vantage of presenting a big difference between the Čerenkov/scintillation pulse ratio for fast charges and for fast monopoles. For example, if a Lucite Čerenkov counter is used, then the ratio for fast monopoles is 2.25 times that for fast charges. The ratio expected from fast charged monopoles is intermediate between the ratio for fast charges and that for fast monopoles. The range of unsaturated response of the scintillation counter will limit the pole strengths for which this type of experiment can be used. These considerations make this type of experiment particularly useful in searching for small-pole-strength monopoles or charged monopoles.

## First-Method Experiment

This type of experiment compares the pulses from two Čerenkov counters with different refractive indices. The difference between the (Čerenkov counter $a$ / Cerenkov counter $b$ ) pulse ratio for fast charges and fast monopoles depends on the permittivity ratio ( $\epsilon_{a}{ }^{\prime} / \epsilon_{b}{ }^{\prime}$ ). For example, if a lucite/water combination is used then
$\left(\epsilon_{a}{ }^{\prime} / \epsilon_{b}{ }^{\prime}\right)=1.27$ and fast monopoles yield a (counter $a /$ counter $b$ ) ratio that is 1.27 times that for fast charges. Fast charged monopoles would yield ratios intermediate between those for fast charges and fast monopoles. In the low-ionization region this type of experiment is limited by two considerations. First, a small pole strength yields a small number of Cerenkovcounter photons which results in poor statistics with which to determine the (counter $a$ /counter b) ratio. Secondly, slow charges can give the same pulse ratio as fast monopoles and, at least for cosmic-ray work, there are a lot of slow charged particles in the low-ionization region. Here "slow" means a particle which is not fast for both counters $a$ and $b$. As pointed out previously, ${ }^{1}$ particles which are slow can be discriminated against by introducing a third Cerenkov counter using a gas.

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# Model Hamiltonians for Local Quantum Field Theory* 

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#### Abstract

We study the problem of renormalization for the interaction of a charged scalar meson field with a fixed two-level point source. Our especial interest is coupling-constant renormalization. We study in particular the problem of obtaining eigenstates and eigenvalues of the Hamiltonian for the fixed-source theory. We propose a model for the fixed-source theory, in which mesons exist only if their momenta ( $k$ ) lie within an infinite set of intervals: $0<k<k_{0}, \frac{1}{2} \Lambda<k<\Lambda, \frac{1}{2} \Lambda^{2}<k<\Lambda^{2}$, etc., where $k_{0}$ is of the order of the meson mass, and $\Lambda$ is much larger. We solve this model by treating the mesons in the $n$th interval (or lower) as a perturbation on mesons in the ( $n+1$ )st interval (and higher). This reduces the problem to the solution of two strongly cut-off Hamiltonians, one of which must be solved for an infinite sequence of coupling constants $\left\{g_{n}\right\}$, one for each momentum interval. We show that even if the low-momentum coupling constants $g_{1}$, $g_{2}$, etc., are small, the sequence goes to infinity as $n \rightarrow \infty$. We analyze the Lee model similarly; here the sequence is undefined above some finite value of $n$. We show a close analogy between our analysis and the analysis of quantum electrodynamics of Gell-Mann and Low. Then we analyze the full fixed-source Hamiltonian qualitatively. We expland the meson field in terms of a complete set of "wave-packet" states, the coefficients being discrete oscillator variables. The states are so chosen that the self-interactions of oscillators dominate the coupling between oscillators. For each order of magnitude for the meson momentum, there is one pair of oscillators coupled to the source; this coupling can be analyzed analogously to our model. The full fixed-source Hamiltonian is thereby reduced to the solution of a Hamiltonian for two oscillators coupled to a two-level source.


## I. INTRODUCTION

T${ }^{\top}$ HE Hamiltonian formulation of quantum mechanics has been essentially abandoned in investigations of the interactions of $\pi$ mesons, nucleons, and strange particles. This is a pity. The Hamiltonian approach has several advantages over the kind of approach (using dispersion relations) presently in use.

[^0]One advantage is that all properties of a system are uniquely determined, given the Hamiltonian of the system. A second advantage is the existence of many approximation methods for solving a given Hamiltonian. The principal approximation methods are perturbation theory, the variational method, and the W.K.B. approximation. A third advantage is that one can often analyze a Hamiltonian intuitively, using the uncertainty principle and classical arguments, to get order of magnitude estimates of its solution; from such
an analysis one can often determine what approximation procedure will work best.

There are two reasons why the Hamiltonian approach was discarded in the study of strong interactions. One reason was that no one knew what Hamiltonian to use, or how to obtain the correct Hamiltonian. The other reason was the problem of renormalization: the problem that whenever one tried to solve a Hamiltonian for a Lorentz-invariant theory, particle self-energies and the like were infinite.

The problem of renormalization has several aspects. One is the practical aspect just mentioned, namely, that whenever one has tried to solve a Lorentz-invariant theory with a standard approximation procedure such as perturbation theory or the Tamm-Dancoff method, one gets infinities. Then there is a more fundamental difficulty. This difficulty is that any estimate of the order of magnitude of the interaction energy gives infinity, for a relativistic quantum-field theory. For example, the simplest relativistically invariant interaction Hamiltonian, for a neutral scalar field $\phi(x)$, is

$$
H_{I}=g_{0} \int \phi^{3}(\mathbf{x}) d^{3} x
$$

where $g_{0}$ is a constant. We might estimate the order of magnitude of $H_{I}$ by computing the square root of the matrix element $\langle\Omega| H_{I}{ }^{2}|\Omega\rangle$ where $|\Omega\rangle$ is the vacuum state for a free field. This matrix element is infinite. This difficulty is a fundamental one, despite the crude nature of such estimates, because any sensible method for solving Hamiltonians requires that one first make just such crude estimates of the order of magnitude of the terms in the Hamiltonian.

A third aspect of the renormalization problem is that there is only one Hamiltonian known that can be solved but requires renormalization. This is the Hamiltonian of the Lee model. ${ }^{1}$ Unfortunately, the Lee model (formulated so that the renormalization is infinite, i.e., without a "cutoff") cannot be solved without introducing a "ghost state"-a state with negative normwhich is undesirable. This example has made many people pessimistic about the possibility of resolving the renormalization problem.

It is not known whether relativistic Hamiltonians can be solved exactly without encountering the same infinities that occur in perturbation theory, and without introducing ghost states such as occur in the Lee model. It is clear that the appearance of infinities in perturbation theory does not imply infinities in the exact theory since one can use perturbation theory only when the interaction energy is small, and this is demonstrably not the case for field-theoretic interactions. There are other approximate methods for solving Hamiltonians besides perturbation theory, but they are equally useless so long as one cannot establish their validity even qualitatively. So to resolve this question one will have

[^1]to learn to make qualitative analyses of relativistic Hamiltonians, using order-of-magnitude estimates, the uncertainty principle, etc.

We do not have the competence to make such an analysis at the present time. To develop the necessary techniques, it is useful to have model Hamiltonians which embody some but not all the difficulties presented by fully relativistic interactions. One such model Hamiltonian is presented in this paper. Our model is an offspring of the theory of scalar $\pi$ mesons interacting with a fixed point source. Our model Hamiltonian involves an interaction Hamiltonian which by any order of magnitude estimate is infinite. In perturbation theory it gives infinities analogous to coupling constant and self-energy divergences in relativistic theories. However, we will be able to analyze it in a sensible way: not completely, but well enough to provide considerable insight into the real problems involved in solving Hamiltonians requiring renormalization (in particular, coupling constant renormalization).

Gell-Mann and Low ${ }^{2}$ have considered the problem of what a solution of a relativistic theory would be like, if a solution exists. They examined quantum electrodynamics in particular. Their principle conclusion was that the "bare" coupling constant $e_{0}$ which appears in the unrenormalized Hamiltonian could not be arbitrary; it could take on only special values, or, in other words, it would have to be the solution of some kind of eigenvalue condition. To quote from their abstract:
"Thus it is shown that the unrenormalized coupling constant $e_{0}{ }^{2} / 4 \pi \hbar c$, which appears in perturbation theory as a power series in the renormalized coupling constant $e_{1}{ }^{2} / 4 \pi \hbar c$ with divergent coefficients, may behave in either of two ways: (a) It may really be infinite as perturbation theory indicates; (b) It may be a finite number independent of $e_{1}^{2} / 4 \pi \hbar c$."
We shall discuss the ideas of Gell-Mann and Low in more detail in Sec. VI.
Now we give a brief outline of the contents of this paper. In Sec. II we formulate our model Hamiltonian. In Sec. III we replace the original model by a sequence of Hamiltonians $H_{n}$ and analyze the individual Hamiltonians in the sequence. In Sec. IV, we show how to define our original model as a limit of the sequence considered in Sec. III. In Sec. V we analyze similarly a model derived from the Lee model. In Sec. VI we compare the results of the analysis of our models with GellMann and Low's analysis of quantum electrodynamics. In Sec. VII we discuss a method for qualitatively analyzing field-theoretic Hamiltonians, which involves expanding the fields in a discrete orthonormal set of functions. We state our conclusions in Sec. VIII.

