

Combining spatiotemporal and particle-number degrees of freedom

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Quadrature bases that incorporate spatiotemporal degrees of freedom are derived as eigenstates of momentum-dependent quadrature operators. The resulting bases are shown to be orthogonal for both the particle-number and the spatiotemporal degrees of freedom. Using functional integration, we also demonstrate the completeness of these quadrature bases.

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I. INTRODUCTION

Quantum information promises various new technologies, including quantum imaging [1–3], quantum metrology [4–6], and quantum communication [7–9]. Many of the systems that implement these quantum information technologies use quantum optics [10]. In such cases, we often find that the physical setup predominantly employs either the spatial degrees of freedom [11–13] or the particle-number degrees of freedom [14–16]. Where the spatial degrees of freedom are used, the implementation is often made in terms of spatial modes, such as orbital angular momentum modes [17,18]. Those implementations that employ the particle-number degrees of freedom often use so-called continuous variables [19,20].

However, one cannot completely exclude the influence of other degrees of freedom from any quantum optics implementation. When spatial modes are used, one usually assumes that the states are single-photon states or two-photon states [21–23]. Nevertheless, the laser sources and nonlinear crystals actually produce coherent states and squeezed states, thus polluting these low-particle-number states with higher-particle-number states [24,25]. The effects of these undesirable contributions are treated as ‘noise’ in the system.

On the other hand, continuous-variable systems always incorporate some spatiotemporal description of the optical field. In these implementations, multiple modes are allowed, but it is often assumed that these are fixed discrete modes [15,19,20,26]. (Obviously, there are many exceptions, for instance, in the context of quantum imaging [27,28].) However, in practical systems, distortions and decoherence can introduce unwanted variability in these modes [29–31] that may not be represented by the finite set of fixed modes.

A comprehensive approach that combines all the degrees of freedom allows one to perform thorough analyses of all such scenarios. It is especially relevant in cases where all these degrees of freedom are affected. Such a comprehensive approach requires the incorporation of both spatiotemporal and particle-number degrees of freedom into the analytic tools. If the current state of the art for implementations that focus on only

one degree of freedom is able to achieve remarkable successes as they do, just imagine how powerful implementations that incorporate all these degrees of freedom would be [32–37]. There is a clear benefit of analytical tools that incorporate both spatiotemporal and particle-number degrees of freedom.

In the endeavor to combine particle-number degrees of freedom and spatiotemporal degrees of freedom, there are different approaches to choose from. One approach is to expand the states under investigation in terms of orthogonal bases. Operators can then be expressed in terms of overlaps among different bases elements. However, such an approach assumes the existence of a basis that is not only complete and orthogonal with respect to the particle-number degrees of freedom but, at the same time, also complete and orthogonal with respect to the spatiotemporal degrees of freedom. To find such a basis is not a simple matter. It is the topic of this paper.

An alternative approach is to define states in terms of operators that would produce those states when they operate on the vacuum state. For normalized states, these would be unitary operators, often expressed as exponential operators, containing Hermitian operators in their exponents. The latter is often expressed as multivariate polynomials of creation and annihilation operators. It is particularly convenient if these polynomials are no higher than second order, thus naturally leading to the notion of Gaussian states [14,20]. Operations on states are represented by products of operators, which can be manipulated via the appropriate commutation relations. Due to the creation and annihilation operators, this approach is often associated with second quantization.

One way to incorporate spatiotemporal degrees of freedom in the latter approach is to include multiple sets of creation and annihilation operators to represent different modes. This necessarily leads to a discrete set of modes representing the spatiotemporal degrees of freedom [19,20]. However, it is not always desirable to perform analyses in terms of discrete modes. The result may require a truncation in the number of modes to allow explicit computations. Such a truncation can lead to large deviations between predictions and experimental results. For example, when discrete modes were used for the analysis of the evolution of biphoton states in turbulence [38], the resulting set of coupled differential equations had to be truncated before they could be solved.

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Moreover, the complexity of the analysis grew rapidly with the size of the truncated set. The predictions obtained from such a truncated set of equations gave a large disagreement compared with experimental results [39]. Only by performing the analysis in terms of the continuous plane-wave basis could this truncation problem be overcome [40,41]. Hence, it is preferable to incorporate spatiotemporal degrees of freedom into a particle-number formalism in such a way that it will allow one to use a continuous parametrization of both the particle-number degrees of freedom and the spatiotemporal degrees of freedom.

Although we focus here mainly on the inclusion of the spatiotemporal degrees of freedom with the particle-number degrees of freedom, with a little more effort, we also incorporate the spin degrees of freedom, thus exhausting all the degrees of freedom associated with the photon field. To do so, we make a slight simplification of the notation to avoid overly complex expressions.

In this paper, we derive the eigenbases associated with so-called fixed-momentum quadrature operators. The resulting bases are called spatiotemporal quadrature bases. Although it makes sense heuristically that these bases would be complete and orthogonal, we proceed to show this explicitly. Since these bases incorporate the spatiotemporal degrees of freedom in terms of continuous degrees of freedom, they naturally lead to a functional formalism—a path-integral approach [42–44]. However, by themselves, these spatiotemporal quadrature bases do not yet represent a fully fledged formalism. They only provide the first step. The next step, which is to use these quadrature bases for the development of a generalized Wigner formalism, is beyond the scope of the current paper. However, we briefly discuss such a generalized Wigner formalism later.

The paper is organized as follows. In Sec. II we provide some background information and define convenient notation. The derivation of the spatiotemporal quadrature bases is provided in Sec. III. Orthogonality and completeness conditions for these bases are considered in Secs. IV and V, respectively. Finally, in Sec. VI a discussion and outlook are provided.

II. PRELIMINARIES

Before we address the main topic of this paper, namely, the derivation of the spatiotemporal quadrature bases, we first consider some preliminary results that would be needed later. It also gives us the opportunity to define convenient notation and to show alternative approaches that do not work.

A. Fixed-momentum Fock states

The quantization of the electromagnetic field in the context of particle physics led to the definition of creation and annihilation operators $a_s^\dagger(\mathbf{k})$ and $a_s(\mathbf{k})$, which carry spatiotemporal degrees of freedom in the form of the wave vector \mathbf{k} , which is proportional to the three-dimensional momentum vector, and spin degrees of freedom, represented by the spin index s . The creation and annihilation operators are assumed to obey a Lorentz covariant commutation relation,

$$[\hat{a}_s(\mathbf{k}_1), \hat{a}_r^\dagger(\mathbf{k}_2)] = (2\pi)^3 \omega_1 \delta_{s,r} \delta(\mathbf{k}_1 - \mathbf{k}_2), \quad (1)$$

where $\delta_{s,r}$ is the Kronecker delta, $\delta(\mathbf{k}_1 - \mathbf{k}_2)$ is a (three-dimensional) Dirac delta function, and the angular frequency ω_1 is related to the wave vector via the dispersion relation $\omega_1 = c|\mathbf{k}_1|$. The creation operators produce the elements of a momentum basis $a_s^\dagger(\mathbf{k})|\text{vac}\rangle = |\mathbf{k}, s\rangle$, which serves as a complete orthogonal basis for all single-photon states. It satisfies a Lorentz covariant orthogonality condition that reads

$$\langle \mathbf{k}_1, r | \mathbf{k}_2, s \rangle = (2\pi)^3 \omega_1 \delta_{r,s} \delta(\mathbf{k}_1 - \mathbf{k}_2). \quad (2)$$

As the first attempt, one can generalize the single-photon momentum basis to a *fixed-momentum* Fock basis, where all photons in each basis element share the same spin and wave vector,

$$\begin{aligned} |n, \mathbf{k}, s\rangle &= \frac{1}{\sqrt{n!}} [a_s^\dagger(\mathbf{k})]^n |\text{vac}\rangle, \\ \langle n, \mathbf{k}, s| &= \frac{1}{\sqrt{n!}} \langle \text{vac} | [a_s(\mathbf{k})]^n, \end{aligned} \quad (3)$$

where n is the occupation number of photons in the state. The appropriate number operator for which these Fock states serve as eigenstates, with the occupation number n being the eigenvalue, is given by

$$\hat{n} = \sum_r \int a_r^\dagger(\mathbf{k}') a_r(\mathbf{k}') \frac{d^3 k'}{(2\pi)^3 \omega'}. \quad (4)$$

At this point, we introduce a simpler notation to alleviate the complexity of subsequent expressions. To this end, integrals over momentum space with spin sums are henceforth denoted

$$\sum_s \int \dots \frac{d^3 k}{(2\pi)^3 \omega} \rightarrow \int \dots \bar{d}_s k. \quad (5)$$

Unfortunately, the attempt to define Fock states with fixed momenta leads to divergences. The inner product between two of these fixed-momentum Fock states with occupation numbers $m, n > 1$ gives a product of Dirac delta functions with the same argument:

$$\begin{aligned} \langle m, \mathbf{k}_1, r | n, \mathbf{k}_2, s \rangle &= \delta_{m,n} \langle \mathbf{k}_1, r | \mathbf{k}_2, s \rangle^n \\ &= \delta_{m,n} [(2\pi)^3 \omega_1 \delta_{r,s} \delta(\mathbf{k}_1 - \mathbf{k}_2)]^n. \end{aligned} \quad (6)$$

The product of Dirac delta functions will inevitably lead to unwanted divergences in any calculation. As a result, such an attempt to incorporate the spatiotemporal degrees of freedom with the particle-number degrees of freedom fails. It follows that any other fixed-momentum basis (such as fixed-momentum quadrature bases or fixed-momentum coherent states) would suffer the same affliction.

