



Bosonic loop diagrams as perturbative
solutions of the classical field equations
in ϕ^4 -theory

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BOSONIC LOOP DIAGRAMS AS PERTURBATIVE SOLUTIONS OF THE CLASSICAL FIELD EQUATIONS IN ϕ^4 -THEORY

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ABSTRACT. Solutions of the classical ϕ^4 -theory in Minkowski space-time are analyzed in a perturbation expansion in the nonlinearity. Using the language of Feynman diagrams, the solution of the Cauchy problem is expressed in terms of tree diagrams which involve the retarded Green's function and have one outgoing leg. In order to obtain general tree diagrams, we set up a “classical measurement process” in which a virtual observer of a scattering experiment modifies the field and detects suitable energy differences. By adding a classical stochastic background field, we even obtain all loop diagrams. The expansions are compared with the standard Feynman diagrams of the corresponding quantum field theory.

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1. INTRODUCTION

The dynamics of interacting quantum fields is most successfully described perturbatively in terms of Feynman diagrams. One distinguishes between tree diagrams,

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which are finite, and loop diagrams, which typically diverge. Performing a suitable “classical limit” $\hbar \rightarrow 0$, the loop diagrams drop out, and only the tree diagrams remain. Therefore, it is a common opinion that tree diagrams are inherent already in the classical theory, whereas loop diagrams describe the particular effects of quantized fields. However, no hints can be found in the literature on how this statement could be made precise. This was our motivation for approaching the question coming from the classical theory: We consider a classical field, being a solution of a nonlinear hyperbolic partial differential equation. Treating the nonlinearity perturbatively, we obtain an expansion which we express in the language of Feynman diagrams. Then the question is, does the resulting expansion contain all tree diagrams of quantum field theory? Are there differences in the tree expansions? Do loop diagrams appear? Is it really impossible to obtain loop diagrams within the realm of classical field theory? Our hope was that analyzing these questions would give a better understanding of what the essence and physical significance of field quantization is. In particular, we wanted to clarify how the nonlinear dynamics of classical fields and the corresponding nonlinear scattering theory fit into the framework of a linear dynamics of quantum fields on the Fock space.

Our results are unexpected in several ways. First, rewriting the solution of the Cauchy problem in terms of Feynman diagrams, we only obtain diagrams with one outgoing leg, in which all lines correspond to retarded Green’s functions. Several outgoing legs are obtained only by considering a scattering experiment and setting up a “classical measurement process”, in which the wave is modified with an inhomogeneity at large times, and the difference of the energies before and after this modification are measured. Furthermore, in order not to distinguish a direction of time, we replace the perturbation expansion involving the retarded Green’s function by the so-called global perturbation expansion. This gives complete agreement with the tree diagrams of quantum field theory, except that, instead of the Feynman propagator, other, necessarily real-valued, Green’s functions appear. In this sense, we can confirm that tree diagrams are in fact inherent already in the classical theory. However, this statement is true only in the special setting of the classical measurement process, and only up to the differences in the choice of the propagators.

As another surprising result, we find that it is indeed possible to obtain loop diagrams in a purely classical context. To this end, we consider the classical field in a stochastic background, as described by a Gaussian ensemble of classical solutions. The resulting expansion in terms of Feynman diagrams includes all loop diagrams of quantum field theory. Our expansion differs from that of QFT in that the Feynman propagators are to be replaced by real-valued Green’s functions and fundamental solutions, with a specific combinatorics. Moreover, there are differences in the combinatorial factors for higher order loop diagrams, as we explain in a simple example.

Working with a stochastic field has some similarity with the approaches to explain quantum effects by adding a stochastic term to the classical equations (see for example Nelson’s stochastic mechanics [14] or [1, 11]). However, in contrast to these approaches, we do not modify the classical equations. We merely add a stochastic background field to the initial data. This additional field participates in the classical interaction as described by the classical field equations. The nonlinearity of the mutual interaction gives rise to the loop diagrams. The physical picture is that the macroscopic field is superimposed by microscopic fluctuations, which can be thought of as a classical

analog of vacuum fluctuations. These microscopic fluctuations can be observed only indirectly via their influence on the dynamics of the macroscopic field.

We point out that our model of a classical field theory in a stochastic background field is clearly not equivalent to a quantized field. In particular, our wave modes are classical, whereas in quantum field theory they correspond to quantum mechanical oscillators. But our point is that on the level of Feynman diagrams, these differences might not be visible. In particular, the radiative corrections might not necessarily be a test for the quantum nature of physical fields.

As is common in perturbative QFT, our treatment is formal in the sense that we do not care about the divergences of diagrams and disregard all convergence issues of the expansions. Thus all our expansions are to be understood as formal power series involving symbolic expressions. Moreover, in order to keep the setting as simple as possible, we always restrict attention to the real massless ϕ^4 -theory in $(3+1)$ -dimensional Minkowski space. But our methods and results immediately extend to any other bosonic field theory in arbitrary dimension.

The paper is organized as follows. In Section 2 we provide the necessary background on the Cauchy problem in classical field theory and on Feynman diagrams in QFT. In Section 3 we consider expansions of the classical field in terms of Feynman diagrams. Taking free fields as the starting point (Section 3.1), we express the solution of the Cauchy problem in terms of Feynman diagrams (Section 3.2). We then set up the “classical measurement process” and derive an expansion in terms of tree diagrams (Section 3.3–3.6).

Section 4 is devoted to the classical field theory in a stochastic background. After introducing the free stochastic background field (Section 4.1), this field is included in the perturbation expansion to obtain loop diagrams (Section 4.2). The comparison to the loop diagrams in QFT is given in Section 4.3.

Finally, in Section 5 we give an outlook on more realistic theories including fermions. We also outline potential applications of our methods to constructive field theory.

2. PRELIMINARIES ON ϕ^4 -THEORY

We introduce classical ϕ^4 -theory in the Lagrangian formulation. The Lagrangian \mathcal{L} is given by

$$\mathcal{L} = \frac{1}{2} (\partial_\mu \phi)(\partial^\mu \phi) - \frac{\lambda}{4!} \phi^4,$$

where ϕ is a real-valued scalar field. Integrating the Lagrangian over Minkowski space-time gives the action \mathcal{S} ,

$$\mathcal{S} = \int \mathcal{L}(\phi, \partial\phi) d^4x.$$

We are working in units where $c = 1 = \hbar$ and denote a chosen length scale by ℓ . Then the field has dimension of inverse length, whereas the coupling constant λ and the classical action are dimensionless,

$$[\phi] = \ell^{-1} \quad \text{and} \quad [\lambda] = \ell^0 = [\mathcal{S}].$$

Considering critical points of the action, one obtains the Euler-Lagrange (EL) equations

$$\square\phi = -\frac{\lambda}{6} \phi^3 \tag{2.1}$$

(where $\square = \partial_t^2 - \Delta_{\mathbb{R}^3}$ is the wave operator).

According to Noether's theorem, the symmetries of the Lagrangian correspond to conserved quantities. In particular, the symmetry under time translations gives rise to the conserved classical energy E ,

$$E(\phi) = \int_{t=T} \left(\frac{1}{2} \dot{\phi}^2 + \frac{1}{2} |\nabla\phi|^2 + \frac{\lambda}{4!} \phi^4 \right) d^3x, \quad (2.2)$$

having the units

$$[E] = \ell^{-1}.$$

2.1. Basics on the Classical Cauchy Problem and Scattering Theory. In the Cauchy problem, one seeks for solutions of the nonlinear hyperbolic partial differential equation (2.1) for initial data $(\phi, \dot{\phi})|_{t_0}$ given at some time t_0 . It is a classical result that the Cauchy problem is well-posed, meaning that there is a unique solution for short times in a suitable Sobolev space. A general method for the proof is provided by the theory of symmetric hyperbolic systems (see [10, 21]). In general, the solution of the Cauchy problem will not exist for all times, as singularities may form. It has been a major topic of mathematical research to obtain global existence results and estimates of the solution for long times (see for example the textbooks [9, 18, 12]). Here we shall disregard all analytical questions by simply assuming that our solution exists for all times.

In scattering theory, one is interested in the asymptotic behavior of the field in the limits $t \rightarrow \pm\infty$. Scattering theory has been developed mainly for linear equations (for good references on linear scattering theory in the mathematical physics literature see [16, 13, 17, 2, 24]). In the standard setting, the dynamics is described by a linear evolution equation of the form of the Schrödinger equation

$$i\partial_t\psi = H(t)\psi, \quad (2.3)$$

where $H(t)$ is a self-adjoint linear operator on a Hilbert space $(\mathcal{H}(t), (\cdot, \cdot))$. Then the corresponding time-evolution operator $U(t_1, t_2)$ defined by

$$U(t_1, t_2) : \mathcal{H}(t_2) \rightarrow \mathcal{H}(t_1) : \psi|_{t_2} \mapsto \psi|_{t_1}$$

is a unitary operator between Hilbert spaces. One wants to compare the interacting dynamics (2.3) with a free dynamics of the form

$$i\partial_t\psi = H_0(t)\psi, \quad (2.4)$$

where H_0 is a self-adjoint linear operator on a Hilbert space $(\mathcal{H}_0, (\cdot, \cdot))$. Usually, the operator H_0 is so simple that the time evolution operator $U_0(t_1, t_2)$ is explicitly known. Then the *wave operators* are introduced by considering the late-time limits

$$\Omega_{\pm}\psi := \lim_{t \rightarrow \pm\infty} U(t_0, t) U_0(t, t_0) \psi : \mathcal{H}_0 \rightarrow \mathcal{H}, \equiv \mathcal{H}(t_0) \quad (2.5)$$

where for large times we identified the Hilbert spaces \mathcal{H}_0 and $\mathcal{H}(t)$. Then the *scattering operator* S is defined by

$$S = (\Omega_-)^* \Omega_+ : \mathcal{H}_0 \rightarrow \mathcal{H}_0; \quad (2.6)$$

it maps the free outgoing field to the corresponding free incoming field.

As the ϕ^4 -theory is non-linear, the formalism of linear scattering does not immediately apply. In order to get a rigorous connection, one would need to prove that the wave dissipates for large times, going over to a solution of the free wave equation. This requires decay estimates in the spirit of [19, 6]; see also [17, Chapter 9] and [20, Section 6]. In this paper, we shall not enter dissipation and decay estimates, but instead

we will restrict attention to a systematic treatment of nonlinear classical scattering on the level of formal perturbation expansions. We will enter this analysis in Section 3. In order to get into the position to compare the different settings, we now briefly summarize the basics on quantum scattering.

2.2. Basics on Quantum Field Theory and Feynman Diagrams. We now review the basics on the corresponding ϕ^4 -quantum field theory. We begin with the formalism of *canonical quantization*. To quantize the free field, we first write the general solution of the classical equation $\square\phi = 0$ as a Fourier integral,

$$\begin{aligned}\phi_0(x) &= \int \frac{d^4k}{(2\pi)^4} \delta(k^2) \phi_0(k) e^{-ikx} \\ &= \frac{1}{(2\pi)^4} \int \frac{d^3k}{2\omega} \left(\phi_0(\omega, \vec{k}) e^{-i\omega t + i\vec{k}\vec{x}} + \phi_0(-\omega, -\vec{k}) e^{i\omega t - i\vec{k}\vec{x}} \right),\end{aligned}\tag{2.7}$$

where we set $\omega = |\vec{k}|$. The fact that our field is real means that the Fourier coefficients satisfy the relation

$$\overline{\phi_0(\omega, \vec{k})} = \phi_0(-\omega, -\vec{k}).\tag{2.8}$$

For the quantization, the Fourier coefficients are replaced by operator-valued distributions according to

$$\phi_0(\omega, \vec{k}) \longrightarrow 2\pi \sqrt{2\omega} a(\vec{k}), \quad \phi_0(-\omega, -\vec{k}) \longrightarrow 2\pi \sqrt{2\omega} a^\dagger(\vec{k}),$$

where $a(\vec{k})$ and $a^\dagger(\vec{k})$ satisfy the commutation relations

$$[a(\vec{k}), a^\dagger(\vec{q})] = (2\pi)^3 \delta^3(\vec{k} - \vec{q}), \quad [a(\vec{k}), a(\vec{q})] = 0 = [a^\dagger(\vec{k}), a^\dagger(\vec{q})].\tag{2.9}$$

We denote the resulting quantized field by $\hat{\phi}_0$,

$$\hat{\phi}_0(x) = \int \frac{d^3k}{(2\pi)^3} \frac{1}{\sqrt{2\omega}} \left(a(\vec{k}) e^{-i\omega t + i\vec{k}\vec{x}} + a^\dagger(\vec{k}) e^{i\omega t - i\vec{k}\vec{x}} \right).\tag{2.10}$$

As an immediate consequence of (2.9) and (2.10), the quantized field has the following properties:

$$\square \hat{\phi}_0(x) = 0\tag{2.11}$$

$$[\hat{\phi}_0(t, \vec{x}), \hat{\phi}_0(t, \vec{y})] = 0 = [\partial_t \hat{\phi}_0(t, \vec{x}), \partial_t \hat{\phi}_0(t, \vec{y})]\tag{2.12}$$

$$[\partial_t \hat{\phi}_0(t, \vec{x}), \hat{\phi}_0(t, \vec{y})] = -i\delta^3(\vec{x} - \vec{y}).\tag{2.13}$$

Noting that $\partial_t \hat{\phi}$ is the canonical momentum, equation (2.13) is the analog of the commutation relation $[p, q] = -i$ of quantum mechanics. For general space-time points x and y , a short calculation yields the commutation relations

$$[\hat{\phi}_0(x), \hat{\phi}_0(y)] = 2\pi K_0(x - y),\tag{2.14}$$

where $K_0(x)$ is the distribution

$$K_0(x) = \int \frac{d^4k}{(2\pi)^4} \delta(k^2) \epsilon(k^0) e^{-ikx}\tag{2.15}$$

$$= \frac{1}{(2\pi)^4} \int \frac{d^3k}{2\omega} \left(e^{-i\omega t + i\vec{k}\vec{x}} - e^{i\omega t - i\vec{k}\vec{x}} \right).\tag{2.16}$$

We remark that the equal-time commutation relations (2.12) and (2.13) immediately follow from (2.14) and (2.16) by differentiating with respect to time and taking the

limit $x^0 - y^0 \rightarrow 0$. The Fourier representation (2.15) shows that $K_0(x)$ is a distributional solution of the wave equation. Moreover, this distribution is causal in the sense that it vanishes for spacelike x . This becomes apparent from the representation

$$K_0 = \frac{1}{2\pi i} (S^\vee - S^\wedge),$$

where S^\vee and S^\wedge are the advanced and retarded Green's function defined by

$$S^\wedge(x) = \lim_{\varepsilon \searrow 0} \int \frac{d^4 k}{(2\pi)^4} \frac{e^{-ikx}}{k^2 + i\varepsilon k^0} \quad (2.17)$$

$$= \begin{cases} -i \int \frac{d^3 k}{(2\pi)^3} \frac{1}{2\omega} \left(e^{-i\omega t + i\vec{k}\vec{x}} - e^{i\omega t + i\vec{k}\vec{x}} \right) & \text{if } t > 0 \\ 0 & \text{if } t \leq 0 \end{cases} \quad (2.18)$$

$$S^\vee(x) = S^\wedge(-x). \quad (2.19)$$

Usually, the operator algebra which is generated by $a(\vec{k})$ and $a^\dagger(\vec{k})$ is realized on a Fock space $(\mathcal{F}, \langle \cdot | \cdot \rangle)$. This is a Hilbert space with a distinguished normalized vector $|0\rangle$ (the ‘‘vacuum state’’) having the following properties:

- (i) For all \vec{k} , the operator $a(\vec{k})$ annihilates the vacuum: $a(\vec{k})|0\rangle = 0$.
- (ii) The vacuum is cyclic in the sense that by iteratively applying the operators $a^\dagger(\vec{k})$ to $|0\rangle$ one generates a dense subspace of \mathcal{F} .
- (iii) For all \vec{k} , the operator $a^\dagger(\vec{k})$ is the formal adjoint of $a(\vec{k})$ with respect to the scalar product $\langle \cdot | \cdot \rangle$.