## II. FORMULATION OF MODEL

In this Section we formulate our model Hamiltonian. We begin by quoting the complete Hamiltonian for

[^2]charged scalar mesons ( $\pi^{ \pm}$) interacting with a fixed point source with two quantum levels ("proton" and "neutron"):
\[

$$
\begin{align*}
& H=\int_{\mathrm{k}} \omega_{\mathrm{k}}\left\{a_{\mathrm{k}}^{\dagger} a_{\mathrm{k}}+b_{\mathrm{k}}^{\dagger} b_{\mathrm{k}}\right\} \\
&+g_{0} \int_{\mathrm{k}}\left(2 \omega_{\mathrm{k}}\right)^{-1 / 2}\left\{\left(a_{\mathrm{k}}+b_{\mathbf{k}}^{\dagger}\right) \tau^{+}+\left(a_{\mathrm{k}}^{\dagger}+b_{\mathrm{k}}\right) \tau^{-}\right\} \tag{1}
\end{align*}
$$
\]

where
$\int_{\mathbf{k}}$ means $(2 \pi)^{-3} \int d^{3} k$,
$\omega_{k}$ means $\left(\mathbf{k}^{2}+\mu^{2}\right)^{1 / 2}$,
$\mu$ is the meson mass, which we assume is of order 1 , $a_{\mathrm{k}}{ }^{\dagger}$ creates $\pi^{+}$,
$b_{\mathrm{k}}{ }^{\dagger}$ creates $\pi^{-}$,
$g_{0}$ is the (unrenormalized) coupling constant,
$\tau^{+}$takes the neutron state into the proton state,
$\tau^{-}$is the Hermitian conjugate of $\tau^{+}$.
The commutation relations of $a_{\mathrm{k}}$ and $a_{\mathbf{k}}{ }^{\dagger}$ are

$$
\begin{equation*}
\left[a_{\mathbf{k}}, a_{\mathbf{k}^{\prime}}+\right]=\left[b_{\mathbf{k}}, b_{\mathbf{k}^{\prime}}+\dagger\right]=(2 \pi)^{3} \delta^{3}\left(\mathbf{k}-\mathbf{k}^{\prime}\right) . \tag{2}
\end{equation*}
$$

This Hamiltonian when solved in perturbation theory leads to two divergences. The ground-state energy is linearly divergent; the renormalized coupling constant (related to a ground-state matrix element of $\tau^{+}$) is logarithmically divergent. It also has the property that an estimate of the interaction energy gives infinity. For example let us estimate the order of magnitude of $\int_{\mathbf{k}} a_{\mathbf{k}}{ }^{\dagger}\left(2 \omega_{\mathbf{k}}\right)^{-1 / 2}$; then we compute

$$
\langle\Omega| \int_{\mathrm{k}} \int_{\mathrm{k}^{\prime}} a_{\mathrm{k}} a_{\mathbf{k}^{\prime}}{ }^{\dagger}\left(2 \omega_{\mathrm{k}}\right)^{-1 / 2}\left(2 \omega_{\mathbf{k}^{\prime}}\right)^{-1 / 2}|\Omega\rangle,
$$

where $|\Omega\rangle$ is the meson vacuum state. This expression is easily computed and yields the quadratically divergent integral

$$
\int_{\mathbf{k}}\left(2 \omega_{\mathbf{k}}\right)^{-1}
$$

Our model Hamiltonian is constructed as follows: We take the Hamiltonian of Eq. (1) and throw away most of it. Specifically, the integration over all $\mathbf{k}$ is replaced by an integral over an infinite set of intervals. The intervals are as follows ( $k$ is the magnitude of $\mathbf{k}$ )

$$
0<k<k_{0}, \frac{1}{2} \Lambda<k<\Lambda, \cdots, \frac{1}{2} \Lambda^{n}<k<\Lambda^{n}, \cdots,
$$

where $k_{0}$ is an arbitrary number of the order of the meson mass $\mu$ and $\Lambda$ is an arbitrary number much larger than either $k_{0}$ or $\mu$. This is the only modification we make; the integrands in Eq. (1) are unchanged. What we have done is to chop out most of momentum space leaving an infinite set of well-separated intervals. We shall do this for both integrals in $H$, so that in our model, mesons with momenta outside our intervals do not exist.

We must ask immediately what are the advantages and disadvantages of our model compared with the original fixed source Hamiltonian. First, we want to make clear that our model Hamiltonian is not supposed to be an approximation to the complete fixed source Hamiltonian: We consider it as a completely new Hamiltonian. As to possible disadvantages, the important question is whether our model Hamiltonian has the same renormalization problems as the original fixed source theory. It is easy to see that it does. To prove this we look at the nature of the divergences for the original fixed source theory. In perturbation theory these divergences take the form of integrals over momentum. A linearly divergent integral will be of the form

$$
\int_{\mathbf{k}} k^{-2}
$$

and a logarithmically divergent integral will be of the form

$$
\int_{\mathrm{k}} k^{-3}
$$

In each case the integrand behaves as $k^{-2}$ or $k^{-3}$ only if $k$ is much larger than the meson mass $\mu$. In our model the integrals over all $k$ are replaced by integrals over our sequence of intervals. For example, the logarithmic divergence becomes

$$
\sum_{n=1}^{\infty} \int_{\mathbf{k}} k^{-3}
$$

with the integrals being over $\frac{1}{2} \Lambda^{n}<k<\Lambda^{n}$. These integrals are independent of $n$, and therefore the infinite sum diverges. So our model Hamiltonian has the divergence difficulties that we wish to investigate.
The advantage of our model Hamiltonian is that we can analyze it. This is because the mesons in different momentum intervals differ in their energies by orders of magnitude. For example a meson of momentum $\Lambda^{n}$ has an energy $\Lambda^{n}$, while a meson of momentum $\Lambda^{n-1}$ has energy $\Lambda^{n-1}$. This suggests that we can treat the mesons of momentum $\Lambda^{n-1}$ or less as a perturbation with respect to mesons of momentum of order $\Lambda^{n}$. This idea is the basis of our analysis.

## III. ANALYSIS OF MODEL HAMILTONIAN

In this section we shall present an analysis of our model Hamiltonian (defined in Sec. II). We use the method just proposed : mesons of momenta $\Lambda^{n-1}$ or less are regarded as a perturbation relative to mesons of momentum of order $\Lambda^{n}$. Evidently, we cannot carry out this program unless we cut off our Hamiltonian at some finite value of $n$. Therefore, instead of working with the model Hamiltonian defined in Sec. II, we now define a sequence of Hamiltonians. We first define " $H_{\text {lab }}$ " to be the Hamiltonian of Eq. (1) with the integral being only
over the interval $0<k<k_{0}$. We define $H_{1}$ to include integration over the first two intervals: $0<k<k_{0}$ and $\frac{1}{2} \Lambda<k<\Lambda$. The remaining Hamiltonians $H_{n}$ are defined analogously: for $H_{n}$ we integrate over all our intervals through $\frac{1}{2} \Lambda^{n}<k<\Lambda^{n}$. We shall analyze each of the Hamiltonians $H_{n}$, and then try to recover our original model Hamiltonian as a limit of the $H_{n}$ as $n$ becomes infinite.

We first study $H_{\text {lab. }}$. It is equivalent to the complete fixed source Hamiltonian with a strong cutoff. ${ }^{3}$ There are no problems of renormalization in $H_{\text {lab }}$. One can solve it when $g_{0}$ is small by perturbation theory. One can solve it when $g_{0}$ is very large by Wentzel's strong coupling approximation. ${ }^{4}$ If $g_{0}$ is of order 1 , one may use variational methods (for example, the Tomonaga approximation $\left.{ }^{5}\right)$. One can also estimate both the freemeson energy and the interaction energy by our crude methods and discover that both are of order $\mu .{ }^{6}$ So $H_{\text {lab }}$ is a Hamiltonian which we understand and can solve to reasonable accuracy.

