Spectral functions in the ϕ^4 -theory from the spectral Dyson-Schwinger equations

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We develop a nonperturbative functional framework for computing real-time correlation functions in strongly correlated systems. The framework is based on the spectral representation of correlation functions and dimensional regularization. Therefore, the nonperturbative spectral renormalization setup here respects all symmetries of the theories at hand. In particular, this includes space-time symmetries, as well as internal symmetries such as chiral symmetry, and gauge symmetries. Spectral renormalization can be applied within general functional approaches such as the functional renormalization group, Dyson-Schwinger equations, and two- or *n*-particle irreducible approaches. As an application, we compute the full, nonperturbative, spectral function of the scalar field in the ϕ^4 -theory in 2 + 1 dimensions from spectral Dyson-Schwinger equations. We also compute the *s*-channel spectral function of the full ϕ^4 -vertex in this theory.

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I. INTRODUCTION

The study of the dynamics and resonance structure of strongly correlated systems requires the knowledge of realtime (timelike) correlation functions. In particular, the evolution of slow modes is dominated by the low-energy regime and is crucial for transport approaches and hydrodynamics. Spectral functions encode the full, nonperturbative, information of the respective degrees of freedom and open the door to additional real-time quantities such as transport coefficients. They are also a particularly useful tool when discussing resonances and bound states, since they give direct access to the spectrum of excitations in a given theory. Emergent composite states are of central interest not just in particle and nuclear physics, but in most physics areas.

The treatment of strongly correlated systems asks for nonperturbative techniques. In recent years, functional and lattice approaches have been applied very successfully to the low-energy regime of theories ranging from QCD to condensed matter systems. The general framework for the quantum-field theoretical description of bound states with functional methods was introduced by Bethe and Salpeter [1,2]. For a recent review, see [3]; for a review of respective lattice results, see [4]. Despite this rapid and very impressive progress in particular in the last decade, the reliable direct nonperturbative calculation of real-time correlation functions is still in its infancy. Functional approaches have matured in recent years and are by now quantitatively competitive in QCD in Euclidean space-time. The extension of such calculations to Minkowski space-time is yet hindered by technical and conceptual complications; for recent advances, see [5–34].

In this work. we develop a novel approach for the direct computation of real-time (Minkowski) observables that is based on spectral representations of correlation functions. The approach comes with the advantage that we can use dimensional regularization for the analytic computation of momentum integrals in fully numerical nonperturbative calculations. Accordingly, the respective renormalization scheme, *spectral renormalization*, is based on standard dimensional regularization and respects the space-time symmetries, internal symmetries such as chiral symmetry, and gauge symmetries of the theory at hand. Clearly, this method is applicable to a broad range of theories, including non-Abelian gauge theories, within a regularization and renormalization scheme which is manifestly gauge invariant.

In the present work, we apply our novel approach to the scalar ϕ^4 -theory in d = 2 + 1 dimensions. This theory is a simple strongly correlated system and serves as a good benchmark for new techniques before applying them more involved theories such as non-Abelian ones. It is also interesting in its own right and has many applications as a model theory for perturbative and nonperturbative phenomena.

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The numerical application in the present work is done within the spectral Dyson-Schwinger approach. We compute the spectral function of the scalar field from the gap equation for the propagator, using its Källén-Lehmann spectral representation. In a first step, all vertices are approximated with the classical ones, and the two-loop diagrams in the Dyson-Schwinger equations (DSEs) are included. In a second step, the use of skeleton expansion of the DSE with a bubble-resummation of the *s*-channel fourpoint function. The nonperturbative *s*-channel spectral function of the vertex is computed and used in the DSE. The setup gives us direct numerical access to Minkowski space-time correlation functions due to the spectral representation. This also allows for a (physical) on-shell renormalization via the spectral renormalization scheme.

The paper is organized as follows: in Sec. II, we introduce spectral renormalization as the example of the Dyson-Schwinger approach along with the necessary technical tools. In particular, we discuss dimensional spectral renormalization and a Bogoliubov-Parasiuk-Hepp-Zimmermann (BPHZ)-type spectral renormalization. Section III works out the calculational details of solving the DSE in the developed scheme, followed by our results in Sec. IV A for classical vertices and in Secs. IV B–IV E for the skeleton expansion. Finally, our conclusion is presented in Sec. V.

II. SPECTRAL RENORMALIZATION

The general real-time renormalization scheme developed here aims at combining a practical numerical implementation in nonperturbative applications while maintaining all underlying symmetries including gauge symmetries. This is achieved by utilizing dimensional regularization, which respects all space-time, internal and gauge symmetries of the theory at hand. We also develop a BPHZ-type subtraction scheme which facilitates the analytical computations significantly. If such a subtraction schemes does not violate any symmetries in the theory at hand, it is the scheme of choice.

A practical implementation of dimensional regularization requires an analytic momentum structure of the propagators and vertices in the given loop integrals. While this allows for its use in perturbation theory, nonperturbative applications, with their necessarily numerical computation of propagators and vertices, usually rely on ultraviolet momentum cutoffs. The latter are neither consistent with space-time symmetries nor with gauge symmetries. It is well known that in gauge theories and supersymmetric theories such a regularization requires symmetry-breaking counterterms. This is not a conceptual problem, but it typically triggers additional power-counting relevant terms that may leads to additional fine-tuning tasks; see e.g., [35,36]. Moreover, the Wick rotation to Minkowski space-times is hampered by the deformation of the momentum integrals, which can lead to additional poles and cuts in the integration contours.

The present renormalization scheme achieves the requirement of analytic momentum integrals by using spectral representations of correlation functions. For the propagators, this is the Källén-Lehmann spectral representation, similar spectral representations also exist for the vertices, though getting increasingly difficult. We call this renormalization scheme spectral renormalization: after inserting the spectral representations in the loop integrals, the momentum integrands take an analytic form. This form is well suited for using dimensional regularization or related analytic computation techniques. The loop-momentum integrations can be performed analytically and we are left with spectral integrals. The whole nonperturbative information is contained in the spectral functions of propagators and vertices. In most nonperturbative applications, the respective spectral integrals can only be computed numerically.

The scheme can be practically applied to any divergent diagram that scales in the UV with loop momentum to some natural power q^m , $m \in \mathbb{N}$ (with $m < n_{\max}$ and n_{\max} given by the renormalizability constraint). This is always the case when using spectral representations for all correlation functions, but also works for classical vertices. This will be detailed in the present work within the example of the Dyson-Schwinger approach introduced in the next section, Sec. II A. The functional DSE for the effective action, (4), is depicted in Fig. 2, that for the inverse propagator is depicted in Fig. 3. For classical vertices, all momentum integrals in functional approaches, see e.g., Fig. 3, are of the standard perturbative form, but with different spectral masses for all lines. Most of these integrals are known from perturbation theory results, e.g., [37]. Note that this reparametrization comes at the cost of spectral integrals for each propagator. Spectral representations of vertices lead to further spectral integrals as well as further classical propagators with spectral masses. In summary, in a spectral functional approach, all momentum integrals are perturbative. Hence, we can implement a symmetry-preserving regularization such as dimensional regularization, leading to a renormalization scheme with symmetry-consistent counterterms.

A relevant example for this important symmetrypreserving property are gauge theories. There, a momentum cutoff or standard subtraction scheme requires explicitly or implicitly a mass counterterm for the gauge field in order to keep the renormalized gauge field massless. In four



FIG. 1. Diagrammatic notation used throughout this work: lines stand for full propagators, small black dots stand for classical vertices, and larger gray dots stand for full vertices.



FIG. 2. Functional DSE for the effective action of the scalar theory under investigation in this work. The notation is given in Fig. 1.

dimensions, this leads to a quadratic fine-tuning task instead of a logarithmic one; for a detailed discussion, see [35,36]. Spectral renormalization with spectral regularization removes the (explicit or implicit) necessity of a mass counterterm for the gauge field, and hence the quadratic fine-tuning task.

After briefly introducing Dyson-Schwinger equations, Sect. II A, and the Källén-Lehmann spectral representation, Sec. II B, we set up spectral renormalization in Secs. II C (spectral dimensional renormalization) II D (spectral BPHZ-renormalization). Further examples and the discussion of the fully nonperturbative setup can be found in Secs. II E and II F. The explicit example used for demonstrating the properties and computational details of the spectral renormalization scheme is the gap equation for scalar ϕ^4 - and ϕ^3 -theories.

A. Dyson-Schwinger equations

The central object in functional approaches such as Dyson-Schwinger equations or functional renormalization group equations is the quantum effective action $\Gamma[\phi]$. It is related to the generating functional Z[J] of correlation functions including their disconnected parts via a Legendre transformation,

$$\Gamma[\phi] = \sup_{J} \left[\int d^d x J(x) \phi(x) - W[J] \right].$$
(1)

Here, $W[J] = \ln Z[J]$ is the Schwinger functional that generated connected correlation functions. The effective action $\Gamma[\phi]$ is the generating functional for one-particle irreducible (1PI) correlation functions. They are obtained from *n*th derivative of the effective action with respect to the fields. In momentum space, this reads

$$\Gamma^{(n)}[\phi](p_1, \dots, p_n) = \frac{\delta^n \Gamma[\phi]}{\delta \phi(p_1) \dots \delta \phi(p_n)}.$$
 (2)

A pivotal role in all functional approaches is played by the full, field-dependent propagator $G[\phi]$. It is simply the inverse of the 1PI two-point function,

$$G[\phi](p,q) = \frac{1}{\Gamma^{(2)}}[\phi](p,q).$$
(3)

With these prerequisites, we straightforwardly arrive at the master DSE, the quantum equation of motion,

$$\frac{\delta\Gamma[\phi]}{\delta\phi} = \frac{\delta S}{\delta\varphi} \left[\varphi = G \cdot \frac{\delta}{\delta\phi} + \phi \right],\tag{4}$$

where $G \cdot \delta/\delta \phi = \int d^d q/(2\pi)^d G(p,q) \delta/\delta \phi(q)$ in momentum space, and $S[\phi]$ is the classical action. A more detailed derivation can be found e.g., in [38]. By taking functional derivatives of (4) with respect to ϕ , all higher-order 1PI correlation functions are generated. For example, the DSE for the two-point function is generated by taking one derivative of (4) and is depicted in Fig. 3.

For our explicit example of a ϕ^4 -theory, the classical action in (4) is given by

$$S[\varphi] = \int d^d x \left[\frac{1}{2} (\partial_\mu \varphi)^2 + \frac{m_{\phi,0}^2}{2} \varphi^2 + \frac{\lambda_{\phi,0}}{4!} \varphi^4 \right].$$
(5)

It depends on two parameters or couplings, the bare fourpoint coupling $\lambda_{\phi,0}$, and the bare mass parameter $m_{\phi,0}$. The third parameter required for renormalization, the wave function renormalization Z_{ϕ} can be scaled out. This is conveniently done by setting it to unity at the renormalization scale.

