

SPACE-TIME TOPOLOGY AS A CONSEQUENCE OF THE DYNAMICS OF CLOSED BOSON STRINGS. II. THE SECOND APPROXIMATION

N. F. Nelipa and M. Yu. Pekar

The possibility of constructing a theory where the space-time topology would be determined by the dynamics of closed boson strings is analyzed. The analysis has the form of an expansion into a series with respect to the coupling constant. The second-order expansion is considered. Some new constraints on the admissible space-time topology are found in addition to those in the first-order approximation.

In the previous paper [1] we described an approach in which the space-time topology was determined by the dynamics of strings, and also considered the zeroth- and first-order approximations with respect to the string coupling constant. The purpose of the present paper is to analyze the second-order approximation. In accordance with our approach, a cohomological complex will be constructed and a topological invariant will be found in the given approximation. It will be shown that the requirement of the differential nilpotency in the second-order approximation strengthens the constraints on the possible choice of the space-time topology.

1. To the second-order approximation there corresponds the one-loop diagram and also a diagram describing the four-string interaction. As before, the expression for the interaction vertex is taken in accordance with formula (1) in [1].

Let us find the system of equations implied by the property of nilpotency of the total differential.

Because in the first-order approximation the differential is expressed by formula (3') in [1],

$$d^1\omega = d^0\omega + F^1(\omega) = d^0\omega + g\Phi * \omega,$$

we shall seek the total differential in the form

$$d\omega = d^0\omega + F^1(\omega) + F^2(\omega) + F^3(\omega) + \dots \quad (1)$$

where $F^i(\omega)$ is a linear functional on the set of differential forms ω , which is proportional to the i th power of the coupling constant. By virtue of (1), we obtain the following expression for the square of the differential:

$$[d]^2\omega = [d^0]^2\omega + \{d^0 F^1(\omega) + F^1(d^0\omega)\} + F^1(F^1(\omega)) \\ + \{d^0 F^2(\omega) + F^2(d^0\omega)\} + F^2(F^1(\omega)) + F^1(F^2(\omega)) + \dots$$

In the first approximation we have

$$d^0(F^1(\omega)) + F^1(d^0\omega) + F^1(F^1(\omega)) = (d^1\Phi) * \omega = 0.$$

i.e., the condition of nilpotency of the differential in the first-order approximation holds on the set of closed forms. This relation corresponds to the covariant equation for the total differential:

$$[d]^2\omega = (d\Phi) * \omega. \quad (2)$$

Using (2) we derive a system of linked equations for the functionals Φ and F^i :

$$[d^0]^2 = 0, \\ d^0(F^i(\omega)) + F^i(d^0\omega) = F^{i+1}(\omega), \\ d^0(F^1(\omega)) + F^1(d^0\omega) + F^1(F^1(\omega)) = (d^1\Phi) * \omega, \\ F^{2i}(\omega) + g(F^2(\Phi * \omega) + \Phi * F^2(\omega)) = gF^2(\Phi) * \omega. \quad (3)$$

In particular, for the functional F^2 we obtain the system

$$\begin{cases} F^{2'}(\omega) = 0, \\ F^2(\Phi * \omega) + \Phi * F^2(\omega) = F^2(\Phi) * \omega, \end{cases} \quad (4)$$

which supplements conditions (1) and (4).

System (3) must be supplemented with a condition following from the property of the differential to raise the order of the form by unity:

$$|F^i(\omega)| = |\omega| + 1. \quad (5)$$

2. Our task is to find the form of the functionals F^i , which are solutions to the system of equations (4), (5).

To satisfy condition (5) it is necessary to introduce the dual differential \bar{d}^0 in the contour space, which does not raise but lowers the order of the form by unity. By definition, this differential operator satisfies the relation

$$d^0 \bar{d}^0 \omega = \omega, \quad \bar{d}^0 d^0 \tilde{\omega} = \tilde{\omega}. \quad (6)$$

Alongside the zeroth-order dual differential \bar{d}^0 , there also exists the total dual differential \bar{d} :

$$\begin{cases} d\bar{d}\omega = \omega, \quad \bar{d}d\tilde{\omega} = \tilde{\omega}, \\ [\bar{d}]^2\omega = 0. \end{cases} \quad (7)$$

for which the conditions

$$\bar{d}\omega \equiv \bar{d}^0\omega + \bar{F}^1(\omega) + \bar{F}^2(\omega) + \dots = \bar{d}^0\omega + \bar{\Phi} \circ \omega + \bar{F}^2(\omega) + \dots$$

hold.