Secondly, we study the Hamiltonian $H_{1}$. The idea here is to divide $H_{1}$ into an unperturbed Hamiltonian $H_{0}$ and a perturbation $H_{I}$. The unperturbed Hamiltonian $H_{0}$ is simply the Hamiltonian $H$ of Eq. (1) with the integrals restricted to the single interval $\frac{1}{2} \Lambda<k<\Lambda$. The perturbation is the same as $H_{\text {lab }}$. Note that our method of dividing $H_{1}$ into an unperturbed part and a perturbation has nothing to do with the size of the coupling constant $g_{0}$; both the free-meson energy and the interaction energy get split up in our procedure. Both terms contribute to the unperturbed Hamiltonian $H_{0}$, and both contribute to the perturbation $H_{I}$.

To analyze $H_{0}$ it is convenient to introduce a scale transformation in order to eliminate (or more correctly, to determine) its dependence on $\Lambda$. First we make a change of variable in the integral, letting

$$
\begin{equation*}
\mathbf{k}=\Lambda \mathbf{p} \tag{3}
\end{equation*}
$$

so that the range of $\mathbf{p}$ is $\frac{1}{2}<p<1$. Next we introduce new creation and destruction operators depending on $\mathbf{p}$ instead of $\mathbf{k}$ :

$$
\begin{array}{ll}
a_{\mathrm{k}}^{\dagger}=\Lambda^{-3 / 2} A_{\mathrm{p}}^{\dagger}, & a_{\mathrm{k}}=\Lambda^{-3 / 2} A_{\mathrm{p}}, \\
b_{\mathrm{k}}^{\dagger}=\Lambda^{-3 / 2} B_{\mathrm{p}}^{\dagger}, & b_{\mathrm{k}}=\Lambda^{-3 / 2} B_{\mathrm{p}} . \tag{5}
\end{array}
$$

The purpose of the factors $\Lambda^{-3 / 2}$ is to ensure that the commutation relations of the $A_{\mathrm{p}}$ etc. will not involve

[^3]$\Lambda$. It follows from Eq. (2) that
\[

$$
\begin{align*}
{\left[A_{\mathrm{p}}, A_{\mathbf{p}^{\prime}}{ }^{\prime}\right]=} & {\left[B_{\mathbf{p}}, B_{\mathbf{p}^{\prime}} \dagger\right] } \\
& =\Lambda^{3}(2 \pi)^{3} \delta^{3}\left(\boldsymbol{\Lambda} \mathbf{p}-\Lambda \mathbf{p}^{\prime}\right)=(2 \pi)^{3} \delta^{3}\left(\mathbf{p}-\mathbf{p}^{\prime}\right) . \tag{6}
\end{align*}
$$
\]

Finally, we have

$$
\begin{equation*}
\omega_{\mathrm{k}} \approx k=\Lambda p . \tag{7}
\end{equation*}
$$

Now one obtains

$$
\begin{equation*}
H_{0}=\Lambda H_{s} \tag{8}
\end{equation*}
$$

with

$$
\begin{align*}
H_{\mathrm{s}}=\int_{\mathrm{p}} p\left\{A_{\mathrm{p}}^{\dagger} A_{\mathrm{p}}\right. & \left.+B_{\mathrm{p}}^{\dagger} B\right\}+g_{0} \int_{\mathrm{p}}(2 p)^{-1 / 2} \\
& \times\left\{\left(A_{\mathrm{p}}+B_{\mathrm{p}}^{\dagger}\right) \tau^{+}+\left(A_{\mathrm{p}}^{\dagger}+B_{\mathrm{p}}\right) \tau^{-}\right\} \tag{9}
\end{align*}
$$

with $\frac{1}{2}<p<1$ in the integrals. The Hamiltonian $H_{s}$ is very similar to $H_{\text {lab }}$; it also has no renormalization problems and can be solved at least roughly for any value of $g_{0}$. It is worth pointing out the importance of the finite lower limit $\frac{1}{2}$ on the momentum $p$ in $H_{s}$. If this limit were replaced by 0 , the charge renormalization integral would be

$$
\int_{0}^{1} p^{-3} d^{3} p
$$

which diverges at the lower limit. There is no mass $\mu$ in $H_{s}$ to provide a low-momentum cutoff for this integral.
The Hamiltonian $H_{s}$ has two symmetries. First of all, it conserves charge, if we think of the proton source level as having charge 1 and the neutron level having charge zero. Secondly, it is invariant to the exchange $\pi^{+} \leftrightarrow \pi^{-}, p \leftrightarrow n$. There is no state of definite charge which is invariant to this transformation, for such a state must have charge $\frac{1}{2}$. Therefore the ground state of $H_{s}$ must be at least doubly degenerate. We shall assume that the ground states have the same quantum numbers as the source levels. This is known to be true for very weak or very strong coupling. We can call these ground states $|P\rangle$ and $|N\rangle$; the source levels with no mesons present will be referred to as $|p\rangle$ and $|n\rangle$.

In addition to the ground states, the Hamiltonian $H_{s}$ has excited states, which are either discrete levels usually called "isobars," or continuum levels which are meson-scattering states. We shall not need to know anything about these states except the energy spacing between the ground state and the first excited state. This spacing is of order 1 unless $g_{0}$ is very large, in which case it is of order $g_{0}{ }^{-2} .{ }^{4}$ The latter case will be ignored for the present.

We have now completed our discussion of $H_{0}$ and can consider the effect of the perturbation $H_{\text {lab }}$. Since the first excited state of $H_{0}$ is of order $\Lambda$ above the ground state, we can consider $H_{\text {lab }}$ as a perturbation provided it is much smaller than $\Lambda$. But according to our previous analysis of $H_{\text {lab }}$ by itself, it is of order $\mu$ and therefore indeed much smaller than $\Lambda$. However $H_{0}$ has a degenerate ground state, so we must use degenerate perturbation theory. In fact, the ground state of $H_{0}$ is
highly degenerate, for to the ground states $|P\rangle$ and $|N\rangle$ of $H_{s}$ we can add any number of "laboratory" mesons (mesons of momenta $<k_{0}$ ) without changing the eigenvalue of $H_{0}$.

Let us now examine the total Hamiltonian $H_{0}+H_{\text {lab }}$ in the lowest order of degenerate perturbation theory. We shall consider only those states which reduce to the ground state of $H_{s}$ in the absence of the perturbation. In lowest order, the energies and eigenstates are determined by an effective Hamiltonian " $H_{\text {eff }}$ " obtained by restricting the full Hamiltonian to the subspace of states spanned by $|P\rangle$ and $|N\rangle$ plus arbitrary numbers of laboratory mesons. The laboratory meson creation and destruction operators are unaffected by this restriction. The unperturbed Hamiltonian $H_{0}$ becomes simply a constant $\Lambda E_{0}$ where $E_{0}$ is the ground-state energy of $H_{8}$. The $\tau$ matrices become operators which act only on the two levels $|P\rangle$ and $|N\rangle$, since they do not affect the laboratory mesons. In particular $\tau^{+}$has only one nonzero matrix element

$$
\begin{equation*}
\alpha=\langle P| \tau^{+}|N\rangle . \tag{10}
\end{equation*}
$$

The other three matrix elements are 0 by charge conservation. We can choose the relative phases of $|P\rangle$ and $|N\rangle$ so that $\alpha$ is positive. We can restate this result by introducing a raising operator $\tau_{R^{+}}$for the states $|P\rangle$ and $|N\rangle$ :

$$
\begin{equation*}
|P\rangle=\tau_{R^{+}}|N\rangle \tag{11}
\end{equation*}
$$

Then, in the restricted Hamiltonian $H_{\text {eff, }} \tau^{+}$becomes $\alpha \tau_{R^{+}}$, likewise $\tau^{-}$becomes $\alpha \tau_{R^{-}}$. Note that if we had not restricted ourselves to the two ground states of $H_{s}$, but allowed excited states as well, then $\tau^{+}$would have been a much more complicated matrix connecting $|P\rangle$ and $|N\rangle$ to all the excited states of $H_{0}$ of appropriate charge.