We close this section with a few comments: to begin with, the underlying Z_2 -symmetry of the theory, $\varphi \rightarrow -\varphi$ implies the same symmetry for the effective action under transformations of the mean field, $\phi \rightarrow -\phi$. Accordingly, the odd vertices vanish at vanishing mean field $\Gamma^{(2n+1)}[\phi = 0] \equiv 0$. Moreover, restricting ourselves to constant background fields ϕ_c , we can formally expand the three-point function in powers of the field,



FIG. 3. DSE of the two-point function for a general background field $\phi \neq 0$. The vacuum polarization and squint diagrams are proportional to $S^{(3)}[\phi] \propto \phi$. They vanish for $\phi = 0$, where the standard form with tadpole and sunset diagrams is obtained. The notation is given in Fig. 1.

$$\Gamma^{(3)}[\phi_c](p_1, p_2, p_3) = \phi_c[\Gamma^{(4)}(p_1, p_2, p_3, 0) + O(\phi_c^2)],$$
(6)

due to the odd vertices vanishing at the expansion point $\phi = 0$. In (6), we used the Fourier transform $\tilde{\phi}_c(p) = (2\pi)^d \phi_c \delta(p)$ of the constant field. We also introduced $\Gamma^{(n)}(p_1, p_2, p_3, 0, ..., 0)$, the *n*-point functions at a vanishing background and n - 3 vanishing momenta. The relation (6) gives rise to the polarization and squint diagram in Fig. 3 for $\phi_c \neq 0$, which are absent for $\phi_c = 0$. In the broken phase, the equation of motion (EoM) ϕ_0 for constant fields is solved for a nonvanishing expectation value of the field, i.e., $\phi_0 \neq 0$. Accordingly, if evaluating the DSEs for correlation functions on the EoM, these diagrams are present.

B. Källén-Lehmann spectral representation

Using the Källén-Lehmann spectral representation [39,40], the propagator can be recast in terms of its spectral function ρ ,

$$G(p_0) = \int_0^\infty \frac{d\lambda}{\pi} \frac{\lambda \rho(\lambda, |\vec{p}|)}{p_0^2 + \lambda^2}.$$
 (7)

For asymptotic states, the spectral function can be understood as a probability density for the transition to an excited state with energy λ . In this way, the spectral function acts as a linear response function of the two-point correlator, encoding the energy spectrum of the theory. The existence of a spectral representation imposes tight restrictions on the analytic structure of the propagator. In turn, the Euclidean propagator also constrains the spectral function; for a rather nontrivial example for the latter constraints, see [41].

From a complex analysis perspective, the spectral function naturally arises as the set of nonanalyticities of the propagator, which are severely restricted in the complex plane by Cauchy's theorem. This results into the following inverse relation between spectral function and the retarded propagator:

$$\rho(\omega, |\vec{p}|) = 2 \operatorname{Im} G(-i(\omega + i0^+), |\vec{p}|), \tag{8}$$

where ω is the real time zero momentum component. This formulation allows us to work only with the frequency argument and set the spatial momentum to zero in practice, since the full phase space can be restored from Lorentz invariance. Hence, for the remainder of this work, $|\vec{p}|$ will be dropped.

The existence of a spectral representation restricts all nonanalyticities of the propagator to lie on the real momentum axis, as is manifest in (7) and (8). This crucial condition as well as the generic structure in the complex plane already mentioned above allows us to recast the spectral function into the form

$$\rho(\lambda) = \frac{\pi}{\lambda} \sum_{i} Z_i \delta(\lambda - m_i) + \tilde{\rho}(\lambda).$$
(9)

The spectral function is split into a set of δ -functions and a continuous scattering part $\tilde{\rho}$, arising from branch cuts in the complex plane of the propagator. For a classical propagator, the spectral function reduces to one mass-shell δ -function with $Z_1 = 1$. There are no further poles, $Z_{i>1} = 0$, and the scattering part is absent, $\tilde{\rho} = 0$. Further details can be found e.g., in [42].

In the scalar ϕ^4 -theory, the spectral function of the scalar field is that of an asymptotic state, and hence is positive semidefinite and has the interpretation of a probability density. Therefore, it is convenient to normalize its integrated weight to unity within an appropriate renormalization scheme. Within this scheme, we have the sum rule

$$\int_{\lambda} \lambda \rho(\lambda) = 1 \quad \text{with} \quad \int_{\lambda} \equiv \int_{0}^{\infty} \frac{d\lambda}{\pi}, \qquad (10)$$

which implies $Z_i \leq 1$. The spectral weight is distributed between poles and cuts, and in the presence of scattering states the weight of the poles is less than 1.

C. Spectral dimensional renormalization

Next, we discuss the spectral renormalization scheme, which is fully based on dimensional regularization. Dimensional regularization renders the loop diagrams finite, and we can swap orders of the initially outer momentum and inner spectral integrals integrals (cf. upper part of Fig. 4).



FIG. 4. Schematic illustration of the spectral renormalization scheme. The function f is some arbitrary divergent integrand. The dependence on the external momentum p is suppressed. The upper two boxes have to be understood as finite by dimensional regularization, but divergent in the limit $\varepsilon \to 0$ in $d - \varepsilon$ dimensions. In a first step, the momentum integrals are analytically evaluated via dimensional regularization. Subsequently, the spectral integrals are renormalized via spectral renormalization, either within the dimensional (Sec. II C) or BPHZ-approach (Sec. II D).

This allows us to first perform the momentum integrals analytically. We are left with finite spectral integrals at finite $\epsilon > 0$, that in general have to be done numerically since the spectral functions may only be known numerically.

Generally, the numerical integration of the spectral integrals can be performed for finite ϵ , and in gauge theories full manifest gauge consistency of spectral renormalization requires that the limit $\epsilon \to 0$ is taken only after performing all integrals. However, it is convenient for the numerical performance to do the spectral integration at $\epsilon = 0$. The same holds for the access to the analytic momentum structure required for extracting Minkowski properties. In this case, for the limit $\epsilon \to 0$, we have to set up a consistent renormalization procedure before performing the spectral integrals, which is worked out in detail in Sec. II D. In particular, it is not sufficient to only remove the $1/\epsilon$ -divergences that originate in the momentum integrations, as the spectral integrations lead to further $1/\epsilon$ -terms. This relates to the swapping of the integration order and performing the limit $\epsilon \to 0$ before evaluating all integrals.

This section is dedicated to the fully gauge-consistent renormalization scheme we call spectral dimensional renormalization. In this scheme, the divergent parts of the spectral integrals are performed analytically before taking the limit $\varepsilon \to 0$. A simple example is the tadpole contribution to the gap equation in Fig. 3 in d = 3 dimensions, which comes with a momentum-independent linearly divergent term. This example is also relevant for our later computation in Secs. III and IV. After the momentum integration is performed, we arrive at a finite result proportional to

$$\int_0^\infty d\lambda \lambda \frac{\mu^{2\epsilon} \lambda^{1-2\epsilon}}{(\lambda^2 + m^2)} = -\frac{\pi}{2} \frac{1}{\cos(\frac{\pi\epsilon}{2})} m \left(\frac{\mu^2}{m^2}\right)^\epsilon.$$
 (11)

In (11), we used a trial spectral function that decays for large spectral values λ according to its momentum or spectral dimension,

$$\rho_{\rm trial}(\lambda,m) = \frac{1}{\lambda^2 + m^2}, \qquad (12)$$

with a positive mass m > 0. The trial spectral function in (12) approximates the correct leading ultraviolet behavior, if we neglect logarithmic corrections. For large spectral values, the UV asymptotic of the spectral function can be extracted from the leading momentum dependence of the respective propagator. In the present example, the latter is assumed to decay quadratically on its branch cut on the real momentum axis.

The finiteness of the result of the momentum integration used on the left-hand side of (11) also points at a specific property of dimensional regularization: in odd dimensions, d = 2n + 1, it already removes all momentum divergences. In even dimensions, d = 2n, it removes subclasses of divergent ones, a prominent being the (one-loop) tadpole in a massless theory such as a gauge theory. If we had put $\epsilon = 0$ before the integration, (11) simply is (linearly) divergent. This is the price to pay for the swapping of integration orders: the divergences are not fully covered by the momentum integrals any more. The example also entails that in odd dimensions including our explicit computation in the ϕ^4 -theory in d = 3, all divergences come via the spectral integrals.

In spectral dimensional renormalization, spectral singularities in even dimensions show up as $1/\epsilon$ -terms. In dimensional regularization in perturbation theory, these divergences are typically removed recursively by introduction of appropriate counterterms. In spectral dimensional renormalization, this procedure is applied too, while keeping a finite ϵ as well as isolating analytically the singular part of the spectral integrals with

$$\rho(\lambda) = \rho_{\rm IR}(\lambda) + \rho_{\rm UV,an}(\lambda). \tag{13}$$

In (13), the numerical "infrared" part $\rho_{\rm IR}(\lambda)$ decays sufficiently fast for large spectral values and renders the respective spectral integrals finite. In turn, the ultraviolet part $\rho_{\rm UV,an}(\lambda)$ carries the ultraviolet asymptotics analytically. Therefore, the respective spectral integrals can be treated analytically with dimensional regularization as done in (11). With (12), we use the IR-UV split,

$$\rho(\lambda) = \rho_{\rm IR}(\lambda, k) + \rho_{\rm trial}(\lambda, k).$$
(14)

We also emphasize that the theory does not depend on the mass parameter *k* that regularizes (in the infrared) the UV part of the spectral function. In particular, we have $\partial_k \rho(\lambda) \equiv 0$, which entails the *k* dependence of the infrared part of the spectral function.

The (leading) ultraviolet behaviour of the spectral function at large spectral values is governed by $\rho_{UV,an} = \rho_{trial}$ and the infrared part decays with the fourth power of the spectral value,

$$\lim_{\lambda \to \infty} \rho_{\rm IR}(\lambda \to \infty) \propto \frac{1}{\lambda^4} \tag{15}$$

Inserting the split (14) into (11) leads us to the final finite result with $\epsilon = 0$,

$$\int_0^\infty d\lambda \mu^{2\epsilon} \lambda^{2-\epsilon} \rho(\lambda) = \int_0^\infty d\lambda \lambda^2 \rho_{\rm IR}(\lambda) - \frac{\pi}{2} k, \quad (16)$$

with a finite (in general numerical) integral over the infrared part of the spectral function due to (15). The numerical convergence of this integral can be further improved systematically, if the UV part $\rho_{\rm UV}(\lambda)$ also includes subleading UV terms of the full spectral function.

The finite result of (16) was obtained without the introduction of any possibly symmetry-breaking counterterms. Since the systematics of this example is general, it can be applied to all divergences of general diagrams. As the demonstrated spectral dimensional renormalization procedure is entirely based on dimensional regularization and the use of spectral representation, it preserves all symmetries of the theory at hand. Especially, for the case of gauge theories, it is manifestly gauge invariant or rather gauge consistent. For example, it reflects the peculiarity of dimensional regularization that integrals without any external scale vanish identically. For k = 0, the integral over the UV part of the spectral function vanishes and we are left with the finite IR part. Accordingly, no mass counterterms are needed in a massless theory such as a gauge theory.

Figure 4 illustrates the general renormalization workflow, including spectral renormalization. The schematic representation holds for the case of spectral dimensional renormalization as well as for the subtraction based and more universal approach of *spectral BPHZ-renormalization*, which is introduced in the following section.

D. Spectral BPHZ-renormalization

In fully symmetry-consistent spectral dimensional renormalization as described in Sec. II C, one has to perform analytic spectral integrals with dimensional regularization on top of the momentum integrations. Already the latter are more complicated than in standard perturbation theory at the same order, since the spectral representations lead to different masses for each line. While the momentum integration has to be necessarily analytic in order to access Euclidean and Minkowski space-time, this is not necessary for the spectral integration, whose nonperturbative infrared part has to be done numerically in most cases anyway.

The need for additional analytic computations of spectral integrals can be circumvented by performing subtractions on the spectral integrals which render the spectral integrals finite. This can be done by subtracting a Taylor expansion of the spectral integrand in momenta according to the BPHZ-scheme with Dyson's formula. The procedure is called spectral BPHZ-renormalization. We shall see, that its work flow is still described within Fig. 4. It is the last "spectral" step from the bottom right to the bottom left which is changed by moving from spectral dimensional to the spectral BPHZ-procedure. We emphasize that the underlying BPHZ-regularization in general does not preserve all symmetries of a given theory and in particular breaks gauge symmetry. This is not a conceptual problem, as the counterterms also break gauge invariance and the final result is gauge consistent; see e.g., [43]. However, in numerical applications to gauge theories the gaugeconsistent spectral dimensional renormalization is arguably worth its price, in particular for investigations of the Gribov problem and the confinement mechanism.