B. Fixed-spectrum Fock states

A slight variation on the theme of fixed-momentum Fock states is the notion of a *fixed-spectrum* Fock state, in which all the photons in the state share the same spectrum of plane waves. Making the presence of the spectrum explicit in the

notation by the subscript F , we express these Fock states

$$\begin{aligned} |n_F\rangle &= \frac{1}{\sqrt{n!}} (\hat{a}_F^\dagger)^n |\text{vac}\rangle = \frac{(|1_F\rangle)^n}{\sqrt{n!}}, \\ \langle n_F| &= \frac{1}{\sqrt{n!}} \langle \text{vac}| (\hat{a}_F)^n = \frac{(\langle 1_F|)^n}{\sqrt{n!}}, \end{aligned} \quad (7)$$

where the fixed-spectrum creation and annihilation operators are defined by

$$\begin{aligned} \hat{a}_F^\dagger &= \int a_s^\dagger(\mathbf{k}) F_s(\mathbf{k}) \tilde{d}_s k, \\ \hat{a}_F &= \int F_s^*(\mathbf{k}) a_s(\mathbf{k}) \tilde{d}_s k. \end{aligned} \quad (8)$$

Fixed-spectrum single-photon states are expressed as

$$\begin{aligned} |1_F\rangle &= \int |\mathbf{k}, s\rangle F_s(\mathbf{k}) \tilde{d}_s k, \\ \langle 1_F| &= \int F_s^*(\mathbf{k}) \langle \mathbf{k}, s| \tilde{d}_s k. \end{aligned} \quad (9)$$

The fixed-spectrum Fock states also act as eigenstates of the number operator in Eq. (4): $\hat{n}|n_F\rangle = |n_F\rangle n$.

In all these expressions, the spectrum is represented by $F_s(\mathbf{k})$, being a function of the three-dimensional wave vector \mathbf{k} and the spin index s . It is normalized according to the expression

$$\int |F_s(\mathbf{k})|^2 \tilde{d}_s k = 1. \quad (10)$$

The normalized spectrum ensures that $\langle 1_F|1_F\rangle = 1$ and $[a_F, a_F^\dagger] = 1$. As a result, the inner product between arbitrary Fock states with different spectra gives

$$\langle m_F|n_G\rangle = \delta_{mn} (\langle F, G\rangle)^n, \quad (11)$$

where

$$\langle F, G\rangle \equiv \int F_s^*(\mathbf{k}) G_s(\mathbf{k}) \tilde{d}_s k. \quad (12)$$

Hence, if we select a discrete set of spatial modes that is orthogonal and complete in terms of the spatiotemporal degrees of freedom to represent the spectra in these Fock states, we would obtain a set that obeys an orthogonality condition both in the spatiotemporal degrees of freedom and in the particle-number degrees of freedom. Such a discrete fixed-spectrum Fock basis is denoted $\{|n_m\rangle\}$, where m is an index for the spatiotemporal modes.

Unfortunately, it would not be complete in both these degrees of freedom. Trying to expand a state consisting of the tensor product of two single-photon elements of the basis $|\psi\rangle = |1_m\rangle|1_n\rangle$, where $m \neq n$, one finds that all the inner products between $|\psi\rangle$ and the discrete fixed-spectrum Fock basis elements are 0. As a result, it cannot be represented in terms of the discrete Fock basis, which means that the discrete fixed-spectrum Fock basis is not complete. Again, our attempt fails.

C. Completeness of the fixed-spectrum Fock basis

For the moment, we ignore the orthogonality requirement and only focus on completeness. In what follows, we

eventually see that it makes sense to include all functions as spectra, and not just those that form an orthogonal basis for the spatial degrees of freedom. A completeness condition over such a space would naturally lead to a path-integral formulation [42–44], because integrals over such a space are functional integrals that run over all the possible spectral functions in the space.

An expansion of an arbitrary pure state in terms of the full set of fixed-spectrum Fock states would have the form of a functional integral,

$$|\psi\rangle = \sum_n \int |n_F\rangle \mathcal{C}_n[F] \mathcal{D}[F], \quad (13)$$

where F represents the complex functions for the spectra that define the fixed-spectrum Fock basis, the coefficient function $\mathcal{C}_n[F]$ is a functional (i.e., a function of functions), and the measure of the integral $\mathcal{D}[F]$ runs over all such functions F .

To investigate the completeness of the set of all fixed-spectrum Fock states, we consider the possibility of resolving the identity operator in terms of this set. For this purpose, we consider the functional integral for an operator given by

$$\hat{L} = \sum_n \int |n_F\rangle \langle n_F| \mathcal{D}[F]. \quad (14)$$

Using Eq. (7), one can write it as

$$\begin{aligned} \hat{L} &= \sum_n \int \frac{1}{n!} (|1_F\rangle \langle 1_F|)^n \mathcal{D}[F] \\ &= \sum_n \int \frac{1}{n!} \left(\int |\mathbf{k}_1, r\rangle F_r(\mathbf{k}_1) F_s^*(\mathbf{k}_2) \langle \mathbf{k}_2, s| \right. \\ &\quad \left. \times \tilde{d}_r k_1 \tilde{d}_s k_2 \right)^n \mathcal{D}[F]. \end{aligned} \quad (15)$$

To evaluate the functional integral, we interpret it as an ensemble averaging process, but with some differences. The majority of functions in the space over which the functional integral runs would be similar to random functions. It is also reasonable to assume that the function values of most of these functions would be normally distributed. Those functions in the space that do not qualify as normally distributed random functions would form a subset of measure 0. To make this statement stronger than a mere assumption, one can specify it as a condition in the definition of the fixed-spectrum Fock basis. As a result, the functional integral of the product of a function with its complex conjugate, evaluated at different arguments, would produce the equivalent of a completeness condition over the space of functions:

$$\int F_r(\mathbf{k}_1) F_s^*(\mathbf{k}_2) \mathcal{D}[F] = (2\pi)^3 \omega_1 \delta_{r,s} \delta(\mathbf{k}_1 - \mathbf{k}_2). \quad (16)$$

Note that there are two differences between what we have here and what is usually implied in delta-correlated random functions $\langle \chi(x_1) \chi^*(x_2) \rangle = \Delta \delta(x_1 - x_2)$. The first difference is that the ensemble average always implies that the sum is divided by the number of elements. In the functional integral, there is no such division process implied. The second difference is that, to produce the singularity at the origin of the Dirac delta function, the random functions must have

amplitudes of infinite magnitude. In the space of spectra that we consider, all functions are of finite energy, implying finite amplitudes. These two differences cancel each other to produce effectively the same result. Since we sum over an infinite number of finite amplitudes but do not divide by the number of functions, we end up with a divergent value at the origin. It is also assumed that

$$\int F_r(\mathbf{k}_1)F_s(\mathbf{k}_2) \mathcal{D}[F] = \int F_r^*(\mathbf{k}_1)F_s^*(\mathbf{k}_2) \mathcal{D}[F] = 0. \quad (17)$$

The functional integrals over products of more than two functions either give 0 for uneven numbers or break up into a sum of products of integrals over just two functions for even numbers of functions, analogous to the ensemble averages of products of normally distributed random functions. For example,

$$\begin{aligned} & \int F_r(\mathbf{k}_1)F_s^*(\mathbf{k}_2)F_u(\mathbf{k}_3)F_v^*(\mathbf{k}_4) \mathcal{D}[F] \\ &= (2\pi)^6 \omega_1 \omega_3 [\delta_{r,s} \delta_{u,v} \delta(\mathbf{k}_1 - \mathbf{k}_2) \delta(\mathbf{k}_3 - \mathbf{k}_4) \\ &+ \delta_{r,v} \delta_{u,s} \delta(\mathbf{k}_1 - \mathbf{k}_4) \delta(\mathbf{k}_3 - \mathbf{k}_2)]. \end{aligned} \quad (18)$$

In general, we can express the completeness condition to all orders as

$$\begin{aligned} & \int \prod_{m=1}^M F(\mathbf{k}_m; r_m) \prod_{n=1}^N F^*(\mathbf{q}_n; s_n) \mathcal{D}[F] \\ &= \delta_{M,N} \sum_{\text{pert}} \prod_{n=1}^N (2\pi)^3 \omega_n \delta[r_n, s_{\mathcal{P}n}] \delta(\mathbf{k}_n - \mathbf{q}_{\mathcal{P}n}), \end{aligned} \quad (19)$$

where the summation runs over all permutations and the subscript $\mathcal{P}n$ represents a permutation of all N indices in the product. The Kronecker delta function for the spin is represented as $\delta[r, s] \equiv \delta_{r,s}$. We see that each function F needs to be matched to a complex conjugate F^* , and vice versa. If any of these functions remains unmatched, the result is 0.