The effective Hamiltonian now reads

$$
\begin{array}{r}
H_{\text {eff }}=\Lambda E_{0}+\int_{\mathrm{k}} \omega_{\mathrm{k}}\left\{a_{\mathrm{k}^{\dagger}} a_{\mathrm{k}}+b_{\mathrm{k}}^{\dagger} b_{\mathrm{k}}\right\}+g_{0} \alpha \int_{\mathbf{k}}\left(2 \omega_{\mathbf{k}}\right)^{-1 / 2} \\
 \tag{12}\\
\times\left\{\left(a_{\mathrm{k}}+b_{\mathbf{k}}^{\dagger}\right) \tau^{+}+\left(a_{\mathbf{k}}^{\dagger}+b_{\mathrm{k}}\right) \tau^{-}\right\}
\end{array}
$$

with the integral over $k$ restricted to $0<k<k_{0}$.
Evidently, $H_{\text {eff }}$ is the same as the original $H_{\text {lab }}$ except for an additive constant $E_{0}$ and, more importantly, a change in coupling constant from $g_{0}$ to $g_{0} \alpha$. This is hardly a surprise. What we have found is that when the mesons of momentum of order $\Lambda$ (" $\Lambda$ mesons") are added to $H_{\text {lab }}$ to form $H_{1}$, the effect on the lowest energy levels (energies small compared with $\Lambda$ ) is simply what one would predict from renormalization theory: the ground-state energy is shifted and the coupling constant is modified.
We can now carry over our previous discussion of $H_{\text {lab }}$ to show that the energy levels of $H_{\text {eff }}$ differ from the ground-state energy of $H_{0}$ by an amount of order $\mu$, as expected, and the perturbation analysis is justified. We must except from this remark the eigenstates of
$H_{\text {eff }}$ which involve a large number of free laboratory mesons and therefore a large energy; I presume that we can restrict our attention to states with not too many mesons.
It is now a trivial matter to analyze the Hamiltonian $H_{n}$. This Hamiltonian involves mesons in all intervals up to the interval $\frac{1}{2} \Lambda^{n}<k<\Lambda^{n}$. Now the mesons with momenta of order $\Lambda^{n}$, interacting with the source, form the unperturbed system, and the mesons of momenta $\Lambda^{n-1}$ or less are treated as a perturbation. The unperturbed Hamiltonian can again be analyzed by means of a scale transformation, which reduces it to the Hamiltonian $H_{s}$ apart from a factor $\Lambda^{n}$. Thus, the full Hamiltonian $H_{n}$ becomes

$$
\begin{equation*}
H_{n}=\Lambda^{n} H_{s}+H_{n-1} . \tag{13}
\end{equation*}
$$

Let us indicate explicitly that these Hamiltonians depend on a coupling constant $g_{0}$ :

$$
\begin{equation*}
H_{n}\left(g_{0}\right)=\Lambda^{n} H_{s}\left(g_{0}\right)+H_{n-1}\left(g_{0}\right) . \tag{14}
\end{equation*}
$$

Now we use degenerate perturbation theory just as we did for $H_{1}$. Here we restrict $H_{n}$ to the subspace of states spanned by $|P\rangle$ and $|N\rangle$ plus arbitrary numbers of mesons of momenta $\Lambda^{n-1}$ or less. The result is an effective Hamiltonian of the form

$$
\begin{equation*}
H_{\text {eff }}=\Lambda^{n} E_{0}\left(g_{0}\right)+H_{n-1}\left[g_{0} \alpha\left(g_{0}\right)\right] \tag{15}
\end{equation*}
$$

The constant $\alpha$ is as before the matrix element $\langle P| \tau^{+}|N\rangle ; \alpha$ depends on $g_{0}$ since the ground states $|P\rangle$ and $|N\rangle$ depend on $g_{0}$. The effective Hamiltonian describes only the ground state of $H_{n}$ and those excited states with energy of order $\Lambda^{n-1}$ or less above the ground state.

What we now have is a recursion formula which we can state as follows: The lowest energy levels of the Hamiltonian $H_{n}\left(g_{0}\right)$ are equivalent to the energy levels of $\Lambda^{n} E_{0}\left(g_{0}\right)+H_{n-1}\left[f\left(g_{0}\right)\right]$, where

$$
\begin{equation*}
f\left(g_{0}\right)=\alpha\left(g_{0}\right) g_{0} \tag{16}
\end{equation*}
$$

By repeated use of this recursion formula, we see that we will generate a sequence of coupling constants:

$$
\begin{align*}
g_{n} & =g_{0}, \\
g_{n-1} & =f\left(g_{n}\right), \\
g_{n-2} & =f\left(g_{n-1}\right),  \tag{17}\\
& \vdots \\
g & =f\left(g_{1}\right) .
\end{align*}
$$

The ground state and lowest excited states of $H_{n}$ are described by $H_{\text {lab }}(g)$ apart from a constant; $H_{\text {lab }}$ describes those excited states with energies of order $\mu$ above the ground state. To get excited states with energies of order $\Lambda$ one must solve $H_{1}\left(g_{1}\right)$. To get excited states of energy of order $\Lambda^{2}$ one must solve $H_{2}\left(g_{2}\right)$, and so forth.

We have now completed our analysis of the cutoff Hamiltonians $H_{n}$. We know the order of magnitudes of
the components of $H_{n}$ : mesons of momenta of order $\Lambda^{m}(m \leq n)$ contribute a free energy and an interaction energy of order $\Lambda^{m}$. This fact allows us to solve $H_{n}$ by perturbation techniques, which reduce the problem of solving $H_{n}$ to the problem of solving $H_{6}$ and $H_{\text {lab }}$, both of which can be handled by conventional methods. Our principal result is the appearance of a sequence of coupling constants $g_{m}$; we must be able to solve $H_{s}$ for each of these coupling constants.

## IV. LIMIT OF THE SEQUENCE OF HAMILTONIANS

Now we consider the problem of solving our original model Hamiltonian, which formally is the limit of $H_{n}$ as $n \rightarrow \infty$. We shall solve this problem by redefining our original Hamiltonian. Namely, we require that the coupling constant $g$ in $H_{\text {lab }}$ stay fixed, as $n \rightarrow \infty$, and permit the original coupling constant $g_{0}$ to vary as necessary to achieve this. This procedure is a logical development from intuitive notions of coupling constant renormalization. To keep the ground-state energy finite we also must subtract from $H_{n}$ its ground-state energy: call it $E_{n}$. Then what we propose, precisely, is the following: We define a sequence of coupling constants, starting with a given value of $g$.

$$
\begin{align*}
& g_{1}=f^{-1}(g), \\
& g_{2}=f^{-1}\left(g_{1}\right),  \tag{18}\\
& \\
& \vdots \\
& g_{n}=f^{-1}\left(g_{n-1}\right), \text { etc. }
\end{align*}
$$

where $f^{-1}$ is the function inverse to $f$. We must think of $g$ as being determined "experimentally," as if our Hamiltonian described some physical system. Since all the properties of the low-lying eigenstates are determined by $H_{\text {lab }}$ which depends on $g$, it is no problem to determine $g$. It does not matter, for our purposes, if the function $f^{-1}$ is multiple valued, for whenever one has several possibilities for a particular $g_{n}$, one can appeal to experiment to make the choice. We now define our model Hamiltonian $H$ as

$$
\begin{equation*}
H=\operatorname{Lim}_{n \rightarrow \infty}\left\{H_{n}\left(g_{n}\right)-E_{n}\left(g_{n}\right)\right\} \tag{19}
\end{equation*}
$$

The limit that we propose is a strange one, since the operators $H_{n}$ do not all act in the same Hilbert space; instead, as $n$ increases the number of eigenstates of $H_{n}$ also increases. (Those who quibble because the number of eigenstates is infinite for any $n$ and therefore cannot increase may observe instead that the number of meson degrees of freedom increases with $n$.) We should discuss what quantities actually have limits as $n \rightarrow \infty$. To simplify our discussion, suppose that $H_{n}$ has a discrete spectrum with eigenvalues denoted by $E_{m n}$ and states $\left|E_{m n}\right\rangle$. Then the following quantities have limits as
$n \rightarrow \infty$ :
(a) the eigenvalues $E_{m n}$ (for fixed $m$ );
(b) the matrix elements $\left\langle E_{m n} \mid E_{m^{\prime} n}\right\rangle=\delta_{m m^{\prime}}$;
(c) the matrix elements $\left\langle E_{m n}\right| a_{\mathbf{k}}\left|E_{m^{\prime} n}\right\rangle$;
(d) the ratio of the matrix elements $\left\langle E_{m n}\right| \tau^{+}\left|E_{m^{\prime} n}\right\rangle$; to the matrix elements $\langle P ; n| \tau^{+}|N: n\rangle$;
where $|P ; n\rangle$ and $|\lambda ; n\rangle$ are the two ground states of $H_{n}$. In fact, every quantity listed will be independent of $n$ (for fixed $m, m^{\prime}$, and $\mathbf{k}$ ) if $n$ is sufficiently large. This can be proved as a consequence of our analysis of the Hamiltonians $H_{n}$. One can easily discuss matrix elements of operators other than $a_{\mathrm{k}}$ or $\tau^{+}$.