The properties (i)–(iii) together with the commutation relations (2.9) determine the scalar product. For example, setting

$$|\vec{k}\rangle := a^\dagger(\vec{k})|0\rangle \quad \text{and} \quad \langle \vec{k}| := \langle | \vec{k} \rangle | \cdot \rangle,$$

we obtain

$$\begin{aligned} \langle \vec{k} | \vec{q} \rangle &= \langle 0 | a(\vec{k}) a^\dagger(\vec{q}) | 0 \rangle = \langle 0 | [a(\vec{k}), a^\dagger(\vec{q})] | 0 \rangle \\ &= (2\pi)^3 \delta^3(\vec{k} - \vec{q}) \langle 0 | 0 \rangle = (2\pi)^3 \delta^3(\vec{k} - \vec{q}). \end{aligned}$$

Thus, introducing for a Schwartz function f the vector

$$|f\rangle = \int \frac{d^3 k}{(2\pi)^3} \frac{1}{\sqrt{2\omega}} f(\vec{k}) |\vec{k}\rangle \in \mathcal{F}, \quad (2.20)$$

one obtains

$$\langle f | g \rangle = \frac{1}{(2\pi)^3} \int \frac{d^3 k}{2\omega} \overline{f(\vec{k})} g(\vec{k}). \quad (2.21)$$

In order to introduce the interaction, one wants to construct new field operators $\hat{\phi}$ on \mathcal{F} which solve the nonlinear classical field equation,

$$\square \hat{\phi}(x) = -\lambda \hat{\phi}(x)^3. \quad (2.22)$$

This can be accomplished by the ansatz

$$\hat{\phi}(t, \vec{x}) = U(t)^\dagger \hat{\phi}_0(t, \vec{x}) U(t), \quad (2.23)$$

where $U(t)$ is a solution of the equation

$$i\partial_t U(t) = -H_{\text{int}}(t) U(t) \quad \text{with} \quad H_{\text{int}}(t) := \frac{\lambda}{4!} \int_{t=\text{const}} \hat{\phi}_0(t, \vec{x})^4 d^3 x. \quad (2.24)$$

Indeed, applying the wave operator to (2.23) and using (2.24) together with the commutation relations (2.12) and (2.13), a straightforward calculation shows that (2.22) is satisfied (for details see for example [22, Section 3.5] in the similar context of a pion-field or [15, Chapter 4]).

Writing the interaction in the form (2.23) and (2.24) is particularly useful for describing a *quantum scattering process*. To this end, we assume that in the limits $t \rightarrow \pm\infty$, the dynamics goes over to that of free fields. Thus the asymptotic states are vectors of the free Fock space $(\mathcal{F}, \langle \cdot | \cdot \rangle)$. We denote the free field operators in the asymptotic regions by $\hat{\phi}_{\text{in}}$ and $\hat{\phi}_{\text{out}}$ (note that they are solutions of the free field equation (2.11)). According to common notation, the Fock space $(\mathcal{F}, \langle \cdot | \cdot \rangle)$ with the basis generated by a_{in}^\dagger and a_{out}^\dagger is denoted by $(\mathcal{F}_{\text{in}}, \langle \cdot | \cdot \rangle_{\text{in}})$ and $(\mathcal{F}_{\text{out}}, \langle \cdot | \cdot \rangle_{\text{out}})$, respectively. By construction, these two Fock spaces are identical, but they come with different orthonormal bases. In order to describe the transformation of one basis to the other, one considers the solution of the above equations (2.23) and (2.24) with $\hat{\phi}_0 = \hat{\phi}_{\text{in}}$ and

$$\lim_{t \rightarrow -\infty} U(t) = \mathbf{1}. \quad (2.25)$$

Then $U(t)$ can be given in terms of a time-ordered exponential,

$$\begin{aligned} U(t) &= \text{Texp} \left(i \int_{-\infty}^t H_{\text{int}}(\tau) d\tau \right) = \text{Texp} \left(\frac{i\lambda}{4!} \int_{\{y^0 < t\}} \hat{\phi}_{\text{in}}(y)^4 d^4y \right) \\ &= \sum_{n=0}^{\infty} \frac{1}{n!} \left(\frac{i\lambda}{4!} \right)^n \int_{\{y_1^0 < t\}} d^4y_1 \cdots \int_{\{y_n^0 < t\}} d^4y_n T \left(\hat{\phi}_{\text{in}}(y_1)^4 \cdots \hat{\phi}_{\text{in}}(y_n)^4 \right). \end{aligned}$$

The scattering operator S is defined by

$$S := \lim_{t \rightarrow \infty} U(t) \quad (2.26)$$

$$= \sum_{n=0}^{\infty} \frac{1}{n!} \left(\frac{i\lambda}{4!} \right)^n \int d^4y_1 \cdots \int d^4y_n T \left(\hat{\phi}_{\text{in}}(y_1)^4 \cdots \hat{\phi}_{\text{in}}(y_n)^4 \right). \quad (2.27)$$

Then, according to (2.23),

$$\hat{\phi}_{\text{out}} = S^\dagger \hat{\phi}_{\text{in}} S.$$

Moreover, noting that H_{int} is a symmetric operator on the Fock space, one sees that S is a unitary operator on \mathcal{F} ,

$$\begin{aligned} S &: \mathcal{F} \rightarrow \mathcal{F} \text{ unitary} \\ |\text{out}\rangle &\mapsto |\text{in}\rangle. \end{aligned}$$

We remark that this construction is very similar to the procedure in classical linear scattering theory. While the scattering operator in the linear classical setting (2.6) is a linear operator on the Hilbert space \mathcal{H}_0 of classical fields, the quantum scattering operator (2.26) is a linear operator on the Fock space \mathcal{F} . Since here we are working in the interaction picture, the free dynamics U_0 is the identity. Then the relations (2.5) and (2.6) reduce to $S = U(-\infty, \infty)$. This is consistent with (2.26) if in view of (2.25) we identify $U(t)$ with $U(-\infty, t)$.

The scattering amplitude of a scattering process can be expressed as a matrix element of the scattering operator with respect to the in-basis,

$$\begin{aligned} \langle \beta_{\text{out}} | \alpha_{\text{in}} \rangle &= \langle \beta_{\text{in}} | S | \alpha_{\text{in}} \rangle \\ &= \sum_{n=0}^{\infty} \frac{1}{n!} \left(-\frac{i\lambda}{4!} \right)^n \int d^4 y_1 \cdots \int d^4 y_n \langle \beta_{\text{in}} | T \left(\hat{\phi}_{\text{in}}(y_1)^4 \cdots \hat{\phi}_{\text{in}}(y_n)^4 \right) | \alpha_{\text{in}} \rangle \\ &= \langle \beta_{\text{in}} | \alpha_{\text{in}} \rangle - \frac{i\lambda}{4!} \int d^4 y \langle \beta_{\text{in}} | \hat{\phi}_{\text{in}}(y)^4 | \alpha_{\text{in}} \rangle + \mathcal{O}(\lambda^2). \end{aligned} \quad (2.28)$$

Generating the in-basis by iteratively applying $\hat{\phi}_{\text{in}}$ to the vacuum (LSZ reduction formalism; see for example [15]), all expectation values can be expressed in terms of the n -point functions defined by

$$G_0(x_1, \dots, x_n) = \langle 0 | T(\hat{\phi}_{\text{in}}(x_1) \cdots \hat{\phi}_{\text{in}}(x_n)) | 0 \rangle.$$

Using the commutation relations, the n -point functions can be written as sums of products of the two-point functions (Wick's theorem). A short computation using (2.9) and (2.10) yields for the two-point function

$$G_0(x, y) = i\Delta_F(x - y), \quad (2.29)$$

where Δ_F is the Feynman propagator,

$$\begin{aligned} \Delta_F(x) &:= \lim_{\varepsilon \searrow 0} \int \frac{d^4 k}{(2\pi)^4} \frac{e^{-ikx}}{k^2 + i\varepsilon} \\ &= -i \int \frac{d^3 k}{(2\pi)^3} \frac{1}{2\omega} \left(\Theta(t) e^{-i\omega t + i\vec{k}\vec{x}} + \Theta(-t) e^{i\omega t - i\vec{k}\vec{x}} \right). \end{aligned} \quad (2.30)$$

The Feynman propagator satisfies the defining equation of a Green's function

$$-\square \Delta_F(x) = \delta^4(x).$$

However, in contrast to the advanced and retarded Green's function, it is complex-valued, and it is non-zero in space-like directions.

For the systematic treatment of perturbation theory, it is most convenient to work with a generating functional in the *path integral formalism*. We thus introduce the generating functional

$$Z_\lambda[j] = \int e^{i(S[\phi] + j \cdot \phi)} \mathcal{D}\phi \quad \text{where} \quad j \cdot \phi := \int j(x) \phi(x) d^4 x.$$

The generating functional is related to the vacuum expectation value of canonical quantization by (see for example [23] and [8])

$$Z_\lambda[j] = \langle 0 | T \exp \left(-\frac{i\lambda}{4!} \int \hat{\phi}_{\text{in}}^4(x) d^4 x + i \int j(x) \hat{\phi}_{\text{in}}(x) d^4 x \right) | 0 \rangle.$$

In order to compute $Z_\lambda[j]$, we rewrite the generating functional as

$$Z_\lambda[j] = \exp \left(-\frac{i\lambda}{4!} \int d^4 x \left(-i \frac{\delta}{\delta j(x)} \right)^4 \right) Z_0[j], \quad (2.31)$$

where $Z_0[j]$ is the generating functional of the free quantum field theory:

$$\begin{aligned} Z_0[j] &:= \lim_{\varepsilon \searrow 0} \int \mathcal{D}\phi \exp \left(i \int \left(-\frac{1}{2} \phi(x) ((\square - i\varepsilon)\phi(x)) + j(x) \phi(x) \right) d^4x \right) \\ &= \lim_{\varepsilon \searrow 0} \int \mathcal{D}\mu_\varepsilon(\phi) e^{i(\mathcal{S}_0[\phi] + j \cdot \phi)}. \end{aligned}$$

Here \mathcal{S}_0 is the free action,

$$\mathcal{S}_0[\phi] = \frac{1}{2} \int (\partial_\mu \phi)(x) (\partial^\mu \phi)(x) d^4x,$$

and $\mathcal{D}\mu_\varepsilon(\phi)$ is the Gaussian measure

$$\mathcal{D}\mu_\varepsilon(\phi) = e^{-\varepsilon \int \phi(x)^2 d^4x} \mathcal{D}\phi.$$

This path integral can be computed explicitly to obtain

$$Z_0[j] = \exp \left(-\frac{i}{2} \iint j(x) \Delta_F(x, y) j(y) d^4x d^4y \right) \quad (2.32)$$

(we use the convention $Z_0[0] = 1$). In particular,

$$G_0(x, y) = -\frac{\delta}{\delta j(x)} \frac{\delta}{\delta j(y)} Z_0[j] \Big|_{j=0} = i \Delta_F(x, y),$$

in agreement with (2.29). Wick's theorem can be expressed as

$$\begin{aligned} G_0(x_1, \dots, x_n) &= (-i)^n \frac{\delta}{\delta j(x_1)} \cdots \frac{\delta}{\delta j(x_n)} Z_0[j] \Big|_{j=0} \\ &= (-i)^n \frac{\delta}{\delta j(x_1)} \cdots \frac{\delta}{\delta j(x_n)} e^{-\frac{i}{2} \iint j(x) \Delta_F(x, y) j(y) d^4x d^4y} \Big|_{j=0}, \end{aligned} \quad (2.33)$$

showing that the n -point functions can indeed be expressed as sums of products of the Feynman propagator.

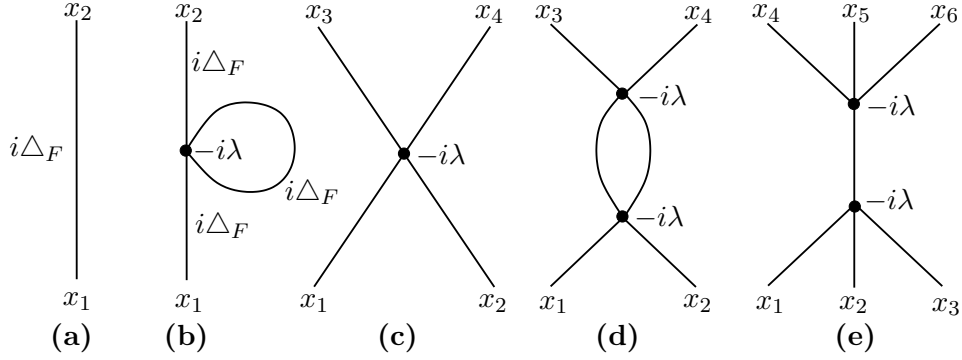
The scattering amplitudes are most conveniently expressed in terms of the *interacting n -point functions* G_λ , which are obtained by taking functional derivatives of the generating functional with interaction,

$$G_\lambda(x_1, \dots, x_n) := (-i)^n \frac{\delta}{\delta j(x_1)} \cdots \frac{\delta}{\delta j(x_n)} Z_\lambda[j] \Big|_{j=0}, \quad (2.34)$$

Their perturbation expansion is again performed by expanding with respect to λ , (2.31), using (2.32) and applying the Wick rules. *Feynman diagrams* are the pictorial representation of the perturbation expansion for the interacting n -point functions. They involve tree and loop diagrams; see Figure 1 for a few examples. A diagram is called *connected* if all outer lines are connected to each other. In the functional calculus, the connected diagrams are represented by the generating functional $W_\lambda(j) = \log Z_\lambda(j)$. With this in mind, in what follows we shall always restrict attention to connected diagrams. For more details and further reading, we refer for example to [23, 25, 22].

