperator A^\times as $A^\times(\cdot) := [A, \cdot]$, \mathcal{H} can be written in an expanded form

$$\mathcal{H} = i \sum_{n=0}^{\infty} f_n H_\alpha^{\times n} (\partial_\alpha H_\alpha), \quad (9)$$

where the coefficient

$$f_n = \frac{(it)^{n+1}}{(n+1)!}. \quad (10)$$

In many real problems, the recursive commutations in Eq. (9) can either repeat or terminate²⁸, indicating an analytical expression of \mathcal{H} . Thus, this representation of quantum Fisher information would be very useful in these problems. For the simplest case that $H_\alpha = \alpha H$, all terms vanish but the first one, then $\mathcal{H} = -tH$. When $[H_\alpha, \partial_\alpha H_\alpha] = C$, with C a constant matrix or proportional to H_α , only the first and second terms remain. In this case, \mathcal{H} reduces to $-t(\partial_\alpha H_\alpha + itC/2)$. A more interesting case is that $[H_\alpha, \partial_\alpha H_\alpha] = c\partial_\alpha H_\alpha$, with c a nonzero constant number, then \mathcal{H} can be written in the form

$$\mathcal{H} = \frac{i}{c} [\exp(itc) - 1] \partial_\alpha H_\alpha. \quad (11)$$

In the following we give an example to exhibit Eq. (9). Consider the interaction Hamiltonian of a collective spin system in a magnetic field

$$H_\theta = B(\cos\theta J_x + \sin\theta J_z) = B\mathbf{J}_{n_\theta}, \quad (12)$$

where $\mathbf{J}_{n_\theta} = \mathbf{n}_\theta \cdot \mathbf{J}$ with $\mathbf{n}_\theta = (\cos\theta, 0, \sin\theta)^T$ and $\mathbf{J} = (J_x, J_y, J_z)^T$. B is the amplitude of the external magnetic field and θ is the angle between the field and the collective spin. Here $J_i = \sum_k \sigma_i^{(k)} / 2$ for $i = x, y, z$ with $\sigma_i^{(k)}$ the Pauli matrix for k th spin. Taking θ as the parameter under estimation, \mathcal{H} can be expressed by

$$\mathcal{H} = 2 \left| \sin\left(\frac{Bt}{2}\right) \right| \mathbf{J}_{n_1}, \quad (13)$$

where $\mathbf{J}_{n_1} = \mathbf{n}_1 \cdot \mathbf{J}$ with the vector

$$\mathbf{n}_1 = \mu \left(\cos\left(\frac{Bt}{2}\right) \sin\theta, -\sin\left(\frac{Bt}{2}\right), -\cos\left(\frac{Bt}{2}\right) \cos\theta \right)^T,$$

where $\mu = \text{sgn}(\sin(Bt/2))$ is the sign function and \mathbf{n}_1 is normalized.

The operator \mathcal{H} for Hamiltonian (12) may be also available to be solved using the procedure in Ref. 27, in the $(2j + 1)$ -dimensional eigenspace of H_θ (j is the total spin). In principle, the eigenstates of H_θ can be found by rotating the Dicke state into the same direction of H_θ . However, even one can analytically obtain all the eigenvalues and eigenvectors, it still requires a large amount of calculations to obtain \mathcal{H} , especially when the spin numbers are tremendous. Comparably, utilizing Eq. (9), it only takes a few steps of calculation, which can be found in the method. This is a major advantage of the expanded form of \mathcal{H} .

