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in Renormalized Perturbation Theory

by



R. Flume

II. Institut für Theoretische Physik, Universität Hamburg

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The Invariance of the S-Matrix under Point Transformations
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R. Flume

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Abstract

We give a simple proof of the invariance of the S-matrix under point transformations of the fields in renormalized perturbation field theory.

I. Introduction

The equivalence theorem of Lagrangian field theory can be stated in the following form: The quantization of two classical Lagrange densities $\mathcal{L}(\phi, \partial_\mu \phi)$ and $\mathcal{L}'(\phi, \partial_\mu \phi)$ related to each other through a point transformation of the fields

$$\phi \rightarrow (\phi + h(\phi)), \quad (h(0) = 0, \quad \frac{dh}{d\phi} = 0)$$

$$\mathcal{L}'(\phi, \partial_\mu \phi) = \mathcal{L}(\phi + h(\phi), \partial_\mu(\phi + h(\phi)))$$

gives rise to the same S-matrix.

There exist several formal proofs for this theorem in the literature^{1,2}. Recently Lam³ proposed a constructive and rigorous proof in renormalized perturbation theory. His method of proof is based on the technique of anisotropically quantized normal products⁴. The purpose of this note is to present a proof of the equivalence theorem, which subsists on a minimum of technicalities and, in particular, avoids the rather involved normal product formalism. Furthermore, in contrast to Lam's approach, we don't need to refer in our proof to the Haag-Ruelle (L.S.Z)-theorem.

II.

We confine ourselves to a Lagrangian $\mathcal{L}(\phi, \partial^\mu \phi)$ which is a function of a single scalar massive field (of mass \bar{m}). We exclude higher than first space-time derivatives. The proof presented below applies modulo a greater amount of book-keeping also to Lagrangians involving several fields of possibly different type. To start with, we construct for a given not necessarily renormalizable Lagrangian $\mathcal{L}(\phi, \partial_\mu \phi)^*$ the S-matrix

$$S(\mathcal{L}) = \lim_{g \rightarrow \text{konst.}} \sum_{n=0}^{\infty} \frac{(-i)^n}{n!} \int T(\bar{\mathcal{L}}(x_1) \dots \bar{\mathcal{L}}(x_n)) g(x_1) \dots g(x_n) dx_1 \dots dx_n, \quad (1)$$

$$\mathcal{L} = \mathcal{L}_{\text{free}} + \mathcal{V},$$

by specifying through some ad hoc prescription, which may or may not correspond to a minimal subtraction scheme, the time ordered operator valued distributions $T(\bar{\mathcal{L}}(x_1) \dots \bar{\mathcal{L}}(x_n))$. One knows from the work of Epstein and Glaser^{5,6} that the adiabatic limit $g \rightarrow \text{konst}$ in (1) exists** and that the theory fulfils in the adiabatic limit all requirements of locality and unitarity in the sense of formal power series.

$\mathcal{L}(\phi, \partial_\mu \phi)$ goes under a point transformation $\phi \rightarrow (\phi + h(\phi))$ over into $\mathcal{L}_h = \mathcal{L}(\phi + h(\phi), \partial_\mu(\phi + h(\phi)))$. h has to be considered as a formal power series in ϕ ($h(0) = 0, dh/d\phi = 0$). We interpolate \mathcal{L} and \mathcal{L}_h by

$$\mathcal{L}_\lambda(\phi) = \mathcal{L}(\phi + \lambda h, \partial_\mu(\phi + \lambda h)) \quad 0 \leq \lambda \leq 1.$$

Specifying the time ordered products $T_\lambda(x_1 \dots x_n) = T(\bar{\mathcal{L}}_\lambda(x) \dots \bar{\mathcal{L}}_\lambda(x))$ we obtain a one parameter family S_λ of S-matrices. We will define the T_λ in such a way that $dS_\lambda/d\lambda$ vanishes and S_0 is given by equ. (1). To make the motivation of our procedure clearer we give first a naive argument, which afterwards will be made rigorous. The argument is taken over with slight modifications from Divakaran².

$dS_\lambda/d\lambda$ is entirely determined by the derivatives with respect to λ of the time ordered products T_λ

* $\mathcal{L}(\phi, \partial_\mu \phi)$ is supposed to have a formal power series expansion in $\phi, \partial_\mu \phi$ around $\phi = \partial_\mu \phi = 0$.