For the later comparison with the classical theory, we now give the combinatorics of the diagrams. We use the short notation

$$\Delta_F(j, j) \equiv \iint \Delta_F(x, y) j(x) j(y) d^4x d^4y. \quad (2.35)$$

FIGURE 1. Feynman diagrams for the interacting n -point functions.

Lemma 2.1. *Suppose that we consider a contribution to $G_\lambda(x_1, \dots, x_n)$ of order k in perturbation theory. Then G_λ vanishes unless n is even. Then the contribution can be represented by a Feynman diagram involving*

$$N = \frac{n}{2} + 2k \quad (2.36)$$

lines. Analytically, the contribution can be written as

$$(-i\lambda)^k i^N \int d^4 y_1 \cdots \int d^4 y_k \prod_{\ell=1}^N \Delta_F(\cdot, \cdot). \quad (2.37)$$

For connected tree diagrams, we have the additional relation

$$k = \frac{n}{2} - 1. \quad (2.38)$$

Proof. According to (2.34), (2.31) and (2.32),

$$\begin{aligned} G_\lambda(x_1, \dots, x_n) &= (-i)^n \frac{\delta}{\delta j(x_1)} \cdots \frac{\delta}{\delta j(x_n)} \\ &\times \frac{1}{k!} \left(-\frac{i\lambda}{4!} \right)^k \int d^4 y_1 \left(-i \frac{\delta}{\delta j(y_1)} \right)^4 \cdots \int d^4 y_k \left(-i \frac{\delta}{\delta j(y_k)} \right)^4 \frac{1}{N!} \left(-\frac{i}{2} \Delta_F(j, j) \right) \Big|_{j=0}. \end{aligned}$$

For this contribution to be non-zero, we must have as many derivatives $\delta/\delta j$ as factors j . This gives (2.36). Carrying out the derivatives, all the factorials and factors $1/2$ are compensated by combinatorial factors. Collecting the factors of $-i$ and using (2.36) gives (2.37).

To prove (2.38) we note that in the case $k = 0$, only the two-point function contributes. Obviously, each vertex increases n by two. \square

We finally review how to take the “classical limit” in which only the tree diagrams remain. As we want to take the limit $\hbar \rightarrow 0$, we clearly need to work in more general units where $\hbar \neq 1$ (but still $c = 1$). Denoting the length scale again by ℓ and the mass scale by m , all objects have dimensions of ℓ and m ; more precisely,

$$[S] = [\hbar] = m \ell, \quad [\phi] = \sqrt{\frac{m}{\ell}}, \quad [\lambda] = \frac{1}{m^2 \ell^2}, \quad [E] = m. \quad (2.39)$$

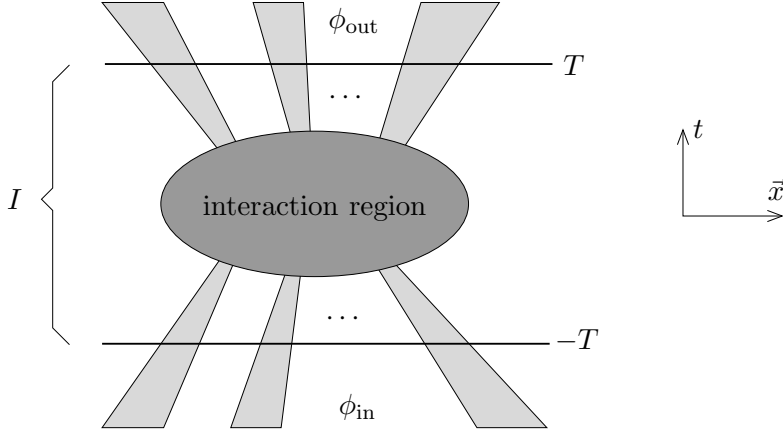


FIGURE 2. A nonlinear classical scattering process.

After inserting factors of \hbar into the n -point functions by

$$G_\lambda(x_1, \dots, x_n) := \frac{\hbar^{-n+1}}{Z_\lambda} \int \phi(x_1) \cdots \phi(x_n) e^{\frac{i}{\hbar} S} \mathfrak{D}\phi,$$

a short calculation using (2.36) and (2.38) shows that in the tree diagrams all factors \hbar drop out. Every loop, however, gives rise to a factor \hbar . Thus taking the limit $\hbar \rightarrow 0$, precisely the tree diagrams remain.

3. A NONLINEAR CLASSICAL SCATTERING PROCESS

In this section we shall analyze a scattering process for a classical field ϕ being a solution of the classical field equation (2.1). The physical situation which we have in mind is shown in Figure 2. We have an incoming wave ϕ_{in} , which may be composed of several wave packets. We assume for convenience that for large negative times, the wave packets become more spread out in space, implying that their amplitude becomes smaller. This has the advantage that in the limit $t \rightarrow -\infty$, ϕ can be treated as a free field (making it unnecessary to “switch off” the interaction adiabatically). Similarly, for large positive times, ϕ should evolve into a free field ϕ_{out} . We assume that the interaction takes place inside the strip $I = [-T, T] \times \mathbb{R}^3$ for a given large time T (in the end, we will take the limit $T \rightarrow \infty$). Inside the interaction region, ϕ satisfies the nonlinear field equation (2.1), whereas outside the interaction region, ϕ is a solution of the linear wave equation.

In the following subsections we first summarize how the Hilbert space of classical fields can be regarded as a subspace of the quantum Fock space. Afterwards, we discuss the perturbation expansion of the classical (non-linear) Cauchy problem. The corresponding pictorial representation of this expansion significantly differs from Feynman diagrams obtained in perturbative quantum field theory. To overcome this difference we motivate and introduce the so-called “nonlinear classical measurement process”. This will make it possible to introduce the *classical scattering operator* and the *classical n -point functions*. The pictorial description of the perturbative expansion of these classical n -point functions is then compared with the tree-diagrams of the corresponding quantum field theory.

3.1. The Free Classical Field. In preparation, we need to describe free fields, being solutions of the linear wave equation

$$\square\phi_0 = 0. \quad (3.1)$$

We assume for simplicity that ϕ_0 is smooth and spatially compact (note that, due to finite propagation speed, the last property is preserved under the time evolution),

$$\phi_0 \in C^\infty(\mathbb{R}, C_0^\infty(\mathbb{R}^3)).$$

For free fields, the coupling constant λ vanishes, and thus the classical energy (2.2) is quadratic. Thus by polarization we can introduce a scalar product,

$$(\phi_0, \psi_0) := \frac{1}{2} \int_{t=\text{const}} \left(\dot{\phi}_0 \dot{\psi}_0 + \vec{\nabla} \phi_0 \vec{\nabla} \psi_0 \right) d^3x. \quad (3.2)$$

Due to energy conservation, this scalar product is independent of t . It makes the space of free solutions to a Hilbert space $(\mathcal{H}_0, (\cdot, \cdot))$.

It will be useful to represent the function ϕ_0 in various ways. First, in view of the unique Cauchy development, we can describe ϕ_0 by its initial values at any given time t ,

$$\Phi(t) := (\phi_0, \partial_t \phi_0)|_t \in C_0^\infty(\mathbb{R}^3) \times C_0^\infty(\mathbb{R}^3). \quad (3.3)$$

Next, we can represent ϕ_0 similar to (2.7) and (2.8) as a Fourier integral supported on the upper and lower mass cone,

$$\phi_0(x) = \frac{1}{(2\pi)^4} \int \frac{d^3k}{2\omega} \left(\phi_0(\vec{k}) e^{-i\omega t + i\vec{k}\vec{x}} + \overline{\phi_0(\vec{k})} e^{i\omega t - i\vec{k}\vec{x}} \right) \quad (3.4)$$

where again $\omega = |\vec{k}|$. Then the energy scalar product (3.2) becomes

$$(\phi_0, \phi_0) = \frac{1}{(2\pi)^4} \int \frac{d^3k}{2\omega} \frac{\omega}{2\pi} |\phi_0(\vec{k})|^2. \quad (3.5)$$

This scalar product has the same units as the classical energy (see after (2.2)),

$$[(\phi_0|\psi_0)] = \ell^{-1}. \quad (3.6)$$

Another scalar product can be obtained by identifying the classical solutions with vectors of the one-particle Fock space: For given ϕ_0 , we seek a Schwartz function f such that

$$\phi_0(x) = 2 \operatorname{Re} \langle 0 | \hat{\phi}_0(x) | f \rangle.$$

Using (2.20) and the commutation relations, we find that f is uniquely determined by

$$f(\hat{k}) = \phi_0(\vec{k}).$$

We introduce the real scalar product $\langle \cdot | \cdot \rangle$ on \mathcal{H}_0 by $\langle \phi_0 | \phi_0 \rangle := \langle f | f \rangle$. Using (2.21), this scalar product has the representation

$$\begin{aligned} \langle \phi_0 | \psi_0 \rangle &= \frac{1}{(2\pi)^3} \operatorname{Re} \int \frac{d^3k}{2\omega} \overline{\phi_0(\vec{k})} \psi_0(\vec{k}) \\ &= \frac{1}{2(2\pi)^3} \int \frac{d^3k}{2\omega} \left(\overline{\phi_0(\vec{k})} \psi_0(\vec{k}) + \phi_0(\vec{k}) \overline{\psi_0(\vec{k})} \right). \end{aligned} \quad (3.7)$$

This scalar product is Lorentz invariant. Moreover, comparing with (3.5) and (3.6), one sees that it is dimensionless,

$$[\langle \phi_0 | \psi_0 \rangle] = \ell^0. \quad (3.8)$$

It is easy to verify that, up to a multiplicative constant, the scalar product (3.7) is indeed the only Lorentz invariant scalar product which can be introduced on the free classical scalar fields. We thus obtain the isometric embedding

$$\begin{aligned} (\mathcal{H}_0, \langle \cdot | \cdot \rangle) &\hookrightarrow (\mathcal{F}, \langle \cdot | \cdot \rangle) \\ \phi_0 &\mapsto |f\rangle. \end{aligned} \quad (3.9)$$

3.2. Perturbative Solution of the Classical Cauchy Problem. We now return to the nonlinear wave equation (2.1). For simplicity, we only consider solutions which are smooth with spatially compactly support,

$$\phi \in C^\infty(\mathbb{R}, C_0^\infty(\mathbb{R}^3)).$$

We denote the set of all such solutions by \mathcal{H} . Note that, since our equation is nonlinear, \mathcal{H} is not a vector space. Moreover, there is no bilinear form or scalar product on \mathcal{H} . The only quantity available is the classical energy (2.2),

$$E : \mathcal{H} \rightarrow \mathbb{R}_0^+.$$

For the description of the Cauchy problem, it is convenient to again combine ϕ and its time derivative to a two-component function Φ , (3.3), and to write the equation as a system of first order in time,

$$\partial_t \Phi = \mathfrak{H}(\Phi) \quad \text{with} \quad \mathfrak{H}(\Phi) = \begin{pmatrix} \dot{\phi} \\ \Delta \phi - \lambda \phi^3/6 \end{pmatrix}. \quad (3.10)$$

Then the initial data at some time t_0 is a vector

$$\Phi(t_0) \in C_0^\infty(\mathbb{R}^3)^2.$$

The time evolution obtained by solving the Cauchy problem gives rise to the mapping

$$U(t, t_0) : C_0^\infty(\mathbb{R}^3)^2 \rightarrow C_0^\infty(\mathbb{R}^3)^2 \quad \text{with} \quad \Phi(t) = U(t, t_0) \star \Phi(t_0).$$

The operator $U(t, t_0)$ is referred to as the *time evolution operator*. Here the star emphasizes that it is a *nonlinear* operator. But clearly, the time evolution has the group property,

$$U(t'', t') \star U(t', t) = U(t'', t) \quad \text{for all } t, t', t'' \in \mathbb{R}. \quad (3.11)$$

Let us rewrite the solution of the Cauchy problem perturbatively. We decompose the operator \mathfrak{H} into its linear and nonlinear parts,

$$\mathfrak{H}(\Phi) = \mathfrak{H}_0 \Phi + \lambda B(\Phi),$$

where

$$\mathfrak{H}_0 = \begin{pmatrix} 0 & 1 \\ \Delta & 0 \end{pmatrix}, \quad B(\Phi) = \frac{1}{6} \begin{pmatrix} 0 \\ -\phi^3 \end{pmatrix}.$$

In the case $\lambda = 0$ without interaction, the equation (3.10) is linear. The corresponding Cauchy problem can be solved formally by an exponential,

$$\Phi(t) = e^{(t-t_0)\mathfrak{H}_0} \Phi(t_0). \quad (3.12)$$

In the next lemma, we express the linear time evolution operator $e^{(t-t_0)\mathfrak{H}_0}$ in terms of the retarded Green's function (2.17).