Utilizing Eq. (13), one can immediately obtain the form of \mathcal{H} for a spin-half system

$$\mathcal{H}_{\text{qubit}} = \left| \sin\left(\frac{Bt}{2}\right) \right| \mathbf{n}_1 \cdot \boldsymbol{\sigma}, \quad (14)$$

with $\boldsymbol{\sigma} = (\sigma_x, \sigma_y, \sigma_z)^T$, which was also discussed in the Hamiltonian eigenbasis in Ref. 27. For any 2-dimensional state, based on Eq. (4), the quantum Fisher information can be expressed by



$$F_\theta = 4 \sin^2\left(\frac{Bt}{2}\right) |\mathbf{r}_{\text{in}}|^2 [1 - (\mathbf{n}_1 \cdot \mathbf{r}_e)^2], \quad (15)$$

where $\mathbf{r}_{\text{in}} = (\langle \sigma_x \rangle, \langle \sigma_y \rangle, \langle \sigma_z \rangle)^T$ is the Bloch vector of the initial state ρ_0 and \mathbf{r}_e is the Bloch vector of any eigenstate of ρ_0 . For pure states, there is $\mathbf{r}_e = \mathbf{r}_{\text{in}}$ and $|\mathbf{r}_{\text{in}}| = 1$. Since the Bloch vector of a 2-dimensional state satisfies $|\mathbf{r}_{\text{in}}| \leq 1$, it can be found that the maximum value of Eq. (15) is

$$F_\theta^{\text{max}} = 4 \sin^2\left(\frac{Bt}{2}\right), \quad (16)$$

which can be saturated when $|\mathbf{r}_{\text{in}}| = 1$ and $\mathbf{n}_1 \cdot \mathbf{r}_{\text{in}} = 0$, namely, the optimal state to access maximum quantum Fisher information here is a pure state perpendicular to \mathbf{n}_1 , as shown in Fig. 1. In this figure, the yellow sphere represents the Bloch sphere and the blue arrow represents the vector \mathbf{n}_1 . It can be found that all states on the joint ring of the green plane and surface of Bloch sphere can access the maximum quantum Fisher information, i.e., all states on this ring are optimal states. One simple example is $\mathbf{r}_{\text{opt}} = \mathbf{n}_0$, and another one is the superposition state of two eigenstates of $\mathcal{H}^{22,27}$.

Alternatively, B could be the parameter that under estimation. In the spin-half case, with respect to B , $\mathcal{H}_B = -t\mathbf{n}_0 \cdot \sigma/2$, then the quantum Fisher information can be expressed by

$$F_B = t^2 |\mathbf{r}_{\text{in}}|^2 [1 - (\mathbf{n}_0 \cdot \mathbf{r}_e)^2]. \quad (17)$$

The optimal states to access the maximum value $F_B^{\text{max}} = t^2$ are the pure states vertical to \mathbf{n}_0 .

Exponential form initial state. For an exponential form initial state $\rho_0 = \exp(G_0)$, the parametrized state reads

$$\rho_x = U \rho_0 U^\dagger = \exp(UG_0 U^\dagger). \quad (18)$$

Recently, Jiang²⁸ studied the quantum Fisher information for exponential states and gave a general form of SLD operator. In his theory, the SLD operator can be expanded as

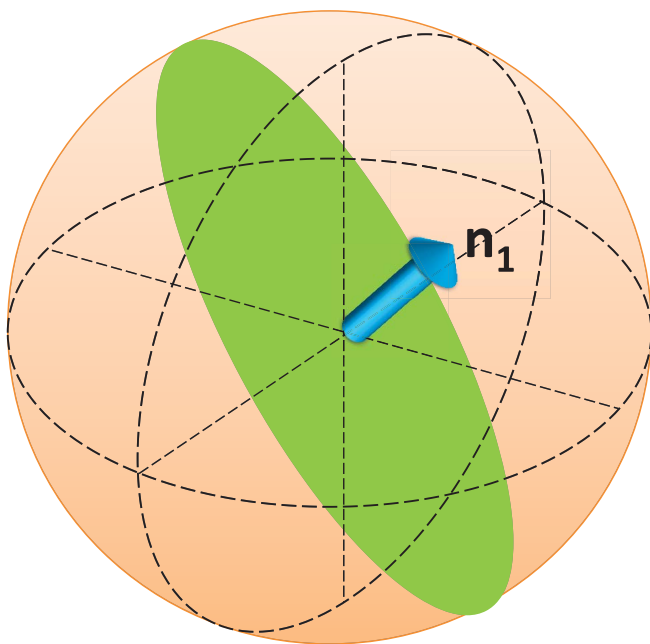


Figure 1 | Optimal states to access maximum quantum Fisher information in a spin-half system. The blue arrow represents the vector \mathbf{n}_1 and all vectors in the green plane are vertical to \mathbf{n}_1 . All the states in the joint ring of green plane and Bloch sphere's surface can access maximum quantum Fisher information.