The relationship in Eq. (19) serves as a generalized completeness condition for the space of all functions. It is now considered as a defining condition for the space of functions that defines the fixed-spectrum Fock basis.

As a result, the expression in Eq. (15) becomes

$$\hat{L} = \sum_n \left(\int |\mathbf{k}, s\rangle \langle \mathbf{k}, s| \mathcal{d}_s k \right)^n = \sum_n \mathcal{I}^n \equiv \mathbb{1}, \quad (20)$$

where \mathcal{I}^n is a projection operator for n -particle states (an identity operator within the subspace of n -particle states) and $\mathbb{1}$ is the identity operator for the entire space, including states with arbitrary numbers of particles. The representation succeeds as a resolution of the identity operator, which shows that, provided that the condition in Eq. (19) is satisfied, the fixed-spectrum Fock basis is a complete basis for both the particle-number degrees of freedom and the spatiotemporal degrees of freedom:

$$\sum_n \int |n_F\rangle \langle n_F| \mathcal{D}[F] = \mathbb{1}. \quad (21)$$

As a result, we obtained a successful completeness condition at the cost of orthogonality of the basis. Based on Eqs. (17) and (21), we also have

$$\int |m_F\rangle \langle n_F| \mathcal{D}[F] = \delta_{mn} \mathcal{I}^n, \quad (22)$$

which we need later.

D. Fixed-spectrum coherent states

In our pursuit for a complete orthogonal basis in spatiotemporal and particle-number degrees of freedom, any attempt to consider coherent states may seem like a waste of time, because one already knows that coherent states are not mutually orthogonal. However, we consider the completeness properties of fixed-spectrum coherent states for the benefit of later use.

The fixed-spectrum coherent states can be defined as an expansion in terms of fixed-spectrum Fock states

$$|\alpha_F\rangle = \exp\left(-\frac{1}{2}|\alpha|^2\right) \sum_{n=0}^{\infty} \frac{\alpha^n}{\sqrt{n!}} |n_F\rangle, \quad (23)$$

where α is a complex constant. In terms of the fixed-spectrum displacement operator, they are given by

$$|\alpha_F\rangle = \hat{D}(\alpha_F)|\text{vac}\rangle = \exp(\hat{a}_\alpha^\dagger - \hat{a}_\alpha)|\text{vac}\rangle, \quad (24)$$

where

$$\begin{aligned} \hat{a}_\alpha^\dagger &= \int \alpha_s(\mathbf{k}) \hat{a}_s^\dagger(\mathbf{k}) \mathcal{d}_s k, \\ \hat{a}_\alpha &= \int \alpha_s^*(\mathbf{k}) \hat{a}_s(\mathbf{k}) \mathcal{d}_s k. \end{aligned} \quad (25)$$

The complex function $\alpha_s(\mathbf{k})$ is composed of the product of the complex parameter α and the normalized complex function $F_s(\mathbf{k})$. Note that, according to Eq. (10),

$$\begin{aligned} \|\alpha_s(\mathbf{k})\|^2 &\equiv \int |\alpha_s(\mathbf{k})|^2 \mathcal{d}_s k \\ &= |\alpha|^2 \int |F_s(\mathbf{k})|^2 \mathcal{d}_s k = |\alpha|^2. \end{aligned} \quad (26)$$

The inner product between different fixed-spectrum coherent states (i.e., where the photons in the respective coherent states in general have completely different spectra) can be calculated with the aid of Eq. (11). It reads

$$\begin{aligned} \langle \alpha_F | \beta_G \rangle &= \exp\left(-\frac{1}{2}|\alpha|^2 - \frac{1}{2}|\beta|^2 + \langle \alpha, \beta \rangle\right) \\ &= \exp\left(-\frac{1}{2}\|\alpha_s(\mathbf{k}) - \beta_s(\mathbf{k})\|^2\right) \exp(i\Theta), \end{aligned} \quad (27)$$

where, from Eq. (26),

$$\|\alpha_s(\mathbf{k}) - \beta_s(\mathbf{k})\|^2 = \int |\alpha_s(\mathbf{k}) - \beta_s(\mathbf{k})|^2 \mathcal{d}_s k \quad (28)$$

gives the metric distance between the two spectra, and

$$\Theta = \text{Im}\{\langle \alpha, \beta \rangle\} = \text{Im}\left\{ \int \alpha_s^*(\mathbf{k}) \beta_s(\mathbf{k}) \mathcal{d}_s k \right\}, \quad (29)$$

with $\text{Im}\{\cdot\}$ giving the imaginary part of the argument.

Therefore, the inner product between different fixed-spectrum coherent states is related to the metric distance between their associated complex spectra in the space of functions. So, even when $\langle \alpha, \beta \rangle = 0$, we still have $\langle \alpha_F | \beta_G \rangle \neq 0$. Not surprisingly, it is not possible to define a basis in terms of such coherent states that are orthogonal. The space of functions for all fixed-spectrum coherent states forms a metric space (or normed vector space). The metric in the space of functions is related to the inner product between the associated fixed-spectrum coherent states, as

$$d\{\alpha, \beta\} \equiv \|\alpha_s(\mathbf{k}) - \beta_s(\mathbf{k})\|^2 = -\ln(|\langle \alpha_F | \beta_G \rangle|^2). \quad (30)$$

To investigate the completeness of the fixed-spectrum coherent bases, we first consider

$$\hat{L} = \int |\alpha_F\rangle \langle \beta_F| \mathcal{D}[F], \quad (31)$$

where the complex constants α and β are allowed to be different and where the functional integral runs only over the normalized functions F . Using the definition of the coherent states in Eq. (23) and applying Eq. (22), we get

$$\hat{L} = \exp\left(-\frac{1}{2}|\alpha|^2 - \frac{1}{2}|\beta|^2 + \alpha^* \beta \mathcal{I}\right). \quad (32)$$

By setting $\alpha = \beta$, we obtain

$$\int |\alpha_F\rangle \langle \alpha_F| \mathcal{D}[F] = \exp(-|\alpha|^2 + |\alpha|^2 \mathcal{I}). \quad (33)$$

Next, we also integrate over α :

$$\begin{aligned} \int |\alpha_F\rangle \langle \alpha_F| \mathcal{D}[\alpha_F] &= \int \exp(-|\alpha|^2 + |\alpha|^2 \mathcal{I}) d\alpha \\ &= \pi \sum_{n=0}^{\infty} \mathcal{I}^n = \pi \mathbb{1}. \end{aligned} \quad (34)$$

Hence, the functional integral over all complex functions $\alpha_F = \alpha F(\mathbf{k})$ provides us with a completeness condition,

$$\mathbb{1} = \frac{1}{\pi} \int |\alpha_F\rangle \langle \alpha_F| \mathcal{D}[\alpha_F], \quad (35)$$

which is analogous to what one obtains without the spatiotemporal degrees of freedom. The fixed-spectrum coherent states represent a (over-) complete basis, even though they are not orthogonal.

Since the identity is idempotent, we must have $\mathbb{1}^2 = \mathbb{1}$. Due to the nonorthogonality of the coherent states, Eq. (27), it leads to the awkward identity

$$\begin{aligned} \mathbb{1} &= \frac{1}{\pi^2} \int |\alpha_F\rangle \langle \alpha_F | \beta_G\rangle \langle \beta_G| \mathcal{D}[\alpha_F] \mathcal{D}[\beta_G] \\ &= \frac{1}{\pi^2} \int |\alpha_F\rangle \exp\left(-\frac{1}{2}|\alpha|^2 - \frac{1}{2}|\beta|^2 + \langle \alpha, \beta \rangle\right) \langle \beta_G| \\ &\quad \times \mathcal{D}[\alpha_F] \mathcal{D}[\beta_G], \end{aligned} \quad (36)$$

which comes in handy later.

E. Fixed-spectrum quadrature bases

One can define fixed-spectrum quadrature operators directly in terms of the fixed-spectrum creation and annihilation

operators

$$\begin{aligned} \hat{q}_F &= \frac{1}{\sqrt{2}}(\hat{a}_F + \hat{a}_F^\dagger), \\ \hat{p}_F &= \frac{-i}{\sqrt{2}}(\hat{a}_F - \hat{a}_F^\dagger). \end{aligned} \quad (37)$$

Their commutation relation is given by

$$[\hat{q}_F, \hat{p}_G] = i\text{Re}\{\langle F, G \rangle\}, \quad (38)$$

where $\text{Re}\{\cdot\}$ represents the real part of the expression.