Lemma 3.1. *For any $t \geq 0$, the operator $e^{t\mathfrak{H}_0}$ can be written as*

$$(e^{t\mathfrak{H}_0}\Phi)(\vec{x}) = \int R_t(\vec{x} - \vec{y}) \Phi(\vec{y}) d^3y, \quad (3.13)$$

where the integral kernel is the distribution

$$R_t(\vec{x}) = \begin{pmatrix} -\partial_t S^\wedge(t, \vec{x}) & -S^\wedge(t, \vec{x}) \\ -\Delta S^\wedge(t, \vec{x}) & -\partial_t S^\wedge(t, \vec{x}) \end{pmatrix}. \quad (3.14)$$

Proof. From the representation (2.18) one sees that for positive times, $S^\wedge(t, \vec{x})$ is a solution of the wave equation. Thus a short calculation yields that the function (3.13) is a solution of the equation $(\partial_t - \mathfrak{H}_0)(e^{t\mathfrak{H}_0}\Phi) = 0$. Next, one easily verifies from (2.18) that

$$\lim_{t \searrow 0} S^\wedge(t, \vec{x}) = 0 \quad \text{and} \quad \lim_{t \searrow 0} \partial_t S^\wedge(t, \vec{x}) = -\delta^3(\vec{x}),$$

showing that (3.13) has the correct initial values at $t = 0$. \square

In order to treat the interaction perturbatively, we write (3.10) as

$$(\partial_t - \mathfrak{H}_0)\Phi = \lambda B(\Phi). \quad (3.15)$$

Similar to the interaction picture, we set

$$\Phi_{\text{int}}(t) = e^{-t\mathfrak{H}_0} \Phi(t), \quad B_{\text{int}}(\Phi_{\text{int}}) = e^{-t\mathfrak{H}_0} B(e^{t\mathfrak{H}_0} \Phi_{\text{int}}).$$

Then (3.15) simplifies to

$$\partial_t \Phi_{\text{int}} = \lambda B_{\text{int}}(\Phi_{\text{int}}). \quad (3.16)$$

Making an ansatz as a formal expansion in powers of λ ,

$$\Phi_{\text{int}}(t) = \Phi_{\text{int}}^{(0)} + \lambda \Phi_{\text{int}}^{(1)}(t) + \lambda^2 \Phi_{\text{int}}^{(2)}(t) + \dots \quad (3.17)$$

(where $\Phi_{\text{int}}^{(0)}$ is a time independent wave with the correct initial data), we obtain for $n = 1, 2, \dots$ the equations

$$\partial_t \Phi_{\text{int}}^{(n)}(x) = - \begin{pmatrix} 0 \\ \rho_{\text{int}}^{(n)}(x) \end{pmatrix}, \quad (3.18)$$

where we set $x = (t, \vec{x})$ and

$$\begin{pmatrix} 0 \\ \rho_{\text{int}}^{(n)}(x) \end{pmatrix} = e^{-t\mathfrak{H}_0} \begin{pmatrix} 0 \\ \rho^{(n)}(x) \end{pmatrix}, \quad (3.19)$$

$$\rho^{(n)}(x) = \sum_{a,b,c \text{ with } a+b+c=n-1} \frac{1}{6} \phi^{(a)}(x) \phi^{(b)}(x) \phi^{(c)}(x). \quad (3.20)$$

Integrating (3.18) on both sides and using that $\Phi_{\text{int}}^{(n)}(t_0, \vec{x}) = 0$ for $n = 1, 2, \dots$, we obtain

$$\Phi_{\text{int}}^{(n)}(t) = - \int_{t_0}^t \begin{pmatrix} 0 \\ \rho_{\text{int}}^{(n)}(\tau) \end{pmatrix} d\tau.$$

Transforming back to Φ gives

$$\Phi^{(n)}(t) = - \int_{t_0}^t e^{(t-\tau)\mathfrak{H}_0} \begin{pmatrix} 0 \\ \rho^{(n)}(\tau) \end{pmatrix} d\tau.$$

Using the explicit form of the free time evolution operator (3.14), we obtain

$$\Phi^{(n)}(t, \vec{x}) = - \int_{t_0}^t d\tau \int d^3y R_{t-\tau}(\vec{x} - \vec{y}) \begin{pmatrix} 0 \\ \rho^{(n)}(\tau, \vec{y}) \end{pmatrix}.$$

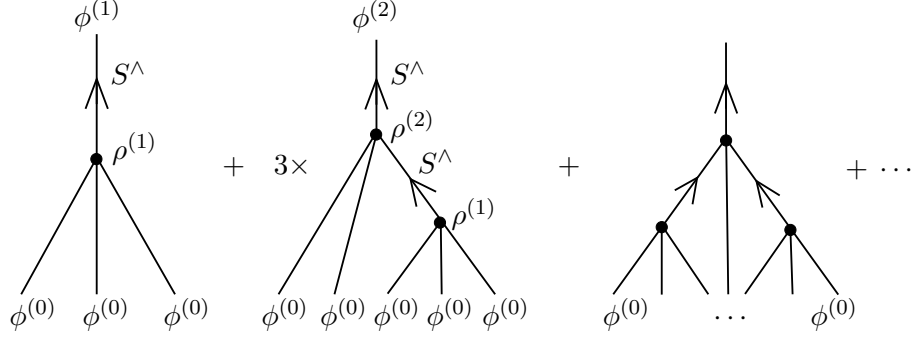


FIGURE 3. The tree diagrams in the perturbation expansion of the classical Cauchy problem.

Setting $y = (\tau, \vec{y})$, the perturbation expansion (3.17) becomes

$$\phi(x) = \sum_{n=0}^{\infty} \lambda^n \phi^{(n)}, \quad (3.21)$$

where $\phi^{(0)} = \phi_0$ is a solution of the free wave equation with the correct initial values, and

$$\phi^{(n)}(x) = \int_{\{y^0 > t_0\}} S^\wedge(x, y) \rho^{(n)}(y) d^4y \quad (n \geq 1). \quad (3.22)$$

We thus obtain an expansion of ϕ in terms of tree diagrams, which can be depicted as shown in Figure 3.

Let us briefly compare the above perturbation expansion for the classical time evolution with the perturbation expansion for the quantum field as outlined in Section 2.2: The equation for the classical evolution equation in the interaction picture (3.16) can be regarded as the analog of the quantum time evolution operator U in (2.24). Although formally similar, they differ in that the classical evolution is non-linear in the classical field, whereas the quantum evolution is linear in the quantum field. The resulting perturbation expansion for the classical field differs from that for the quantum field in that only non-loop diagrams appear (see Figures 1 and 3). Moreover, the diagrams corresponding to the perturbative expansion of the classical field always involve the retarded Green's function, instead of the Feynman propagator, and they all have exactly one outgoing leg.

3.3. A Nonlinear Classical Measurement Process. Let us try to mimic the construction in linear scattering theory as outlined in Section 2.1 (see (2.3)–(2.6)). Denoting the free classical dynamics (3.12) by $U_0(t, t') = e^{(t-t')\mathfrak{H}_0}$, we can introduce in analogy to (2.5)

$$\Omega_{\pm} \psi := U(t_0, \pm T) \star U_0(\pm T, t_0) \psi : \mathcal{H}_0 \rightarrow \mathcal{H},$$

where at time T we identify free and interacting solutions. Note that the operators Ω_{\pm} are nonlinear. Moreover, as \mathcal{H} is not a linear space, we cannot take their adjoints. Rewriting (2.6) with the inverse, the obvious idea is to define the nonlinear scattering operator by

$$(\Omega_-)^{-1} \star \Omega_+ = U_0(t_0, -T) \star U(-T, T) \star U_0(T, t_0),$$

where in the last step we used the group property (3.11). As the operator $U_0(T, t_0)$ maps a free solution at time t_0 to the same solution at time T , it is the identity on \mathcal{H}_0 . Thus the naive ansatz for the scattering operator is

$$S : \mathcal{H}_0 \rightarrow \mathcal{H}_0, \quad S \phi_{\text{out}} = U(-T, T) \star \phi_{\text{out}}.$$

Similar to (2.28), a matrix element of the scattering operator would be given by

$$(\beta_{\text{out}}, \alpha_{\text{in}}) = (S^{-1} \star \beta_{\text{in}}, \alpha_{\text{in}}) \quad (3.23)$$

(alternatively, one could work with the scalar product (3.7); this would make no difference for the following consideration). However, this naive approach does not work for the following reasons. The first problem is that the construction manifestly distinguishes a direction of time. This can be seen from the fact that, similar to (3.22), the perturbation expansion for $U(-T, T)$ will involve only advanced but no retarded Green's functions. Moreover, as $U(-T, T)$ is a nonlinear operator, the expression (3.23) is nonlinear in β_{in} , but it is linear in α_{in} . A related problem is that S is not unitary,

$$(S^{-1} \star \beta_{\text{in}}, \alpha_{\text{in}}) \neq (\beta_{\text{in}}, S \star \alpha_{\text{in}}).$$

This inequality is obvious because the left side is linear in α_{in} , whereas the right side is not.

In order to find a more appropriate description, let us discuss the physical situation that an observer at some large time T wants to detect the result of the scattering process. The only well-defined quantity which we can use for his observation is the classical energy (2.2). The most obvious method to make observations is to modify the wave ϕ by some ‘‘test wave’’ $\delta\phi$ and to consider how the energy changes. If the test wave is taken into account linearly, the energy is perturbed by

$$\delta E = \int_{t=T} \left(\dot{\phi} \delta \dot{\phi} + \nabla \phi \nabla (\delta \phi) + \frac{\lambda}{6} \phi^3 \delta \phi \right) d^3 x.$$

Using that at time T , the outgoing wave packets should be so spread out that the term involving λ can be dropped, we obtain the simple expression

$$\delta E = (\phi, \delta \phi), \quad (3.24)$$

where (\cdot, \cdot) is the scalar product obtained by polarizing the free energy (3.2) at time T . Despite the fact that we are considering purely classical fields, the resulting situation has similarity with a measurement process in quantum mechanics. Namely, the computation of the expectation value in (3.24) can be interpreted that a measurement of the wave ϕ is performed with a prepared ‘‘end-state’’ $\delta\phi$. By modifying $\delta\phi$, one can determine ϕ completely. In particular, one can measure the distribution of ϕ in momentum space.

The important point for what follows is that the relation (3.24) only holds in the linear approximation. If the amplitude of $\delta\phi$ is increased, it has a nonlinear effect on ϕ , which influences the result of the measurement process. In order to analyze such nonlinear effects, we consider the following idealized situation (see Figure 4). It is useful to describe the incoming field by an inhomogeneity ρ_{in} which lies in the past of the interaction region and generates a wave moving to the future. Similarly, the observer is described by an inhomogeneity ρ_{out} , which is located in the future of the interaction region and generates a wave moving to the past. The resulting wave ϕ will be a solution of the nonlinear inhomogeneous wave equation

$$-\square \phi - \frac{\lambda}{6} \phi^3 = \frac{1}{2} (\rho_{\text{in}} + \rho_{\text{out}}). \quad (3.25)$$

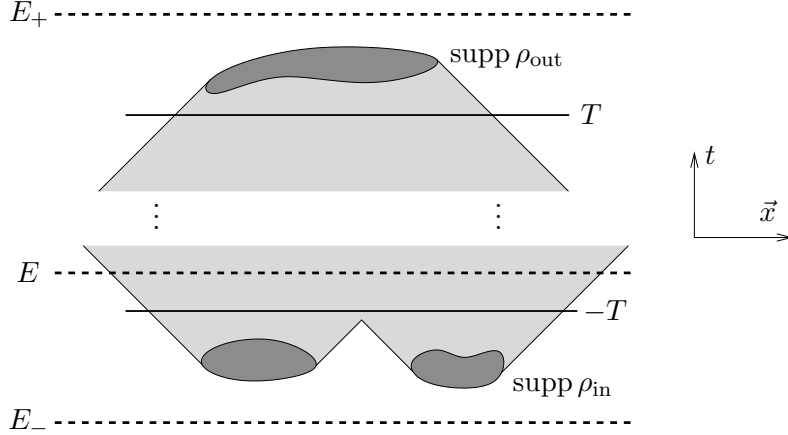


FIGURE 4. A nonlinear classical measurement process.

Due to the inhomogeneity, the classical energy (2.2) is not conserved. But it is conserved in the regions of space-time where both ρ_{in} and ρ_{out} vanish. We can thus distinguish three different energies: the energy E in the interaction region, the energy E_- in the past of ρ_{in} , and the energy E_+ in the future of ρ_{out} .

We take the point of view that the observer can only measure *energy differences*, because a constant offset to the classical energy can be interpreted as being the contribution of a background field which cannot be detected. Alternatively, one can say that a constant offset to the energy can always be removed by redefining the zero point of the energy. Moreover, as we are interested in the effect of the interaction, it is convenient to also subtract the corresponding energies of the free fields, which we denote by the additional superscript “free”. This leads us to introduce the quantity

$$\Delta E := (E - E^{\text{free}}) - (E_+ - E_+^{\text{free}}). \quad (3.26)$$

It gives the energy shift caused by the interaction as being detected by the observer at late times. We interpret ΔE as the result of our classical measurement process.

3.4. The Global Perturbation Expansion. In what follows, similar to (2.35) we use the short notations

$$S^\wedge(x, \rho) \equiv \int S^\wedge(x, y) \rho(y) d^4y, \quad S^\wedge(\rho_1, \rho_2) \equiv \iint S^\wedge(x, y) \rho_1(x) \rho_2(y) d^4x d^4y,$$

also analogously for other two-point distributions. In order to model the classical measurement process (see (3.25) and Figure 4), we choose the free solution $\phi^{(0)}$ as

$$\begin{aligned} \phi^{(0)} &= \phi_{\text{in}} + \phi_{\text{out}} \quad \text{with} \\ \phi_{\text{in}}(x) &:= \frac{1}{2} S^\wedge(x, \rho_{\text{in}}), \quad \phi_{\text{out}}(x) := \frac{1}{2} S^\vee(x, \rho_{\text{out}}). \end{aligned} \quad (3.27)$$

It is a solution of the inhomogeneous linear equation $-\square\phi_0 = (\rho_{\text{in}} + \rho_{\text{out}})/2$. In order to construct a corresponding solution of the nonlinear equation (3.25), we make a perturbation ansatz as a formal power series in λ ,

$$\phi = \sum_{n=0}^{\infty} \lambda^n \phi^{(n)}. \quad (3.28)$$

By direct calculation one verifies that the functions $\phi^{(1)}, \phi^{(2)}, \dots$ can be written iteratively as

$$\begin{aligned} \phi^{(n)}(x) &= \int S^{(n)}(x, y) \rho^{(n)}(y) d^4y \quad \text{with} \\ \rho^{(n)}(y) &:= \chi_I \sum_{a,b,c \text{ with } a+b+c=n-1} \frac{1}{6} (\phi^{(a)} \phi^{(b)} \phi^{(c)})(y), \end{aligned} \quad (3.29)$$

where $S^{(n)}$ is a general Green's function, meaning that it is a distributional solution of the equation

$$-\square S^{(n)}(x) = \delta^4(x).$$

Here χ_I is the characteristic function defined by $\chi_I(x) = 1$ if $x \in I$ and $\chi_I(x) = 0$ otherwise. It has the effect that in (3.29) we only integrate over the interaction region. This is a technical simplification which will be used in the proof of Lemma 3.3. It has no physical significance because in the scattering process under consideration, we assumed that the dynamics outside the interaction region is linear, so that cubic expressions in ϕ can be omitted.

