

C031-060-080
 [N23 080]

Interaction operators
 which will be called
 the third class.

Properties of operators
 III. ~~Matrix~~

Among the operators of the second class, there are a particular family of operators, which can be expressed as sums of products of nonlocal field operators such U, U^* and operators of the first class. Such operators will be called A_{III} . Any one of them can be written in the form

$$A_{III} = A_0 + A_1 U + A_2 U^* + A_3 U^* A_4 U + \dots$$

$$+ B_1 \psi^* \gamma^4 \psi + \dots$$

$$+ C_1 U \psi^* \gamma^4 \psi + \dots$$

where A_0, A_1, A_2 etc., $B_1, B_2, \dots, C_1, C_2, \dots$ are all operators of the first class, A of the third class,

as will be shown in the following case, the space-time average of any operator which is defined

$$A = \int \int (dx_i^{\mu}) (dx_j^{\nu}) (x_i^{\mu} | A | x_j^{\nu}),$$

which is still a submatrix, of any operator A of the third class, is of

(4)

Obviously,

particular importance in connection with the physical interpretation of the formalism of nonlocal field theory, \bar{A} is ~~still~~ a ^{kind of} submatrix of A with rows and columns characterized by the distinct number of particles in each of all possible quantum states. It can be shown very generally that the space-time average \bar{A} of arbitrary operator A of the third class is an ~~submatrix~~ ~~with~~ ~~the~~ ~~following~~ property: ~~the~~ ~~submatrix~~ ~~elements~~ ~~of~~ ~~\bar{A}~~ ~~the~~ ~~submatrix~~ ~~\bar{A}~~ ~~which~~ ~~form~~ (Thus, we can write \bar{A} more precise in the

$$\bar{A}_{(n)(k, l, m)} = \sum_{n'} \bar{A}_{(n')(k, l, m)}$$

where $n'(k, l, m)$ are the numbers of particles in the state k, l, m characterizing the row and column respectively of the submatrix.

Now it can be shown very generally that the matrix element of \bar{A} is different

(5)

from zero, only if the row and the column are ^{represented} characterized by the states of the ^{total} system of particles with the same total energy and total momentum. It should be notice that we mean by the total energy and the total momentum of the system the ^{total} sum of the energies of motions of ⁿ centers of mass of particles present and that of the momenta of motions of centers of mass of particles present, respectively

In other word the matrix element $(\dots | A | \dots)$ is different from zero, only if the

$$\sum_{\underline{k}, l, m} k_{\mu} n'(\underline{k}, l, m) = \sum_{\underline{k}, l, m} k_{\mu} n''(\underline{k}, l, m),$$

where the summation covers all the possible states of particles of all types, in consideration.
(of particles)

(6)

In order to ~~show~~^{see} that this is generally true, however, it is sufficient to any of simple ~~examples~~^{cases}, because look at the ~~derivation~~ way of method of proof can easily be extended to more general cases. For example, in ~~the~~ case A has a particular form

$$A_3 U^* A_4 U,$$

where A_3, A_4 are arbitrary operators of the ~~the~~ first class and U, U^* are nonlocal scalar field operators, it can ^{always} be decomposed into ~~the form~~ terms such as

$$L_I = \exp(i \int d^4x \mathcal{L}(x)) \prod_{k,l} A_{k,l}^* \exp(i \int d^4x \mathcal{L}(x)) \exp(i \int d^4x \mathcal{L}(x)),$$

$\prod_{k,l} A_{k,l}^*$ and $\prod_{k,l} A_{k,l}$ are ~~or similar~~ apart from the purely numerical $\exp(i \int d^4x \mathcal{L}(x))$ factors, the space-time average of ~~each of~~

$\exp(i \int d^4x \mathcal{L}(x))$ the matrix elements of $(\alpha_j | U | \alpha_i)$, $(\alpha_j | U | \alpha_i)$ etc. is evidently equivalent to the integration ~~each~~

(8)