$$L = \sum_{n=0}^{\infty} g_n G^{\times n} (\partial_x G), \quad (19)$$

where the coefficient

$$g_n = \frac{4(2^{n+2} - 1) \mathcal{B}_{n+2}}{(n+2)!} \quad (20)$$

for even n and g_n vanishes for odd n . Here \mathcal{B}_{n+2} is the $(n+2)$ th Bernoulli number and in our case, $G = UG_0U^\dagger$. Through some straightforward calculation, the derivative of G on α reads

$$\partial_x G = -iU(G_0^\times \mathcal{H})U^\dagger. \quad (21)$$

Based on this equation, the n th order term in Eq. (19) is

$$G^{\times n} (\partial_x G) = -iU\left((G_0^\times)^{n+1} \mathcal{H}\right)U^\dagger, \quad (22)$$

where \mathcal{H} is given by Eq. (9). Generally, it is known that the quantum Fisher information reads

$$F = \text{Tr}(U \rho_0 U^\dagger L^2) = \text{Tr}(\rho_0 L_{\text{eff}}^2), \quad (23)$$

where the effective SLD operator $L_{\text{eff}} = U^\dagger L U$. The effective SLD operator for pure states is already shown in Eq. (6). Substituting Eq. (22) into Eq. (19), the effective SLD operator can be expanded as

$$L_{\text{eff}} = -i \sum_{n=0}^{\infty} g_n (G_0^\times)^{n+1} \mathcal{H}. \quad (24)$$

In most mixed states cases, to obtain quantum Fisher information, the diagonalization of initial state is inevitable, which is the reason why the usual form of quantum Fisher information is expressed in the eigenbasis of density matrix. Thus, it is worth to study the expression of effective SLD operator and quantum Fisher information in the eigenbasis of G_0 . We denote the i th eigenvalue and eigenstate of G_0 as a_i and $|\phi_i\rangle$, and in the eigenbasis of G_0 , the element of $G_0^{\times n} \mathcal{H}$ satisfies the recursion relation

$$[G_0^{\times n} \mathcal{H}]_{ij} = (a_i - a_j) [(G_0^{\times n})^{n-1} \mathcal{H}]_{ij}, \quad (25)$$

where $[\cdot]_{ij} := \langle \phi_i | \cdot | \phi_j \rangle$. Utilizing this recursive equation, a general formula of n th order term can be obtained,

$$[G_0^{\times n} \mathcal{H}]_{ij} = (a_i - a_j)^n \mathcal{H}_{ij}. \quad (26)$$

Substituting above equation into the expression of L_{eff} and being aware of the equality

$$\sum_{n=0}^{\infty} g_n (a_i - a_j)^{n+1} = 2 \tanh\left(\frac{a_i - a_j}{2}\right), \quad (27)$$

the element of effective SLD operator in Eq. (24) can be written as

$$[L_{\text{eff}}]_{ij} = -i2 \tanh\left(\frac{a_i - a_j}{2}\right) \mathcal{H}_{ij}. \quad (28)$$

Based on the equation $F = \text{Tr}(e^{G_0} L_{\text{eff}}^2)$, the quantum Fisher information in the eigenbasis of G_0 can finally be expressed by

$$F = \sum_{i>j} 4(e^{a_i} + e^{a_j}) \tanh^2\left(\frac{a_i - a_j}{2}\right) |\mathcal{H}_{ij}|^2. \quad (29)$$

This is one of the main results in this paper. In some real problems, the eigenspace of G_0 could be found easily. For instance, the eigenspace of a bosonic thermal state is the Fock space. Thus, as long as the formula of \mathcal{H} in Fock space is established, the quantum Fisher information can be obtained from Eq. (29).