Since we want ϕ to be a real-valued function, the Green's function should be real. This gives rise to the general ansatz

$$S^{(n)}(x, y) = S_0(x, y) + c_n P_0(x, y) + id_n K_0(x, y),$$

where S_0 is the causal Green's function

$$S_0(x, y) = \frac{1}{2} (S^\vee + S^\wedge)(x, y) = \int \frac{d^4k}{(2\pi)^4} \frac{\text{PP}}{k^2} e^{-ik(x-y)}, \quad (3.30)$$

whereas K_0, P_0 are the fundamental solutions of the free wave equation (2.15) and

$$P_0(x, y) = \int \frac{d^4k}{(2\pi)^4} \delta(k^2) e^{-ik(x-y)}, \quad (3.31)$$

and the constants c_n and d_n are free real parameters (note that P_0 is real-valued, whereas K_0 is purely imaginary). In order to specify the Green's function, we can use that our expansion should be time symmetric, meaning that all our constructions should be symmetric under the transformations

$$t \rightarrow -t \quad \text{and} \quad \phi_{\text{in}} \leftrightarrow \phi_{\text{out}}. \quad (3.32)$$

Since the distributions S_0 and P_0 are symmetric in its two arguments, whereas K_0 is anti-symmetric, this leads us to impose that $d_n = 0$. This leaves us with the relations

$$S^{(n)}(x, y) = S_0(x, y) + c_n P_0(x, y), \quad (3.33)$$

which involve one parameter c_n which can be chosen freely to every order in perturbation theory.

At this stage, there seems no reason why the parameters c_n should depend on the order n of the perturbation expansion (note that this is different for the perturbation expansion for fermions, where the concept of the Dirac sea and normalization conditions give rise to a non-trivial combinatorics of the operator products [3, 5]). Therefore, in the remainder of this paper we always make the simplest choice where all c_n vanish,

$$S^{(n)}(x, y) = S_0(x, y). \quad (3.34)$$

But more general choices of the coefficients c_n are also possible and might be worth considering in the future. The resulting perturbation expansion (3.29) and (3.34) is

neither advanced nor retarded and should thus be regarded as a global expansion in space-time. Therefore, we refer to it as the *global perturbation expansion*. It gives rise to a nonlinear mapping \mathfrak{p} which to every $\phi^{(0)} \in \mathcal{H}^{(0)}$ associates a corresponding solution ϕ of the nonlinear equation,

$$\mathfrak{p} : \mathcal{H}^{(0)} \rightarrow \mathcal{H}. \quad (3.35)$$

We remark that the Feynman propagator (2.30) can be written similar to (3.33) as

$$\Delta_F = S_0 - i\pi P_0.$$

This shows once more that the Feynman propagator is complex-valued (indeed, it is imaginary in spacelike directions). Therefore, it cannot be used for the perturbation expansion of real-valued classical fields.

3.5. A Perturbation Expansion of the Classical Energy. We now consider the classical energy (2.2) of our solution ϕ ,

$$E(t) = \int_{t=\text{const}} \left(\frac{1}{2} \dot{\phi}^2 + \frac{1}{2} |\nabla\phi|^2 + \frac{\lambda}{4!} \phi^4 \right) d^3x. \quad (3.36)$$

As in Figure 4, we denote the energy in the interaction region by E , whereas E_{\pm} are the energies for large positive and negative times, respectively.

Proposition 3.2. *The classical energies E and E_{\pm} have the expansions*

$$E_+ = \frac{i\pi}{4} \dot{K}_0(\rho_{in} + \tilde{\rho}, \rho_{in} + \tilde{\rho}) + \lim_{t \rightarrow \infty} \frac{\lambda}{4!} \int_{t=\text{const}} \phi^4 d^3x \quad (3.37)$$

$$E = \frac{i\pi}{4} \dot{K}_0(\rho_{in} - \rho_{out} + \tilde{\rho}, \rho_{in} - \rho_{out} + \tilde{\rho}) + \frac{\lambda}{4!} \int_{t=T} \phi^4 d^3x \quad (3.38)$$

$$= \frac{i\pi}{4} \dot{K}_0(\rho_{in} - \rho_{out} - \tilde{\rho}, \rho_{in} - \rho_{out} - \tilde{\rho}) + \frac{\lambda}{4!} \int_{t=-T} \phi^4 d^3x \quad (3.39)$$

$$E_- = \frac{i\pi}{4} \dot{K}_0(\rho_{out} + \tilde{\rho}, \rho_{out} + \tilde{\rho}) + \lim_{t \rightarrow -\infty} \frac{\lambda}{4!} \int_{t=\text{const}} \phi^4 d^3x, \quad (3.40)$$

where $\dot{K}_0(x, y) = \partial_{x^0} K_0(x, y)$ is the time derivative of the distribution (2.15). Here the function ϕ is defined inductively by (3.27)–(3.29), and $\tilde{\rho}$ is given by

$$\tilde{\rho} = \sum_{n=1}^{\infty} \rho^{(n)}. \quad (3.41)$$

Proof. Due to energy conservation, the energy E can be computed at any time $t \in [-T, T]$. We first compute it at time $t = T$ to derive (3.38). We again polarize the free energy to obtain the scalar product (\cdot, \cdot) introduced in (3.2). We now substitute the perturbation expansion (3.28) into (3.36). Multiplying out the free part of the energy, we need to compute expressions of the form

$$(S^{\vee}(\cdot, x), S_0(\cdot, y)), \quad (S^{\wedge}(\cdot, x), S_0(\cdot, y)), \quad (S_0(\cdot, x), S_0(\cdot, y)), \dots$$

Furthermore, we know that the argument x of the factors $S^{\wedge}(\cdot, x)$ and $S_0(\cdot, x)$ always lies to the past of time T (because $x \in I \cup \text{supp } \rho_{in}$), whereas the argument of the factors $S^{\vee}(\cdot, x)$ lies to the future of the time T (because in this case, $x \in \text{supp } \rho_{out}$). According to (2.16), (2.18), (3.30) and (2.15), we may thus exchange the Green's functions with the replacement rules

$$S^{\wedge}(\cdot, x) \rightarrow -2\pi i K_0(\cdot, x), \quad S^{\vee}(\cdot, x) \rightarrow 2\pi i K_0(\cdot, x), \quad S_0(\cdot, x) \rightarrow -i\pi K_0(\cdot, x) \quad (3.42)$$

by corresponding fundamental solutions K_0 . After these replacements, we can apply Lemma 3.3 below. In particular, we obtain for any $n, n' \geq 1$,

$$\begin{aligned} (\phi^{(n)}, \phi^{(n')}) &= (-i\pi)^2 \iint (K_0(\cdot, x), K_0(\cdot, y)) \rho^{(n)}(x) \rho^{(n')}(y) d^4x d^4y \\ &= -\pi^2 \left(-\frac{i}{4\pi} \right) \iint \dot{K}_0(x, y) \rho^{(n)}(x) \rho^{(n')}(y) d^4x d^4y = \frac{i\pi}{4} \dot{K}_0(\rho^{(n)}, \rho^{(n')}). \end{aligned}$$

Adding all the terms of the perturbation expansion gives (3.38).

To derive the expansion (3.39), we compute the energy similarly at time $t = -T$. Then the interaction region lies in the future of t , and therefore the replacement rule for S_0 in (3.42) is to be replaced by

$$S_0(\cdot, x) \rightarrow i\pi K_0(\cdot, x).$$

The expansion for E_+ is derived similar to (3.38). However, as the argument x of the factors $S^\vee(\cdot, x)$ now lies in the past of t , all the term involving a scalar product with $S^\vee(\cdot, x)$ drop out. Then energy conservation allows us to take the limit $t \rightarrow \infty$. Similarly, the expansion for E_- is obtained from (3.39) by omitting the terms involving $S^\wedge(\cdot, x)$ and taking the limit $t \rightarrow -\infty$. \square

It remains compute the energy scalar product of two factors K_0 .

Lemma 3.3. *The following relation holds in the sense of distributions,*

$$(K_0(\cdot, x), K_0(\cdot, y)) = -\frac{i}{4\pi} \dot{K}_0(y, x).$$

Proof. We write the scalar product in momentum space as

$$\begin{aligned} (K_0(\cdot, x), K_0(\cdot, y)) &= \lim_{\tau \rightarrow \infty} \frac{1}{2} \int d^3z \int \frac{d^4k}{(2\pi)^4} \epsilon(k^0) \delta(k^2) e^{-ik^0(\tau-x^0)+i\vec{k}(\vec{z}-\vec{x})} \\ &\quad \times \int \frac{d^4q}{(2\pi)^4} \epsilon(q^0) \delta(q^2) e^{-iq^0(\tau-y^0)+i\vec{q}(\vec{z}-\vec{y})} \left(-k^0 q^0 - \vec{k}\vec{q} \right). \end{aligned}$$

Carrying out the spatial integral, we only get a contribution if $\vec{q} = -\vec{k}$. Setting $\Omega = q^0$, we thus obtain

$$\begin{aligned} (K_0(\cdot, x), K_0(\cdot, y)) &= \lim_{\tau \rightarrow \infty} \frac{1}{4\pi} \int \frac{d^4k}{(2\pi)^4} \int_{-\infty}^{\infty} d\Omega \epsilon(k^0) \delta(k^2) \epsilon(\Omega) \delta(\Omega^2 - |\vec{k}|^2) \\ &\quad \times e^{-ik^0(\tau-x^0)-i\Omega(\tau-y^0)+i\vec{k}(\vec{y}-\vec{x})} \left(-k^0\Omega + |\vec{k}|^2 \right). \end{aligned}$$

Due to the two delta distributions, the integrand vanishes unless $\Omega = \pm k^0$. In the case $\Omega = +k^0$, the last factor vanishes, because

$$-k^0\Omega + |\vec{k}|^2 = (k^0)^2 - |\vec{k}|^2 = k^2 = 0.$$

Thus we may set $\Omega = -k^0$ to obtain

$$\begin{aligned} (K_0(\cdot, x), K_0(\cdot, y)) &= -\frac{1}{4\pi} \int \frac{d^4k}{(2\pi)^4} \frac{1}{2|\vec{k}|} \delta(k^2) e^{-ik^0(y^0-x^0)+i\vec{k}(\vec{y}-\vec{x})} 2|\vec{k}|^2 \\ &= -\frac{1}{4\pi} \int \frac{d^4k}{(2\pi)^4} \epsilon(\omega) \omega \delta(k^2) e^{-ik(y-x)}. \end{aligned}$$

Comparing with (2.15) gives the result. \square

We now analyze the result of Proposition 3.2. First, if we assume that the solution ϕ is dissipative and for large times goes over to a solution of the free wave equation, the last integrals in (3.37) and (3.40) vanish. This dissipation could be made precise using the techniques described in [19, 6, 17, 20]. Here we take the decay property for granted and simply drop the last integrals in (3.37) and (3.40). Similarly, if we choose the interaction time T sufficiently large, we can also drop the last integrals in (3.38) and (3.39). Then taking the difference of (3.38) and (3.39) gives the identity

$$\dot{K}_0(\rho_{\text{in}} - \rho_{\text{out}}, \tilde{\rho}) = 0. \quad (3.43)$$

Using this identity, we obtain

$$\begin{aligned} E_+ &= \frac{i\pi}{4} \dot{K}_0(\rho_{\text{in}} + \tilde{\rho}, \rho_{\text{in}} + \tilde{\rho}) \\ E &= \frac{i\pi}{4} \dot{K}_0(\rho_{\text{in}} - \rho_{\text{out}}, \rho_{\text{in}} - \rho_{\text{out}}) + \frac{i\pi}{4} \dot{K}_0(\tilde{\rho}, \tilde{\rho}) \\ E_- &= \frac{i\pi}{4} \dot{K}_0(\rho_{\text{out}} + \tilde{\rho}, \rho_{\text{out}} + \tilde{\rho}). \end{aligned}$$

Here the function $\tilde{\rho}$ describes the component of the wave generated by the interaction, whereas ρ_{in} and ρ_{out} describe the free incoming and outgoing components. In order to analyze how the energies are effected by the interaction, we subtract the free energies as given by

$$\begin{aligned} E_+^{\text{free}} &= \frac{i\pi}{4} \dot{K}_0(\rho_{\text{in}}, \rho_{\text{in}}) \\ E^{\text{free}} &= \frac{i\pi}{4} \dot{K}_0(\rho_{\text{in}} - \rho_{\text{out}}, \rho_{\text{in}} - \rho_{\text{out}}) \\ E_-^{\text{free}} &= \frac{i\pi}{4} \dot{K}_0(\rho_{\text{out}}, \rho_{\text{out}}). \end{aligned}$$

We thus obtain the compact formulas

$$\begin{aligned} E - E^{\text{free}} &= \frac{i\pi}{4} \dot{K}_0(\tilde{\rho}, \tilde{\rho}) \\ E_+ - E_+^{\text{free}} = E_- - E_-^{\text{free}} &= \frac{i\pi}{2} \dot{K}_0(\rho_{\text{out}}, \tilde{\rho}) + \frac{i\pi}{4} \dot{K}_0(\tilde{\rho}, \tilde{\rho}). \end{aligned}$$

Let us discuss these formulas. First, the formulas for the free energies can be understood immediately from the fact that for large positive (negative) times, only the wave generated by ρ_{in} (resp. ρ_{out}) is present, whereas in the interaction region, both waves are present. The interaction affects the energies by two different contributions: The term

$$\frac{i\pi}{4} \dot{K}_0(\tilde{\rho}, \tilde{\rho})$$

is delocalized in space-time in the sense that it adds to the energies E_{\pm} and E in exactly the same way. As a consequence, this contribution is not accessible to the observer, because he can only measure the energy difference $E - E_+$, but a constant offset to the energy is not detectable. The situation is different for the contribution

$$\frac{i\pi}{2} \dot{K}_0(\rho_{\text{out}}, \tilde{\rho}),$$

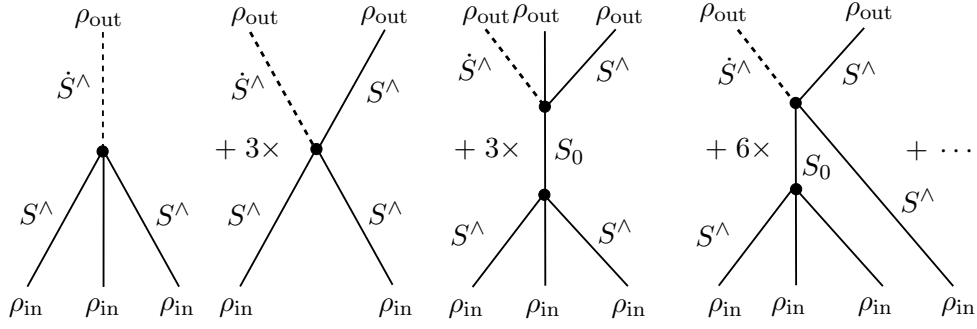


FIGURE 5. The perturbation expansion of the energy shift ΔE in the classical measurement process.

which only effects the energy in the interaction region. The observer at time T can detect this contribution by measuring the classical field energy before and after modifying the field by ρ_{out} . We now identify this contribution with (3.26) and bring it into a more convenient form.