In the present example of a scalar ϕ^4 -theory, the BPHZscheme is consistent with both space-time and the internal Z_2 symmetry. We will therefore utilise it for explicit computation. We introduce and explain the setup within a specific example, the sunset graph in the gap equation of the scalar ϕ^4 -theory in *d* dimensions; see Fig. 3. This diagram also carries subdivergences while still being relatively simple. In Sec. II E, we additionally analyze an explicit one-loop example.

Due to the spectral representation of the sunset graph, the fully perturbative momentum integrals including the sub-traction can be solved analytically in both d = 3 and d = 4:

In d = 4, the sunset graph is superficially quadratically divergent with logarithmic divergences in the subdiagrams. The respective spectral power counting follows from the momentum dimension of the spectral value, $[\lambda] = 1$, where $[\mathcal{O}]$ counts the momentum dimension of \mathcal{O} . In particular, that of the spectral function is $[\rho(\lambda)] = -2$. This can be read off from the classical spectral function of a field with mass *m* with $\rho_{\rm cl}(\lambda) = 2\pi\delta(\lambda^2 - m^2)$. Trivial examples for such a power counting of spectral integrals are (11) and the propagator itself with $\left[\int d\lambda \lambda \rho(\lambda)\right)/(\omega^2 + \lambda^2) = -2$. We subtract the zeroth and first orders of the Taylor expansion in external momentum about $p^2 = \mu^2$ as well as the zeroth term in Taylor expansions about the external momenta of the subdiagrams. These counterterms contribute to the mass renormalization as well as the wave function renormalization of the scalar field. The subtractions remove the leading and also the next-to-leading order (NLO) contributions in the spectral parameters λ_i in the integrand of the spectral integral for $\lambda_i \to \infty$. Consequently, the integrand decays faster by two powers of λ_i^2 and the spectral integrals over λ_i with i = 1, 2, 3 are UV-finite.

In d = 3, there are no divergent subdiagrams and the sunset is superficially logarithmically divergent. We only subtract the zeroth order of the Taylor expansion, which contributes to the mass renormalization.

We are left with finite spectral integrations for the sunset graph in d = 3, 4 at $\epsilon = 0$. The integrand depends analytically on the external momentum, the spectral values λ_i with i = 1, 2, 3 of the three internal lines as well the respective spectral functions $\rho(\lambda_i)$. In general, the remaining spectral integrals have to be performed numerically.

In summary, spectral BPHZ-renormalization, as described in detail above, has the same work flow as in the last section and is depicted in Fig. 4: first, we apply dimensional regularization to the momentum integrations and swap the order of momentum and spectral integrals. Then we perform the momentum integration analytically, right bottom corner in Fig. 4. Finally, we apply the spectral BPHZ-step: the Taylor expansion in all momenta about the renormalization scale μ , which allows us to take the limit $\epsilon \rightarrow 0$. This leaves us with the task to perform the finite spectral integrals either analytically or numerically, depending on the application.

E. One-loop example: ϕ^3 -theory in d = 6

For further illustration, we now apply spectral BPHZrenormalization within the simple perturbative example of the one-loop DSE for the two-point function of the (renormalizable) ϕ^3 -theory in d = 6 dimensions with a coupling $g/(3!) \int_x \phi^3$. The advantage of this example is that already at one loop it requires both a mass renormalization and a wave function renormalization. Hence, both can be discussed within a simple one-loop computation. In contrast, in the ϕ^4 -theory, the wave function renormalization only arises at two-loop from the sunset diagram (even for $\phi_c \neq 0$).

In the ϕ^3 -theory, the only diagrams in the gap equation Fig. 3 are the polarization diagram and the squint diagram. At one loop, we only have to consider the polarization diagram. In d = 6, this diagram is quadratically divergent. The respective DSE for the (inverse) propagator reads schematically,

$$\Gamma^{(2)}(p) = Z_{\phi,0}(\mu)(p^2 + m_{\phi,0}^2(\mu)) + g^2 \int_{\lambda_1,\lambda_2} \lambda_1 \lambda_2 \rho(\lambda_1) \rho(\lambda_2) F(p,\mu;\lambda_1,\lambda_2).$$
(17)

The integrand *F* results from the momentum integration and depends on the renormalization group scale μ due to dimensional regularization. $Z_{\phi,0}$ is the wave function renormalization and $m_{\phi,0}^2$ is the bare mass squared. Both bare parameters contain counterterms that remove the divergences in the diagrams within an expansion about $p^2 = \mu^2$: the constant quadratic divergence as well as the logarithmic divergence proportional to p^2 . This amounts to the choices

$$Z_{\phi,0}(\mu) = 1 - g^2 \int_{\lambda_1,\lambda_2} \lambda_1 \lambda_2 \rho(\lambda_1) \rho(\lambda_2) \frac{F(p,\mu;\lambda_1,\lambda_2)}{\partial p^2} \Big|_{p^2 = \mu^2},$$

$$Z_{\phi,0}(\mu) m_{\phi,0}^2(\mu) = m_{\phi}^2 - g^2 \int_{\lambda_1,\lambda_2} \lambda_1 \lambda_2 \rho(\lambda_1) \rho(\lambda_2) \left[F(p,\mu;\lambda_1,\lambda_2) - \mu^2 \frac{F(p,\mu;\lambda_1,\lambda_2)}{\partial p^2} \Big|_{p^2 = \mu^2} \right]$$
(18)

for wave function and mass renormalization, respectively. The counterterms proportional to g^2 in (18) provide the first two terms of the Taylor expansion about $p^2 = \mu^2$ of the diagram. To see this, we insert (18) in DSE (17),

$$\Gamma^{(2)}(p) = p^2 + m_{\phi}^2 + g^2 \int_{\lambda_1, \lambda_2} \lambda_1 \lambda_2 \rho(\lambda_1) \rho(\lambda_2) \left[F(p, \mu; \lambda_1, \lambda_2) - F(\mu, \mu; \lambda_1, \lambda_2) - (p^2 - \mu^2) \frac{F(p, \mu; \lambda_1, \lambda_2)}{\partial p^2} \Big|_{p^2 = \mu^2} \right].$$
(19)

Equation (19) is depicted in Fig. 5. Accordingly, (18) implements the standard renormalization conditions, that the quantum corrections vanish at $p^2 = \mu^2$,

$$\Gamma^{(2)}(p^2 = \mu^2) = Z_{\phi}(\mu^2 + m_{\phi}^2)$$

$$\partial_{p^2}\Gamma^{(2)}(p^2 = \mu^2) = Z_{\phi}$$
(20)

for the two-point function with $Z_{\phi} = 1$. In the present ϕ^3 example, these two renormalization conditions are complemented by that for the coupling g, which is also logarithmically divergent in d = 6. As we have introduced this example only for illustration of spectral BPHZrenormalization, we refrain from discussing this any further. Mode details on the spectral renormalization conditions can be found in Sec. III B. Continuing with the discussion of spectral renormalization for the two-point function, the subtraction of *F* in (19) by its own Taylor expansion at $p^2 = \mu^2$ leads to finite spectral integrals. We emphasize that this is not achieved by simply subtracting the $1/\epsilon$ -terms before performing the spectral integration: since *F* scales with λ^2 for large λ with $\lambda = \lambda_1$, λ_2 , the spectral integrals in (17) are quadratically divergent. After spectral BPHZ-renormalization, however, the subtracted scalar integrand in (19) scales as $1/\lambda^2$. The subtraction scheme cancels the leading and subleading contributions in λ to *F* and leads to finite spectral integrals.

F. Nonperturbative spectral renormalization

The discussions of the last three sections, Secs. II C–II E, entail that spectral renormalization leads to two different



FIG. 5. Schematic spectral BPHZ-renormalization procedure at the example of the one-loop scalar propagator DSE for the ϕ^3 -theory. The diagram is quadratically divergent in d = 6. First, the loop-momentum divergences are discarded by the usual momentum renormalization part of dimensional regularization (i.e., implicitly assumed on the lhs). By introducing mass and wave function counterterms, the diagram is subtracted by the first two terms of its own Taylor expansion around the RG scale μ . This cancels the leading order quadratic and subleading logarithmic divergences of the spectral integrals.

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parts in the counterterms: the first part is related to the momentum divergences and has all the properties of the counterterms in dimensional regularization. The second part comes from the spectral divergences. The counterterms in spectral dimensional renormalization respect all symmetries including gauge symmetries and is tantamount to dimensional regularization and renormalization. The counterterms in spectral BPHZ-renormalization lack the full symmetries. In particular, in gauge theories, the BPHZcounterterms are necessarily not gauge invariant, precisely for restoring gauge consistency of the full renormalized result.

Importantly, in both spectral renormalization schemes, the spectral counterterms follow the same recursive relations known from standard perturbation theory. This makes it a consistent renormalization scheme to all orders of perturbation theory.

In nonperturbative applications of the present spectral approach, the nonperturbative information is solely present in the spectral functions of propagators and vertices. Moreover, within the DSE, we only deal with one- and two-loop diagrams with *n* vertices derived from the master DSE in (4); see also Fig. 2. We have one classical (bare) vertex and n - 1 full vertices as well as full propagators. If recursively written in terms of loops, both, the *finite* full vertices and propagators, carry subtractions of the divergences in these diagrams that render the diagrams finite. The leftover subtractions from these redistribution renormalize the one or two explicit loops in the DSE diagrams.

In summary, nonperturbative spectral renormalization only concerns the counterterms for the explicit loops in the DSE, while the rest of the renormalization is carried by the finite full vertices and propagators that have to obey the renormalization conditions. This is a consistent numerical nonperturbative renormalization scheme.

III. ϕ^4 -THEORY IN 2+1 DIMENSIONS

The ϕ^4 -theory in 2 + 1 dimensions is super-renormalizable, and the initial two renormalization conditions for the two-point function in Fig. 3 reduce to the first one for the mass. Moreover, we do not need to renormalize the coupling. This entails that in the DSE for the scalar twopoint function spectral BPHZ-renormalization as discussed in Sec. II D simply amounts to subtracting the zeroth order term in the Taylor expansion about $p^2 = \mu^2$. After the momentum integrals are computed analytically within dimensional regularization, we are left with the finite spectral integrals.

After completing renormalization, the iterative solution procedure in the DSE is briefly described as follows: with a given input spectral function, the renormalized DSE is evaluated in Minkowski space-time. The input spectral function is either the initial guess or the result of the last iteration step. Then, an updated retarded two-point function is computed from the result. This allows us to extract an updated spectral function, which is fed back as input into the next iteration step. In this section, we discuss the calculation sketched above in a detailed way, step by step.

A. Momentum integration and spectral renormalization

The DSE can be expressed as the sum of the bare twopoint function and the loop diagrams D_j ,

$$\Gamma^{(2)}(p) = p^2 + m_{\phi,0}^2 + \sum_{\{j\}} D_j(p), \qquad (21)$$

with j = tad, pol, sun, squint. In (21), we have used that the ϕ^4 -theory in d = 3 is super-renormalizable and the only divergent term is the mass term. From now on, we drop the μ dependence of the spectral integrands for notational simplicity. With the Källén-Lehmann spectral representation for the full propagator (7), as well as momentum-independent vertices, an arbitrary loop diagram D_j in the DSE takes the form

$$D_j(p) = g_j \prod_i^{N_j} \left(\int_{\lambda_i} \lambda_i \rho(\lambda_i) \right) I_j(p; \lambda_1, \dots, \lambda_{N_j}).$$
(22)

The prefactors g_j are the products of the combinatorial prefactors in the DSE and the vertices of the corresponding diagram. In Table I, we provide the prefactors for the Minkowski version of (22); see (25).