The fixed-spectrum quadrature basis elements are the eigenstates of the fixed-spectrum quadrature operators

$$\begin{aligned} \hat{q}_F |q_F\rangle &= |q_F\rangle q, \\ \hat{p}_F |p_F\rangle &= |p_F\rangle p \end{aligned} \quad (39)$$

and are assumed to be given by the expansions

$$\begin{aligned} |q_F\rangle &= \sum_n |n_F\rangle \Theta_n(q_F), \\ |p_F\rangle &= \sum_n |n_F\rangle \Phi_n(p_F) \end{aligned} \quad (40)$$

in terms of the fixed-spectrum Fock states. To find the expressions for the coefficient functions, we compute the overlaps

$$\begin{aligned} \Theta_n(q_F) &= \langle n_F | q_F \rangle = \frac{1}{\sqrt{n!}} \langle \text{vac} | (\hat{a}_F)^n | q_F \rangle, \\ \Phi_n(p_F) &= \langle n_F | p_F \rangle = \frac{1}{\sqrt{n!}} \langle \text{vac} | (\hat{a}_F)^n | p_F \rangle, \end{aligned} \quad (41)$$

where we have used Eq. (7) to express the fixed-spectrum Fock states in terms of fixed-spectrum annihilation operators. We now convert the products of annihilation operators into generating functions for such products,

$$\begin{aligned} \sum_n \frac{\eta^n}{\sqrt{2^n n!}} \Theta_n(q_F) &= \langle \text{vac} | \exp\left(\frac{\eta}{\sqrt{2}} \hat{a}_F\right) | q_F \rangle, \\ \sum_n \frac{(-i)^n \eta^n}{\sqrt{2^n n!}} \Phi_n(p_F) &= \langle \text{vac} | \exp\left(\frac{-i\eta}{\sqrt{2}} \hat{a}_F\right) | p_F \rangle, \end{aligned} \quad (42)$$

where η is the generating parameter and where we have introduced convenient constants in anticipation of our goal. Next, we exploit the fact that

$$\langle \text{vac} | \exp(K \hat{a}_F^\dagger) = \langle \text{vac} | \quad (43)$$

to insert exponentiated creation operators,

$$\begin{aligned} \sum_n \frac{\eta^n}{\sqrt{2^n n!}} \Theta_n(q_F) &= \langle \text{vac} | \exp\left(\frac{\eta}{\sqrt{2}} \hat{a}_F^\dagger\right) \exp\left(\frac{\eta}{\sqrt{2}} \hat{a}_F\right) | q_F \rangle, \\ \sum_n \frac{(-i)^n \eta^n}{\sqrt{2^n n!}} \Phi_n(p_F) &= \langle \text{vac} | \exp\left(\frac{i\eta}{\sqrt{2}} \hat{a}_F^\dagger\right) \exp\left(\frac{-i\eta}{\sqrt{2}} \hat{a}_F\right) | p_F \rangle. \end{aligned} \quad (44)$$

The exponential operators are combined using the Baker-Campbell-Hausdorff formula. The resulting combined exponential operators can be expressed in terms of fixed-spectrum quadrature operators. Hence,

$$\begin{aligned} \sum_n \frac{\eta^n}{\sqrt{2^n n!}} \Theta_n(q_F) &= \langle \text{vac} | \exp\left(\eta \hat{q}_F - \frac{\eta^2}{4}\right) | q_F \rangle \\ &= \Theta_0(q_F) \exp\left(\eta q - \frac{\eta^2}{4}\right), \\ \sum_n \frac{(-i)^n \eta^n}{\sqrt{2^n n!}} \Phi_n(p_F) &= \langle \text{vac} | \exp\left(\eta \hat{p}_F - \frac{\eta^2}{4}\right) | p_F \rangle \\ &= \Phi_0(p_F) \exp\left(\eta p - \frac{\eta^2}{4}\right), \end{aligned} \quad (45)$$

where we have used the eigenvalue equations in Eq. (39) to pull the exponentiated quadrature operators through the basis elements. These results resemble the generating function for Hermite polynomials, given by

$$\exp(2xv - v^2) = \sum_{n=0}^{\infty} \frac{v^n}{n!} H_n(x), \quad (46)$$

where v is the generating parameter. The coefficient functions are therefore given by

$$\begin{aligned} \Theta_n(q_F) &= \frac{\Theta_0(q_F)}{\sqrt{2^n n!}} H_n(q), \\ \Phi_n(p_F) &= \frac{(i)^n \Phi_0(p_F)}{\sqrt{2^n n!}} H_n(p) \end{aligned} \quad (47)$$

in terms of Hermite polynomials, up to the zeroth-order coefficient functions $\Theta_0(q)$ and $\Phi_0(p)$, which are yet to be determined. At this point, we note that the possible dependences on the normalized spectrum F appear only in the zeroth-order coefficient functions. So, if they depend only on the q and p parameters, and not on the normalized spectrum, then the same will be true for all coefficient functions. Therefore, we assume that one can replace $q_F \rightarrow q$ and $p_F \rightarrow p$ in these expressions.

To find the expressions for the zeroth-order coefficient functions, we consider the inner product between coefficient functions of different orders. Comparing the resulting expressions with the orthogonality condition for Hermite polynomials

$$\int H_m(x) H_n(x) \exp(-x^2) dx = \sqrt{\pi} 2^n n! \delta_{mn}, \quad (48)$$

we see that, if

$$\begin{aligned} |\Theta_0(q)|^2 &= \frac{1}{\sqrt{\pi}} \exp(-q^2), \\ |\Phi_0(p)|^2 &= 2\sqrt{\pi} \exp(-p^2), \end{aligned} \quad (49)$$

then the coefficient functions will obey the orthogonality conditions

$$\begin{aligned} \int \Theta_m(q) \Theta_n^*(q) dq &= \delta_{mn}, \\ \int \Phi_m(p) \Phi_n^*(p) dp &= 2\pi \delta_{mn}. \end{aligned} \quad (50)$$

Assuming that the zeroth-order coefficient functions are real-valued functions, we obtain expressions for the coefficient functions, given by

$$\begin{aligned} \Theta_n(q_F) &= \frac{1}{\pi^{1/4} \sqrt{2^n n!}} H_n(q) \exp\left(-\frac{q^2}{2}\right), \\ \Phi_n(p_F) &= \frac{(i)^n \sqrt{2} \pi^{1/4}}{\sqrt{2^n n!}} H_n(p) \exp\left(-\frac{p^2}{2}\right). \end{aligned} \quad (51)$$

These coefficient functions are the same as those one has without the spatiotemporal degrees of freedom. In addition to the orthogonality conditions given in Eq. (50), the coefficient functions of Eq. (51) also obey completeness conditions, given by

$$\begin{aligned} \sum_n \Theta_n^*(q) \Theta_n(q') &= \delta(q - q'), \\ \sum_n \Phi_n^*(p) \Phi_n(p') &= 2\pi \delta(p - p'). \end{aligned} \quad (52)$$

What is the effect of the spectra on the inner products among quadrature basis elements? Based on Eq. (11), the inner product between q states with different spectra becomes

$$\langle q_F | q'_G \rangle = \sum_n \mu^n \Theta_n^*(q) \Theta_n(q'), \quad (53)$$

where $\mu \equiv \langle F, G \rangle$. Using Eq. (51) and Mehler's formula [45], which is given by

$$\begin{aligned} \sum_{n=0}^{\infty} \frac{\rho^n}{2^n n!} H_n(x) H_n(y) \\ = \frac{1}{\sqrt{1 - \rho^2}} \exp\left[\frac{2\rho xy}{1 - \rho^2} - \frac{(x^2 + y^2)\rho^2}{1 - \rho^2}\right], \end{aligned} \quad (54)$$

we obtain an expression for the inner product that reads

$$\begin{aligned} \langle q_F | q'_G \rangle &= \frac{1}{\sqrt{\pi(1 - \mu^2)}} \exp\left[-\frac{\mu(q - q')^2}{1 - \mu^2}\right] \\ &\times \exp\left[-\frac{(1 - \mu)(q^2 + q'^2)}{2(1 + \mu)}\right]. \end{aligned} \quad (55)$$

For $G \rightarrow F$, but keeping the q 's different, we have $\mu \rightarrow 1$. We see that Eq. (55) is singular for $\mu \rightarrow 1$. So we replace $\mu = 1 + \epsilon/2$ and consider the limit where $\epsilon \rightarrow 0$. It gives

$$\langle q_F | q'_F \rangle = \lim_{\epsilon \rightarrow 0} \frac{1}{\sqrt{\pi\epsilon}} \exp\left[-\frac{(q - q')^2}{\epsilon}\right]. \quad (56)$$

One can show that the right-hand side represents a limit process for the Dirac delta function

$$\lim_{\epsilon \rightarrow 0} \frac{1}{\sqrt{\pi\epsilon}} \exp\left[-\frac{(q - q')^2}{\epsilon}\right] = \delta(q - q'). \quad (57)$$

Hence, we obtain the orthogonality condition

$$\langle q_F | q'_F \rangle = \delta(q - q_0). \quad (58)$$

A similar condition applies for the p basis.