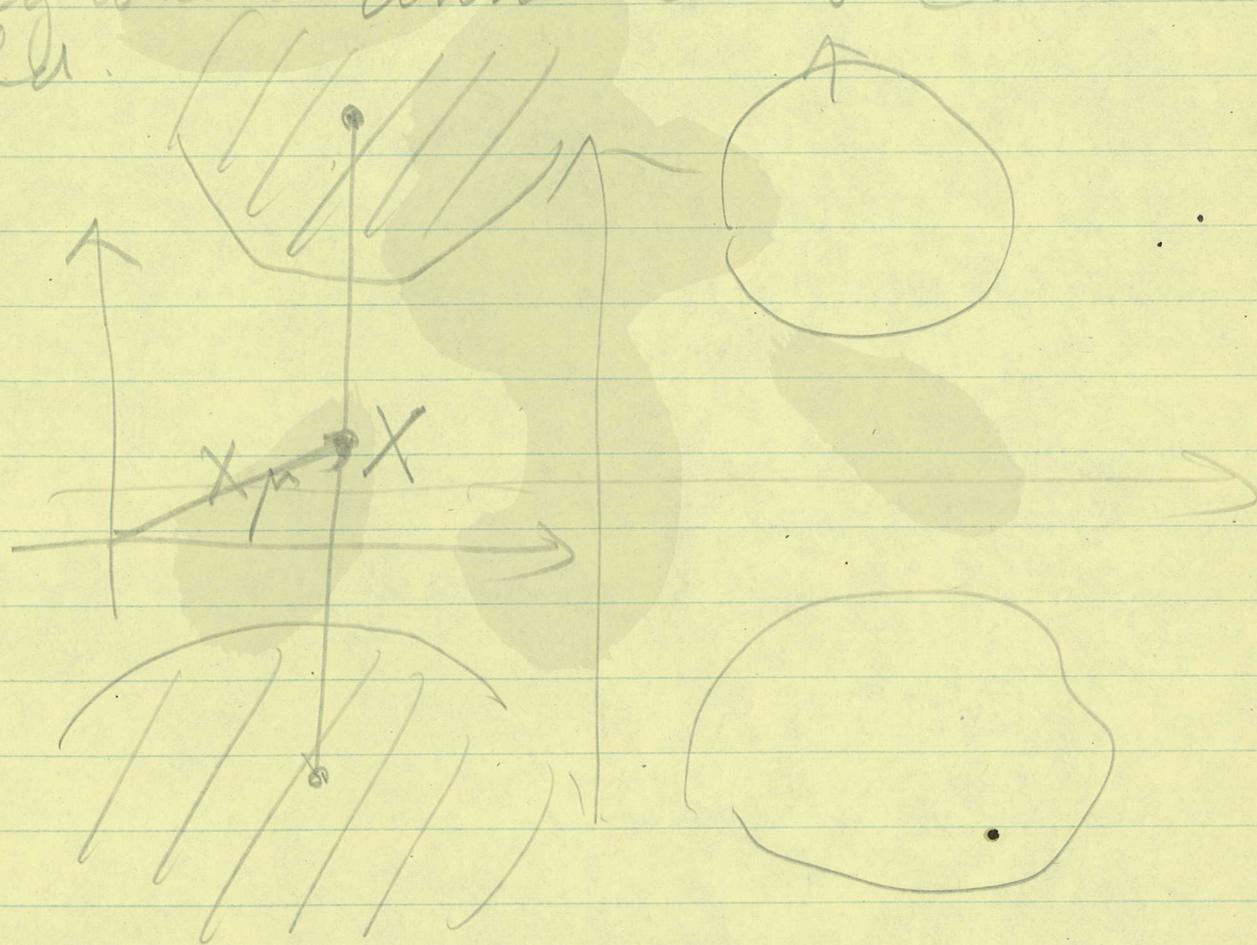
if we

It is true even more generally that for many nonlocal operators A , \bar{A}

$$\bar{A} = \int \int A(x, v) \rho(v) dx (dv)^4$$

is a submatrix with rows and columns can be regarded

characterized by distribution in number of particles and the matrix elements are different from zero only when conservation laws hold.



(9)

So $\rho(x)$ can be so chosen that
 $\rho(x)$ is different from zero
only for

$$x_4' - x_4'' \geq cT,$$

where T is a time interval ^{of order}
compared with the collision time
in question.