One is tempted, of course, to ask what happens if we keep $g_{0}$ fixed, in the limiting process, instead of $g$. But this is a pointless question, for, once we have determined $g$ experimentally, we know that the Hamiltonian $H$ must be given by the limit of Eq. (19), and no new information is gained by seeing whether another limiting process gives the same result.
Let us look in more detail at the sequence of coupling constants $g$, $g_{1}$, etc. First, we show that the quantity $\alpha$ is less than one. This is because

$$
\begin{equation*}
\tau^{+} \tau^{-}+\tau^{-} \tau^{+}=1 \tag{20}
\end{equation*}
$$

Let $|x\rangle$ be a complete set of eigenstates of $H_{s}$. Then

$$
\begin{equation*}
\sum_{x}\left\{\langle P| \tau^{+}|x\rangle\langle x| \tau^{-}|P\rangle+\langle P| \tau^{-}|x\rangle\langle x| \tau^{+}|P\rangle\right\}=1 \tag{21}
\end{equation*}
$$

Every term in the sum over $x$ is positive, and one of the terms is $\alpha^{2}$, namely the term with $|x\rangle=|N\rangle$. So $\alpha^{2}$ is less than or equal to one. It is trivial to show that $\alpha$ cannot be equal to one (except when $g_{0}$ is 0 ).

Because $\alpha$ is less than one, we have

$$
\begin{equation*}
f\left(g_{0}\right)<g_{0} . \tag{22}
\end{equation*}
$$

Secondly, we note that for $g_{0} \gg 1, f\left(g_{0}\right) \approx \frac{1}{2} g_{0}$. This is a consequence of Wentzel's strong coupling theory ${ }^{4}$ and will not be proved here. For small $g_{0}$ we have $f\left(g_{0}\right)$ $=g_{0}+$ order $g_{0}{ }^{3}$. If $f\left(g_{0}\right)$ is a continuous function of $g_{0}$, it follows from these results that the equation $g_{n-1}=f\left(g_{n}\right)$ has at least one solution $g_{n}$ for any value of $g_{n-1}$. Hence there exists a complete sequence $\left\{g_{n}\right\}$ for any value of $g$. Furthermore, this sequence is increasing. The sequence cannot have a finite upper bound for if so it would approach a limit point, which we can call $g_{0}$; and we would have

$$
g_{0}=f\left(g_{0}\right)
$$

which is impossible.
So whatever the value of $g$, the sequence of coupling constants $g_{n}$ is an increasing sequence which approaches infinity as $n$ becomes large.

This is unfortunate, for when $n$ is so large that $g_{n}{ }^{2}$ is of order $\Lambda$, our perturbation analysis ceases to be valid. This is because the first excited (isobar) state of $H_{s}$ has an energy only $g_{n}{ }^{-2}$ above the ground state when $g_{n}$ is large; when multiplied by $\Lambda^{n}$, this gives an energy of
order $\Lambda^{n-1}$, the same order as the terms we treated as a perturbation. So to solve $H_{n}$ for very large $g_{n}$, the whole set of isobar states of $H_{s}$ must be taken into account in the perturbation analysis, rather than just the two ground states. This is an involved task; we do not attempt it here. Because of this difficulty, the analysis of our model is incomplete. Our analysis would also be incorrect if, for values of $g_{0}$ of order 1 , the ground states of $H_{s}$ did not have the same charge as the two source levels.

## V. AN OFFSPRING OF THE LEE MODEL

Another Hamiltonian that can be analyzed in the manner of Secs. III and IV is derived from the Lee model. ${ }^{1}$ The Lee model Hamiltonian can be obtained from the fixed source Hamiltonian of Eq. (1) by dropping the $\pi^{-}$meson terms (i.e., we suppress $b_{\mathbf{k}}$ and $b_{\mathbf{k}^{\dagger}}{ }^{\dagger}$ ). We can then construct a model Hamiltonian by replacing the full momentum integration by an integral over our sequences of intervals. We can then perform an analysis analogous to Secs. III and IV. We describe this analysis briefly.

First of all we have charge conservation, but we no longer have charge symmetry. To ensure that the physical proton state has the same energy as the physical neutron state, we let the bare proton state have an undetermined energy. This means adding a term $E \tau^{++}$to the Hamiltonian, where $\tau^{++}$has the properties:

$$
\begin{align*}
& \tau^{++}|p\rangle=|p\rangle,  \tag{23}\\
& \tau^{++}|n\rangle=0 .
\end{align*}
$$

( $|p\rangle$ and $|n\rangle$ are the two source levels.)
Now we define, as before, a sequence of Hamiltonians $H_{n}$. The added term for $H_{n}$ will be denoted by $E_{n} \tau^{++}$. We break $H_{n}$ into an unperturbed and perturbed Hamiltonian, as before. The added term must likewise be split, say

$$
\begin{equation*}
E_{n}=\Lambda^{n} E_{0}+\epsilon_{n} \tag{24}
\end{equation*}
$$

Then after performing the scaling as in Sec. III, we obtain

$$
\begin{equation*}
H_{n}=\Lambda^{n} H_{s}+H_{n-1} \tag{25}
\end{equation*}
$$

where the added term in $H_{n-1}$ is $\epsilon_{n} \tau^{++}$, and

$$
\begin{align*}
H_{s}=E_{0} \tau^{++}+\int_{\mathrm{p}} p A_{\mathrm{p}}^{\dagger} A_{\mathrm{p}}+g_{0} \int_{\mathrm{p}} & (2 p)^{-1 / 2} \\
& \times\left\{A_{\mathrm{p}} \tau^{+}+A_{\mathrm{p}}^{\dagger} \tau^{-}\right\} \tag{26}
\end{align*}
$$

with $\frac{1}{2}<p<1$. We choose $E_{0}$ so that the ground states $|P\rangle$ and $|N\rangle$ of $H_{s}$ are degenerate. These states and $E_{0}$ are easily derived; we obtain

$$
\begin{align*}
& |P\rangle=Z\left\{|p\rangle-g_{0} \int_{\mathrm{p}}(2)^{-1 / 2} p^{-3 / 2} A_{\mathrm{p}}^{\dagger}|N\rangle\right\},  \tag{27}\\
& |\cdot\rangle\rangle=|n\rangle \tag{28}
\end{align*}
$$

$$
\begin{align*}
& E_{0}=\frac{1}{2} g_{0} \int_{\mathrm{p}} p^{-2},  \tag{29}\\
& Z^{-2}=1+\frac{1}{2} g_{0} \int_{\mathrm{p}} p^{-3}, \tag{30}
\end{align*}
$$

with $\frac{1}{2}<p<1$ in all integrals.
We now solve $H_{n}$ by degenerate perturbation theory. We restrict $H_{n}$ to the subspace of states spanned by $|P\rangle$ and $|N\rangle$ plus mesons of momentum $\Lambda^{n-1}$ or less. Under this restriction

$$
\begin{align*}
\tau^{+} & \rightarrow \alpha \tau_{R^{+}}  \tag{31}\\
\tau^{++} & \rightarrow \beta \tau_{R^{++}} \tag{32}
\end{align*}
$$

where $\tau_{R}{ }^{++}|P\rangle=|P\rangle, \tau_{R^{++}}|\lambda\rangle=0$, and

$$
\begin{align*}
& \alpha=\langle P| \tau^{+}|N\rangle=Z,  \tag{33}\\
& \beta=\langle P| \tau^{++}|P\rangle=Z^{3} . \tag{34}
\end{align*}
$$

Thus, we obtain an effective Hamiltonian which is just $H_{n-1}\left(\alpha g_{0}\right)$; the added term is $E_{n-1} \tau_{R}{ }^{++}$with $E_{n-1}$ set equal to $Z^{2} \epsilon_{n}$.
It is now easy to see that exactly as in Sec. IV, we want to define a sequence of coupling constants $g_{n}$. In the present problem we will have

$$
\begin{equation*}
g_{n}=f^{-1}\left(g_{n-1}\right), \tag{35}
\end{equation*}
$$

where $f^{-1}$ is the inverse function for the function $f$ given by

$$
\begin{equation*}
f\left(g_{0}\right)=Z\left(g_{0}\right) g_{0} \tag{36}
\end{equation*}
$$

But now for large $g_{0}, Z\left(g_{0}\right)$ behaves as $g_{0}{ }^{-1}$. Hence $f\left(g_{0}\right)$ has a finite upper bound. If $g_{n-1}$ is larger than this upper bound, $g_{n}$ does not exist. But since $Z\left(g_{0}\right)$ is always smaller than one, the sequence $\left\{g_{n}\right\}$ is increasing; there will inevitably be a finite value of $n$, above which $g_{n}$ does not exist.