Now we exhibit Eq. (29) with a spin-half thermal state. The initial state is taken as



$$\begin{aligned}\rho_0 &= \frac{1}{Z} \exp(-\beta\sigma_z), \\ &= \exp(-\beta\sigma_z - \ln Z),\end{aligned}\quad (30)$$

where $\beta = 1/(k_b T)$ with k_b the Boltzmann constant and T the temperature. In Planck unit, $k_b = 1$. The partition function reads $Z = \text{Tr}[\exp(-\beta\sigma_z)] = 2 \cosh \beta$. In this case, $G_0 = -\beta\sigma_z - \ln z$. Denoting the eigenstates of σ_z as $|0\rangle$ and $|1\rangle$, i.e., $\sigma_z = |0\rangle\langle 0| - |1\rangle\langle 1|$, the eigenvalues of G_0 read $a_1 = -\beta\sigma_z - \ln z$ and $a_2 = \beta\sigma_z - \ln z$. The parametrization process is still taken as $H_\theta = B\mathbf{n}_\theta \cdot \boldsymbol{\sigma}/2$ with θ the parameter under estimation, indicating that $\mathcal{H} = \left| \sin\left(\frac{Bt}{2}\right) \right| \mathbf{n}_1 \cdot \boldsymbol{\sigma}$, then the squared norm of the off-diagonal element of \mathcal{H} in the eigenbasis of σ_z reads

$$|\mathcal{H}_{01}|^2 = \sin^2\left(\frac{Bt}{2}\right) \left[1 - \cos^2\theta \cos^2\left(\frac{Bt}{2}\right) \right]. \quad (31)$$

Immediately, the quantum Fisher information can be obtained from Eq. (29) as

$$F_T = 4 \tanh^2\beta \sin^2\left(\frac{Bt}{2}\right) \left[1 - \cos^2\theta \cos^2\left(\frac{Bt}{2}\right) \right]. \quad (32)$$

The maximum value of above expression is obtained at $Bt = (4k + 1)\pi$ for $k = 0, 1, \dots$ and

$$F_T^{\max} = 4 \tanh^2(\beta). \quad (33)$$

From this equation, one can see that the value of maximum quantum Fisher information is only affected by the temperature. With the increase of temperature, the maximum value reduces. In the other hand, quantum Fisher information in Eq. (32) is related to Bt and θ . Fig. 2 shows the quantum Fisher information as a function of Bt and θ . The values of Bt and θ are both within $[0, 2\pi]$ in the plot. The temperature is set as $T = 1$ here. From this figure, it can be found that the maximum quantum Fisher information is robust for θ since it is always obtained at $Bt = \pi$ for any value of θ . Furthermore, this optimal condition of Bt is independent of temperature. With respect to Bt , there is a large regime near $Bt = \pi$ in which the quantum Fisher information's value can surpass 2, indicating that the quantum Fisher information can be still very robust and near its maximum value even when Bt is hard to set exactly at π .

Multiparameter processes. For a multiparameter system, the element of quantum Fisher information matrix in Ref. 30 can also be written with \mathcal{H} operator,

$$\begin{aligned}\mathcal{F}_{\alpha\beta} &= \sum_{i=1}^M 4p_i \text{cov}_i(\mathcal{H}_\alpha, \mathcal{H}_\beta) \\ &\quad - \sum_{i \neq j} \frac{8p_i p_j}{p_i + p_j} \text{Re}\left(\langle \psi_i | \mathcal{H}_\alpha | \psi_j \rangle \langle \psi_j | \mathcal{H}_\beta | \psi_i \rangle\right),\end{aligned}\quad (34)$$

where U is dependent on a series of parameters α, β and so on, and

$$\mathcal{H}_m = i(\partial_m U^\dagger)U, \quad (35)$$

with the index $m = \alpha, \beta, \dots$. The covariance matrix on the i th eigenstate of initial state is defined as