On the other hand, if F and G are orthogonal, we have $\mu \rightarrow 0$, so that

$$\langle q_F | q'_G \rangle \rightarrow \Theta_0^*(q) \Theta_0(q') = \frac{1}{\sqrt{\pi}} \exp \left[\frac{-(q^2 + q'^2)}{2} \right]. \quad (59)$$

As a result, the fixed spectrum quadrature basis elements lose their orthogonality. Even if we use a discrete modal basis for the spatiotemporal degrees of freedom, we still do not obtain an orthogonal basis for both particle-number and spatiotemporal degrees of freedom. So, if we remain within the subspace associated with a specific spectral function, the fixed-spectrum quadrature bases are orthogonal bases, but beyond that they are not orthogonal.

F. Completeness of fixed-spectrum quadrature bases

Considering the completeness of the fixed-spectrum quadrature bases, we first restrict the bases to the subspace of a specific spectral function and integrate over the q parameter. We obtain

$$\begin{aligned} \int |q_F\rangle \langle q_F| dq &= \sum_{m,n} |m_F\rangle \langle n_F| \Theta_m(q) \Theta_n^*(q) dq \\ &= \sum_n |n_F\rangle \langle n_F| = \mathcal{I}_F, \end{aligned} \quad (60)$$

thanks to the orthogonality of the coefficient functions. The result represents a projection operator for states with arbitrary numbers of photons, but where their spatiotemporal degrees of freedom are defined by a specific spectral function F .

To investigate the completeness of the fixed-spectrum quadrature bases for the entire space of spectral functions, we need to employ functional integrals, as done in Sec. II C. First, we consider the case where the spectrum is the same, but the q parameters are different:

$$\hat{L} = \int |q_F\rangle \langle q'_F| \mathcal{D}[F]. \quad (61)$$

Substituting Eq. (40), we obtain

$$\begin{aligned} \hat{L} &= \sum_{mn} \int |n_F\rangle \Theta_m(q) \Theta_n^*(q') \langle m_F| \mathcal{D}[F], \\ &= \sum_n \Theta_n(q) \Theta_n^*(q') \mathcal{I}^n, \end{aligned} \quad (62)$$

where we have used Eq. (22). It cannot be simplified further. However, if we also integrate over q and use the orthogonality of the coefficient functions, Eq. (50), we get

$$\begin{aligned} \int \hat{L} dq &= \sum_n \int \Theta_n(q) \Theta_n^*(q) dq \mathcal{I}^n \\ &= \sum_n \mathcal{I}^n = \mathbb{1}. \end{aligned} \quad (63)$$

So, the entire space, including all normalized functions and all q parameters, provides a complete basis. The same applies for the p basis.

We are almost there. The fixed-spectrum quadrature bases are complete, but not orthogonal, unless we restrict them to subspaces for specific spectral functions.

III. EIGENSTATES OF FIXED-MOMENTUM QUADRATURE OPERATORS

The stage is set to derive bases that are both complete and orthogonal with respect to the entire space of quantum states that can be defined in terms of both spatiotemporal degrees of freedom and particle-number degrees of freedom. These bases are what we call *spatiotemporal* quadrature bases.

For derivation of the spatiotemporal quadrature bases, we start with the notion of *fixed-momentum* quadrature operators. These operators are directly defined in terms of the creation and annihilation operators for photonic states, which obey the Lorentz covariant commutation relation given in Eq. (1). The fixed-momentum quadrature operators are given by

$$\begin{aligned} \hat{q}_s(\mathbf{k}) &= \frac{1}{\sqrt{2}} [\hat{a}_s(\mathbf{k}) + \hat{a}_s^\dagger(\mathbf{k})], \\ \hat{p}_s(\mathbf{k}) &= \frac{-i}{\sqrt{2}} [\hat{a}_s(\mathbf{k}) - \hat{a}_s^\dagger(\mathbf{k})]. \end{aligned} \quad (64)$$

They obey a Lorentz covariant commutation relation that reads

$$[\hat{q}_r(\mathbf{k}_1), \hat{p}_s(\mathbf{k}_2)] = i(2\pi)^3 \omega_1 \delta_{r,s} \delta(\mathbf{k}_1 - \mathbf{k}_2). \quad (65)$$

The term *fixed-momentum* quadrature operator follows from the fact that they explicitly depend on the value of the wave vector.

It is now assumed that the fixed-momentum quadrature operators give rise to eigenstates and eigenvalue functions according to the following eigenvalue equations:

$$\begin{aligned} \hat{q}_s(\mathbf{k})|q\rangle &= |q\rangle q_s(\mathbf{k}), \\ \hat{p}_s(\mathbf{k})|p\rangle &= |p\rangle p_s(\mathbf{k}). \end{aligned} \quad (66)$$

The eigenvalue functions are real-valued functions because the quadrature operators are Hermitian. They can be separated into a real-valued magnitude times real-valued normalized functions $q_s(\mathbf{k}) = q F_s(\mathbf{k})$ and $p_s(\mathbf{k}) = p G_s(\mathbf{k})$. For the sake of notational simplicity, we generally distinguish between the whole eigenvalue function and the magnitude simply by the presence or absence of the argument. The functions $F_s(\mathbf{k})$ and $G_s(\mathbf{k})$ are normalized in the sense

$$\int F_s^2(\mathbf{k}) d_s k = \int G_s^2(\mathbf{k}) d_s k = 1. \quad (67)$$

In the derivation that follows, we ignore the spin index to alleviate the complexity of the expressions. After the derivation, we reintroduce the spin indices.

It has been shown in Sec. II A that, when the eigenstates depend explicitly on the momentum, they will not be well defined. Therefore, we assume that all the momentum dependences are integrated out. As a result, $|q\rangle \neq |q(\mathbf{k})\rangle$ and $|p\rangle \neq |p(\mathbf{k})\rangle$. Focusing on the first equation in Eq. (66), we therefore start with an ansatz for $|q\rangle$ of the form

$$\begin{aligned} |q\rangle &= |\text{vac}\rangle V_0 + \int |\mathbf{k}\rangle V_1(\mathbf{k}) d_k \\ &+ \frac{1}{2!} \int |\mathbf{k}_a\rangle |\mathbf{k}_b\rangle V_2(\mathbf{k}_a, \mathbf{k}_b) d_k a d_k b \\ &+ \frac{1}{3!} \int |\mathbf{k}_a\rangle |\mathbf{k}_b\rangle |\mathbf{k}_c\rangle V_3(\mathbf{k}_a, \mathbf{k}_b, \mathbf{k}_c) d_k a d_k b d_k c + \dots, \end{aligned} \quad (68)$$

where $V_n(\cdot)$ denotes coefficient functions to be determined. The subscript n is not the spin; it represents the order of the term in which it appears. Applying the eigenvalue equation to the ansatz and solving the coefficient functions order by order, one obtains expressions for all coefficient functions with $n > 0$ in terms of V_0 . The first few are given by

$$\begin{aligned} V_1(\mathbf{k}) &= \sqrt{2}V_0q(\mathbf{k}), \\ V_2(\mathbf{k}_a, \mathbf{k}_b) &= 2V_0q(\mathbf{k}_a)q(\mathbf{k}_b) - V_0\delta(\mathbf{k}_a - \mathbf{k}_b), \\ V_3(\mathbf{k}_a, \mathbf{k}_b, \mathbf{k}_c) &= 2\sqrt{2}V_0q(\mathbf{k}_a)q(\mathbf{k}_b)q(\mathbf{k}_c) \\ &\quad - 3\sqrt{2}V_0q(\mathbf{k}_a)\delta(\mathbf{k}_b - \mathbf{k}_c), \end{aligned} \quad (69)$$

and so forth. Substituting these coefficient functions back into the ansatz in Eq. (68), we obtain an expression of the form

$$\begin{aligned} |q\rangle &= |\text{vac}\rangle V_0 + (|Q\rangle - |R\rangle)V_0 + \frac{1}{2!}(|Q\rangle - |R\rangle)^2 V_0 \\ &\quad + \frac{1}{3!}(|Q\rangle - |R\rangle)^3 V_0 + \dots \\ &= V_0 \sum_{m=0}^{\infty} \frac{1}{m!} (|Q\rangle - |R\rangle)^m \\ &= V_0 \exp(|Q\rangle - |R\rangle), \end{aligned} \quad (70)$$

where V_0 is a global constant and

$$\begin{aligned} |Q\rangle &\equiv \sqrt{2} \int |\mathbf{k}, s\rangle q_s(\mathbf{k}) \tilde{d}_s k, \\ |R\rangle &\equiv \frac{1}{2} \int |\mathbf{k}, s\rangle |\mathbf{k}, s\rangle \tilde{d}_s k, \end{aligned} \quad (71)$$

with the spin indices being reintroduced. Note that

$$\begin{aligned} \hat{a}_s(\mathbf{k}')|Q\rangle &= |\text{vac}\rangle \sqrt{2}q_s(\mathbf{k}'), \\ \hat{a}_s(\mathbf{k}')|R\rangle &= |\mathbf{k}', s\rangle, \end{aligned} \quad (72)$$

which implies that

$$\begin{aligned} \hat{a}_s(\mathbf{k}') \exp(|Q\rangle) &= \exp(|Q\rangle) \sqrt{2}q_s(\mathbf{k}'), \\ \hat{a}_s(\mathbf{k}') \exp(-|R\rangle) &= -\exp(-|R\rangle) |\mathbf{k}', s\rangle. \end{aligned} \quad (73)$$