## VI. COMPARISON WITH QUANTUM ELECTRODYNAMICS

We want to compare the results of our analysis of two model Hamiltonians with the analysis of quantum electrodynamics by Gell-Mann and Low. First we shall restate the ideas of Gell-Mann and Low in a form suitable for our purposes. For background the reader is referred to their paper. ${ }^{7}$ Let $e$ be the physical electron charge, and let $m$ be the physical electron mass. ${ }^{8}$ Let $d_{C}\left(k^{2} / m^{2}, e^{2}\right)$ be the photon propagator apart from a factor $k^{-2}$; we require the customary renormalization condition

$$
\begin{equation*}
d_{C}\left(0, e^{2}\right)=1 \tag{37}
\end{equation*}
$$

[^4]Now define a one-parameter family of coupling constants $e_{\lambda}$ as

$$
\begin{equation*}
e_{\lambda}^{2}=e^{2} d_{C}\left(\lambda^{2} / m^{2}, e^{2}\right) \tag{38}
\end{equation*}
$$

These constants $e_{\lambda}$ depend on $e$, but we shall not need to indicate this explicitly. Gell-Mann and Low introduce a fancy subtraction scheme for renormalizing perturbation theory, in which the subtractions are performed at momentum $\lambda$ instead of momentum 0 . The result of their scheme is
(1) their renormalized coupling constant is $e_{\lambda}$, not $e$;
(2) their photon propagator is normalized to 1 at $k^{2}=\lambda^{2}$ instead of 0 ;
(3) their photon propagator is a function which may be written $d\left(k^{2} / \lambda^{2}, m^{2} / \lambda^{2}, e_{\lambda}^{2}\right)$;
(4) their photon propagator is related to the usual one according to the equation

$$
\begin{equation*}
e^{2} d_{C}\left(k^{2} / m^{2}, e^{2}\right)=e_{\lambda}^{2} d\left(k^{2} / \lambda^{2}, m^{2} / \lambda^{2}, e_{\lambda}^{2}\right) \tag{39}
\end{equation*}
$$

Alternatively, one can use Eq. (38) to express $e^{2}$ as a function of $e_{\lambda}{ }^{2}$, and then define the function

$$
d\left(k^{2} / \lambda^{2}, m^{2} / \lambda^{2}, e_{\lambda}{ }^{2}\right)
$$

by Eq. (39).
(5) They argue that the function $d\left(k^{2} / \lambda^{2}, m^{2} / \lambda^{2}, e_{\lambda}^{2}\right)$ does not depend on $m$ when $m^{2} / \lambda^{2}$ is small.

This result is crucial to their discussion. They illustrate it with an example from fourth-order perturbation theory and propose a reason for it to be true to all orders [see, e.g., preceding Eq. (5.1) of Ref. 2]. Some people tend to regard this result skeptically, and to use this skepticism as an excuse to disregard the consequences which Gell-Mann and Low deduce from it. None of these people have succeeded (to my knowledge) in finding a specific counterexample to this claim of Gell-Mann and Low, although the paper it was made in is now over ten years old.

The specific equation we need is obtained as follows. We rewrite Eq. (39) as

$$
\begin{equation*}
e_{k}^{2}=e_{\lambda}^{2} d\left(k^{2} / \lambda^{2}, m^{2} / \lambda^{2}, e_{\lambda}^{2}\right) \tag{40}
\end{equation*}
$$

Differentiate with respect to $k$, holding $\lambda$ and $e_{\lambda}$ fixed (which means that $e$, on which $e_{k}$ also depends, is fixed), and then set $k^{2}=\lambda^{2}$ : we obtain

$$
\begin{equation*}
d\left(e_{\lambda}^{2}\right) / d \lambda=2 \lambda^{-1} \psi\left(m^{2} / \lambda^{2}, e_{\lambda}^{2}\right) \tag{41}
\end{equation*}
$$

where

$$
\begin{equation*}
\psi\left(\frac{m^{2}}{\lambda^{2}}, e_{\lambda}^{2}\right)=\left.e_{\lambda}^{2} \frac{\partial d}{\partial x}\left(x, \frac{m^{2}}{\lambda^{2}}, e_{\lambda}^{2}\right)\right|_{x=1} \tag{42}
\end{equation*}
$$

Now, according to result 5 , for $\lambda \gg m$, the dependence of $\psi$ on $m^{2} / \lambda^{2}$ is negligible, so that to a good approximation $\psi$ depends only on $e_{\lambda}{ }^{2}$. Thus we have

$$
\begin{equation*}
d\left(e_{\lambda}^{2}\right) / d \lambda=2 \lambda^{-1} \psi\left(e_{\lambda}{ }^{2}\right) \tag{43}
\end{equation*}
$$

We can understand this equation by setting up an analogy with the results of the analysis of our model Hamiltonians. The function $e_{\lambda}{ }^{2}$ is analogous to our sequences $\left\{g_{n}\right\}$, and has the same significance: The constant $g_{n}$ determines the solution of the unperturbed part of the Hamiltonian $H_{n}$, which determines the essential features of the behavior of mesons of momentum $\Lambda^{n}$. The constant $e_{\lambda}$ determines the essential features of the behavior of propagators and vertex functions when the momenta are of order $\lambda$; in particular these functions can be computed by a perturbation expansion in $e_{\lambda}$ if $e_{\lambda}$ is small.

We can illustrate this point by studying the photon propagator in perturbation theory. Let $\lambda$ be very large, and let $k$ be of order $\lambda$. Then $d\left(k^{2} / \lambda^{2}, m^{2} / \lambda^{2}, e_{\lambda}^{2}\right)$ takes the form of a double power series in $e_{\lambda}{ }^{2}$ and $\ln \left(k^{2} / \lambda^{2}\right) .{ }^{2}$ For order of magnitude purposes, the coefficients can be set equal to one and we obtain

$$
\begin{align*}
d\left(k^{2} / \lambda^{2}, m^{2} / \lambda^{2}, e_{\lambda}^{2}\right)= & 1+e_{\lambda}^{2}\left[\ln \left(k^{2} / \lambda^{2}\right)+1\right] \\
& +e_{\lambda}^{4}\left[\ln ^{2}\left(k^{2} / \lambda^{2}\right)+\cdots\right]+\cdots . \tag{44}
\end{align*}
$$

Since $\ln \left(k^{2} / \lambda^{2}\right)$ is, by assumption, of order 1 , this series in $e_{\lambda}^{2}$ converges rapidly if $e_{\lambda}^{2} \ll 1$; we can compute $d$ using the perturbation expansion in $e_{\lambda}$ if $e_{\lambda} \ll 1$. This is not true of the function $d_{C}$ expanded in powers of $e$; for large $k$ we have (in order of magnitude)

$$
\begin{align*}
d_{C}\left(k^{2} / m^{2}, e^{2}\right)=1+e^{2}[ & \left.\ln \left(k^{2} / m^{2}\right)+1\right] \\
& +e^{4}\left[\ln ^{2}\left(k^{2} / m^{2}\right)+\cdots\right]+\cdots \tag{45}
\end{align*}
$$

Clearly, this power series converges rapidly only if $e^{2} \ln \left(k^{2} / m^{2}\right) \ll 1$. So the criterion for whether the photon propagator can be computed in perturbation theory, when $k$ is of order $\lambda$, is not $e \ll 1$ but $e_{\lambda} \ll 1$.

Equation (43) for $e_{\lambda}$ is analogous to our recursion formula for $g_{n}$ :

$$
\begin{equation*}
g_{n-1}=f\left(g_{n}\right) \tag{17}
\end{equation*}
$$

This equation determines how $g_{n}$ changes with $n$, given $g_{n}$, while Eq. (43) for $e_{\lambda}$ determines the rate of change of $e_{\lambda}$, given $e_{\lambda}$. We note that there is no question that the function $f\left(g_{n}\right)$ is independent of the meson mass $\mu$; this mass disappeared when we approximated $\omega_{\mathbf{k}}$ by $k$, for $k$ large. Thus, when we go from quantum electrodynamics to a theory which we understand, the question of mass dependence becomes trivial. I suspect the same will be true of quantum electrodynamics when a decent formulation of it becomes available.

Gell-Mann and Low discussed the possible behavior of $e_{\lambda}$ for very large $\lambda$, using Eq. (43). There is a close analogy between this question and the problem of how the sequence $\left\{g_{n}\right\}$ behaves for large $n$. First let us discuss how the sequence $g_{n}$ behaves, in terms of possible types of functions $f$. We show in Fig. 1 three types of functions $f$. The curve labeled A is the function that results from our Lee-model-type Hamiltonian discussed in Sec. V. The curve labeled B results from our fixed source model. The curve labeled C has not been derived

from any model but it is instructive to discuss the behavior of the sequence $\left\{g_{n}\right\}$ for such a function. We have discussed curves A and B previously. If $f$ is given by curve $A$, then the sequence $\left\{g_{n}\right\}$ stops after a finite number of terms. If $f$ is given by curve $B$, then the sequence $\left\{g_{n}\right\}$ exists for all $n$, it is increasing, and it approaches infinity as $n \rightarrow \infty$. Now consider the curve C. We have not shown the whole of this curve because of the enormous number of possibilities for the remaining part of the curve. The part shown has the following properties:
(a) The inverse function $f^{-1}(y)$ is uniquely defined for not too large $y$.
(b) The equation $y=f^{-1}(y)$ has a solution $y=g_{0}$.
(c) If $y<g_{0}$, then $y<f^{-1}(y)<g_{0}$; if $y>g_{0}$, then $y>f^{-1}(y)>g_{0}$.