Note that conditions (6) and (7) are understood in the sense of cohomological equivalence classes, i.e., accurate to the addition of an arbitrary exact form to ω .

Writing in full Eq. (7) we find a system of linked equations for the functionals \bar{F}^i , which is obtained by means of the transformation $F \rightarrow \bar{F}$, $\Phi \rightarrow \bar{\Phi}$, $* \rightarrow \circ$ in (3). Besides, there appears an additional system of equations:

$$\begin{cases} d^0(\bar{\Phi} \circ \omega) = -\Phi * \bar{d}^0\omega, \\ \bar{d}(\Phi * \omega) = -\bar{\Phi} \circ d^0\omega. \end{cases} \quad (8)$$

$$\begin{cases} \bar{\Phi} \circ (\Phi * \omega) + \bar{d}^0 F^2(\omega) + \bar{F}^2(d^0\omega) = 0, \\ \Phi * (\bar{\Phi} \circ \omega) + d^0 \bar{F}^2(\omega) + F^2(\bar{d}^0\omega) = 0. \end{cases} \quad (9)$$

Equations (8) express additional operation rules on the functionals Φ and $\bar{\Phi}$, and system (9) supplements the system of equations (4) for the functionals F^2 and \bar{F}^2 . Analyzing system (4), (7), (9), common for F^2 and \bar{F}^2 , with consideration for rules (8) and rules (4) in [1], we find the following form of the desired functionals in the second-order approximation with respect to the coupling constant:

$$\begin{aligned} F^2 &= g^2(d^0[\bar{\Phi} \circ ((\bar{d}^0\Phi) * d^0\omega)] - \Phi * \bar{d}^0(\Phi * \omega)), \\ \bar{F}^2 &= g^2(\bar{d}^0[(\bar{d}^0\Phi) * d^0(\Phi \circ \omega)] - \bar{\Phi} \circ d^0(\bar{\Phi} \circ \omega)). \end{aligned} \quad (10)$$

The composition $*$ means the transition of a single string z into two strings x and y . Therefore the dual composition \circ means the reverse transition $z' = x, y$. The first terms in expressions (10) describe the one-loop interaction, and the second terms describe a four-string vertex.

3. We now find the topological invariant in the second-order approximation with respect to the coupling constant, which determines the effective action.

We first determine the integration measure on U^2 . Since the first-order approximation with respect to the coupling constant means a higher order of contact as compared to the first-order approximation, the neighborhood over which the integration extends in the expressions for the topological invariants contains

U^1 . Therefore in the integral measure on the neighborhood of U^2 there appear two new terms as compared to U^1 :

$$\begin{aligned} \int_2^a &= \frac{1}{2} \left\{ \int_{-\tau_1}^{\tau_1} d\tau' d\tau'' \oint Dx(\sigma, \tau') \oint Dy(\sigma, \tau'') \delta(\tau' - \tau'') v(xyz) \right. \\ &= \int_{\tau'}^{\tau_1} d\tau \oint Dz(\sigma, \tau) \bar{v}(zxy) \int_{\tau}^{\tau_1} d\tilde{\tau} d\tilde{\tau}' \oint Dx(\sigma, \tilde{\tau}) \oint Dy(\sigma, \tilde{\tau}') \\ \int_2^b &= \frac{1}{2} \int_{\tau_0}^{\tau_1} d\tau \oint Dz(\sigma, \tau) v(xyz) \int_{\tau}^{\tau_1} d\tau' d\tau'' \oint Dx(\sigma, \tau') \\ &= \oint Dy(\sigma, \tau'') \delta(\tau' - \tau'') \bar{v}(zxy) \int_{\tau'}^{\tau_1} d\tilde{\tau} \oint Dz(\sigma, \tilde{\tau}). \end{aligned}$$