$$\text{cov}_i(\mathcal{H}_\alpha, \mathcal{H}_\beta) = \frac{1}{2} \langle \psi_i | \{ \mathcal{H}_\alpha, \mathcal{H}_\beta \} | \psi_i \rangle - \langle \psi_i | \mathcal{H}_\alpha | \psi_i \rangle \langle \psi_i | \mathcal{H}_\beta | \psi_i \rangle,$$

with $\{ \cdot, \cdot \}$ the anti-commutation. For a single qubit system, Eq. (34) reduces to

$$\mathcal{F}_{\text{qubit}, \alpha\beta} = 4(2\text{Tr}\rho^2 - 1) \text{cov}_{1(2)}(\mathcal{H}_\alpha, \mathcal{H}_\beta). \quad (36)$$

Similarly with the single-parameter scenario, the subscript in Eq. (36)

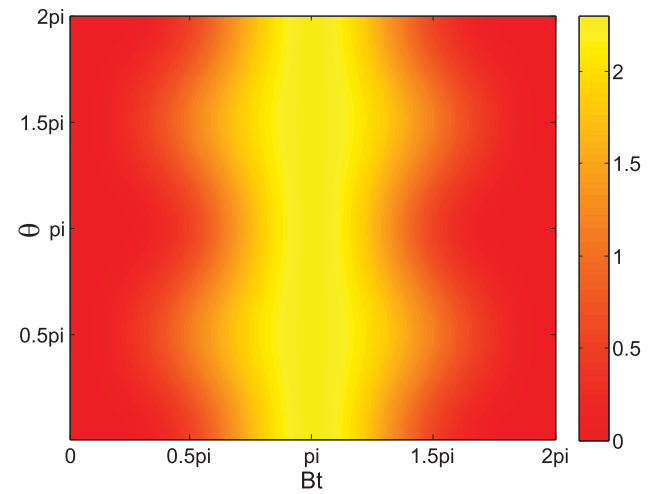


Figure 2 | Quantum Fisher information as a function of Bt and θ . The initial state is a spin-half thermal state and the temperature is set as $T = 1$ here.

can be chosen as 1 or 2 since the covariance for two Hermitian operators are the same on two orthonormal states in 2-dimensional Hilbert space. From this equation, the element of quantum Fisher information matrix for pure states can be immediately obtained as

$$\mathcal{F}_{\text{pure}, \alpha\beta} = 4\text{cov}_{\text{in}}(\mathcal{H}_\alpha, \mathcal{H}_\beta), \quad (37)$$

namely, for pure states, the element of quantum Fisher information matrix is actually the covariance between two \mathcal{H} operators on the initial state. When the total Hamiltonian can be written as $\sum_i \alpha_i H_i$ and $[H_i, H_j] = 0$ for any i, j , above equation can reduce to the covariance between H_i and H_j ³¹. For the diagonal elements, they are exactly the quantum Fisher information for the corresponding parameters.

For multiparameter estimations, the Cramér-Rao bound cannot always be achieved. In the scenario of pure states, the condition of this bound to be tight is $\text{Im}\langle \psi_{\text{out}} | L_\alpha L_\beta | \psi_{\text{out}} \rangle = 0, \forall \alpha, \beta$ ^{32,33}. Here $|\psi_{\text{out}}\rangle$ is dependent on the parameter under estimation. In the unitary parametrization, $|\psi_{\text{out}}\rangle = U|\psi_0\rangle$ and this condition can be rewritten into $\text{Im}\langle \psi_0 | L_{\text{eff}}^\alpha L_{\text{eff}}^\beta | \psi_0 \rangle = 0, \forall \alpha, \beta$. Here $L_{\text{eff}}^{\alpha(\beta)} = U^\dagger L_{\alpha(\beta)} U$ is the effective SLD operator for parameter $\alpha(\beta)$. Utilizing Eq. (6), this condition can be expressed in the form of \mathcal{H} operator,

$$\langle \psi_0 | [\mathcal{H}_\alpha, \mathcal{H}_\beta] | \psi_0 \rangle = 0, \quad \forall \alpha, \beta. \quad (38)$$

In other word, $\langle \psi_0 | \mathcal{H}_\alpha, \mathcal{H}_\beta | \psi_0 \rangle$ needs to be a real number for any α and β . When \mathcal{H}_α commutes with \mathcal{H}_β for any α and β , above condition can always be satisfied for any initial state.