From these properties it follows that if the low-momentum coupling constant $g$ is smaller than $g_{0}$, then the sequence $\left\{g_{n}\right\}$ will be increasing, and approach $g_{0}$ in the limit $n \rightarrow \infty$. If $g$ is greater than $g_{0}$, but not above the range of the curve C, then the sequence $\left\{g_{n}\right\}$ will be decreasing and again approach the limit $g_{0}$ as $n \rightarrow \infty$.

For both the curves B and C, we see that the sequence $\left\{g_{n}\right\}$ approaches a limit as $n \rightarrow \infty$ which is independent of the low-momentum coupling constant $g$ that we start with. For the curve B , the limit of $g_{n}$ is infinite; for curve C the limit is the finite number $g_{0}$ which solves the equation

$$
\begin{equation*}
g_{0}=f\left(g_{0}\right) \tag{46}
\end{equation*}
$$

We can invent possible forms for the function $\psi(x)$ which are analogous to the three curves $\mathrm{A}, \mathrm{B}$, and C for $f$. These are shown in Fig. 2. To set up the analogy, we note that $g_{n}-g_{n-1}$ constitutes the change in $g_{n}$ as the momentum changes by a factor $\Lambda$; therefore $g_{n}-g_{n-1}$ is roughly analogous to the derivative of $e_{\lambda}{ }^{2}$ with respect to $\ln \lambda$, which is just the function $\psi$. Now for curve A for $f$, the difference $g_{n}-g_{n-1}$ becomes very large as $g_{n} \rightarrow \infty$. By analogy we obtain the curve A of Fig. 2. For the curve B for $f$, the difference $g_{n}-g_{n-1}$ is just a constant times $g_{n}$ when $g_{n}$ is large. Thus we obtain the curve B of Fig. 2 which is linear in $x$ for large $x$. Finally, when $f$ is
of type C, the difference $g_{n}-g_{n-1}$ passes through 0 as $g_{n}$ passes through $g_{0}$; hence the curve C of Fig. 2.

To find $e_{\lambda}{ }^{2}$ as a function of $\lambda$ from Eq. (43), we simply integrate this equation. To simplify matters let $\psi$ be independent of $m$ if $\lambda>m$, and let $e_{\lambda}$ be $e_{m}$ at $\lambda=m$. Then

$$
\begin{equation*}
\ln \left(\lambda^{2} / m^{2}\right)=\int_{e_{m}{ }^{2}}^{e^{2}} d x / \psi(x) \tag{47}
\end{equation*}
$$

For curve A, the integral is finite for $e_{\lambda}^{2} \rightarrow \infty$ which means we get a finite upper bound for $\lambda$, beyond which $e_{\lambda}$ is undefined. For curve $B, \lambda$ increases as $e_{\lambda}^{2}$ increases and $\lambda \rightarrow \infty$ as $e_{\lambda}{ }^{2} \rightarrow \infty$. In this case we can invert our equation to get $e_{\lambda}$ as a function of $\lambda$; we see that $e_{\lambda}{ }^{2}$ increases as $\lambda$ increases and $e_{\lambda}{ }^{2} \rightarrow \infty$ as $\lambda \rightarrow \infty$. For curve C, the integral diverges as $e_{\lambda}^{2} \rightarrow e_{0}{ }^{2}$. For $e_{\lambda}{ }^{2}<e_{0}{ }^{2}$, $\lambda$ increases as $e_{\lambda}{ }^{2}$ increases. Hence, if $e_{m}{ }^{2}<e_{0}{ }^{2}$, the function $e_{\lambda}{ }^{2}$ will be increasing, and approach the limit $e_{0}{ }^{2}$ as $\lambda \rightarrow \infty$.
The curve A will never occur in practice. For if it did occur then $e_{\lambda}$ would not exist for large $\lambda$. But because of our definition of $e_{\lambda}$ [Eq. (38)], the photon propagator also would not exist for large photon momentum $k$. But quantum electrodynamics cannot exist without a photon propagator, and in particular the function $\psi$ will then not exist either. So, instead of the curve A being a possibility, we have the possibility that no consistent quantum electrodynamics exists.
The curves $B$ and $C$ were the specific alternatives considered by Gell-Mann and Low, and led to their alternatives (a) and (b) which we quoted in the Introduction.

## VII. QUALITATIVE ANALYSIS OF COMPLETE FIXED SOURCE HAMILTONIAN

In the Introduction we emphasized the importance of making qualitative estimates of various terms in a Hamiltonian before trying to solve it quantitatively. For this purpose, the creation and destruction operators $a_{k}$ and $a_{k}{ }^{\dagger}$ are a nuisance. In particular, the operator $a_{\mathbf{k}}{ }^{\dagger}$, since it creates a meson in a plane wave state, creates states of infinite norm. Thus, if one tries to estimate the "order of magnitude" of $a_{\mathbf{k}}{ }^{\dagger}$ in the manner

Fig. 2. Plots of the function $\psi(x)$ versus $x$. Three cases are shown: the curves labeled A, B, and C.



Fig. 3. Our division of phase space into cells of unit volume shown in two dimensions (one coordinate and one momentum dimension). Each cell is labeled with its indices $(n, m)$.
we have used, which is to calculate the norm of the state created by $a_{k}{ }^{\dagger}$, we get infinity. So far we have not had to estimate $a_{\mathbf{k}}{ }^{\dagger}$ itself, but only integrals involving $a_{\mathbf{k}}{ }^{\dagger}$ of the form $\int a_{\mathbf{k}}{ }^{\dagger} u(\mathbf{k}) d^{3} k$. If the function $u(\mathbf{k})$ in the integral is normalizable, then the resulting operator creates a normalizable state and estimates are possible. However, in relativistic Hamiltonians, the operators analogous to $a_{\mathbf{k}}{ }^{\dagger}$ are integrated with $\delta$ functions (to ensure momentum conservation) rather than normalizable functions.

What we shall propose here is that the creation and destruction operators, which depend on the continuous variable $\mathbf{k}$, be expanded in terms of a discrete, complete, orthonormal set of functions $u_{l}(\mathbf{k})$. The coefficients of this expansion are themselves creation and destruction operators for mesons in normalizable "wave-packet" states. We obtain

$$
\begin{gather*}
a_{\mathbf{k}}=\sum_{l} u_{l}(\mathbf{k}) a_{l}  \tag{48}\\
a_{\mathbf{k}^{\dagger}}=\sum_{l} u_{l}^{*}(\mathbf{k}) a_{l}^{\dagger} . \tag{49}
\end{gather*}
$$

The operators $a_{l}$ and $a_{l}{ }^{\dagger}$ are operators with the same commutation rules as an infinite set of harmonic oscillator variables:

$$
\begin{align*}
{\left[a_{l}, a_{m}\right] } & =0, \\
{\left[a_{l}, a_{m}^{\dagger}\right] } & =\delta_{l m},  \tag{50}\\
{\left[a_{l} \dagger, a_{m}^{\dagger}\right] } & =0
\end{align*}
$$

where $\delta_{l m}$ is the Kronecker delta.
When a Hamiltonian involving $a_{\mathrm{k}}$ and $a_{\mathbf{k}}{ }^{\dagger}$ is reexpressed in terms of the operators $a_{l}$ and $a_{l}{ }^{\dagger}$, we obtain a Hamiltonian for an infinite set of coupled, perhaps anharmonic, oscillators. It seems to be easier to analyze such Hamiltonians qualitatively than to analyze Hamiltonians containing the operators $a_{\mathbf{k}}$ and $a_{\mathbf{k}}{ }^{\dagger}$. We shall illustrate this shortly by discussing the full fixed source Hamiltonian of Eq. (1).

There are very many complete orthonormal sets of functions $u_{l}(\mathbf{k})$. However, the complete sets of interest to us will have to satisfy some restrictions, which we shall now discuss. The restrictions will be imposed with the aim of making the self-interactions of the individual oscillators more important than interactions between
different oscillators. We shall not try to demonstrate that our choice of restrictions accomplishes this until we discuss a particular Hamiltonian.