 D_j has N_j internal lines, each of them coming with one spectral integral and a corresponding spectral function. The I_j are nothing but a product of (momentum) loop integrals over N_j classical propagators with different spectral masses λ_i ,

$$I_{j}(p;\lambda_{1},...,\lambda_{N_{j}}) = \prod_{k}^{N_{j}^{\text{loops}}} \int \frac{d^{3}q_{k}}{(2\pi)^{3}} \prod_{i}^{N_{j}} \frac{1}{\lambda_{i}^{2} + l_{i}^{2}}.$$
 (23)

TABLE I. Prefactors of the propagator DSE diagrams in the different approximation schemes. The prefactors are obtained by the standard DSE prefactors and the loop-momentum and spectral parameter-independent parts of the vertices. The tadpole factor in the approximation with classical vertices is set to zero, as the tadpole is absorbed completely in the mass renormalization. The sunset prefactor in the skeleton expansion compensates the two-loop contributions of the tadpole with full four-vertex. In the self-consistent approximation of Sec. IV E, the skeleton prefactors apply with exception of the polarization diagram, which is given by (48).

	Classical vertices	Skeleton
$g_{\rm pol}$	$-\frac{3}{2}\Gamma^{(2)}(\omega=0)\lambda_{\phi}$	$-\frac{3}{2}\Gamma^{(2)}(\omega=0)\Gamma^{(4)}(\omega=0)$
g_{sunset}	$-\frac{1}{6}\lambda_{\phi}^2$	$\frac{1}{12}\Gamma^{(4)}(\omega=0)^2$
$g_{\rm squint}$	$\frac{3}{2}\lambda_{\phi}^{2}\Gamma^{(2)}(\omega=0)$	0
$g_{\rm tad}$	0	$\frac{1}{2}$

In (23) the momenta l_i are linear combinations of the loopmomenta q_k and the external momentum p. The number of loops is denoted by N_j^{loops} in (23). The analytic solutions of these integrals in d = 3 are known from perturbation theory, e.g., [37], and can be used here.

With the analytic expressions for I_j , we apply spectral BPHZ-renormalization, cf. Sec. II D. Please note that only the logarithmically divergent sunset diagram D_{sun} explicitly requires renormalization, since the bare tadpole simply can be absorbed in the definition of the bare mass squared. To cancel the logarithmic divergence of this diagram, within the BPHZ-approach, we subtract the zeroth order term in the Taylor expansion of D_{sun} about $p^2 = \mu^2$. The renormalized diagram reads

$$D_{\rm sun}^{\rm ren}(\omega) = g_{\rm sun} \int_{\lambda_1, \lambda_2} \lambda_1 \lambda_2 \rho(\lambda_1) \rho(\lambda_2) \\ \times [I_{\rm sun}(\omega; \lambda_1, \lambda_2) - I_{\rm sun}(\mu; \lambda_1, \lambda_2)].$$
(24)

B. Analytic continuation

The $I_j(p)$ are evaluated in Minkowski space-time for the retarded two-point function. This is done by parametrizing the complex (Euclidean) frequency as $p_0 = -i(\omega + i\varepsilon)$ and taking the limit $\varepsilon \to 0$. In a slight abuse of notation, we denote the continued expression as $I_j(\omega)$. They are given explicitly in the Appendix C. The DSE in Minkowski space-time reads

$$\Gamma^{(2)}(\omega) = -\omega^2 + m_{\phi,0}^2 + \sum_{\{j\}} g_j \prod_i^{N_j} \left(\int_{\lambda_i} \lambda_i \rho(\lambda_i) \right) I_j(\omega,\mu;\lambda_1,\dots,\lambda_{N_j}).$$
(25)

The prefactors of the diagrams of the DSE (25) with the classical vertex approximation and the skeleton expansion can be found in Table I. It can be deduced from the analytic structure of the spectral integrands $I_j(\omega)$, that the support in ω of the imaginary part of the polarization diagram starts at $\omega = \lambda_1 + \lambda_2$, as expected. Similarly, we find that the support of the sunset diagram starts at $\omega = \lambda_1 + \lambda_2 + \lambda_3$. Hence, the support of the expressions in the two-point function is given by multiples of the pole mass. This shows how the Cutkosky cutting rules can be easily extracted from the present spectral approach.

It is clear from (25) that the spectral renormalization approach allows for the implementation of physical onshell renormalization conditions. This is in contrast to Euclidean computations, where on-shell renormalization can only be implemented for massless modes. On-shell renormalization has the advantage that it minimizes the quantum corrections in a study of the resonance spectrum

$$\Gamma^{(2)}[p^2 = -m_{\rm pole}^2] = 0, \tag{26}$$

since both the coupling and the kinetic term do not require renormalization. The triviality of the wave function renormalization or rather the vanishing anomalous dimension entail, that the leading large momentum behavior of the propagator is given by $1/p^2$. Accordingly, the canonical commutation relations of the scalar field are unchanged, as is the normalization of dynamical states. The sum rule (10) is hence formally always satisfied in 2 + 1 dimensions, and thus provides a nontrivially benchmark test of our results.

C. Spectral integration and iteration

The remaining multidimensional integrals over the spectral parameters λ_i in (25) are solved numerically. Details on the numerical part of the calculation can be found in Appendix B. Subsequently, the spectral function is extracted from the updated two-point function via (8) and fed back into the DSE. In this way, the DSE is solved iteratively by successively integrating the right-hand side of the DSE with the updated spectral function from the last step until the solution converges, comp. Figure 6. Note that all the dynamical information is stored in the spectral functions, and the integrands I_j do not change within the iterations. Hence, each iteration only involves the numerical solution of the respective multidimensional spectral integral of each diagram.

More details on the convergence test can be found in Appendix B. There, the rapid convergence is illustrated as an exemplary case; see in particular Fig. 15.

As a starting point for the iterative procedure, an initial guess for the spectral function has to be made that is close enough to the solution (in the attraction basin of the solution in terms of the iteration). In the present case with the on-shell renormalization (26), the spectral function of the classical theory carries already the correct pole position by definition. In general, this will improve the convergence properties of the iteration. This is yet another property that



FIG. 6. Iteration procedure for computing the spectral function. With the initial guess ρ_0 for the spectral function, the two-point function $\Gamma^{(2)}$ is computed via DSE. The resulting spectral function is fed back into the DSE for the two-point function. This procedure is iterated until the convergence for the spectral function is reached.

singles out on-shell renormalization. The classical spectral function is given by

$$\rho_0(\omega) = \delta(\omega^2 - m_{\text{pole}}^2)$$

= $\frac{\pi}{m_{\text{pole}}} [\delta(\omega - m_{\text{pole}}) - \delta(\omega + m_{\text{pole}})].$ (27)

The delta-function peaks are located at the physical pole mass squared, and the delta-functions at $\pm m_{\text{pole}}$, are related by antisymmetry. With the initial guess in (27), the spectral integrals in the propagator DSE (25) reduce to the perturbative expressions. Further iterations then lead to resummation of the respective diagram classes.

The branch cuts of the loop corrections generate a continuous tail in the spectral function when inserting the initial spectral function ρ_0 on the right-hand side of the DSE. This entails via the sum rule in (10), that the residue of the mass pole decreases to Z < 1 due to the positive weight of the tail. The mass pole residue of the updated spectral function is obtained via the relation

$$Z = -\frac{2m_{\text{pole}}}{\partial_{\omega}\Gamma^{(2)}(\omega)}\Big|_{\omega=m_{\text{pole}}}.$$
(28)

The counterterm for the mass is conveniently extracted from $\operatorname{Re} \Gamma^{(2)}(\omega = m_{\text{pole}}) = 0.$

The three-point function $\Gamma^{(3)}[\phi_c]$ in the gap equation is evaluated at the constant field value $\phi_c = \phi_0$, which solves the equation of motion, $\partial_{\phi} V_{\text{eff}}[\phi_0] = 0$, at each step of the iteration. For fields in the vicinity of ϕ_0 , we can expand the effective potential in powers of $\phi^2 - \phi_0^2$ in the broken phase, leading to

$$V_{\rm eff}[\phi] = \sum_{n=2}^{\infty} \frac{v_n}{2n!} (\phi^2 - \phi_0^2)^n.$$
(29)

Accordingly, the two-, three-, and four-point functions at vanishing momentum are given by

$$\Gamma^{(2)}[\phi_0] = \frac{1}{3} v_2 \phi_0^2,$$

$$\Gamma^{(3)}[\phi_0] = v_2 \phi_0 \left(1 + \frac{1}{15} \frac{v_3}{v_2} \phi_0^2 \right),$$

$$\Gamma^{(4)}[\phi_0] = v_2 \left(1 + \frac{1}{5} \frac{v_3}{v_2} \phi_0^2 + \frac{1}{105} \frac{v_4}{v_2} \phi_0^4 \right).$$
 (30)

Dropping the higher-order terms $O((\phi^2 - \phi_0^2)^3)$ in (29), the terms in parentheses of (30) all reduce to unity. We introduce the curvature mass $m_{\text{cur}}^2(m_{\text{pole}}, \lambda_{\phi})$ as the value of the two-point function at vanishing momentum, i.e.,

$$m_{\rm cur}^2(m_{\rm pole},\lambda_{\phi}) \equiv \Gamma^{(2)}[\phi_0](p=0).$$
 (31)

Note that m_{cur} is determined by the two free parameters of the theory, m_{pole} and λ_{ϕ} . It is no new parameter. Using (31) in (30) and dropping the higher-order contributions, the three- and four-point functions at vanishing momentum are given by

$$\Gamma^{(3)}[\phi_0] = \sqrt{3} v_2 m_{\rm cur},$$

$$\Gamma^{(4)}[\phi_0] = v_2.$$
(32)

 v_2 is nothing but the full four-vertex at vanishing momentum, $v_2 = \Gamma^{(4)}[\phi_0](p = 0)$. For a general, momentumdependent four-point function in the *s*-channel, we find for the momentum-dependent three-point function,

$$\Gamma^{(3)}(p) = \Gamma^{(4)}(p) \sqrt{\frac{3}{v_2}} m_{\rm cur}, \tag{33}$$

dropping the ϕ_0 dependence of the correlators from now on. Note that by choosing $\Gamma^{(3)}$ consistently with $\Gamma^{(4)}$ as done above, the three-point function becomes dynamical and is hence updated through each iteration by its dependence on the two- and four-point functions.

IV. RESULTS

In this section, we compute and discuss the solution to the DSE for the propagator with two different approximations for the vertices. At first, we solve the DSE with the classical four-point vertex and the related three-point vertex as derived in Sec. III C. Subsequently, the DSE is considered in a skeleton expansion. Both, the full three- and four-point functions $\Gamma^{(3)}$, $\Gamma^{(4)}$, are based on the bubble resummed *s*-channel approximation to the four-point vertex derived from its Bethe-Salpeter equation. We also use the DSE for $\Gamma^{(4)}$ to compute results within a self-consistent version of this setup.

A. DSE with classical four-point vertex

In the present section, we approximate the full vertices in the gap equation in Fig. 3 with their classical counterparts while keeping the two-loop terms. This is depicted in Fig. 7. The two-loop terms constitute vertex corrections to the classical four-point function in the tadpole diagram (sunset) and for the classical three-point function in the polarization diagram (squint). In the latter case, this is but half of the vertex correction; the other half has been dropped in the current approximation when approximating the full three-point function in the polarization diagram by its classical counterpart. This approximation resums the propagator but is expected to fail in a regime where vertex corrections grow large.