The creation operator just adds a factor of $|\mathbf{k}'\rangle$. As an aside, it is interesting that $\exp(|R\rangle)$ acts like an ‘antivacuum’ state in the sense that the application of an annihilation operator creates a ket vector in the same way that it would normally be created by a creation operator acting on the vacuum state.

One can use the results in Eq. (73) to test whether our solution satisfies the first eigenvalue equation in Eq. (66):

$$\begin{aligned} \hat{q}_s(\mathbf{k})|q\rangle &= \frac{V_0}{\sqrt{2}} [\exp(|Q\rangle) \exp(-|R\rangle) \sqrt{2}q_s(\mathbf{k}) \\ &\quad - \exp(|Q\rangle) \exp(-|R\rangle) |\mathbf{k}, s\rangle \\ &\quad + \exp(|Q\rangle) \exp(-|R\rangle) |\mathbf{k}, s\rangle] \\ &= |q\rangle q_s(\mathbf{k}). \end{aligned} \quad (74)$$

It confirms that the expression for the eigenstates, given in Eq. (70), satisfies the eigenvalue equation. However, the constant V_0 is still unspecified.

A similar procedure can be followed to obtain an expression for the eigenstate $|p\rangle$. It is given by

$$|p\rangle = W_0 \exp(i|P\rangle + |R\rangle), \quad (75)$$

where W_0 is a global constant, $|R\rangle$ is given in Eq. (71), and

$$|P\rangle \equiv \sqrt{2} \int |\mathbf{k}, s\rangle p_s(\mathbf{k}) \tilde{d}_s k. \quad (76)$$

IV. ORTHOGONALITY

Since the fixed-momentum quadrature operators are Hermitian, it makes sense that the spatiotemporal quadrature bases would be orthogonal bases. However, what does it mean in the context of an infinite functional space? To investigate the orthogonality of these bases and to find expressions for V_0 and W_0 , we need to compute $\langle q|q'\rangle$ and $\langle p|p'\rangle$. At the same time, it is helpful to compute $\langle q|p\rangle$. For this purpose, we follow an operator approach, defining annihilation operators for $|Q\rangle$, $|P\rangle$, and $|R\rangle$, given by

$$\begin{aligned} \hat{a}_Q &= \sqrt{2} \int \hat{a}_s(\mathbf{k}) q_s(\mathbf{k}) \tilde{d}_s k, \\ \hat{a}_P &= \sqrt{2} \int \hat{a}_s(\mathbf{k}) p_s(\mathbf{k}) \tilde{d}_s k, \\ \hat{a}_R &= \frac{1}{2} \int \hat{a}_s(\mathbf{k}) \hat{a}_s(\mathbf{k}) \tilde{d}_s k. \end{aligned} \quad (77)$$

The corresponding creation operators are obtained as the adjoint operators, so that $|Q\rangle = \hat{a}_Q^\dagger |\text{vac}\rangle$, $|P\rangle = \hat{a}_P^\dagger |\text{vac}\rangle$, and $|R\rangle = \hat{a}_R^\dagger |\text{vac}\rangle$. Using these operators, we then define creation operators for the eigenstates

$$\begin{aligned} \hat{a}_q^\dagger &= V_0 \exp(\hat{a}_Q^\dagger - \hat{a}_R^\dagger), \\ \hat{a}_p^\dagger &= W_0 \exp(i\hat{a}_P^\dagger + \hat{a}_R^\dagger), \end{aligned} \quad (78)$$

so that $|q\rangle = \hat{a}_q^\dagger |\text{vac}\rangle$ and $|p\rangle = \hat{a}_p^\dagger |\text{vac}\rangle$. The overlaps among the eigenstates are then given by

$$\begin{aligned} \langle q|q'\rangle &= \langle \text{vac} | \hat{a}_q \hat{a}_q^\dagger | \text{vac} \rangle, \\ \langle p|p'\rangle &= \langle \text{vac} | \hat{a}_p \hat{a}_p^\dagger | \text{vac} \rangle, \\ \langle q|p\rangle &= \langle \text{vac} | \hat{a}_q \hat{a}_p^\dagger | \text{vac} \rangle. \end{aligned} \quad (79)$$

Starting with the last expression in Eq. (79), we obtain

$$\begin{aligned} \langle q|p\rangle &= V_0 W_0 \langle \text{vac} | \exp(\hat{a}_Q) \exp(-\hat{a}_R) \\ &\quad \times \exp(i\hat{a}_P^\dagger) \exp(\hat{a}_R^\dagger) | \text{vac} \rangle. \end{aligned} \quad (80)$$

To evaluate this expression, one can rearrange the exponential operators in normal order. However, in the process new operators are generated in addition to the current ones. As a result, we expect to get an expression of the form

$$\begin{aligned} &\exp(k_1 \hat{a}_Q) \exp(k_2 \hat{a}_R) \exp(k_3 \hat{a}_P^\dagger) \exp(k_4 \hat{a}_R^\dagger) \\ &= \exp(h_0) \exp(h_1 \hat{a}_Q^\dagger) \exp(h_2 \hat{a}_P^\dagger) \exp(h_3 \hat{a}_R^\dagger) \\ &\quad \times \exp(h_4 \hat{s}) \exp(h_5 \hat{a}_Q) \exp(h_6 \hat{a}_P) \exp(h_7 \hat{a}_R), \end{aligned} \quad (81)$$

where k_1, k_2, k_3 , and k_4 are assumed to be known constants; $h_0, h_1, h_2, h_3, h_4, h_5, h_6$, and h_7 are unknown constants; and

$$\hat{s} \equiv \frac{1}{2} \int [\hat{a}_s^\dagger(\mathbf{k})\hat{a}_s(\mathbf{k}) + \hat{a}_s(\mathbf{k})\hat{a}_s^\dagger(\mathbf{k})] d_s k. \quad (82)$$

To obtain the required relationship, we follow a standard procedure where one introduces an auxiliary variable t into the exponents on the left-hand side and converts the unknown constants into unknown functions of t on the right-hand side:

$$\begin{aligned} & \exp(tk_1\hat{a}_Q) \exp(tk_2\hat{a}_R) \exp(tk_3\hat{a}_P^\dagger) \exp(tk_4\hat{a}_R^\dagger) \\ &= \exp[h_0(t)] \exp[h_1(t)\hat{a}_Q^\dagger] \exp[h_2(t)\hat{a}_P^\dagger] \exp[h_3(t)\hat{a}_R^\dagger] \\ & \quad \times \exp[h_4(t)\hat{s}] \exp[h_5(t)\hat{a}_Q] \exp[h_6(t)\hat{a}_P] \exp[h_7(t)\hat{a}_R]. \end{aligned} \quad (83)$$

For consistency, the unknown functions must go to 0 for $t = 0$.