The transformation rule for functionals in the transition from one neighborhood to another in the second-order approximation is written thus:

$$\delta^2 \Phi = d^2 \Psi = d^0 \Psi + g \Phi * \Psi + F^2(\Psi). \quad (11)$$

As can easily be shown by means of direct verification, the topological invariant in the second-order approximation with respect to the coupling constant, i.e., an expression invariant relative to (11), takes the form

$$\Gamma^2 = \frac{\Gamma^1}{U^2} + \frac{2}{5} \int_{U^2} \langle \Phi, F_a^2(\Phi, \bar{\Phi}) \rangle + \frac{1}{2} \int_{U^2} \langle \Phi, F_b^2(\Phi, \bar{\Phi}) \rangle, \quad (12)$$

where

$$\begin{aligned} F_a^2 &= g^2 d^0 [\bar{\Phi} \circ ((\bar{d}^0 \Phi) * d^0 \Phi)], \\ F_b^2 &= -g^2 \Phi * \bar{d}^0 (\Phi * \Phi). \end{aligned}$$

Using (3) these expressions can be brought to the form not involving $\bar{\Phi}$:

$$F_a^2(\Phi) = g^2 \Phi * \bar{d}^0 [(\bar{d}^0 \Phi) * d^0 \Phi].$$

The variation of (12) results in equations of motion:

$$d^2 \Phi = d^0 \Phi + g \Phi * \Phi + F^2(\Phi) = 0.$$

One can see that the equations coincide with the closure condition for the form Φ . Thus, the invariant Γ^2 , which describes the topological properties of the contour space, assumes a discrete set of values on the set of closed forms.

The quantity Γ^2 is the one-loop approximation to the effective action, which immediately results in the second-quantized theory.

4. To illustrate the mechanism of the appearance of additional constraints on the space-time topology due to the second-order approximation, we consider a simple example of a topology admissible in the first-order approximation.

As an example we take a torus $T = R^{(D-1,1)}/l$, where l is the lattice $\{e_i\}$ on the space $R^{(D-1,1)}$, $(e_i, e_j) \equiv \sum_{IJ} e_i^I e_j^J \eta_{IJ} \equiv g_{ij}$. Consider expression (10) for F_a^2 . From the physical standpoint, it describes the transition of a single string into two strings and the subsequent reverse transition of the two strings into the single string. If before the interaction the indices of the mapping of the string z on the torus were zero, then after the interaction the indices of the mapping of the strings x, y become equal to $\pm k_i \neq 0$, respectively,

on a lattice vector e_i . Upon repeated interaction all mapping indices vanish again. This interaction can be realized on a torus whose lattice contains at least two vectors e_i, e_j ($\pi_1(T) \neq Z$, where Z is the set of integers). Moreover, the dual operator \circ involved in (8) is a composition $x, y \rightarrow \bar{x}$ on the dual lattice \bar{l} :

$$(\bar{e}_i, e_j) = \delta_{ij}.$$

This means that a lattice on the space-time must be self-dual.

In the approximation order with respect to the coupling constant we are dealing with, there appears one more constraint on the possible space-time topology in the case of a torus. Consider the graph of the lattice l , which means that the elements e_i are denoted by points and the i th and j th elements are joined by a line if $(e_i, e_j) \neq 0$. Taking into account that we parametrized the contour space by passing consecutively from one lattice vector to another, from the requirement that the parametrization should be single-valued we conclude that any two points in the graph can be joined by only a single path. This is equivalent to the property that the graph possesses a tree structure, i. e., it contains no closed cycles.

Hence, taking account of the second-order approximation leads to additional constraints on the possible space-time topology as compared to the first-order approximation. It should be expected that the inclusion of higher-order approximations will result in a further concretization of the space-time topology and at the same time will elucidate some questions related to compactification.

REFERENCES

1. N. F. Nelipa and M. Yu. Pekar. *Vest. Mosk. Univ. Fiz. Astron.*, vol. 32, no. 6, p. 18, 1991.

15 May 1991

Research Institute of Nuclear Physics