Generally, for the unitary parametrization process, the element of quantum Fisher information matrix can be expressed by $F = \text{Tr}(\rho\{L_\alpha, L_\beta\}) = \text{Tr}\left(\rho_0\left\{L_{\text{eff}}^\alpha, L_{\text{eff}}^\beta\right\}\right)$. From the definition equation of SLD, one can see that L_{eff}^α satisfies the equation $\partial_\alpha \rho = U\{\rho_0, L_{\text{eff}}^\alpha\}U^\dagger/2$. The quantum Fisher information matrix has more than one definitions. One alternative candidate is using the so-called Right Logarithmic Derivative (RLD)^{24,34,35}, which is defined as $\partial_\alpha \rho = \rho R_\alpha$ with R_α the RLD. The element of RLD quantum Fisher information matrix can be written as

$$\mathcal{J}_{\alpha\beta} = \text{Tr}\left(\rho R_\alpha R_\beta^\dagger\right) = \text{Tr}\left(\rho_0 R_{\text{eff}}^\alpha R_{\text{eff}}^{\beta\dagger}\right), \quad (39)$$

where the effective RLD reads $R_{\text{eff}}^{\alpha(\beta)} = U^\dagger R_{\alpha(\beta)} U$. For a unitary parametrization process, assuming the initial state has nonzero determin-



ant, R_{eff}^z can be expressed by \mathcal{H}_z and the initial state ρ_0 , i.e.,

$$R_{\text{eff}}^z = i(\rho_0^{-1} \mathcal{H}_z \rho_0 - \mathcal{H}_z). \quad (40)$$

With this equation, the element of RLD quantum Fisher information matrix can be expressed by

$$\mathcal{J}_{\alpha\beta} = \text{Tr}(\mathcal{H}_\alpha \rho_0^2 \mathcal{H}_\beta \rho_0^{-1} - 2\mathcal{H}_\beta \mathcal{H}_\alpha \rho_0 + \mathcal{H}_\alpha \mathcal{H}_\beta \rho_0). \quad (41)$$

When the parametrization process is displacement, this equation can reduce to the corresponding form in Ref. 35. For pure states, the element reads $\mathcal{J}_{\text{pure},\alpha\beta} = \text{Tr}[(\partial_\alpha \rho)(\partial_\beta \rho)] = \mathcal{F}_{\text{pure},\alpha\beta}/2$. Recently, Genoni *et al.*³⁵ proposed a most informative Cramér-Rao bound for the total variance of all parameters under estimation. From the relation between $\mathcal{J}_{\text{pure},\alpha\beta}$ and $\mathcal{F}_{\text{pure},\alpha\beta}$, one can see that $\text{Tr}\mathcal{F}_{\text{pure}}^{-1}$ is always larger than $\text{Tr}\mathcal{F}_{\text{pure}}^{-1}$, namely, the SLD Cramér-Rao bound is always more informative than the RLD counterpart in this scenario.

We still consider the spin-half system with the Hamiltonian $H = B\mathbf{n}_0 \cdot \boldsymbol{\sigma}/2$. Take both B and θ as the parameters under estimations. First, based on aforementioned calculation, the \mathcal{H} operator for B and θ read

$$\mathcal{H}_B = -\frac{t}{2} \mathbf{n}_0 \cdot \boldsymbol{\sigma}, \quad (42)$$

$$\mathcal{H}_\theta = \left| \sin\left(\frac{Bt}{2}\right) \right| \mathbf{n}_1 \cdot \boldsymbol{\sigma}. \quad (43)$$