For the four-point function, the classical vertex approximation amounts to

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FIG. 7. DSE of the two-point function with the classical vertex approximation as used in Sec. IVA. The tadpole is not present because its contribution is absorbed into the mass renormalization in the bare inverse propagator; for more, details can be found in the main text. The two-loop terms constitute vertex corrections of the classical vertices in the one-loop diagrams. The DSE is not two-loop complete as further vertex corrections have been dropped due to the classical vertex approximation. The notation is given in Fig. 1.

$$\Gamma^{(4)} = S^{(4)} = \lambda_{\phi}.$$
 (34)

With (34) and (32), the three-point function is given by

$$\Gamma^{(3)} = S^{(3)}[\phi_0] = \sqrt{3\lambda_{\phi}}m_{\rm cur},$$
(35)

see (32). The resulting prefactors g_j of the diagrams are listed in Table I. The tadpole diagram only contributes a momentum-independent term that shifts the mass and is absorbed into the mass renormalization.

All units will be given in the value of the pole mass. Put differently, we introduce dimensionless units

$$\lambda_{\phi} \to \frac{\lambda_{\phi}}{m_{\text{pole}}}, \qquad \omega \to \frac{\omega}{m_{\text{pole}}}, \qquad p \to \frac{p}{m_{\text{pole}}}.$$
 (36)

This also entails $m_{\text{pole}} = 1$, and the massless limit is taken with $\lambda_{\phi} \rightarrow \infty$. The dimensionless units also emphasize the well-known fact that the physics of the ϕ^4 -theory in d = 2 + 1 is specified by one dimensionless parameter, the ratio of coupling and (pole) mass.

We solve the DSE for three different values of the classical coupling constant $\lambda_{\phi} = 5$, 10, 20. The renormalized mass is fixed by the on-shell RG conditions (26) with $m_{\text{pole}} = 1$, and the quantum corrections to the mass vanish on-shell. For the smallest classical coupling used here,

 $\lambda_{\phi} = 5$, we take the classical spectral function as initial choice. For the further couplings, we use as the initial choice the full quantum spectral function of the closest coupling value available. This stabilizes the iterative procedure when successively moving further to larger couplings inducing larger quantum corrections.

The resulting spectral functions, as well as the corresponding propagators, are shown in Fig. 8 for different values of the classical four-point coupling, $\lambda_{\phi} = 5$, 10, 20. The mass pole as well as the onset of the two-particle threshold, $\phi \rightarrow \phi \phi$, at twice the pole mass is clearly visible. The three-particle threshold, $\phi \rightarrow \phi \phi \phi$, at $3m_{\text{pole}}$ becomes visible for large enough coupling. We emphasize that also all higher *n*-particle thresholds are present in the result and can be resolved numerically. Since they are suppressed by the inverse of the respective threshold energy squared, they are not visible in the plots. When further zooming in on the higher onsets which were marked by the dashed lines, the respective onsets would become visible.

The main effect of a stronger coupling (or small pole mass) can be understood intuitively very well. The residue of the mass pole becomes smaller, while the scattering cut gets larger contributions, i.e., scatterings get enhanced due to the large coupling or the small pole mass. As mentioned before, the only parameter present is $\lambda_{\phi}/m_{\text{pole}}$, and in the present units with $m_{\text{pole}} = 1$, this simply is λ_{ϕ} ; see (36).



FIG. 8. Spectral function (left) and propagator (right) in the scalar theory for the coupling choices $\lambda_{\phi} = 5$, 10, 20 from the full DSE with classical vertices using on-shell renormalization (26). All dimensionful quantities were rescaled in units of the respective mass pole result. The different height of the delta peaks encodes the magnitude of the residue relative to the other spectral functions. The gray dashed lines mark the n-particle onsets. For large enough coupling, the three-particle onset becomes visible. The propagators were computed by the Källén-Lehmann spectral representation.



FIG. 9. Truncated DSE of the two-point function as used in Sec. IV D. Notation is given in Fig. 1. The full *s*-channel four-vertex in the tadpole diagram enters via its spectral representation (39). The full vertices in polarization and sunset are approximated at zero frequency; see Sec. IV B 3. The different prefactor in front of the sunset diagram as compared to Fig. 7 is due to the tadpoles contribution to the sunset topology, cf. Sec. IV B 2. Squint, kite, and double-bubble topology are dropped, as motivated in Sec. IV B.

At the largest coupling value considered here, $\lambda_{\phi} = 20$, also the higher scattering processes start contributing significantly: the threshold at $3m_{\text{pole}}$ is clearly visible in the spectral function for this coupling, shown in the left panel of Fig. 8. For large couplings also, the higher thresholds kick in. This scattering physics also leaves its traces in the Euclidean propagator, shown in the right panel of Fig. 8: while first the propagator drops for small momenta with the increasing coupling (measured in the respective pole masses), it increases again for even larger ones due to the more pronounced scattering physics present in the spectral function.

The resolution of the higher particle thresholds is nearly impossible within reconstruction methods due to their exponential suppression in the Euclidean data. This emphasizes once again the strength of our semianalytic approach of calculating spectral functions directly from real-time correlators, where all these thresholds are incorporated explicitly.

B. Fully nonperturbative DSE

The practical applicability of the spectral renormalization scheme has been shown in the last section within the DSE for the two-point function with classical vertices; see Fig. 7. This approximation implements a full resummation of the propagator. The approximation also includes some corrections in the higher-order diagrams in the DSE as already discussed in the last section. While these diagrams contribute to the higher-order scattering thresholds, this may not be sufficient in the limit of asymptotically large couplings $\lambda_{\phi}/m_{\text{pole}} \rightarrow \infty$, also tantamount to small pole masses. In this regime, the vertex corrections should be taken into account consistently.

In the present section, we discuss nonperturbative expansion schemes of the DSE as well as resummations of the vertices. This allows us to study the strongly correlated regime of the theory. A full quantitative study is beyond the scope of the present contribution and is deferred to future work.

A nonperturbative expansion scheme for the DSE is given by the skeleton expansion. In this expansion, all vertices are full-dressed, and higher loop order diagrams with dressed propagators and vertices have to be introduced successively. Instead of an expansion in classical vertices, it is an expansion in fully dressed ones. This expansion is closely related to nPI-resummation schemes; in the ϕ^4 -theory, it is related to a 4-PI scheme.

Here we consider the two-loop order of the skeleton expansion in the broken phase. A first observation is that the prefactor of the squint diagram vanishes: it is fully contained in the polarization diagram with two dressed three-point functions $\Gamma^{(3)}$. At perturbative two-loop level, the expansion involves a kite diagram as well as a doublebubble diagram. Both topologies are only generated from the polarization diagram in the DSE, more precisely from vertex corrections of the dressed three-point function. These contributions have to be subtracted in terms of explicit kite and double-bubble diagram in the skeleton expansion. For small field expectation values $\phi_0 \ll 1$, it is reasonable to neglect the kite diagram, since it scales like ϕ_0^4 due to its four three-point functions. We will also drop the double-bubble diagram which is of order ϕ_0^2 . These approximations are discussed again later. The remaining diagrams are the polarization, sunset, and tadpole. The present approximation of the two-loop skeleton expansion of the gap equation is depicted in Fig. 9.

1. Bubble-resummed s-channel four-point function

It is left to specify the approximations for the threepoint and four-point vertices. To begin with, we still use the relation (33) for the three-point function with the assumption of small field values ϕ_0^2 . This leaves us with the four-point function, for which we resort to a bubbleresummed *s*-channel expansion, e.g., [42,44], shown graphically in Fig. 10. Algebraically, the momentum dependence of the *s*-channel in the four-point function can be expressed as

$$\Gamma^{(4)}(p) = \frac{\lambda_{\phi}}{1 + \lambda_{\phi} \Pi_{\text{fish}}(p)}.$$
(37)

Here, Π_{fish} is the one-loop part of the *s*-channel self-energy (apart from λ_{ϕ}^2),

$$\Pi_{\text{fish}}(p) = \frac{1}{2} \int_{\lambda_1, \lambda_2} \lambda_1 \lambda_2 \rho(\lambda_1) \rho(\lambda_2) I_{\text{pol}}(p; \lambda_1, \lambda_2). \quad (38)$$



FIG. 10. *s*-channel expansion of the momentum-dependent four-point function $\Gamma^{(4)}$. With a bubble resummation, one arrives at (37). The notation is defined in Fig. 1.

This is exactly the polarization diagram also appearing in the DSE with prefactor $g_{\text{pol}} = 1$, cf. Fig. 5 and Eq. (17). The analytically continued expression for $\Gamma^{(4)}(\omega)$ is obtained by simply replacing $I_{\text{pol}}(p)$ in (38) with $I_{\text{pol}}(\omega)$.

The resulting $\Gamma^{(4)}(\omega)$ depends on the full propagator through the spectral functions in (38), and the equations for $\Gamma^{(2)}$ and $\Gamma^{(4)}$ are coupled. We also note that the *s*-channel resummation used here is obtained in the NLO-expansion of the 1/N-expansion of an O(N)-theory with *N* real scalar fields. In the present N = 1 case, *s*-channel dominance is merely an assumption. For a full quantitative study, this assumption would have to be validated numerically. Otherwise, other scattering channels would have to be taken into account as well.

The iteration procedure does not change for such a coupled system, even though coupled systems generically show worse convergence properties. For a given input pair $\Gamma^{(2)}$ and $\Gamma^{(4)}$, we compute the next iteration from the right-hand side of the DSE for $\Gamma^{(2)}$, Fig. 9, and the resummed representation of $\Gamma^{(4)}$, (37). This is repeated until convergence is reached.

2. Tadpole contribution to the sunset topology

Before turning to the explicit approximation used in the skeleton scheme, we emphasize again that the fully dressed tadpole diagram and the fully dressed sunset diagram are related. They both carry the *s*-channel of the four-point vertex. While the tadpole simply is proportional to the *s*-channel four-point vertex, the sunset includes the fish diagram as a subdiagram. Indeed, the perturbative two-loop sunset graph is a combination of the respective contributions, and the prefactor g_{sun} of the sunset diagram in the skeleton expansion is such that the perturbative prefactor, cf. Table I.

The tadpole diagram is proportional to the *s*-channel four-point vertex and we use the full momentum-dependent four-point vertex obtained from the bubble resummation (37). Inserting the diagrammatic vertex expansion explicitly into the diagram, one sees that the dressed tadpole contributes to the sunset topology on the perturbative two-loop level. However, this contribution does not account for the full prefactor of the latter. To arrive at the correct perturbative prefactor of the sunset diagram, the prefactor of the fully dressed sunset diagram in the skeleton expansion needs to be adjusted accordingly; see Fig. 9 and Table I.

3. Vertex approximation in the skeleton expansion

In the sunset diagram, the 2 four-point vertices are averaged due to the two-loop momenta that run through both vertices. This averaging holds true for both the Euclidean branch and the Minkowski one. For this reason, we approximate the full momentum-dependent four-point vertices by that at vanishing momentum, $\Gamma^{(4)}[\phi_0]$.

In our approximation with $\Gamma_{\text{pol}}^{(3)}(p) = \phi_0 \Gamma^{(4)}(p)$ with external momentum p, the vertices in the polarization diagram are, as in the tadpole, proportional to the *s*-channel four-point vertex. For the sake of simplicity, we also use $\Gamma_{\text{pol}}^{(3)} = \phi_0 \Gamma^{(4)}(\omega = 0)$. In any case, the spectral integrands I_{pol} and I_{sun} of polarization and sunset diagram remain the same as in Sec. IVA. However, their prefactors g_{pol} and g_{sun} are modified by the skeleton expansion, cf. Table I.