** After correct mass and wave function renormalization.

$$\frac{d}{d\lambda} T_\lambda(x_1 \dots x_n) = \sum_i T(\bar{\mathcal{L}}_\lambda(x_1) \dots \bar{\mathcal{L}}_\lambda(x_{i-1}) \cdot \frac{d}{d\lambda} \mathcal{L}_\lambda(x_i) \dots \mathcal{L}_\lambda(x_n)) .$$

We can manipulate $d\mathcal{L}_\lambda/d\lambda$ as follows:

$$\frac{d\mathcal{L}_\lambda}{d\lambda} = \frac{d}{d\lambda} \mathcal{L}(\phi + \lambda h, \partial_\mu(\phi + \lambda h)) = \frac{\delta \mathcal{L}_\lambda}{\delta \psi} h + \frac{\delta \mathcal{L}_\lambda}{\delta(\partial^\mu \psi)} \partial_\mu \phi h'$$

$$\psi = (\phi + \lambda h), \quad h' = \frac{dh}{d\phi}$$

We set:

$$f = \frac{h}{1 + \lambda h'}$$

$$\begin{aligned} \frac{d\mathcal{L}_\lambda}{d\lambda} &= \frac{\delta \mathcal{L}_\lambda}{\delta \phi} f + \frac{\delta \mathcal{L}_\lambda}{\delta(\partial^\mu \phi)} \partial^\mu f = \\ &= f \left(\frac{\delta \mathcal{L}_\lambda}{\delta \phi} - \partial_\mu \left(\frac{\delta \mathcal{L}_\lambda}{\delta(\partial^\mu \phi)} \right) \right) + \partial_\mu \left(\frac{\delta \mathcal{L}_\lambda}{\delta(\partial^\mu \phi)} \right) f \end{aligned} \quad (2)$$

One expects that the first term in (2) - f multiplied with the equation of motion - does not contribute on the mass shell. The second term should drop out also off mass shell in the adiabatic limit because of the total derivative in front.

We model the inductive construction à la Epstein and Glaser^{5,6} of the time ordered products closely after the naive argument given above.

Assume that all time ordered products with less than n points have already been defined. We include in our induction hypothesis the following assumptions (whose consistency is easily verified in the case of the two point function):

$$\begin{aligned} \frac{d}{d\lambda} T_\lambda^{(-)}(x_1, \dots, x_m) &= \frac{d}{d\lambda} T_{\lambda, m}^{(-)} = \\ &= \sum_{i=1}^m \left\{ (\partial_\mu^{x_i}) T_{\lambda, m}^{(-)1, i, \mu} + T_{\lambda, m}^{(-)2, i} + \right. \\ &\quad \left. (-i) \sum_{j \neq i} \delta(x_i - x_j) T_{\lambda, m-1}^{(-)2, j}(x_1 \dots x_{i-1} x_{i+1} \dots x_m) + \right. \\ &\quad \left. \dots (m^2 + \square) \phi : T_\lambda^{(-)3} \right\} \quad (m < n) \end{aligned} \quad (3)$$

$T^{(-)}$ denotes alternatively chronological or antichronological ordering.

$$T_{\lambda,m}^{1,i,\mu} = T(:\bar{\Psi}_\lambda(x_1): \dots : f \frac{\delta \bar{\Psi}_\lambda}{\delta(\partial^\mu \phi)}(x_1) : \dots : \bar{\Psi}'_\lambda(x_m):) \quad (4)$$

$$T_{\lambda,m}^{2,i} = T(:\bar{\Psi}_\lambda(x_1): \dots : f \mathcal{L}_{\lambda,\phi}(x_i) : \dots : \bar{\Psi}_\lambda(x_m):) \quad (5)$$

$$\mathcal{L}_{\lambda,\phi} = \frac{\delta \bar{\Psi}_\lambda}{\delta \phi} - \partial_\mu \frac{\delta \bar{\Psi}_\lambda}{\delta(\partial^\mu \phi)} - (\square + \bar{m}^2)\phi$$

$:(\square + \bar{m}^2)\phi \dots : T_{\lambda,m}^3$ denotes a sum of expressions with at least one operator $(\square + \bar{m}^2)\phi$ in their Wick ordered operator part. $T_{\lambda,m}$, $T_{\lambda,m}^{1,i,\mu}$, $T_{\lambda,m}^{2,i}$, $T_{\lambda,m}^3$ can be represented as formal power series in λ .

$$T_{\lambda,m} = \sum_n T_{\lambda,m}^{(n)} \lambda^n$$

$$\frac{d}{d\lambda} T_{\lambda,m} = \sum_n T_{\lambda,m}^{(n)} \lambda^{n-1} n$$

$T_{\lambda,m}^{(0)}$ is equal T_m of equ. (1). Equ. (3) has to be understood as equality in the sense of formal power series. Furthermore we assume that the c-number kernels, which appear in a Wick product expansion of $T_{\lambda,m}$, $T_{\lambda,m}^{i,j,(\mu)}$ have all properties necessary for establishing the adiabatic limits

$$\int T_{\lambda,m}^{(i,j),(\mu)}(x_1, \dots, x_m) g(x_1) \dots g(x_m) dx_1 \dots dx_m, \quad g \rightarrow \text{konst.}$$

Equation (3) implies $dS_\lambda/d\lambda = 0$: the contributions to $dS_\lambda/d\lambda$ from $T_{\lambda,m}^{2,i}$ are cancelled by those from

$$(-i) \sum_{j \neq i} \delta(x_i - x_j) T_{\lambda,m}^{2,i}.$$

The second term picks up from the perturbation expansion a factor $(-i)/(m+1)$ relative to the first one. $1/m$ cancels against the sum over j and $(-i)$ produces a relative minus sign.