Based on the property of Pauli matrices $\{\mathbf{n}_0 \cdot \boldsymbol{\sigma}, \mathbf{n}_1 \cdot \boldsymbol{\sigma}\} = 2\mathbf{n}_0 \cdot \mathbf{n}_1$, the anti-commutation in the covariance reads

$$\{\mathcal{H}_B, \mathcal{H}_\theta\} = -t \left| \sin\left(\frac{Bt}{2}\right) \right| \mathbf{n}_0 \cdot \mathbf{n}_1. \quad (44)$$

For a pure initial state, the off-diagonal element of the quantum Fisher information matrix is expressed by

$$\mathcal{F}_{B\theta} = 2t \left| \sin\left(\frac{Bt}{2}\right) \right| (\mathbf{n}_0 \cdot \mathbf{r}_{\text{in}})(\mathbf{n}_1 \cdot \mathbf{r}_{\text{in}}), \quad (45)$$

where \mathbf{r}_{in} is the Bloch vector of the initial pure state and the equality $\mathbf{n}_0 \cdot \mathbf{n}_1 = 0$ has been used. When the initial pure state is vertical to \mathbf{n}_0 or \mathbf{n}_1 , this off-diagonal element vanishes. Compared with the optimal condition for maximum quantum Fisher information for B and θ individually, the Bloch vector $\mathbf{n}_2 = \mathbf{n}_0 \times \mathbf{n}_1$ can optimize both the diagonal elements of quantum Fisher information matrix and vanish the off-diagonal elements. However, all above is only necessary conditions for the achievement of Cramér-Rao bound. To find out if the bound can be really achieved, the condition (38) needs to be checked. In this case,

$$[\mathcal{H}_B, \mathcal{H}_\theta] = -it \left| \sin\left(\frac{Bt}{2}\right) \right| \mathbf{n}_2 \cdot \boldsymbol{\sigma}. \quad (46)$$

With this equation, condition (38) reduces to $\mathbf{n}_2 \cdot \mathbf{r}_{\text{in}} = 0$, i.e., to make the Cramér-Rao bound achievable, the Bloch vector of the initial state needs to be in the plane of \mathbf{n}_0 and \mathbf{n}_1 . Unfortunately, \mathbf{n}_2 is not in this plane. Thus, B and θ cannot be optimally joint measured simultaneously.

In the plane constructed by \mathbf{n}_0 and \mathbf{n}_1 , any Bloch vector of pure state can be written as $\mathbf{r}_{\text{in}} = \mathbf{n}_0 \cos \phi + \mathbf{n}_1 \sin \phi$, then we have $\mathcal{F}_{BB} = t^2 \sin^2 \phi$, $\mathcal{F}_{\theta\theta} = 4 \sin^2(Bt/2) \cos^2 \phi$, and $\mathcal{F}_{B\theta} = 2t |\sin(Bt/2)| \cos \phi \sin \phi$. From these expressions, one can see that the determinant of quantum Fisher information matrix is zero, i.e., $\det \mathcal{F} = 0$. This fact indicates that, utilizing any pure state in this plane, the variances of B and θ cannot be estimated simultaneously through the Cramér-Rao theory.

Discussion

We have discussed the quantum Fisher information with unitary parametrization utilizing an alternative representation. The total information of the parametrization process is involved in a \mathcal{H} operator in this representation. This operator is totally determined by the parameter and parametrization transformation U . As long as the parameter and transformation are taken, \mathcal{H} is a settled operator and independent of the initial state. More interestingly, \mathcal{H} can be expressed in an expanded form. For the Hamiltonians owning recursive commutations with their partial derivative on the parameter under estimation, this expanded form shows a huge advantage. Utilizing this representation, we give a general analytical expression of quantum Fisher information for an exponential form initial state. Moreover, we have also studied the \mathcal{H} representation in multiparameter processes. The condition of Cramér-Rao bound to be achievable for pure states are also presented in the form of \mathcal{H} operator. In addition, we give the \mathcal{H} representation of Right Logarithmic Derivative and the corresponding quantum Fisher information matrix.