Proposition 3.4. *In the classical measurement process described in Section 3.3, the energy shift (3.26) as measured by the observer at time T is given by*

$$\Delta E = \frac{1}{4} \dot{S}^\vee(\tilde{\rho}, \rho_{\text{out}}) = -\frac{1}{4} \dot{S}^\wedge(\tilde{\rho}, \rho_{\text{in}}). \quad (3.44)$$

Proof. Comparing the above formulas for the classical energies with (3.26), we find

$$\Delta E = \frac{i\pi}{2} \dot{K}_0(\rho_{\text{out}}, \tilde{\rho}) = \frac{i\pi}{2} \dot{K}_0(\rho_{\text{in}}, \tilde{\rho}).$$

The function $\tilde{\rho}$ is localized in the interaction region. Therefore, we can use (3.42) to replace \dot{K} by the time derivatives of the causal Green's functions. \square

We remark that (3.44) shows that our constructions are symmetric under time reversals (note that a time reversal corresponds to the replacements $\rho_{\text{out}} \leftrightarrow \rho_{\text{in}}$, $S^\vee \leftrightarrow S^\wedge$ and $\partial_t \leftrightarrow -\partial_t$).

Expressing $\tilde{\rho}$ by (3.41) and the perturbation expansion (3.27)–(3.29), we obtain an expansion of ΔE in terms of tree diagrams. The resulting diagrams are shown in Figure 5. This perturbation expansion differs from that for the solution of the Cauchy problem (see Figure 3) in several ways: First, we can now have several outgoing legs. Second, the inner lines now involve the Green's function S_0 (instead of the retarded Green's function S^\wedge); this is a direct consequence of our choice of the Green's function (3.34) in the global perturbation expansion. Finally, the time derivative in (3.44) has the effect that in each diagram, exactly one of the outgoing legs is differentiated with respect to t . This leg is denoted by a dotted line. This time derivative obviously destroys the Lorentz invariance. This can be understood from the fact that as an energy, ΔE is not a Lorentz scalar but the zero component of a Lorentz vector.

3.6. The Classical Scattering Operator and Classical n -Point Functions. Let us reconsider the previous constructions. The result of the classical measurement process was associated to a suitable difference of classical energies ΔE as given by (3.26). To compute these energies, we started from the free solution (3.27), performed the

global perturbation expansion, and could then evaluate the energy integrals (see Proposition 3.2). Describing the perturbation expansion by the nonlinear operation (3.35), we thus obtain the nonlinear mapping

$$\Delta E : \mathcal{H}_0 \times \mathcal{H}_0 \rightarrow \mathbb{R}, \quad \Delta E(\phi_{\text{out}}, \phi_{\text{in}}) := \Delta E(\mathfrak{p} \star (\phi_{\text{out}} + \phi_{\text{in}})). \quad (3.45)$$

Plugging in the perturbation expansion (3.29) and multiplying out, we get an infinite sum of multilinear mappings,

$$\Delta E(\phi_{\text{out}}, \phi_{\text{in}}) = \sum_{p,q=0}^{\infty} \Delta E_{p,q}(\underbrace{\phi_{\text{out}}, \dots, \phi_{\text{out}}}_p, \underbrace{\phi_{\text{in}}, \dots, \phi_{\text{in}}}_q). \quad (3.46)$$

At this point, it is helpful to apply the standard procedure of rewriting the multilinear operators as linear operators on the tensor product. To this end, we introduce the free p -particle classic bosonic Fock space

$$\mathcal{F}_{\text{class}}^n = \underbrace{(\mathcal{H}_0 \otimes \dots \otimes \mathcal{H}_0)}_{n \text{ factors}}_{\text{symm}},$$

where ‘‘symm’’ means total symmetrization (and \mathcal{F}^0 are simply the real numbers). On $\mathcal{F}_{\text{class}}^n$ we let $\langle \cdot | \cdot \rangle$ be the scalar product induced from the Lorentz invariant scalar product (3.7) on \mathcal{H}_0 . Taking the direct sum gives the

$$\text{classical Fock space} \quad \mathcal{F}_{\text{class}} = \bigoplus_{n=0}^{\infty} \mathcal{F}_{\text{class}}^n.$$

We remark that lifting the isometric embedding (3.9) to the tensor product yields a canonical isometric embedding of the Fock spaces,

$$(\mathcal{F}_{\text{class}}, \langle \cdot | \cdot \rangle) \hookrightarrow (\mathcal{F}, \langle \cdot | \cdot \rangle).$$

The two Fock spaces differ by the fact that $\mathcal{F}_{\text{class}}$ is a real, whereas \mathcal{F} is a complex vector space. In fact, \mathcal{F} is the complexification of the image of $\mathcal{F}_{\text{class}}$ under the above embedding.

As the multilinear mappings in (3.46) give rise to unique linear mappings on the corresponding symmetric tensor products, we obtain a unique bilinear operator

$$\Delta E : \mathcal{F}_{\text{class}} \times \mathcal{F}_{\text{class}} \rightarrow \mathbb{R}. \quad (3.47)$$

The classical scattering operator S is introduced by

$$S : \mathcal{F}_{\text{class}} \rightarrow \mathcal{F}_{\text{class}} \quad \text{with} \quad \Delta E(\Phi_{\text{out}}, \Phi_{\text{in}}) = \frac{1}{2} \langle \partial_t \Phi_{\text{out}} | S^\dagger \Phi_{\text{in}} \rangle. \quad (3.48)$$

In the next lemma we bring the classical scattering operator into a more explicit form. Our proof will also explain why the factor ∂_t in (3.48) is needed in order to make the definition Lorentz invariant.

Proposition 3.5. *The scattering operator has the matrix elements*

$$\langle \phi_{\text{out}}^1 \otimes \dots \otimes \phi_{\text{out}}^p | S^\dagger \phi_{\text{in}}^1 \otimes \dots \otimes \phi_{\text{in}}^q \rangle = S_{p+q}(\phi_{\text{out}}^1, \dots, \phi_{\text{out}}^p, \phi_{\text{in}}^1, \dots, \phi_{\text{in}}^q). \quad (3.49)$$

Here the functions S_{p+q} have a perturbation expansion in terms of Feynman tree diagrams. The order k and the total number of lines N are given by (2.36) and (2.38),

where $n = p + q$. The contributions to S_n without the outer lines are given analytically by

$$\lambda^k \int d^4 y_1 \cdots \int d^4 y_k \prod_{\ell=1}^{N-n} S_0(\cdot, \cdot) \quad (3.50)$$

(where S_0 is the Green's function (3.30)).

The functions $S_n(x_1, \dots, x_n)$ are referred to as the **classical n -point functions**.

Proof of Proposition 3.5. Combining (3.44) with (3.27), we obtain

$$\Delta E = \frac{1}{2} \int \tilde{\rho}(x) \dot{\phi}_{\text{out}}(x) d^4 x. \quad (3.51)$$

Expressing $\tilde{\rho}$ by the perturbation series (3.41) and (3.29), one gets an expansion in terms of tree diagrams. This explains (2.36) and (2.38). In order to get hold of the combinatorics, we consider the summand $\Delta E_{p,q}$ in (3.46) and number the arguments by $\phi_{\text{out}}^1, \dots, \phi_{\text{out}}^p$ and $\phi_{\text{in}}^1, \dots, \phi_{\text{in}}^q$. Moreover, we restrict attention to the contributions to $\Delta E_{p,q}$ corresponding to a fixed tree diagram, whose outer lines are labeled by the outer arguments. According to (3.51), exactly one of the outer arguments $\phi_{\text{in}}^1, \dots, \phi_{\text{in}}^p$ involves a time derivative. If we fix this dotted outer leg, the only combinatorial factor comes from the freedom to permute the three incoming legs at each vertex. This leads to a factor $3!$ at each vertex, which cancels the factor $1/6$ in (3.29). We thus obtain

$$\begin{aligned} & \Delta E_{p,q}(\phi_{\text{out}}^1, \dots, \phi_{\text{out}}^p, \phi_{\text{in}}^1, \dots, \phi_{\text{in}}^q) \\ &= \frac{1}{2} \left(S_n(\dot{\phi}_{\text{out}}^1, \dots, \phi_{\text{out}}^p, \phi_{\text{in}}^1, \dots, \phi_{\text{in}}^q) + \cdots + S_n(\phi_{\text{out}}^1, \dots, \dot{\phi}_{\text{out}}^p, \phi_{\text{in}}^1, \dots, \phi_{\text{in}}^q) \right) \\ &= \frac{1}{2} \langle \partial_t(\phi_{\text{out}}^1 \otimes \cdots \otimes \phi_{\text{out}}^p) | S^\dagger \phi_{\text{in}}^1 \otimes \cdots \otimes \phi_{\text{in}}^q \rangle, \end{aligned} \quad (3.52)$$

where S_n is formed according to (3.50), and S^\dagger is given by (3.49). Comparing with (3.48) shows that S^\dagger indeed agrees with the classical scattering operator. This concludes the proof. \square

Let us compare the classical n -point functions with the n -point functions of quantum field theory as computed in Lemma 2.1. One obvious difference is that the classical n -point functions involve *only tree diagrams*, whereas the loop diagrams of quantum field theory are missing. Before comparing the combinatorics (3.50) with (2.37), one should note that in (3.49) we inserted free fields ϕ_{in}^ℓ and ϕ_{out}^ℓ as arguments. This is why in (3.50) there are no propagators for the outer lines (this corresponds to the usual procedure of considering the “truncated” diagrams). This difference also accounts for the additional factors of i in (2.37). Namely, according to (2.36) and (2.38), for tree diagrams the identity $n = N - k - 1$ holds, and thus the factor i^{N-k} in (2.37) can be generated in the classical scattering operator if to every outer line we associate a factor of i . Then the combinatorial factors in (3.50) coincide with those in (2.37), up to an irrelevant prefactor which could be absorbed into the definition of the classical scattering operator (3.48) (the reason for the convention (3.48) is that we want the classical scattering operator to be real). We conclude that on the level of tree diagrams, the classical and the quantum scattering agree, including the combinatorics of the expansion. The only difference is that in the classical n -point functions, the inner lines are described by the Green's function S_0 , whereas in the quantum scattering process the Feynman propagator Δ_F appears. This difference can be understood from the fact

that the classical field is real, and therefore its dynamics cannot be described with the complex-valued Feynman propagator (see the end of Section 3.4).

We note that the classical scattering operator also differs from the quantum scattering operator in that it only involves the connected diagrams. As a consequence, the classical scattering operator $S : \mathcal{F}_{\text{class}} \rightarrow \mathcal{F}_{\text{class}}$ is not unitary (this is obvious already without interaction, because in the case $\lambda = 0$, the operator S is the identity on the one-particle subspace $\mathcal{F}_{\text{class}}^1$ but vanishes on the many-particle subspaces). A possible strategy to obtain a unitary classical scattering operator would be to take the sum of tensor products of S , like for example the expression

$$\sum_{n=0}^{\infty} \frac{1}{n!} \underbrace{S \otimes \cdots \otimes S}_{n \text{ factors}} .$$

This would give a unitary operator in the case $\lambda = 0$. However, it is not clear whether this operator would also be unitary in the case with interaction. If not, one could take its polar decomposition and absorb the non-unitary part into the scalar product of the classical Fock spaces. Entering the details of this analysis or discussing its potential physical significance goes beyond the scope of the present paper.

We finally remark how the identity (3.43) can be understood in terms of Feynman diagrams. In the proof of Proposition 3.5 we computed ΔE with the first equation in (3.44). The identity (3.43) makes it possible to compute ΔE alternatively by the second equation in (3.44). In analogy to (3.51) and (3.52), we thus obtain

$$\Delta E = -\frac{1}{2} \int \tilde{\rho}(x) \dot{\phi}_{\text{in}}(x) d^4x \quad (3.53)$$

$$\Delta E_{p,q} = -\frac{1}{2} \langle \phi_{\text{out}}^1 \otimes \cdots \otimes \phi_{\text{out}}^p | S^\dagger \partial_t (\phi_{\text{in}}^1 \otimes \cdots \otimes \phi_{\text{in}}^q) \rangle \quad (3.54)$$

Subtracting the last equation from (3.52), we find that (3.43) is equivalent to the identity for the S -matrix elements

$$\langle \partial_t \Phi_{\text{out}} | S^\dagger \Phi_{\text{in}} \rangle + \langle \Phi_{\text{out}} | S^\dagger \partial_t \Phi_{\text{in}} \rangle = 0 .$$

This means that the value of the Feynman diagram remains unchanged if both the incoming and the outgoing waves are translated in time. It corresponds to the usual concept of *conservation of quantum mechanical energy* in a Feynman diagram, stating that the incoming energy, defined as the sum of the frequencies of the incoming waves, coincides with the outgoing energy.

4. NONLINEAR CLASSICAL SCATTERING IN A STOCHASTIC BACKGROUND FIELD

4.1. A Free Stochastic Field. We begin by analyzing a free stochastic field ξ . By “free” we mean that $\xi \in \mathcal{H}_0$ is a solution of the linear wave equation. Similar to (3.3), we can characterize ξ by its initial values for example some fixed time t_0 ,

$$\Xi := (\xi, \dot{\xi})|_{t_0} \in C_0^\infty(\mathbb{R}^3) \times C_0^\infty(\mathbb{R}^3) .$$

“Stochastic” means that we prescribe a probability distribution for ξ which we assume to be *Gaussian*. We write the corresponding probability measure $\mathfrak{D}\Xi$ on the initial data as

$$\mathfrak{D}\Xi = \frac{1}{Z} e^{-\frac{1}{2} F(\Xi, \Xi)} d\xi \wedge d\dot{\xi} , \quad (4.1)$$

where Z is a normalization constant, and F is a positive bilinear form, i.e.

$$F(\Xi, \Xi) = \int \langle \Xi(\vec{x}), F \Xi(\vec{x}) \rangle_{\mathbb{R}^2} d^3x$$

with F being a linear operator on the initial data. Next, it is natural to assume that our probability measure is *translation invariant*, meaning that the operator F is diagonal in momentum space. Thus setting

$$\Xi(\vec{x}) = \int \frac{d^3k}{(2\pi)^3} \Xi(\vec{k}) e^{i\vec{k}\vec{x}}, \quad (4.2)$$

the operator F can be written as

$$\widehat{(F\Xi)}(\vec{k}) = F(\vec{k}) \Xi(\vec{k}),$$

where $F(\vec{k})$ is a 2×2 -matrix. The fact that Ξ is real-valued and that F acts on a space of real-valued functions implies that

$$\Xi(-\vec{k}) = \overline{\Xi(\vec{k})} \quad \text{and} \quad F(-\vec{k}) = \overline{F(\vec{k})}, \quad (4.3)$$

where the bar denotes complex conjugation of each vector or matrix component. Using Plancherel's theorem, we can write the bilinear form F as

$$F(\Xi, \Xi) = \int \frac{d^3k}{(2\pi)^3} \langle \Xi(\vec{k}), F(\vec{k}) \Xi(\vec{k}) \rangle_{\mathbb{C}^2}. \quad (4.4)$$

A Gaussian probability measure is completely characterized by the *covariance* \mathcal{C} defined as the expectation value of a quadratic polynomial, i.e. in our setting

$$\mathcal{C}(x, y) := \int \xi(x) \xi(y) \mathfrak{D}\Xi. \quad (4.5)$$

The expectation values of higher order polynomials can then be computed using Wick's theorem (similar as outlined in Section 2.2 in the context of path integrals or see for example [7]).