As outlined, our solution method requires an analytic solution of the momentum integrals for all diagrams. In the current approximation, this holds true for the polarization and sunset diagram. It is not the case for the tadpole diagram because the loop momentum is probing the non-trivial momentum structure of the resummed four-point function (37). This problem can be resolved if we can use a spectral representation for the resummed four-point function, which is discussed in the following section.

C. Spectral representation for the four-point function

While the existence and practical form of spectral representations for full four-point functions pose an intricate problem, spectral representations for (approximations of) single exchange channels of the four-point function can be derived. From a practical perspective, we may treat such a channel similarly to a propagator. This is well-motivated by considering that the resonant channels of a four-point function correspond to particle exchange interactions. Technically, this can be made explicit by means of an Hubbard-Stratonovich transformation. In analogy to a propagator, we can make the same ansatz for a spectral representation for a single channel of the resummed vertex

$$\Gamma^{(4)}(p) = \lambda_{\phi} + \int_{\lambda} \frac{\lambda \rho_4(\lambda)}{p^2 + \lambda^2}, \qquad (39)$$

with

$$\rho_4(\omega) = 2 \operatorname{Im} \Gamma^{(4)}(-i(\omega + i0^+)). \tag{40}$$



FIG. 11. Left: comparison of the momentum-dependent (Euclidean) four-vertex $\Gamma^{(4)}(p)$ in its initial form (37) with its spectral representation (39) for the coupling choices $\lambda_{\phi} = \{10, 15, 20\}$ using a classical propagator, i.e., a delta pole spectral function peaked at $m_{\text{pole}} = 1$. All dimensionful quantities were rescaled in units of the respective mass pole result. Right: corresponding spectral functions of the four-point functions. The different height of the delta peaks encodes the magnitude of the residue relative to the other spectral functions. Also here, all dimensionful quantities were rescaled in units of the respective mass pole result.

In (39), the constant classical part λ_{ϕ} has to be separated. It has no spectral representation and does not need one for the present purpose. Indeed, classical vertices have been already considered in Sec. IVA.

Our results confirm that the analytic structure of the resummed vertex is compatible with (39) and works well. This can be seen in Fig. 11, and the computational details can be found in the next section, Sec. IV D. The spectral function ρ_4 displayed in the right panel exhibits a continuous tail corresponding to the $\phi\phi \rightarrow \phi\phi$ scattering continuum for $\omega \ge 2m_{\text{pole}}$. The spectral representation (39) of the four-point function is used in the tadpole diagram in complete analogy to that of propagators.

Importantly, the spectral representation of the vertex effectively just leads to another classical propagator with spectral mass λ to the loop-momentum integral. The momentum flowing through the four-vertex is the sum of the loop and external momentum. Thus, the momentum integral of the tadpole is identical to that of the polarization diagram, since the internal line of the four-vertex carries p + q and the initial internal line just the loop momentum q. The tadpole diagram can therefore be expressed as

$$D_{\rm tad}(\omega) = g_{\rm tad} \int_{\lambda_1, \lambda_2} \lambda_1 \lambda_2 \rho(\lambda_1) \rho_4(\lambda_2) I_{\rm pol}(\omega; \lambda_1, \lambda_2).$$
(41)

The spectral integral is logarithmically divergent, since the vertex spectral function ρ_4 drops off in the UV as λ^{-1} (as opposed to $\rho \sim \lambda^{-2}$ in the UV). Again, we employ spectral BPHZ-renormalization to subtract the zeroth order term of the Taylor expansion of I_{pol} . Finally, the renormalized diagram reads

$$D_{\text{tad}}^{\text{ren}}(\omega) = g_{\text{tad}} \int_{\lambda_1, \lambda_2} \lambda_1 \lambda_2 \rho(\lambda_1) \rho_4(\lambda_2) \\ \times [I_{\text{pol}}(\omega; \lambda_1, \lambda_2) - I_{\text{pol}}(\mu; \lambda_1, \lambda_2)]. \quad (42)$$

D. Results for the coupled system of propagator and vertices

The DSE in the skeleton expansion is solved for the couplings also used in Sec. IVA, $\lambda_{\phi} = 5$, 10, 20, measured in the pole mass $m_{\text{pole}} = 1$.

For the first iteration, initial choices ρ_0 , $\rho_{4,0}$ for the spectral function of the propagator and that the four-point vertex are required. For ρ_0 , we use the classical spectral function (27), as already done in Sec. IVA. For the spectral function of the four-point function $\rho_{4,0}$, compute it from the resummed representation of the four-point function, (37) with the initial choice of the spectral function of the propagator, ρ_0 . This results in

$$\rho_{4,0}(\omega) = 2 \operatorname{Im} \frac{\lambda_{\phi}}{1 + \lambda_{\phi} \Pi_{\text{fish},0}(\omega)}, \qquad (43)$$

with

$$\Pi_{\text{fish},0} = \frac{1}{2} \int_{\lambda_1,\lambda_2} \lambda_1 \lambda_2 \rho_0(\lambda_1) \rho_0(\lambda_2) I_{\text{pol}}(\omega;\lambda_1,\lambda_2)$$
$$= \frac{1}{2} I_{\text{pol}}(\omega;m_{\text{pole}},m_{\text{pole}}), \qquad (44)$$

with $m_{\text{pole}} = 1$. While one could also simply take the classical vertex, the convergence speed and potentially also the convergence radius (coupling range) are increased by the improved choice (43). For further couplings λ_{ϕ} , we take as initial choices ρ_0 and $\rho_{4,0}$, the full solutions ρ and ρ_4 of the closest coupling value already computed. This procedure has already been used in Sec. IV A and speeds up the convergence.

The spectral function and propagator obtained from the coupled system of resummed four-point function and DSE are displayed in the top panels of Fig. 12. As for the case of bare vertices, shown in Sec. IVA, we find a distinct one-particle mass pole as well as a scattering tail. The $\phi \rightarrow \phi \phi \phi$



FIG. 12. Results from the skeleton expanded DSE (comp. Fig. 9) with a bubble resummed *s*-channel expansion of the four-point function for coupling choices $\lambda_{\phi} = 5$, 10, 20 using on-shell renormalization (26). The curves were rescaled by the respective mass poles. All vertices except for the tadpole one were approximated at $\omega = 0$. TOP: spectral function (left) and propagator (right). The weight of the continuous tail increases with coupling, and the mass pole residue decreases. Higher *n*-particle onsets are not visible in the spectral function. The different height of the delta peaks encodes the magnitude of the residue relative to the other spectral functions. The propagators were computed by the Källén-Lehmann spectral representation. Increasing coupling makes the propagators deviate more from the classical propagator, which approaches 1 at momentum $p \rightarrow 0$. Bottom: four-vertex spectral function (left) and four-vertex (right). Weight of the vertex spectral functions. The four-vertices were computed by their spectral-like representation [comp. (39)]. The quantum corrections in the IR increase with coupling. All vertices approach their classical value in the UV.

onset is not visible in the spectral function for any of the coupling configurations. It can be seen that for increasing coupling λ_{ϕ} , the tail of the spectral function becomes more enhanced, since the higher-scattering states are more accessible. In turn, the mass pole residue decreases. The corresponding propagators in the right panel show similar behavior as for the DSE with bare vertices. The larger the coupling gets, the further the propagators deviate from the classical behavior $G(p) \rightarrow 1$ for $p \rightarrow 0$.

In the bottom panels of Fig. 12, we display the spectral function of the *s*-channel four-point function and the Euclidean four-point function itself. The consistency of the spectral representation has been discussed already in the previous section and is confirmed numerically; see Fig. 11. Technically, the negativity of the spectral function displayed in Fig. 12 can be understood from the dominant quantum correction to the classical vertex, which is negative. On the conceptual side, within the Hubbard-Stratonovich transformation related to minus the spectral function of the exchange particle.

We find a continuous $2 \rightarrow 2$ scattering tail, starting at $2m_{\text{pole}}$ for all coupling choices. The spectral function is strongly enhanced with increasing coupling. Additionally, for larger coupling, the spectral functions also clearly show the $1 \rightarrow 3$ scattering onsets starting at $3m_{\text{pole}}$, which was not visible in the propagator spectral functions (cf. top left panel of Fig. 12). By simple dimensional analysis, it becomes clear that the higher *n*-particle thresholds in the propagator spectral function are suppressed by their respective energy threshold squared. This is not the case for the vertex spectral function: it decays with λ^{-1} , making the higher onsets less suppressed. In turn, the invisibility of four, five, and higher particle onsets is due to their decreasing amplitude, as every next higher onset comes with one additional loop. Further, we note that the visible size of the $1 \rightarrow 3$ scattering onset has its sole origin in the tadpole diagram. This diagram contributes to the $1 \rightarrow 3$ scattering process due to the s-channel resummed fourpoint function, cf. Sec. IV B 2. The contribution of the sunset itself is very suppressed in comparison. This points toward a general feature of our approximation: In the large coupling (massless) limit, the tadpole becomes the dominating diagram in the spectral function (for $\omega > 3m_{\text{pole}}$). We elaborate more on the massless limit in Appendix A.

While presenting the spectral and correlation functions in units of the fixed mass pole $m_{\text{pole}} = 1$ allows for a comparison of the relative strength of the different contributions, the approach to the massless limit is better studied if the results are presented in units of a uniform interaction strength. This is achieved by measuring all results in units of the coupling λ_{ϕ} ; see Appendix A. For the parameters $\lambda_{\phi}/m_{\text{pole}} = 5$, 10, 20 studied here, this entails that we consider theories with the coupling $\lambda_{\phi} = 1$ with pole masses $m_{\text{pole}} = 1/5$, 1/10, 1/20. Evidently, within these units, the spectral functions pole position moves toward zero and the onset of the scattering states gets more pronounced; see Fig. 14.

E. Self-consistent skeleton expansion

Within the current approximation, we have explicitly dropped the kite and double-bubble diagrams. Implicitly, we have also dropped the squint diagram that corresponds to vertex corrections to the three-point function in the vacuum polarization: we have used two dressed three-point vertices that are derived from the bubble resummation of the four-point function. Such an approximation of the three-point function does not contain the squint topology.

An alternative approximation of the Dyson-Schwinger equation is derived as follows: we start from the initial, full propagator DSE (Fig. 3) and consider it in a diagrammatic expansion in orders of the constant field ϕ_0 . In our approximation, these higher orders come via the three-vertices $\Gamma^{(3)}(p) = \phi_0 \Gamma^{(4)}(p) + O(\phi_0^3)$. Acting on the DSE with two derivatives with respect to the constant field ϕ_0 and multiplying by $\phi_0^2/2$ afterward, one finds the schematic relation

$$O_{\Gamma^{(2)},\text{diag}}[\phi_0^2](p) = \frac{1}{2}\phi_0^2[\Gamma^{(4)}(p,-p,0,0) - \lambda_{\phi}], \quad (45)$$

where $O_{\Gamma^{(2)},\text{diag}}[\phi_0^2]$ represents all diagrams in the propagator DSE with two external constant field legs, including their initial prefactors. This includes the squint diagram and the vacuum polarization.

On the right-hand side of (45), we have the full fourpoint function subtracted by its classical value. We also made explicit the specific momentum dependence of the four-point function. By differentiating twice with respect to the momentum-independent field ϕ_0 , $\Gamma^{(4)}$ only depends on one external momentum p.

This entails that we can reexpress all diagrams of $O_{\Gamma^{(2)},\text{diag}}[\phi_0^2]$ through an *s*-channel four-point function. What is missing is the classical part of the vertex DSE. It is included by adding and subtracting the classical vertex

contribution multiplied by an appropriate prefactor involving the constant field, $\frac{1}{2}\phi_0^2\lambda_{\phi}$ to the propagator DSE. This leads us to (45). Evidently, the additional constant part $-\frac{1}{2}\phi_0^2\lambda_{\phi}$ in (45) is absorbed in the mass renormalization.