Next, we apply a derivative with respect to t on both sides and then remove as many of the exponential operators as possible by operating with the respective inverse operators on the right-hand sides of both sides of the equation. The rather complicated expression that is obtained can be simplified with the aid of the identity

$$\begin{aligned} \exp(\hat{X})\hat{Y} \exp(-\hat{X}) &= \hat{Y} + [\hat{X}, \hat{Y}] + \frac{1}{2!}[\hat{X}, [\hat{X}, \hat{Y}]] \\ & \quad + \frac{1}{3!}[\hat{X}, [\hat{X}, [\hat{X}, \hat{Y}]]] + \dots, \end{aligned} \quad (84)$$

where \hat{X} and \hat{Y} are two arbitrary operators. In Appendix A, we provide all the necessary commutation relations to perform this task. The simplified expression can then be separated into different differential equations for all the unknown functions. Upon solving these differential equations, we obtain

$$\begin{aligned} h_0(t) &= \frac{(k_1^2 k_4 q^2 t + k_2 k_3^2 p^2 t + 2k_1 k_3 \mu) t^2}{1 - k_2 k_4 t^2}, \\ h_1(t) &= \frac{k_1 k_4 t^2}{1 - k_2 k_4 t^2}, \\ h_2(t) &= \frac{k_3 t}{1 - k_2 k_4 t^2}, \\ h_3(t) &= \frac{k_4 t}{1 - k_2 k_4 t^2}, \\ h_4(t) &= -\ln(1 - k_2 k_4 t^2), \\ h_5(t) &= \frac{k_1 t}{1 - k_2 k_4 t^2}, \\ h_6(t) &= \frac{k_2 k_3 t^2}{1 - k_2 k_4 t^2}, \\ h_7(t) &= \frac{k_2 t}{1 - k_2 k_4 t^2}, \end{aligned} \quad (85)$$

where $\mu = \langle q, p \rangle$ is the inner product between the eigenvalue functions. One can now substitute the functions back into Eq. (83). For the case under consideration, as given in Eq. (80), we substitute $k_1 = 1, k_2 = -1, k_3 = i$, and $k_4 = 1$. Then we set $t = 1$ to obtain the expression for the product of

exponential operators in normal order:

$$\begin{aligned} \hat{a}_q \hat{a}_p^\dagger &= V_0 W_0 \exp\left(\frac{q^2}{2} + \frac{p^2}{2} + i\mu\right) \exp\left(\frac{\hat{a}_Q^\dagger}{2}\right) \\ & \quad \times \exp\left(\frac{i\hat{a}_P^\dagger}{2}\right) \exp\left(\frac{\hat{a}_R^\dagger}{2}\right) \exp[-\ln(2)\hat{s}] \\ & \quad \times \exp\left(\frac{\hat{a}_Q}{2}\right) \exp\left(\frac{-i\hat{a}_P}{2}\right) \exp\left(\frac{-\hat{a}_R}{2}\right). \end{aligned} \quad (86)$$

The expression for the overlap follows by contracting the vacuum state on both sides. For this purpose, we note that $\exp(K\hat{a}_Q)|\text{vac}\rangle = |\text{vac}\rangle$, regardless of the value of K . The same applies for \hat{a}_P and \hat{a}_R . Using the normal-ordered form for \hat{s} , as given in Eq. (A6), we obtain

$$\langle \text{vac} | \exp[-\ln(2)\hat{s}] | \text{vac} \rangle = \exp[-\ln(2)\Omega] = \frac{1}{2^\Omega}, \quad (87)$$

where Ω is defined in Eq. (A6). So, the overlap becomes

$$\langle q | p \rangle = \frac{V_0 W_0}{2^\Omega} \exp\left(\frac{q^2}{2} + \frac{p^2}{2} + i\mu\right). \quad (88)$$

We now define the global constants as

$$\begin{aligned} V_0 &= 2^{\Omega/2} \exp\left(-\frac{q^2}{2}\right), \\ W_0 &= 2^{\Omega/2} \exp\left(-\frac{p^2}{2}\right). \end{aligned} \quad (89)$$

As a result, the overlap becomes

$$\langle q | p \rangle = \exp(i\mu) = \exp\left[i \int q_s(\mathbf{k}) p_s(\mathbf{k}) d_s k\right]. \quad (90)$$

To consider the other two overlaps in Eq. (80), we note that they can be obtained from the same result with the appropriate substitutions. For $\langle q | q' \rangle$, we need to replace $P \rightarrow Q'$ and $p \rightarrow q'$. The inner product now represents $\mu \rightarrow \langle q, q' \rangle$. Using these replacements, together with the appropriate assignment for the constants — $k_1 = 1, k_2 = -1, k_3 = 1$, and $k_4 = -1$ — we obtain

$$\begin{aligned} h_0(t) &= \frac{-[q^2 t + (q')^2 t - 2\mu] t^2}{1 - t^2}, \\ h_1(t) = h_6(t) &= \frac{-t^2}{1 - t^2}, \\ h_2(t) = -h_3(t) = h_5(t) = -h_7(t) &= \frac{t}{1 - t^2}, \\ h_4(t) &= -\ln(1 - t^2). \end{aligned} \quad (91)$$

These functions are all singular at $t = 1$. Therefore, we need to consider the overlap as a limit. For this purpose, we substitute $t = 1 - \epsilon$ and Eq. (89). The resulting overlap then becomes

$$\langle q | q' \rangle = \lim_{\epsilon \rightarrow 0} \frac{1}{\epsilon^\Omega} \exp\left\{\frac{-[q^2 + (q')^2 - 2\mu]}{2\epsilon}\right\}. \quad (92)$$

Note that

$$q^2 + (q')^2 - 2\mu = \|q_s(\mathbf{k}) - q'_s(\mathbf{k})\|^2 \geq 0, \quad (93)$$

where

$$||f_s(\mathbf{k})||^2 \equiv \int f_s^2(\mathbf{k}) d_s k. \quad (94)$$

For $||q_s(\mathbf{k}) - q'_s(\mathbf{k})||^2 > 0$ the limit in Eq. (92) gives 0, and for $||q_s(\mathbf{k}) - q'_s(\mathbf{k})||^2 = 0$ the limit gives ∞ . Hence, the result behaves as a Dirac delta functional, which runs over an ∞ number of degrees of freedom. It enforces the equality of the eigenvalue functions $q(\mathbf{k})$ and $q'(\mathbf{k})$. We can thus express the overlap as

$$\langle q|q' \rangle = \delta[q_s(\mathbf{k}) - q'_s(\mathbf{k})], \quad (95)$$

where we have ignored some unspecified constant factor. A similar expression applies for the remaining overlap:

$$\langle p|p' \rangle = \delta[p_s(\mathbf{k}) - p'_s(\mathbf{k})]. \quad (96)$$

The expressions in Eqs. (95) and (96) represent the required orthogonality conditions for the spatiotemporal quadrature bases.

V. COMPLETENESS

To investigate the completeness of the spatiotemporal quadrature bases, we consider an operator defined by

$$\hat{B} = \int |q\rangle\langle q| \mathcal{D}[q]. \quad (97)$$

Having shown in Eq. (35) that the fixed-spectrum coherent states can be used to resolve the identity, we operate on both side with identities resolved in terms of coherent states:

$$\begin{aligned} \hat{B} &= \int \mathbb{1}|q\rangle\langle q|\mathbb{1} \mathcal{D}[q] \\ &= \frac{1}{\pi^2} \int |\alpha_F\rangle\langle\alpha_F|q\rangle\langle q|\beta_G\rangle\langle\beta_G| \mathcal{D}[q] \mathcal{D}[\alpha_F] \mathcal{D}[\beta_G]. \end{aligned} \quad (98)$$

To develop the expression further, we need the overlap between the quadrature basis and the fixed-spectrum coherent states. It is computed in Appendix B. After substituting the expressions for these overlaps, Eq. (B5), into the above expression, we evaluate the functional integration over q . The result is

$$\begin{aligned} \hat{B} &= \frac{\kappa}{\pi^2} \int |\alpha_F\rangle \exp \left\{ - \int \frac{1}{2} |\alpha_s(\mathbf{k})|^2 - \alpha_s^*(\mathbf{k})\beta_s(\mathbf{k}) \right. \\ &\quad \left. + \frac{1}{2} |\beta_s(\mathbf{k})|^2 dk \right\} \langle\beta_G| \mathcal{D}[\alpha_F] \mathcal{D}[\beta_G], \end{aligned} \quad (99)$$

where κ as an unspecified constant. Comparing this result with Eq. (36), we see that it is proportional to an identity operator. Hence,

$$\frac{1}{\kappa} \int |q\rangle\langle q| \mathcal{D}[q] = \mathbb{1}. \quad (100)$$

A similar procedure gives

$$\frac{1}{\kappa} \int |p\rangle\langle p| \mathcal{D}[p] = \mathbb{1}. \quad (101)$$

The expressions in Eqs. (100) and (101) imply that the spatiotemporal quadrature bases do indeed obey the completeness conditions.

VI. DISCUSSION AND OUTLOOK

In our endeavor to find bases that incorporate both particle-number degrees of freedom and spatiotemporal degrees of freedom, we have found that the eigenstates of fixed-momentum quadrature operators satisfy our requirements. These quadrature bases are complete and orthogonal with respect to both particle-number degrees of freedom and spatiotemporal degrees of freedom. We refer to these as spatiotemporal quadrature bases, because the term ‘quadrature’ already gives reference to the particle-number degrees of freedom.

We have derived expressions for the spatiotemporal quadrature basis elements, both in terms of integrals over the single-photon momentum basis and in terms of operators that create these basis elements from the vacuum. The latter allows us to compute the overlaps among elements of these bases to show that they give rise to the notion of Dirac delta functionals, which represent orthogonality conditions for the bases. Using functional integrals, we have also shown that the spatiotemporal quadrature bases satisfy the completeness conditions.