As a demonstration, we apply this representation in a collective spin system and show the expression of \mathcal{H} . Furthermore, we provide an analytical expression of quantum Fisher information in a spin-half system. If we consider this system as a multiparameter system, the corresponding quantum Fisher information matrix can also be straight-forwardly obtained by this representation. From these expressions, one can find the optimal states to access the maximum quantum Fisher information. For the parameter B , the optimal state is a pure state vertical to \mathbf{n}_0 , and for the parameter θ , the optimal one is also a pure state, but vertical to \mathbf{n}_1 . By analyzing the off-diagonal element of quantum Fisher information matrix, the states to optimize the diagonal elements and make the off-diagonal elements vanish are found. However, these states fail to satisfy the condition of achievement. Thus, B and θ cannot be optimally joint measured.

Methods

Collective spin system in a magnetic field. For the Hamiltonian (12), its derivative on parameter θ is $\partial_\theta H_\theta = \mathbf{n}'_0 \cdot \mathbf{J} = J_{n'_0} = -iH_\theta^\times J_y$ with the vector $\mathbf{n}'_0 = d\mathbf{n}_0/d\theta = (-\sin\theta, 0, \cos\theta)^\top$. Based on Eq. (9), \mathcal{H} can be written as

$$\mathcal{H} = [\exp(itH_\theta^\times) - 1] J_y. \quad (47)$$

It is worth to notice that $H_\theta^\times = BJ_{n'_0}^\times$, then \mathcal{H} is

$$\mathcal{H} = [\exp(iBtJ_{n'_0}^\times) - 1] J_y. \quad (48)$$

Being aware of the commutation relations

$$[J_{n_0}, J_y] = iJ_{n'_0}, \quad (49)$$

$$[J_{n_0}, J_{n'_0}] = -iJ_y, \quad (50)$$

one can straightforwardly obtain the n th order term as below

$$J_{n_0}^{\times n} J_y = \begin{cases} J_y & \text{for even } n; \\ iJ_{n'_0} & \text{for odd } n. \end{cases} \quad (51)$$

With this equation, \mathcal{H} can be expressed by

$$\mathcal{H} = [\cos(Bt) - 1] J_y - \sin(Bt) J_{n'_0}, \quad (52)$$

equivalently, it can be written in an inner product form: $\mathcal{H} = \mathbf{r} \cdot \mathbf{J}$, where the elements of \mathbf{r} read $r_x = \sin(Bt) \sin \theta$, $r_y = \cos(Bt) - 1$ and $r_z = -\sin(Bt) \cos \theta$. After the normalization process, \mathcal{H} is rewritten into the form of Eq. (13).

For a spin-half system, the quantum Fisher information can be expressed by

$$F = 4 \sin^2\left(\frac{Bt}{2}\right) |\mathbf{r}_{\text{in}}|^2 \left[1 - (\mathbf{n}_1 \cdot \langle \boldsymbol{\sigma} \rangle_{1(2)})^2 \right], \quad (53)$$

where \mathbf{r}_{in} is the Bloch vector of ρ_0 and can be obtained through the equation

$$\rho_0 = \frac{1}{2} \mathbb{1} + \frac{1}{2} \sum_{i=x,y,z} r_{\text{in},i} \sigma_i, \quad (54)$$

with $\mathbb{1}$ the identity matrix. $\langle \boldsymbol{\sigma} \rangle_i = (\langle \sigma_x \rangle_i, \langle \sigma_y \rangle_i, \langle \sigma_z \rangle_i)^\top$ is the vector of expected values on



the i th ($i = 1, 2$) eigenstate of ρ_0 . It can also be treated as the Bloch vector of the eigenstates. In previous sections, we denote $r_c := \langle \sigma \rangle_i$.

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Author contributions

X.W. and J.L. contributed the idea. J.L. performed the calculations and prepared the figures. X.J. checked the calculations. J.L. wrote the main manuscript and X.W. made an improvement. All authors contributed to discussion and reviewed the manuscript.

Additional information

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