The expectation value ϕ_0^2 can be expressed in terms of $\Gamma^{(2)}(0) = m_{\text{cur}}^2$ and $\Gamma^{(4)}[0]$. This leads us to

$$\frac{1}{2}\phi_0^2\Gamma^{(4)}(p) = \frac{3}{2}m_{\rm cur}^2\frac{\Gamma^{(4)}(p)}{\Gamma^{(4)}(0)},\tag{46}$$

which has the scaling of a two-point function and reduces to $3/2m_{\rm cur}^2$ at vanishing momentum. Making use of (45) and (46), we are led to the DSE for the propagator with Minkowski frequencies ω ,

$$\Gamma^{(2)}(\omega) = -\omega^{2} + m_{\text{pole}}^{2} + D_{\text{tad}}^{\text{ren}}(\omega) + D_{\text{sun}}^{\text{ren}}(\omega) + \frac{3}{2} m_{\text{cur}}^{2} \left[\frac{\Gamma^{(4)}(\omega) - \Gamma^{(4)}(m_{\text{pole}})}{\Gamma^{(4)}(0)} \right],$$
(47)

and i = tad, sun. Note that the polarization and squint diagram have been absorbed into the last term of (47) proportional to m_{cur}^2 as a result of (45). Due to the onshell renormalization condition (26), all renormalized diagrams vanish at $\omega = m_{\text{pole}}$, i.e., $D_i^{\text{ren}}(m_{\text{pole}}) = 0$. In summary, (47) is exact up to higher orders of ϕ_0^2 , leaving us with a self-consistent systematic expansion scheme. The self-consistency refers to the fact that in the present order in ϕ_0^2 , the polarization diagram is given exactly in terms of the four-point vertex. Therefore, approximations to the latter are transported to the former.

1. Vertex-approximation in the self-consistent skeleton expansion

As discussed above, within the self-consistent DSE in (47), it suffices to specify the approximation for the fourpoint function. Here we again resort to the bubbleresummed four-point function of (37), already used in the previous section. This approximation of the four-point function neglects in particular contributions in the DSE of the four-point function that originate in the squint diagram. The self-consistency of (47) is reflected in the fact that the contribution of the polarization diagram is given by its bubble-resummed $1/2\phi_0^2\Gamma^{(4)}$,

$$\frac{1}{2}\phi_0^2\Gamma^{(4)}(p) = \frac{1}{2}\phi_0^2[\lambda_\phi - \lambda_\phi \Pi_{\text{fish}}\Gamma^{(4)}(p)].$$
(48)

The four-point function, (48), lacks the contributions from the squint diagram. Note also that these contributions are related to the u, t-channel. Hence, with (48), we consistently neglect the squint topology in the DSE if assuming dominance of the s-channel vertex corrections for the four-point vertex. We emphasize that the assumption of s-channel dominance is well supported in the large-N limit, but less so in the present N = 1 case. Diagrammatically, the DSE is still represented by Fig. 9, with the polarization diagram given by the last term in (47). The prefactors of tadpole and sunset diagram are identical to that in the standard skeleton scheme used in Sec. IV D. They are listed in Table I. In the tadpole diagram, the four-point function again enters via its spectral representation (39).

2. Results

Numerical results for the self-consistent skeleton scheme are displayed in Fig. 13. The propagator spectral function is similar in shape to that obtained in the standard skeleton approximation; see Fig. 12. However, the magnitude of the scattering tail close to the threshold is roughly a third for all coupling choices. Comparing the two skeleton schemes, one realizes that the correct momentum scaling behavior of the standard skeleton expansion came at the price of two dressed vertices. In turn, the self-consistent skeleton expansion has the correct momentum scaling as well as the correct vertex strength. Arguably, this property is particularly important in the vicinity of *s*-channel resonances or for asymptotically large couplings.

Higher *n*-particle onsets are not visible in the propagator spectral function similarly to the results in the standard skeleton expansion. We emphasize again that they are present nevertheless, as well as easily accessible in the present spectral approach. The three-particle onset can again be seen in the vertex spectral function (bottom left panel of Fig. 13), although it is less pronounced as in Fig. 12. The magnitude of the vertex spectral functions matches very well in the two different schemes however.

The corresponding Euclidean correlation functions are shown in the top and bottom right panels. Both receive slightly less quantum corrections as for the plain skeleton expansion. For the propagator, this is quite clear from the much smaller spectral functions in the self-consistent approximation. For the four-point vertex, the differences



FIG. 13. Results from the self-consistent skeleton expanded DSE (comp. Fig. 9) with a bubble resummed *s*-channel expansion of the four-point function for coupling choices $\lambda_{\phi} = 5$, 10, 20 using on-shell renormalization (26). The polarization diagram is here expressed through the *s*-channel four-vertex, cf. (47). The curves were rescaled by the respective mass poles. All vertices except for the tadpole one were approximated at $\omega = 0$. Top: spectral function (left) and propagator (right). The weight of the continuous tail increases with coupling, and the mass pole residue decreases. Higher *n*-particle onsets are not visible. The different height of the delta peaks encodes the magnitude of the residue relative to the other spectral functions. The propagators were computed by the Källén-Lehmann spectral representation. Increasing coupling makes the propagators deviate more from the classical propagator, which approaches 1 at momentum $p \rightarrow 0$. Bottom: four-vertex spectral function (left) and four-vertex (right). Weight of the vertex spectral functions continuous tail increases with coupling. For the largest coupling choice, the three-particle onset is visible in the spectral function. The four-vertices were computed by their spectral-like representation [comp. (39)]. The quantum corrections in the IR increase with coupling. All vertices approach their classical value in the UV.

are less pronounced and we refrain from discussing them. For a more detailed discussion of the general features, see Sec. IV D.

3. Low-lying bound state close to phase transition

Lattice calculations [45–47] show an additional lowlying excitation in the spectrum of the scalar ϕ^4 -theory in d = 2 + 1 close to the phase transition with a mass of $m \approx 1.8m_{\text{pole}}$. This state has been interpreted as a bound state of the fundamental excitation in [47]. It is also been observed within the recent functional RG study [48]. The approximation scheme underlying the Euclidean computation has been done in [48]. It is close in spirit to the *s*channel approximation scheme used in the present work in Minkowski space-time. In [48], the spectral function of the propagator, numerically reconstructed from Euclidean data, shows a bound state close to the phase transition at a mass ratio consistent with that found in the lattice studies. However, further away from the phase transition, the clear signal of the bound state is lost in [48].

The present direct computation of spectral functions is not performed close to the phase transition, i.e., for $m_{\text{pole}} \rightarrow 0$. This regime will be studied elsewhere within the selfconsistent skeleton expansion developed in the present section. Our results for $m_{\text{pole}} \neq 0$ indicate that the bound state may indeed only exist close to the phase transition. However, the present *s*-channel approximation does not allow for a fully conclusive statement, as the latter requires a multichannel analysis. Yet, the *s*-channel resummation typically captures the dominant resonances and is trustworthy as next-to-leading order in a 1/N-expansion in the large-N limit. The spectral properties of O(N) models as well as the phase transition regime will be studied elsewhere.

V. CONCLUSION

In this work, we developed a spectral functional approach for the direct nonperturbative computation of real-time (timelike) correlation functions. This approach is based on a novel renormalization scheme called spectral renormalization: this renormalization scheme is based on the use of spectral representations and dimensional regularization. The spectral representation allows to perform the momentum integrals within dimensional regularization. This leaves us with the spectral integrals, whose spectral divergences can be renormalized within dimensional regularization as well; see Sec. II C. This scheme is called spectral dimensional renormalization and respects all symmetries of the theory at hand that are maintained within dimensional regularization. The latter set also includes gauge symmetries, and hence spectral dimensional renormalization as developed here is a manifestly gauge-invariant renormalization scheme.

The renormalization step can also be done with a standard subtraction procedure within a Taylor expansion in momenta as done in BPHZ-renormalization. This scheme is called spectral BPHZ-renormalization. It maintains less symmetries than spectral dimensional renormalization and in particular requires counterterms in gauge theories that break gauge symmetry (or rather BRST symmetry). The appeal of spectral BPHZ-renormalization lies in its relative numerical simplicity.

In summary, spectral renormalization allows for direct access to the real momentum axis by analytic continuation as a result of the fully analytic solution of all momentum integrals, while maintaining all symmetries of the theory at hand.

We performed explicit, nonperturbative computations within the spectral Dyson-Schwinger approach to the scalar ϕ^4 -theory in 2 + 1 space-time dimensions. Our results include the spectral function of the scalar propagator and that of the four-point function (*s*-channel), and are obtained by solving the spectral DSE recursively. We have first considered the approximation with classical vertices and full propagators, but including the two-loop diagrams in the DSE. The resulting spectral functions show a distinct one-particle pole and a clear scattering tail with onset at twice the pole mass. The spectral function contains all higher-scattering thresholds, which are easily accessible due to the analytic nature of the momentum integrations. For increasing couplings, the 1 \rightarrow 3 scattering onset becomes more pronounced.

We also considered an approximation with nonperturbative vertices, based on a skeleton expansion scheme. The respective four-point function was given by an *s*-channel bubble resummation. The propagator spectral function again shows a distinct one-particle pole and a continuous scattering tail. The spectral function of the resummed fourvertex features a scattering tail as well and additionally shows a distinct onset for the $1 \rightarrow 3$ scattering process. For larger couplings, the fully nonperturbative nature of the approximation leads to large, though only quantitative differences compared to the classical vertex computation.

In the last part of this work, Sec. IV E, we have developed a self-consistent skeleton expansion scheme. The self-consistency was obtained by relating a class of diagrams with three-point functions to the *s*-channel four-point function. This entails that the approximation used in the computation of the four-point function is also used within the three-point function diagrams. The results are qualitatively similar to that of the standard skeleton scheme. The magnitude of the scattering tail of the propagator spectral function turns out much smaller in the upgraded scheme and indicates an overestimation of the polarization diagram before. We expect that the use of such a self-consistent scheme is important close to the phase transition of the theory or, more generally, in the presence of resonant *s*-channel interactions.

In conclusion, we have developed and put to work a fully nonperturbative spectral functional approach to the computation of real-time correlation functions. The approach was successfully tested within the scalar ϕ^4 -theory in 2 + 1 dimensions. It is specifically attractive in gauge theories, as spectral dimensional renormalization preserves gauge invariance. We hope to report on respective results soon.

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APPENDIX A: MASSLESS LIMIT

The scalar ϕ^4 -theory in 2 + 1 dimensions depends on one dimensionless parameter, $\lambda_{\phi}/m_{\text{pole}}$. In the present work, we have shown all results in terms of the respective pole mass. Hence, the above parameter simply relates to different couplings λ_{ϕ} , measured in units of the pole mass $m_{\text{pole}} = 1$. If interested in the massless limit of the theory, it is more convenient to keep the coupling fixed $\lambda_{\phi} = 1$ and to depict all results for different pole masses. For example, in the work we have used $\lambda_{\phi} = 5$, 10, 20 with $m_{\text{pole}} = 1$, which can be read as $\lambda_{\phi} = 1$ and $m_{\text{pole}} = 1/5$, 1/10, 1/20. This leads us to the spectral functions for propagator and four-vertex as well as the respective Euclidean correlators itself of the scalar field, depicted or rather redrawn in Fig. 14. Rescaling the results in this way, it gets clear the massless limit is readily investigated through the limit $\lambda_{\phi} \rightarrow \infty$. For a consistent treatment of the DSE, the vertices need to be well defined and have the appropriate scaling properties in this limit. This is given for the resummed *s*-channel four-point function that was introduced in Sec. IV B and used in calculation of Secs. IV D and IV E.