One may get the impression that the spatiotemporal quadrature bases are very similar to the fixed-spectrum quadrature bases. The only real difference is the appearance of $|R\rangle$ in the expressions for the spatiotemporal quadrature bases. Yet, it is quite easy to show that the fixed-spectrum quadrature basis elements are not able to serve as eigenstates of the fixed-momentum quadrature operators, according to the eigenvalue equations in Eq. (66). What is the role of $|R\rangle$ that could lead to such a significant difference between these bases? To demonstrate the role of $|R\rangle$, one can try to expand a spatiotemporal quadrature basis element in terms of a fixed-spectrum Fock basis for a spectral function given by the eigenvalue function $q(\mathbf{k})$ for that element. Such an expansion reproduces only the equivalent fixed-spectrum quadrature basis element (without the $|R\rangle$), and not the spatiotemporal quadrature basis element (with the $|R\rangle$). Hence, we conclude that the fixed-spectrum quadrature basis elements are the projections of the spatiotemporal quadrature basis elements onto the subspaces associated with the spectral functions given by $q(\mathbf{k})$. The presence of $|R\rangle$ in the expressions of the spatiotemporal quadrature bases imply that their elements extend beyond the subspaces defined by their spectral functions $q(\mathbf{k})$.

With the spatiotemporal quadrature bases in hand, one can formulate powerful analytical tools to investigate the evolution of quantum states that incorporate all the degrees of freedom that photonic states can possess. The idea is to generalize Wigner functions, which represent the particle-number degrees of freedom, to the notion of Wigner functionals, which incorporate both the particle-number degrees of freedom and the spatiotemporal degrees of freedom. Formally, such a Wigner functional would be defined by a functional integral,

$$\begin{aligned} W[q, p] &\equiv \int \left\langle q + \frac{x}{2} \left| \hat{\rho} \right| q - \frac{x}{2} \right\rangle \\ &\quad \times \exp \left[-i \int p(\mathbf{k})x(\mathbf{k})dk \right] \mathcal{D}[x], \end{aligned} \quad (102)$$

based on the spatiotemporal q basis. Such Wigner functionals would then be able to represent not only quantum states that depend on all these degrees of freedom, but also operators that incorporate all these degrees of freedom. As such, the formalism would be useful for measurements of states with arbitrary numbers of photons that incorporate spatiotemporal degrees of freedom [32–37]. One would compute the predicted outcomes from such measurements with a functional integral,

$$\langle \hat{A} \rangle = \int W_{\hat{A}}[q, p] W_{\hat{\rho}}[q, p] \mathcal{D}[q, p], \quad (103)$$

where \hat{A} represents the Hermitian operator for the measurement.

Apart from the Wigner functionals, one can also consider the generalization of other quasidistributions, such as the Glauber-Sudarshan P distribution [46,47] or the Husimi Q distribution [48]. For such cases, one may employ the fixed-spectrum coherent states, leading to functional integral expressions. For instance, a density operator would be represented in terms of the functional P distribution by

$$\hat{\rho} = \int |\alpha_F\rangle P[\alpha_F] \langle \alpha_F| \mathcal{D}[\alpha_F]. \quad (104)$$

These quasidistributions would be related to the Wigner functionals via functional integral expressions.

Such a formalism based on Wigner functionals (or functional quasidistributions) is currently still a work in progress and is therefore beyond the scope of the current paper. The hope is eventually to develop a tool that would enable one to study physical situations in quantum optics where multiple degrees of freedom are playing significant roles in what is being observed.

APPENDIX A: COMMUTATION RELATIONS

Here, we derive the commutation relations that are associated with the operators in Eq. (77). Our starting point is the commutation relation for $\hat{a}(\mathbf{k})$ and $\hat{a}^\dagger(\mathbf{k})$, given in Eq. (1).

If $[\hat{X}, \hat{Y}] = \hat{Z}$, then $[\hat{Y}^\dagger, \hat{X}^\dagger] = \hat{Z}^\dagger$. So, we do not provide commutation relations that can be obtained directly from others by performing an adjoint operation.

We group the commutation relations in batches. First, all annihilation operators commute among themselves:

$$\begin{aligned} [\hat{a}_s(\mathbf{k}), \hat{a}_Q] &= [\hat{a}_s(\mathbf{k}), \hat{a}_P] = [\hat{a}_s(\mathbf{k}), \hat{a}_R] = 0, \\ [\hat{a}_Q, \hat{a}_Q] &= [\hat{a}_P, \hat{a}_P] = [\hat{a}_R, \hat{a}_R] = 0, \\ [\hat{a}_Q, \hat{a}_P] &= [\hat{a}_Q, \hat{a}_R] = [\hat{a}_P, \hat{a}_R] = 0. \end{aligned} \quad (A1)$$

Unless the relation contains both \hat{a}_Q and \hat{a}_P , as in

$$[\hat{a}_Q, \hat{a}_P^\dagger] = 2 \int q_s(\mathbf{k}) p_s(\mathbf{k}) \vec{d}_s k \equiv 2\mu, \quad (A2)$$

we do not, henceforth, show the commutation relation for them both since they are the same. In Eq. (A2), μ represents the inner product between the functions

$$\mu = \langle q, p \rangle \equiv \int q_s(\mathbf{k}) p_s(\mathbf{k}) \vec{d}_s k. \quad (A3)$$

The next batch involves $\hat{a}^\dagger(\mathbf{k})$:

$$\begin{aligned} [\hat{a}_Q, \hat{a}_s^\dagger(\mathbf{k})] &= \sqrt{2} q_s(\mathbf{k}), \\ [\hat{a}_P, \hat{a}_s^\dagger(\mathbf{k})] &= \sqrt{2} p_s(\mathbf{k}), \\ [\hat{a}_R, \hat{a}_s^\dagger(\mathbf{k})] &= \hat{a}_s(\mathbf{k}). \end{aligned} \quad (A4)$$

Then we have relations that include the R operators,

$$\begin{aligned} [\hat{a}_R, \hat{a}_Q^\dagger] &= \hat{a}_Q, \\ [\hat{a}_R, \hat{a}_R^\dagger] &= \hat{s}, \end{aligned} \quad (A5)$$

where \hat{s} is the symmetrized number operator, given in Eq. (82). It can also be written in normal-ordered form,

$$\hat{s} = \int \hat{a}_s^\dagger(\mathbf{k}) \hat{a}_s(\mathbf{k}) \vec{d}_s k + \frac{1}{2} \int \delta(0) \vec{d}_s k = \hat{n} + \Omega, \quad (A6)$$

where Ω is a divergent constant. The commutation relations involving the symmetrized number operator are

$$\begin{aligned} [\hat{a}_s(\mathbf{k}), \hat{s}] &= \hat{a}_s(\mathbf{k}), \\ [\hat{a}_Q, \hat{s}] &= \hat{a}_Q, \\ [\hat{a}_R, \hat{s}] &= 2\hat{a}_R. \end{aligned} \quad (A7)$$

APPENDIX B: QUADRATURE REPRESENTATIONS OF FIXED-SPECTRUM COHERENT STATES

Here, we consider how to expand fixed-spectrum coherent states in terms of the spatiotemporal quadrature bases. For this purpose, we employ the eigenstate property of the coherent states,

$$\begin{aligned} \hat{a}_Q |\alpha_F\rangle &= |\alpha_F\rangle \sqrt{2} \langle q, \alpha \rangle, \\ \hat{a}_R |\alpha_F\rangle &= |\alpha_F\rangle \frac{1}{2} \mathcal{A}_2, \end{aligned} \quad (B1)$$

where

$$\mathcal{A}_2 \equiv \int \alpha_s^2(\mathbf{k}) \vec{d}_s k. \quad (B2)$$

Therefore,

$$\begin{aligned} \langle q | \alpha_F \rangle &= V_0 \langle \text{vac} | \exp(\hat{a}_Q - \hat{a}_R) | \alpha_F \rangle \\ &= V_0 \langle \text{vac} | \alpha_F \rangle \exp \left[\sqrt{2} \langle q, \alpha \rangle - \frac{1}{2} \mathcal{A}_2 \right]. \end{aligned} \quad (B3)$$

If we express $\alpha_s(\mathbf{k})$ in its real and imaginary parts,

$$\alpha_s(\mathbf{k}) \rightarrow \frac{1}{\sqrt{2}} [q_{s0}(\mathbf{k}) + i p_{s0}(\mathbf{k})], \quad (B4)$$

and use Eq. (89), we obtain

$$\begin{aligned} \langle q | \alpha_F \rangle &= 2^{\Omega/2} \exp \left\{ - \int \frac{1}{2} [q_s(\mathbf{k}) - q_{s0}(\mathbf{k})]^2 \right. \\ &\quad \left. - i p_{s0}(\mathbf{k}) \left[q_s(\mathbf{k}) - \frac{1}{2} q_{s0}(\mathbf{k}) \right] \vec{d}_s k \right\}. \end{aligned} \quad (B5)$$

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