Lemma 4.1. *For any space-time points $x, y \in I$, the covariance (4.5) is given by*

$$\begin{aligned} \mathcal{C}(x, y) &= \int \frac{d^4k}{(2\pi)^4} \delta(k^2) \frac{\pi}{|\omega|} e^{-i\vec{k}(\vec{x}-\vec{y})} \\ &\times \left[e^{i\omega(x^0-y^0)} \left\langle \begin{pmatrix} -i\omega \\ 1 \end{pmatrix}, F(\vec{k})^{-1} \begin{pmatrix} -i\omega \\ 1 \end{pmatrix} \right\rangle_{\mathbb{C}^2} \right. \\ &\quad \left. - e^{i\omega(x^0+y^0+2t)} \left\langle \begin{pmatrix} -i\omega \\ 1 \end{pmatrix}, F(\vec{k})^{-1} \begin{pmatrix} i\omega \\ 1 \end{pmatrix} \right\rangle_{\mathbb{C}^2} \right]. \end{aligned} \quad (4.6)$$

$$- e^{i\omega(x^0+y^0+2t)} \left\langle \begin{pmatrix} -i\omega \\ 1 \end{pmatrix}, F(\vec{k})^{-1} \begin{pmatrix} i\omega \\ 1 \end{pmatrix} \right\rangle_{\mathbb{C}^2}. \quad (4.7)$$

Proof. In momentum space, the Gaussian measure (4.1) becomes

$$\mathfrak{D}\Xi = \frac{1}{Z} \exp \left(-\frac{1}{2} \int \frac{d^3k}{(2\pi)^3} \langle \Xi(\vec{k}), F(\vec{k}) \Xi(\vec{k}) \rangle_{\mathbb{C}^2} \right) d\Xi_1(\vec{k}) \wedge d\Xi_2(\vec{k}). \quad (4.8)$$

By the standard calculation rules for Gaussian measures (see for example [7]), we obtain

$$\int \Xi(\vec{k}) \otimes \Xi(\vec{k}')^\dagger \mathfrak{D}\Xi = (2\pi)^3 \delta^3(\vec{k} - \vec{k}') F(\vec{k})^{-1}. \quad (4.9)$$

Next, we know from (3.13) that

$$\xi(x) = \int \left(-\dot{S}_0(x, (t, \vec{y})) \quad -S_0(x, (t, \vec{y})) \right) \Xi(\vec{y}) d^3y$$

and thus

$$\xi(x) \xi(y) = \int_{\mathbb{R}^3 \times \mathbb{R}^3} d^3 z_1 d^3 z_2 \left\langle \begin{pmatrix} -\dot{S}^0(x, (t, \vec{z}_1)) \\ -S^0(x, (t, \vec{z}_1)) \end{pmatrix}, \Xi(\vec{z}_1) \otimes \Xi(\vec{z}_2)^\dagger \begin{pmatrix} -\dot{S}^0(y, (t, \vec{z}_2)) \\ -S^0(y, (t, \vec{z}_2)) \end{pmatrix} \right\rangle_{\mathbb{C}^2}$$

Similar to (3.42), we may apply the replacement rule

$$S_0^\wedge(x, y) \rightarrow -2\pi i K_0(x, y).$$

As in Lemma 3.3, for the momenta we use the notation $k = (\omega, \vec{k})$ and $q = (\Omega, \vec{q})$. We then obtain

$$\begin{aligned} \xi(x) \xi(y) &= 4\pi^2 \int \frac{d^4 k}{(2\pi)^4} e^{ikx} \epsilon(\omega) \delta(k^2) \int \frac{d^4 q}{(2\pi)^4} e^{-iqy} \epsilon(\Omega) \delta(q^2) \\ &\quad \times e^{i(\omega-\Omega)t} \left\langle \begin{pmatrix} i\omega \\ -1 \end{pmatrix}, \Xi(\vec{k}) \otimes \Xi(\vec{q})^\dagger \begin{pmatrix} i\Omega \\ -1 \end{pmatrix} \right\rangle_{\mathbb{C}^2}. \end{aligned}$$

Now we can carry out the Gaussian integral with (4.9),

$$\begin{aligned} \int \xi(x) \xi(y) \mathfrak{D}\Xi &= 4\pi^2 \int \frac{d^4 k}{(2\pi)^4} e^{ikx} \epsilon(\omega) \delta(k^2) \int \frac{d\Omega}{2\pi} e^{-i\Omega y^0 + i\vec{k}\vec{y}} \epsilon(\Omega) \delta(\Omega - |\vec{k}|^2) \\ &\quad \times e^{i(\omega-\Omega)t} \left\langle \begin{pmatrix} i\omega \\ -1 \end{pmatrix}, F(\vec{k})^{-1} \begin{pmatrix} i\Omega \\ -1 \end{pmatrix} \right\rangle_{\mathbb{C}^2} \\ &= \pi \int \frac{d^4 k}{(2\pi)^4} e^{ikx} \epsilon(\omega) \delta(k^2) \\ &\quad \times \frac{1}{\Omega} e^{-i\Omega y^0 + i\vec{k}\vec{y}} e^{i(\omega-\Omega)t} \left\langle \begin{pmatrix} i\omega \\ -1 \end{pmatrix}, F(\vec{k})^{-1} \begin{pmatrix} i\Omega \\ -1 \end{pmatrix} \right\rangle_{\mathbb{C}^2} \Big|_{\Omega=\pm\omega}. \end{aligned}$$

Rearranging this expression gives the result. \square

Let us discuss the result of this lemma. We first point out that the covariance is obviously real-valued and symmetric under permutations and x and y . Moreover, it satisfies the free wave equation,

$$\mathcal{C}(x, y) = \mathcal{C}(y, x) \quad \text{and} \quad \square_x \mathcal{C}(x, y) = 0.$$

The last equation is obvious from (4.5) because the factors $\xi(x)$ and $\xi(y)$ are solutions of the wave equation. Alternatively, it can also be verified in the expression of Lemma 4.1 by applying the wave operator and using the factor $\delta(k^2)$ in the integrand on the right. In particular, it is *impossible to arrange* that $\mathcal{C}(x, y)$ is a *Green's function*.

In order to obtain a covariant theory, we would like the covariance \mathcal{C} to be Lorentz invariant. This raises the questions for which choices of F the covariance is Lorentz invariant, and which Lorentz invariant distributions \mathcal{C} can be realized. We first explain what the naive answer to these questions is, and then prove that this naive guess is indeed correct. The general Lorentz invariant bi-solution of the scalar wave equation is a linear combination of the distributions P_0 and K_0 introduced in (3.31) and (2.15). As $K_0(x, y)$ is anti-symmetric under permutations of x and y and $P_0(x, y)$ is real-valued, we conclude that \mathcal{C} should be a real multiple of $P_0(x, y)$.

Proposition 4.2. *Assume that the covariance (4.5) of the Gaussian measure (4.1) is Lorentz invariant. Then there is a parameter $\beta > 0$ such that*

$$\mathcal{C}(x, y) = \frac{1}{\beta} P_0(x, y) \quad (4.10)$$

$$F(\Xi, \Xi) = \frac{\beta}{\pi} \langle \xi | \xi \rangle, \quad (4.11)$$

where $\langle \xi | \xi \rangle$ denotes the scalar product (3.7).

Proof. Lorentz invariance clearly implies that $\mathcal{C}(x, y)$ must be rotationally symmetric, meaning that $F(\vec{k})$ must depend only on $|\vec{k}|$. Combining this fact with the second equation in (4.3), we conclude that $F(\vec{k})$ must have real matrix entries. Thus its inverse can be written as

$$F(\vec{k})^{-1} = \begin{pmatrix} a & b \\ b & c \end{pmatrix}. \quad (4.12)$$

with real functions a, b, c depending on $|\vec{k}|$. Next, Lorentz invariance implies that $\mathcal{C}(x, y)$ should be independent of F , meaning that the contribution (4.7) must vanish, i.e.

$$\left\langle \begin{pmatrix} -i\omega \\ 1 \end{pmatrix}, F(\vec{k})^{-1} \begin{pmatrix} i\omega \\ 1 \end{pmatrix} \right\rangle_{\mathbb{C}^2} = 0 \quad \text{for } \omega = \pm |\vec{k}|.$$

Combining this condition with (4.12) we conclude that $b = 0$ and that $c = \omega^2 a$. Then the expectation value in (4.6) becomes

$$\left\langle \begin{pmatrix} -i\omega \\ 1 \end{pmatrix}, F(\vec{k})^{-1} \begin{pmatrix} -i\omega \\ 1 \end{pmatrix} \right\rangle_{\mathbb{C}^2} = 2\omega^2 a(|\vec{k}|).$$

Substituting these results into the formula of Lemma 4.1, we obtain

$$\mathcal{C}(x, y) = \int \frac{d^4 k}{(2\pi)^3} \delta(k^2) e^{ik(x-y)} |\omega| a(|\vec{k}|).$$

This distribution is Lorentz covariant only if $a(|\vec{k}|) = (2\pi\beta |\vec{k}|)^{-1}$ for a suitable real parameter β . This gives (4.10). Moreover,

$$F(\vec{k})^{-1} = \frac{1}{2\pi\beta |\vec{k}|} \begin{pmatrix} 1 & 0 \\ 0 & \omega^2 \end{pmatrix} \quad \text{and thus} \quad F(\vec{k}) = 2\pi\beta \begin{pmatrix} |\vec{k}| & 0 \\ 0 & 1/|\vec{k}| \end{pmatrix}. \quad (4.13)$$

Next, we write ξ similar to (3.4) as a Fourier integral over the upper and lower mass cone, i.e.

$$\xi(x) = \frac{1}{(2\pi)^4} \int \frac{d^3 k}{2\omega} \left(\xi(\vec{k}) e^{-i\omega t + i\vec{k}\vec{x}} + \overline{\xi(\vec{k})} e^{i\omega t - i\vec{k}\vec{x}} \right).$$

Taking the time derivative, evaluating for simplicity at $t = 0$ and comparing with (4.2), we find that

$$\Xi_1(\vec{k}) = \frac{1}{4\pi\omega} \left(\xi(\vec{k}) + \overline{\xi(-\vec{k})} \right) \quad \text{and} \quad \Xi_2(\vec{k}) = -\frac{i}{4\pi} \left(\xi(\vec{k}) - \overline{\xi(-\vec{k})} \right).$$

Hence the functional F as given by (4.4) and (4.13) becomes

$$F(\Xi, \Xi) = \frac{\beta}{2\pi} \int \frac{d^3 k}{(2\pi)^3} \frac{1}{\omega} |\xi(\vec{k})|^2. \quad (4.14)$$

Comparing with (3.7) gives the result. \square

Recall from (3.8) that the scalar product $\langle \xi | \xi \rangle$ is dimensionless. Moreover, the functional F , being the argument of the exponential in (4.1), must clearly be dimensionless. Hence the parameter β in (4.11) is also dimensionless,

$$[\beta] = \ell^0. \quad (4.15)$$

This parameter describes the strength of stochastic background field. In the limit $\beta \rightarrow \infty$, the covariance (4.10) vanishes, so that no stochastic field is present. If β is decreased, the amplitude of the stochastic field gets larger.

4.2. The Perturbation Expansion and Classical Scattering. We now want to include the stochastic background field in the interaction. To this end, we first add the free stochastic field ξ to the free in- and outgoing fields in (3.27),

$$\phi_0 = \phi_{\text{in}} + \phi_{\text{out}} + \xi. \quad (4.16)$$

The corresponding interacting field is then introduced exactly as in Section 3.4 by the perturbation series (3.28) and (3.29). Expanding the classical energy exactly as in Section 3.5, we obtain in analogy to (3.45) the nonlinear mapping

$$\Delta E : \mathcal{H}_0 \times \mathcal{H}_0 \times \mathcal{H}_0 \rightarrow \mathbb{R}, \quad \Delta E(\phi_{\text{out}}, \phi_{\text{in}}, \xi) := \Delta E(\mathfrak{p} \star (\phi_{\text{out}} + \phi_{\text{in}} + \xi)).$$

After expanding similar to (3.46) in a series of multilinear mappings,

$$\Delta E = \sum_{p,q,r=0}^{\infty} \Delta E_{p,q,r}(\underbrace{\phi_{\text{out}}, \dots, \phi_{\text{out}}}_{p \text{ arguments}}, \underbrace{\phi_{\text{in}}, \dots, \phi_{\text{in}}}_{q \text{ arguments}}, \underbrace{\xi, \dots, \xi}_{r \text{ arguments}}), \quad (4.17)$$

we carry out the Gaussian integral with the probability distribution (4.1),

$$\begin{aligned} \Delta E(\phi_{\text{out}}, \phi_{\text{in}}) &= \sum_{p,q=0}^{\infty} \Delta E_{p,q}(\phi_{\text{out}}^p, \phi_{\text{in}}^q) \quad \text{with} \\ \Delta E_{p,q}(\phi_{\text{out}}^p, \phi_{\text{in}}^q) &:= \sum_{r=0}^{\infty} \int \Delta E_{p,q,r}(\phi_{\text{out}}^p, \phi_{\text{in}}^q, \xi^r) d\Xi. \end{aligned}$$

Now the classical scattering operator is introduced similar as in Section 3.6 by (3.47) and (3.48).