The contribution from $\partial_\mu^{x_i} T_{m,\lambda}^{1,i,\mu}$ drops out in the adiabatic limit and the contributions from the terms $:(\square + \bar{m}^2)\phi \dots : T_{\lambda,m}^3$ vanish on the mass shell.

In order to construct the time ordered n-point product $T_\lambda(x_1 \dots x_n)$ (along Epstein and Glaser's lines) we consider first

$$D_\lambda^{(n)}(x_1 \dots x_{n-1}; x_n) = \sum_{\substack{J \cup J' = \{x_1 \dots x_{n-1}\} \\ J \cap J' = \emptyset \\ J \neq \emptyset}} \left[T_\lambda(J', x_n), \bar{T}_\lambda(J) \right] (-1)^{|J|}$$

Applying the induction hypothesis (3) to the factors $T_\lambda, \bar{T}_\lambda$ (with less than n points) one obtains

$$\begin{aligned} \frac{d}{d\lambda} D_{\lambda,n} &= \sum_{k=1}^n \left\{ \partial_\mu^{x_k} D_{\lambda,n}^{1,\mu,k} + D_{\lambda,n}^{2,k} + (-i) \sum_{j \neq k} \delta(x_j - x_k) \cdot \right. \\ &\quad \left. D_{\lambda,n-1}^{2,j}(x_1 \dots x_{k-1} x_{k+1} \dots x_n) + :(\square + \bar{m}^2)\phi : D_{\lambda,n}^3 \right\} \end{aligned} \quad (6)$$

$D_{\lambda,n}^{1,\mu,k}, D_{\lambda,n}^{2,k}$ are similarly defined as $T_{\lambda,n}^{1,\mu,k}, T_{\lambda,n}^{2,k}$, that is, the operator $:\bar{\Psi}_\lambda^*(x_k):$ of $D_{\lambda,n}$ is substituted by $:f_{\frac{\delta \mathcal{L}}{\delta (\partial^\mu \phi)}}(x_k):$ and $:f_{\mathcal{L},\phi}(x_k):$ respectively. The only feature of the term $:(\square + \bar{m}^2)\phi \dots : D_{\lambda,n}^3$ to be noted is the explicit occurrence of an operator $(\square + \bar{m}^2)\phi$.

A proper definition of the time ordered n-point product is obtained by 'splitting' $D_{\lambda,n}$, that is, one looks for a 'retarded' product $R_{\lambda,n}(x_1 \dots x_n)$, whose support is contained in $V^- = \{x = \{x_1 \dots x_n\} \in R^{4n}, (x_i - x_n)^2 \geq 0, x_i^0 - x_n^0 > 0, i < n\}$ such that the support of $(D_{\lambda,n} + R_{\lambda,n})$ is in $V^+ = -V^-$. (The support of D is contained in $V^+ \cup V^-$).

Every term on the right side of equ. (6) has according to the induction hypothesis the same structure as D_λ . This fact enables us to split every term individually. Afterwards we integrate formally with respect to λ

$(\lambda^n \rightarrow \frac{\lambda^{n+1}}{n+1})$ and add $R_{n,\lambda=0}$. This procedure renders an acceptable definition of $R_\lambda(x_1 \dots x_{n-1}; x_n)$ and thereby also a proper definition of the time ordered product

$$T_\lambda(x_1 \dots x_n) = R_\lambda(x_1 \dots x_{n-1}; x_n) - \sum \bar{T}_\lambda(J) T_\lambda(J', x_n) (-1)^{|J|} = R_{\lambda,n} - R'_{\lambda,n}$$

$$J \cup J' = \{x_1 \dots x_{n-1}\}$$

$$J \cap J' = \emptyset$$

$$J \neq \emptyset$$

By a computation analogous to the computation of $\frac{d}{d\lambda} D_{\lambda,n}$ (equ. 6) one can easily verify that $R'_{\lambda,n}$ satisfies by itself the induction hypothesis.

We have to build in all splittings the correct mass and wave function renormalisations in order to guarantee the existence of the adiabatic limit. The term

$$\sum_{j \neq k} \delta(x_k - x_j) D_{\lambda,n-1}^{2,k}$$

of equ. (6) is handled in the same manner as $D_{\lambda,n-1}^{2,k}$ in the preceding induction step (the δ -functions are unessential). $D_{n,\lambda} = 0$ is treated according to some ad hoc prescription as it was fixed after equ. (1). Concerning the term $\partial_\mu^{x_k} D_{n,\lambda}^{1,k,\mu}$ we first split $D_{n,\lambda}^{1,k,\mu}$ and apply only afterwards the derivative $\partial_\mu^{x_k}$, that is, we leave the derivative outside the T-product. Following these prescriptions we reproduce the induction hypothesis for $T_{\lambda,n}$. One should note that apart from the restrictions for the selfenergy kernels one can choose the splittings of $D_{\lambda,n}^{1,k,\mu}$, $D_{\lambda,n}^{2,k}$ and $D_{\lambda,n}^3$ arbitrarily.

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