Alternatively, we can write

$$\bar{A} = \text{Trace } A D_{\frac{1}{2}T}$$

$$= \sum_{x', x''} (x' | A | x'') (x'' | D_{\frac{1}{2}T} | x')$$

where $D_{\frac{1}{2}T}$ is an invariant, (field
independent) operator with the matrix
elements $(x' | D_{\frac{1}{2}T} | x'')$, which are
different from zero only when
 $x_4' + x_4''$ is in time direction and
 $x_4' - x_4'' \leq cT,$

(10)

IV. S-Matrix in N.L.F.T.

Now we can go into the question of the interaction of fields. The first question, with which we meet, is whether there is the Schrödinger equation or any substitute for it in N.L.F.T. It is difficult to answer this question directly. However, irrespective of the ~~existence~~ ^{presence} or absence of such an equation, it is possible ~~in~~ for our case to define the S-matrix in the sense of Heisenberg and Møller.^(?) Suppose that a system ~~there exist at~~ ^{exists at the beginning} a number of particles with ^{the} internal structure in various quantum states. After a long time, ^{both} the number ^{of particles and} the ~~states~~ ^{quantum states} of particle their distribution in various quantum states will be changed due to the interaction between particles. The question now is to determine the probability that ^{any} ~~any~~ ^{arbitrary} distribution of in number of particles in different quantum states, provided that the initial distribution is known,

(41)

It is clear that the distribution will not be changed, if the particles are independent of each other. This can be expressed mathematically by the statement that there exists a matrix diagonal matrix with rows and columns, each being characterized by the distribution in number of particles of various types in various states, ~~and~~ the absolute square of absolute value of these the diagonal element must be 1 each if there is no interaction between

be equal to 1, corresponding to the fact that the probability of finding the ~~system~~ original distribution ^{after a long time} is 1. One can simplify the state ^{vector} further by assuming that this diagonal matrix is the unit matrix (submatrix)

The effect of the interaction is then amount to change this matrix from the unit matrix to some nondiagonal matrix S . This matrix ~~must~~ satisfy, at least, the following two conditions

(12)

The square of the absolute value of the matrix element of \bar{S} is supposed to represent the probability that the distribution corresponding to the row will be realized, when the system is initially in the state characterized by the column.

In order that such an interpretation is permissible, the matrix \bar{S} must satisfy, at least, the following three conditions:

$$(i) \quad \sum_j \left(\bar{S}_{ij}^* \bar{S}_{ji} \right) = 1$$

independent of i .

(ii) The matrix element of \bar{S} must be zero, unless the total energy and momentum of the state of the system corresponding to the row is equal to those corresponding to the column.

The latter condition is satisfied for any \bar{S} defined by

$$(i) \quad \bar{S}_{ij} = \int \int (x' | S_{ij} | x'') \rho(x' - x'') \frac{dx'}{(dx')^4} \frac{dx''}{(dx'')^4}$$

$$= \int \int S_{ij}(X, r) \rho(r) (dX)^4 (dr)^4$$

(iii) $\bar{S}_{ij}^* \bar{S}_{ji}$ must be rel. inv.

(13)

where S is an arbitrary invariant operator, which can be expressed as a sum of products of field operators, and $\rho(r_\mu)$ is any invariant function of r_μ .

\bar{S} can alternatively be written as

$$\bar{S} = \text{Trace } S \mathcal{D}_T$$

where \mathcal{D}_T is an operator depending only on p_μ , so that it can be written in the form

$$\mathcal{D}_T = \int \rho_T(i p_\mu) \exp(i l^\mu p_\mu / \hbar) (d p_\mu)$$

or

$$\langle x' | \mathcal{D}_T | x'' \rangle = \mathcal{D}_T(x, r) = \rho(\pm r_\mu)$$

with $\rho(r_\mu)$ being an invariant fn of r_μ .

$$\begin{aligned} \text{Trace } S \mathcal{D}_T &= \int \int (x' | S_{ij} | x'') (x'' | \rho | x') (d x')^4 (d x'')^4 \\ &= \int \int S_{ij}(X, r) \rho(-r) (d X)^4 (d r)^4 \end{aligned}$$

(14)

The condition (iii) is also satisfied, if ρ is an invariant operator depending only on ψ , because the form of $\bar{\psi}$ as the function of ψ and ψ' does not change by Lorentz transformation.

The condition (i) is not identical with the more familiar condition

$$\bar{\psi}^* \bar{\psi} = 1,$$

the latter being a much stronger requirement to $\bar{\psi}$.

Our main task is to find an operator $\bar{\psi}$, so as to satisfy that $\bar{\psi}$ satisfies the condition (iii) and reduces to the usual probability amplitude or wave matrix in Moller's sense in the limit of local fields.