We now outline a general procedure to express the classical scattering operator in terms of Feynman diagrams. In order to get hold on the combinatorics, it is preferable to first derive the expansion for fixed ξ , and to integrate over Ξ afterwards. Similar to (3.47), we rewrite the multilinear operators in (4.17) as linear operators on classical Fock spaces,

$$\begin{aligned} \Delta E : \mathcal{F}_{\text{class}} \times \mathcal{F}_{\text{class}} \times \mathcal{F}_{\text{class}} &\rightarrow \mathbb{R}, \\ \Delta E(\underbrace{\phi_{\text{out}} \otimes \dots \otimes \phi_{\text{out}}}_{p \text{ factors}}, \underbrace{\phi_{\text{in}} \otimes \dots \otimes \phi_{\text{in}}}_{q \text{ factors}}, \underbrace{\xi \otimes \dots \otimes \xi}_{r \text{ factors}}) &= \Delta E_{p,q,r}(\phi_{\text{out}}^p, \phi_{\text{in}}^q, \xi^r). \end{aligned}$$

For a fixed stochastic field $\Xi_{\text{stoch}} \in \mathcal{F}$, we now introduce the scattering operator $S[\Xi_{\text{stoch}}]$ in analogy to (3.48) by

$$S[\Xi_{\text{stoch}}] : \mathcal{F}_{\text{class}} \rightarrow \mathcal{F}_{\text{class}} \quad \text{with} \quad \Delta E(\Phi_{\text{out}}, \Phi_{\text{in}}, \Xi_{\text{stoch}}) = \frac{1}{2} \langle \partial_t \Phi_{\text{out}} | S[\Xi_{\text{stoch}}]^\dagger \Phi_{\text{in}} \rangle.$$

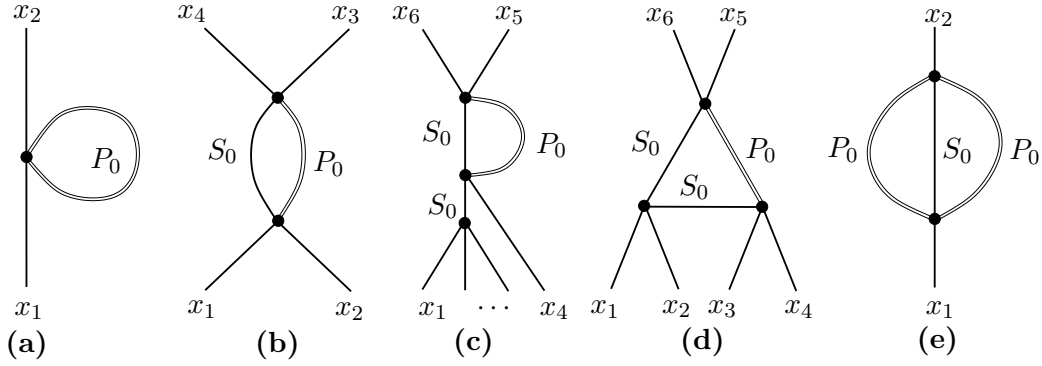


FIGURE 6. Examples of loop diagrams for the classical measurement process in a stochastic background field.

We next consider the matrix elements of $S[\Xi_{\text{stoch}}]$ for product states. Following the procedure in Proposition 3.5, we obtain

$$\begin{aligned} & \langle \phi_{\text{out}}^1 \otimes \cdots \otimes \phi_{\text{out}}^p \mid S[\xi_1 \otimes \cdots \otimes \xi_r]^\dagger \phi_{\text{in}}^1 \otimes \cdots \otimes \phi_{\text{in}}^q \rangle \\ & = S_{p+q+r}(\phi_{\text{out}}^1, \dots, \phi_{\text{out}}^p, \phi_{\text{in}}^1, \dots, \phi_{\text{in}}^q, \xi_1, \dots, \xi_r), \end{aligned}$$

where S_{p+q+r} are again the classical n -point functions. We thus obtain an expansion in terms of tree diagrams, but now with additional external lines labeled by ξ_1, \dots, ξ_r . The analytic expressions for the Feynman diagrams including the combinatorial factors are again given by (3.50).

In order to carry out the integral over Ξ , we first replace the factors ξ_1, \dots, ξ_r by our stochastic field ξ to obtain an r -multilinear function in ξ . Then the Ξ -integration gives a Gaussian integral, which can be carried out with the Wick rules (see (2.33)). We thus obtain the sum over all pairings between the factors ξ , where every pair $\xi(x)\xi(y)$ is replaced by the covariance $\mathcal{C}(x, y)$ (see (4.5)). In view of (4.10), every pairing generates an additional line described by the distribution $P_0(x, y)/\beta$ (where P_0 is the fundamental solution (3.31)). We thus obtain loop diagrams as shown in Figure 6.

4.3. Comparison of Classical and Quantum Loop Diagrams. Let us compare this expansion to the perturbation expansion of quantum field theory as shown in Figure 1. First of all, the expansions are very similar in that they both involve tree as well as loop diagrams. The only difference between the diagrams is that the lines of the quantum diagrams are formed of the Feynman propagator, whereas the lines in the classical diagrams are formed of either the Green's function S_0 or the fundamental solution P_0 , with a specific combinatorics.

Let us explain the significance of the parameter β . We already noted in (4.15) that it is a dimensionless parameter. In the limit $\beta \rightarrow \infty$, the covariance and thus the stochastic background field vanish. This corresponds to the classical limit where no loop diagrams are present. As a diagram with l loops involves a factor β^{-l} , by decreasing β we can make the loop diagrams larger, giving a more “quantum-like behavior.” Thus the parameter β^{-1} plays a similar role as Planck's constant \hbar in quantum theory. In order to make this connection precise, we again need to consider the more general units (2.39) where $\hbar \neq 1$. Comparing (3.2) with (3.7), it follows that the scalar product $\langle \xi | \xi \rangle$ has dimensions $m\ell$. Thus in order for the functional F

in (4.11) to be dimensionless, we need to choose

$$\beta = \frac{\text{const}}{\hbar}$$

with a numerical constant. In this way, that the free parameter β^{-1} which describes the strength of the stochastic background field can indeed be identified with the parameter \hbar in quantum theory.

From the analytic point of view, replacing the Feynman propagator Δ_F by S_0 or P_0 has no influence on the qualitative behavior of the loops. In particular, the classical loop diagrams have the same divergent behavior as the corresponding quantum diagrams. The following lemma even gives a quantitative connection between the classical and the quantum loops.

Lemma 4.3. *Assume that the momenta k_1, \dots, k_r with $r \geq 1$ are all different. Then for almost all momenta \vec{q} ,*

$$\int_{-\infty}^{\infty} dq^0 \Delta_F(k_1 + q) \cdots \Delta_F(k_r + q) = -i\pi \int_{-\infty}^{\infty} dq^0 \sum_{a=1}^r P_0(k_a + q) \prod_{b \neq a} S_0(k_b + q).$$

Proof. We set $\omega = q^0$, $\omega_a = k_a^0$ and $p_a = |\vec{k}_a + \vec{q}| > 0$. Then, according to (2.30),

$$\Delta_F(k_a + q) = \lim_{\varepsilon \searrow 0} \frac{1}{(\omega + \omega_a)^2 - p_a^2 + i\varepsilon} \quad (4.18)$$

$$= \lim_{\varepsilon \searrow 0} \frac{1}{2p_a} \left(\frac{1}{(\omega + \omega_a) - p_a + i\varepsilon} - \frac{1}{(\omega + \omega_a) + p_a - i\varepsilon} \right). \quad (4.19)$$

We take the product of r such factors for fixed $\varepsilon > 0$. Since the momenta k_1, \dots, k_r are pairwise different, the resulting function has simple poles for almost all \vec{q} . In the case $r \geq 2$, we can close the contour either in the upper or in the lower half plane. Computing the mean value of these two contour integrals and taking the limit $\varepsilon \searrow 0$, we obtain

$$\begin{aligned} \int_{-\infty}^{\infty} dq^0 \Delta_F(k_1 + q) \cdots \Delta_F(k_r + q) &= -i\pi \sum_{a=1}^r \frac{1}{2p_a} \prod_{b \neq a} S_0(k_b + q) \Big|_{\omega = -\omega_a \pm p_a} \\ &= -i\pi \int_{-\infty}^{\infty} d\omega \sum_{a=1}^r \frac{1}{2p_a} \left(\delta(\omega + \omega_a + p_a) + \delta(\omega + \omega_a - p_a) \right) \prod_{b \neq a} S_0(k_b + q) \\ &= -i\pi \int_{-\infty}^{\infty} d\omega \sum_{a=1}^r \delta((k_a + q)^2) \prod_{b \neq a} S_0(k_b + q). \end{aligned}$$

Using (3.31) gives the result. In the case $r = 1$, the two summands in (4.19) do not decay fast enough for closing the contours. However, one can compute the contour integrals using (4.18), to again obtain the result. \square

This lemma shows that in a closed loop, a product of r Feynman propagators may be replaced by the product of one factor P_0 and $r - 1$ factors S_0 , if we take the sum over all such combinations. The condition that the momenta k_1, \dots, k_r should all be different can be justified as follows: When computing scattering amplitudes, the momenta k_1, \dots, k_r are expressed in terms of the incoming and outgoing momenta. Then the momenta k_a are all different except at the poles of the scattering amplitudes. Such poles correspond to resonances; they need to be described non-perturbatively

with a resummation technique. If the loop under consideration is contained inside another loop, one also needs to integrate over one or several of the momenta k_a . In this case, the poles of the Feynman diagrams if some of the momenta k_1, \dots, k_r coincide are of no relevance, as only the divergences for large momenta are of interest.

Thus, in order to decide whether the classical and the quantum loop diagrams agree quantitatively, it remains to study the combinatorics. Let us begin with the tadpole diagram (see Figure 1 **(b)**). It arises when integrating the simple tree diagram (see the left of Figure 5) over Ξ , if two of the incoming fields are the stochastic field ξ . As there are three possibilities for the choice of the two ξ -legs, and taking into account the factor $1/6$ in (3.29), we get a total combinatorial factor of $1/2$. Hence Lemma 4.3 allows us to replace the factor P_0 in the classical tadpole by

$$P_0 \rightarrow -\frac{1}{2\pi i} \Delta_F.$$

As the factor $(-i)$ is already taken care of in (2.37), we get complete agreement between the quantum and the classical tadpole if we choose (see (4.5) and (4.10))

$$\beta = \frac{1}{2\pi}. \quad (4.20)$$

We next consider the simple loop diagram in Figure 1 **(d)**. In the classical setting, this diagram arises by performing the stochastic integral of a second order tree diagram with two ξ -legs (see Figure 6 **(b)**). As the tree diagram has a combinatorial factor one (see Proposition 3.5), the inner lines are formed by the product $\lambda^2 S_0 P_0$. Due to the symmetry in the momenta of the two inner lines, Lemma 4.3 allows us to replace the inner lines by $\Delta_F \Delta_F / (-2\pi i)$. Hence we again get complete agreement between the quantum and the classical diagram if we choose β according to (4.20). This argument applies just as well to simple loops inside larger diagrams, as is exemplified in Figure 6 **(c)**.

The situation changes with longer loops, as we now discuss in the example of the triangle loop in Figure 6 **(d)**. The shown diagram is obtained in a unique way as the stochastic integral of a third order tree diagram with two ξ -legs. The combinatorial factor is again one. Similarly, one gets the two loop diagrams where the line P_0 is one of the other sides of the triangle. Thus we can again apply Lemma 4.3. In order to obtain complete agreement to the corresponding quantum diagram we need to choose

$$\beta = \frac{1}{\pi}. \quad (4.21)$$

We point out that this differs from (4.20) by a factor of two. This shows that the expansion in terms of classical loop diagrams is not mathematically equivalent to quantum field theory. One difference is that longer loops have a weight which is double that of a tadpole or a simple loop.

For higher loops (see Figure 6 **(e)** for the simplest example), one can apply inductively Lemma 4.3 (or a generalization where on the left factors Δ_F are replaced by S_0) to get a connection between the classical and the corresponding quantum diagrams. However, the combinatorial details of this analysis go beyond the scope of the present paper.

5. OUTLOOK

The previous analysis shows that by considering the classical ϕ^4 -theory in a stochastic background field, one gets a perturbative description which involves all the

tree and loop diagrams of the corresponding quantum field theory. The main result of the paper is to demonstrate that loop diagrams appear in a purely classical context. By adapting our methods, this result immediately carries over to any other bosonic field theory in any space-time dimension.

We also observed that for ϕ^4 -theory, the combinatorics of the tadpole and of the simple loops differs from that of the longer loops (see Section 4.3). This result is quite sensitive on the theory under consideration. It might well be that for other theories, these combinatorial differences between the classical and the quantum loop diagrams disappear. In particular, it was crucial for our combinatorial considerations that there are distinct lines which begin at the same vertex i and both end at the same vertex j . As such lines do not exist in QED, is conceivable that for QED there might be full agreement between classical and the quantum loop diagrams.

In any case, working out the combinatorial details for a realistic model to higher order seems an important project for the future. Here we only make three remarks:

- (a) We note that for complex fields (like a complex Klein-Gordon field), the charge is an additional conserved quantity. This makes it possible to work in the classical measurement process instead of the energy with the charge. An advantage of this procedure is that the charge is a Lorentz scalar, so that the time derivative in (3.48) could be omitted. The reason why we preferred to work with a real field is that the components of the electromagnetic field are also real-valued.
- (b) We remark that in order to extend our methods, one could replace the Gaussian measure (4.1) by a more general probability distribution on the classical solutions. This would change the detailed form of the higher order loop diagrams. The non-Gaussian probability distribution would have to be justified as arising from the nonlinear classical dynamics prior to the scattering process.
- (c) We point out that all our methods only work for *bosonic* fields. The reason is that the Pauli exclusion principle and pair creation seem to make it impossible to interpret the Dirac equation as a classical field equation. Thus in order to extend our methods to models involving fermions, one possibility is to couple a classical bosonic field to a second-quantized fermion field. An alternative is to describe the fermions in the framework of the fermionic projector [4].

We finally outline how our constructions could be helpful for the rigorous construction of interacting quantum fields. To this end, let us assume that there is a physical theory for which there is complete agreement between the classical and the quantum Feynman diagrams. Then one could prove the existence of the quantum theory by rigorously constructing the classical theory in the presence of a stochastic background field. More precisely, one could first introduce an ultraviolet regularization by replacing the covariance of the Gaussian measure (4.10) by a regularized covariance P^ε/β (for example with a cutoff in momentum space; clearly, the regularization would violate Lorentz invariance). Then the initial values $\phi_0|_{t=-T}$ (with ϕ_0 according to (4.16)) would be in a suitable Sobolev space. If one could prove global existence of solutions uniformly in ξ , integrating the solutions over Ξ would give all interacting n -point functions. If one also succeeded in controlling the behavior of the classical solutions in the limit when the regularization is removed (this would probably require some kind of “non-perturbative renormalization method”), one would have proved non-perturbatively that the quantum field theory in the Lorentzian setting is well-defined and finite. In this way, it seems possible at least in principle to make the

powerful methods of nonlinear partial differential equations applicable to one of the most outstanding problems of quantum field theory.

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