Note that in the skeleton expansion as approximated in this work, the only other appearing vertex, which is the three-point function, is obtained directly from $\Gamma^{(4)}$ and hence has the same property. The resummation is hence suitable for studying the massless case in our skeleton expansion. In all diagrams except for the tadpole, all vertices are approximated at frequency zero. In consequence, they do not carry loop



FIG. 14. Results from the skeleton expanded DSE (comp. Fig. 9) with a bubble resummed *s*-channel expansion of the four-point function for coupling choices $\lambda_{\phi} = 5$, 10, 20 using on-shell renormalization (26). All vertices except for the tadpole one were approximated at $\omega = 0$. The curves were rescaled by their respective coupling parameters, i.e., $m_{\text{pole},i} = \frac{1}{\lambda_{\phi,i}}$ and consequently $\lambda_{\phi} = \frac{\lambda_{\phi,i}}{\lambda_{\phi,i}} = 1$ for all curves. Top: spectral function (left) and propagator (right). The weight of the continuous tail decreases with larger mass pole, and the mass pole residues increase. The different height of the delta peaks encodes the magnitude of the residue relative to the other spectral functions. The propagators were computed by the Källén-Lehmann spectral representation. The pole contributions are dominant, since for smaller mass pole the propagator is strongly enhanced.



FIG. 15. Example of a convergent iteration of the scalar DSE with classical vertices. The curves were *not* rescaled by the respective mass poles to demonstrate convergence of the iteration also for the pole mass. Instead, renormalization was done at $\mu = 0$. Units thus given by the input parameters. The spectral functions are alternating, approaching the final orange curve (left). Twenty iterations have been performed. The curves from iterations 10 to 19 were left out as they were graphically indistinguishable from the final curve. An iterative behavior as displayed is taken taken to be convergent, i.e., signaling a solution to the DSE where left- and right-hand sides of the equation coincide. The corresponding pole positions also converge very quickly (right). Left of the convergence boundary, assuming a (relative) precision of 10^{-4} , all points are identical.

momentum and merely enter as multiplicative factor, which makes them well under control in the large coupling limit. For the tadpole however, this is not the case. Here, the four-point function enters via its spectral representation (39). The additive classical contribution drops out by renormalization, as it contributes momentum independently. What is left is the vertex spectral function contributing to the tadpole loop integral, cf. (42). With increasing coupling, this spectral function gets larger; see Fig. 12 or 13. Since the multidimensional spectral integrals need to be evaluated numerically, a UV cutoff for the integrals needs to be chosen that minimizes the error caused by doing so. This results in increasingly long-range integrals for the large coupling limit. The appropriate treatment of these integrals hence results in a technical obstacle, which will be the subject of a follow-up project. Apart from the technical aspect, this suggests that the tadpole diagram is the dominant contribution to the DSE and thus to the spectral function for $\omega > 3m_{\text{pole}}$ in the *s*-channel approximation. As a result, one is left solely with polarization and tadpole diagram, further simplifying the setup.

APPENDIX B: NUMERICS

In order to compute the spectral integrals in (25), the integrands I_j are discretized on suitable, evenly spaced momentum grids of usually around 100 points. The grids are chosen differently for each diagram such that peaked or

Polarization:

discontinuous structures like the onset jumps in the imaginary parts are ideally resolved. The spectral integrations are performed numerically in *Mathematica* with standard global adaptive integration strategies using a relative precision goal of 10^{-3} . All diagrams are interpolated separately in real and imaginary parts in order to treat the sharp onset of the imaginary parts properly. All interpolations are performed using B-splines up to order 2 as all interpolants are real due to the split of real and imaginary parts. The spectral function is then computed from the interpolated diagrams.

For a given set of parameters, convergence was usually reached within less than ten iterations. Figure 15 demonstrates convergent behavior. We estimate the relative precision of our routine to be $\geq 10^{-4}$. Based on that, all mass poles shown in the right panel of Fig. 15 left of the convergence boundary are indistinguishable. Normalizing the scattering onsets of all spectral functions in the left panel to be identical, the continuous tail also converges pointwise beyond the convergence boundary, based on above precision estimate.

APPENDIX C: ANALYTIC EXPRESSIONS OF THE DIAGRAMS

Analytic expressions for the integrands in (25) before and after analytic continuation.

$$I_{\text{pol}}(p;\lambda_1,\lambda_2) = \frac{1}{4\pi p} \arctan\left(\frac{p}{\lambda_1 + \lambda_2}\right),$$

$$I_{\text{pol}}(\omega;\lambda_1,\lambda_2) = \frac{1}{4\pi \omega} \left[\operatorname{artanh}\left(\frac{\omega}{\lambda_1 + \lambda_2}\right) + i \arg\left(1 - \frac{\omega}{\lambda_1 + \lambda_2}\right) \right].$$
(C1)

Sunset:

$$I_{sun}(p;\lambda_{1},\lambda_{2},\lambda_{3}) = \frac{1}{(4\pi)^{2}} \left(\frac{1}{2} \log\left(\frac{1}{(\lambda_{1}+\lambda_{2}+\lambda_{3})^{2}+p^{2}}\right) - \frac{\lambda_{1}+\lambda_{2}+\lambda_{3}}{p} \arctan\left(\frac{p}{\lambda_{1}+\lambda_{2}+\lambda_{3}}\right) \right),$$

$$I_{sun}(\omega;\lambda_{1},\lambda_{2},\lambda_{3}) = \frac{1}{(4\pi)^{2}} \left(\frac{1}{2} \log\left(\frac{1}{(\lambda_{1}+\lambda_{2}+\lambda_{3})^{2}-\omega^{2}}\right) - \frac{\lambda_{1}+\lambda_{2}+\lambda_{3}}{\omega} \left[\operatorname{artanh}\left(\frac{\omega}{\lambda_{1}+\lambda_{2}+\lambda_{3}}\right) + \operatorname{i} \operatorname{arg}\left(1-\frac{\omega}{\lambda_{1}+\lambda_{2}+\lambda_{3}}\right) \right] \right).$$
(C2)

Squint:

$$I_{\text{squint}}(p;\lambda_{1},\lambda_{2},\lambda_{3},\lambda_{4}) = \frac{1}{(8\pi)^{2}\omega\lambda_{4}} \left(2\log\left(\frac{\lambda_{2}+\lambda_{3}+\lambda_{4}}{\lambda_{2}+\lambda_{3}-\lambda_{4}}\right) \operatorname{artanh}\left(\frac{\omega}{\lambda_{1}+\lambda_{4}}\right) + \mathrm{i}\left[\operatorname{Li}_{2}\left(\frac{\mathrm{i}p-\lambda_{1}-\lambda_{4}}{\lambda_{2}+\lambda_{3}-\lambda_{4}}\right) - \operatorname{Li}_{2}\left(\frac{\mathrm{i}p-\lambda_{1}-\lambda_{4}}{\lambda_{2}+\lambda_{3}-\lambda_{4}}\right)\right]\right),$$

$$I_{\text{squint}}(\omega;\lambda_{1},\lambda_{2},\lambda_{3},\lambda_{4}) = \frac{1}{(8\pi)^{2}\omega\lambda_{4}} \left(\operatorname{Re}\left[2\log\left(\frac{\lambda_{2}+\lambda_{3}+\lambda_{4}}{\lambda_{2}+\lambda_{3}-\lambda_{4}}\right) \operatorname{artanh}\left(\frac{\omega}{\lambda_{1}+\lambda_{4}}\right) - \operatorname{Li}_{2}\left(\frac{\omega-\lambda_{1}-\lambda_{4}}{\lambda_{2}+\lambda_{3}-\lambda_{4}}\right) + \operatorname{Li}_{2}\left(\frac{-\omega-\lambda_{1}-\lambda_{4}}{\lambda_{2}+\lambda_{3}-\lambda_{4}}\right) - \operatorname{Li}_{2}\left(\frac{-\omega-\lambda_{1}+\lambda_{4}}{\lambda_{2}+\lambda_{3}-\lambda_{4}}\right) + \operatorname{Li}_{2}\left(\frac{-\omega-\lambda_{1}+\lambda_{4}}{\lambda_{2}+\lambda_{3}+\lambda_{4}}\right) - \operatorname{Li}_{2}\left(\frac{-\omega-\lambda_{1}+\lambda_{4}}{\lambda_{2}+\lambda_{3}-\lambda_{4}}\right) + \operatorname{Li}_{2}\left(\frac{\omega-\lambda_{1}+\lambda_{4}}{\lambda_{2}+\lambda_{3}-\lambda_{4}}\right) - \operatorname{Li}_{2}\left(\frac{-\omega-\lambda_{1}-\lambda_{4}}{\lambda_{2}+\lambda_{3}-\lambda_{4}}\right) + \operatorname{Li}_{2}\left(\frac{-\omega-\lambda_{1}-\lambda_{4}}{\lambda_{2}+\lambda_{3}-\lambda_{4}}\right) + \operatorname{Li}_{2}\left(\frac{-\omega-\lambda_{1}+\lambda_{4}}{\lambda_{2}+\lambda_{3}-\lambda_{4}}\right) - \operatorname{Li}_{2}\left(\frac{-\omega-\lambda_{1}-\lambda_{4}}{\lambda_{2}+\lambda_{3}-\lambda_{4}}\right) + \operatorname{Li}_{2}\left(\frac{-\omega-\lambda_{1}+\lambda_{4}}{\lambda_{2}+\lambda_{3}-\lambda_{4}}\right) - \operatorname{Li}_{2}\left(\frac{-\omega-\lambda_{1}+\lambda_{4}}{\lambda_{2}+\lambda_{3}-\lambda_{4}}\right) - \operatorname{Li}_{2}\left(\frac{-\omega-\lambda_{1}-\lambda_{4}}{\lambda_{2}+\lambda_{3}-\lambda_{4}}\right) + \operatorname{Li}_{2}\left(\frac{-\omega-\lambda_{1}+\lambda_{4}}{\lambda_{2}+\lambda_{3}-\lambda_{4}}\right) - \operatorname{Li}_{2}\left(\frac{-\omega-\lambda_{1}+\lambda_{4}}{\lambda_{2}+\lambda_{3}-\lambda_{4}}\right) - \operatorname{Li}_{2}\left(\frac{-\omega-\lambda_{1}+\lambda_{4}}{\lambda_{2}+\lambda_{3}-\lambda_{4}}\right) - \operatorname{Li}_{2}\left(\frac{-\omega-\lambda_{1}+\lambda_{4}}{\lambda_{2}+\lambda_{3}-\lambda_{4}}\right) + \operatorname{Li}_{2}\left(\frac{-\omega-\lambda_{1}+\lambda_{4}}{\lambda_{2}+\lambda_{3}+\lambda_{4}}\right) - \operatorname{Li}_{2}\left(\frac{-\omega-\lambda_{1}+\lambda_{4}}{\lambda_{2}+\lambda_{3}-\lambda_{4}}\right) + \operatorname{Li}_{2}\left(\frac{-\omega-\lambda_{1}+\lambda_{4}}{\lambda_{2}+\lambda_{3}-\lambda_{4}}\right) - \operatorname{Li}_{2}\left(\frac{-\omega-\lambda_{1}+\lambda_{4}}{\lambda_{2}+\lambda_{3}-\lambda_{4}}\right) + \operatorname{Li}_{2}\left(\frac{-\omega-\lambda_{1}+\lambda_{4}}{\lambda_{2}+\lambda_{3}+\lambda_{4}}\right) - \operatorname{Li}_{2}\left(\frac{-\omega-\lambda_{1}+\lambda_{4}}{\lambda_{2}+\lambda_{3}-\lambda_{4}}\right) + \operatorname{Li}_{2}\left(\frac{-\omega$$

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