Our first requirement is that all the functions $u_{l}(\mathbf{k})$ be good wave packet functions. By this we mean that, if $\Delta k_{l}$ is the momentum width of the function $u_{l}(\mathbf{k})$, and $\Delta x_{l}$ is the width of the Fourier transform $v_{l}(\mathbf{x})$ of $u_{l}$, then the product $\Delta x_{l} \Delta k_{l}$ is near the lower limit set by the Heisenberg uncertainty principle

$$
\begin{equation*}
\Delta x_{l} \Delta k_{l} \approx \frac{1}{2} \tag{51}
\end{equation*}
$$

For order of magnitude purposes, each function $u_{l}$ can now be thought of as occupying a "cell" in phase space, of unit volume. If the set of functions $u_{l}$ is complete, then the total volume occupied by the functions $u_{l}$ must fill all of phase space. The orthogonality of the different $u_{l}$ means that regions occupied by different $u_{l}$ should not overlap.
These conditions are supposed to be understood qualitatively. That is, the function $u_{l}(\mathbf{k})$ and $v_{l}(\mathbf{x})$ are certainly not 0 outside their assigned regions in momentum space and position space, respectively but, hopefully, their "tails" outside these regions are small and go rapidly to zero as one goes away from the assigned region. I have been unable to construct a specific example of a complete orthonormal set of functions satisfying these requirements: the functions I construct always have long tails in position space and no tails in momentum space, or vice versa. But the conditions proposed represent, I think, the physics of a complete orthonormal set of wave functions, and if so it is reasonable to proceed as if such a set of functions did exist.

We still have to specify how phase space is to be divided into unit cells; we shall presume that we may do this arbitrarily. The particular division that we shall adopt is one in which momentum space is divided evenly on a logarithmic scale. To be precise, we divide momentum space into an infinite number of nested spherical shells, the $n$th shell being

$$
\begin{equation*}
2^{n-1}<k<2^{n} \tag{52}
\end{equation*}
$$

Now for each shell separately we divide position space linearly into cubes of the appropriate size. For the $n$th shell the volume in momentum space is of order $2^{3 n}$ so the volume of a cell in position space will be of order $2^{-3 n}$. So in order of magnitude the cubes may be defined as

$$
\begin{equation*}
m_{i}-\frac{1}{2} \leq 2^{n} x_{i}<m_{i}+\frac{1}{2}, \tag{53}
\end{equation*}
$$

where the $x_{i}$ are position space coordinates and $m_{1}, m_{2}$, and $m_{3}$ are arbitrary integers. Our division of phase space is illustrated in Fig. 3.

Now we shall assume that there exists a complete orthonnormal set of functions $u_{l}(\mathbf{k})$ corresponding to the division of phase space just described. To be precise, $u_{l}(\mathbf{k})$ will be labeled by four indices ( $n, m_{1}, m_{2}, m_{3}$ ) which specify the location of the corresponding phase space cell; let us use the single symbol $l$ to stand for
these four indices. From the normalization requirement

$$
\begin{equation*}
\int_{\mathbf{k}}\left|u_{l}(\mathbf{k})\right|^{2}=\int\left|v_{l}(\mathbf{x})\right|^{2} d^{3} x=1 \tag{54}
\end{equation*}
$$

we can compute the order of magnitude of $u_{l}$ and $v_{l}$. The volume of the $n$th shell in momentum space is of order $2^{3 n}$, and the volume of the corresponding cell in position space is of order $2^{-3 n}$. Hence, inside these regions we have

$$
\begin{align*}
& \left|u_{l}(\mathbf{k})\right| \sim 2^{-3 n / 2} \\
& \left|v_{l}(\mathbf{x})\right| \sim 2^{+3 n / 2} \tag{55}
\end{align*}
$$

Outside these regions, $u_{l}$ and $v_{l}$ are smaller.
To illustrate the usefulness of expanding the creation and destruction operators in terms of oscillator variables, we shall examine the Hamiltonian of Eq. (1) (with no limitations on $k$ in the integrals). First we must express $H$ in terms of the operators $a_{l}$ and $a_{l}{ }^{\dagger}$. We must introduce corresponding oscillator variables also for $b_{k}$ and $b_{\mathbf{k}}{ }^{\dagger}$. Then $H$ has the form
$H=\sum_{j} \sum_{l} C_{j l}\left(a_{j}^{\dagger} a_{l}+b_{j}{ }^{\dagger} b_{l}\right)+g_{0} \sum_{l} D_{l}\left(a_{l} \tau^{+}+b_{l} \tau^{-}\right)$

$$
\begin{equation*}
+g_{0} \sum_{l} D_{l}^{*}\left(b_{l}{ }^{\dagger} \tau^{+}+a_{l}^{\dagger} \tau^{-}\right) \tag{56}
\end{equation*}
$$

where

$$
\begin{align*}
C_{j i} & =\int_{\mathbf{k}} \omega_{\mathbf{k}} u_{j} *(\mathbf{k}) u_{l}(\mathbf{k}),  \tag{57}\\
D_{l} & =\int_{\mathbf{k}}\left(2 \omega_{\mathbf{k}}\right)^{-1 / 2} u_{l}(\mathbf{k}) . \tag{58}
\end{align*}
$$

The Hamiltonian $H$ is now the Hamiltonian of an infinite number of oscillators coupled to each other and to a two-level source.

To simplify our Hamiltonian, we examine the coefficients $C_{j l}$ and $D_{l}$. Except for the factor $\omega_{\mathbf{k}}$, the matrix $C$ would be diagonal because the functions $u_{l}$ are orthogonal. Even with the factor $\omega_{\mathbf{k}}$, the off-diagonal terms should be small: If the momentum shells $n_{j}$ and $n_{l}$ are distinct, then the functions $u_{j}$ and $u_{l}$ do not overlap very much. If $u_{j}$ and $u_{l}$ are in the same momentum shell but different spatial cells, this fact will be reflected in a rapid variation in the phase of $u_{j}^{*} u_{l}$ as a function of $\mathbf{k}$, which, again, makes the integral small. So, for a first approximation, we keep only the diagonal terms $C_{l l}$ of the matrix $C$. These we can estimate to be a mean energy times the normalization integral. In the $n$th shell the mean energy is either about $\mu$, or about $2^{n}$, whichever is larger. Let $\mu$ be 1 for convenience. So

$$
\begin{array}{cc}
C_{l l} \sim 2^{n} &  \tag{59}\\
\sim 1 & (n>0) \\
\sim 1 & \\
(n<0) .
\end{array}
$$

Now consider $D_{l}$. The function $u_{l}(\mathbf{k})$ carries a $\mathbf{k}$ dependent phase factor associated with the center of the appropriate cell in position space, unless this center is at the origin. Thus $D_{l}$ will be small unless the $m_{i}$ in $l$ are
all 0: $l=(n, 0,0,0)$. These latter $D_{l}$ we denote by $D_{n}$. They are easily estimated from the known volume of the $n$th shell and the known order of magnitude of $u_{l}(\mathbf{k})$; we get

$$
\begin{align*}
D_{n} & \sim 2^{n} & & (n>0) \\
& \sim 2^{3 n / 2} & & (n<0) . \tag{60}
\end{align*}
$$

We have now as a rough approximation

$$
\begin{align*}
H \sim \sum_{l} C_{l l}\left(a_{l}^{\dagger} a_{l}+b_{l}{ }^{\dagger} b_{l}\right) & +g_{0} \sum_{n} D_{n}\left(a_{n} \tau^{+}+b_{n} \tau^{-}\right) \\
& +g_{0} \sum_{n} D_{n}^{*}\left(b_{n}^{\dagger} \tau^{+}+a_{n}^{\dagger} \tau^{-}\right), \tag{61}
\end{align*}
$$

where the operators $a_{n}$, etc. are the operators $a_{l}$, etc. with $m_{i}=0$. In this approximation the free-meson field has been replaced by independent harmonic oscillators for each phase space cell, with a frequency depending only on the mean momentum of the cell. The interaction of the meson field with the source has been replaced by an interaction of those oscillators located at the origin (where the source is) with the source. The remaining terms of the original Hamiltonian are to be considered as a perturbation.

To get orders of magnitude, we first neglect the interaction with the source, in Eq. (61). Then we have independent harmonic oscillators. In this case the orders of magnitude are evidently

$$
\begin{equation*}
a_{l} \sim 1, \quad a_{l}^{\dagger} \sim 1 \tag{62}
\end{equation*}
$$

for all $l$. The energy contributed by the $l$ th oscillator is of order $2^{n}$ or 1 whichever is larger. As in our model, the oscillators associated with large momentum in phase space contribute most to the energy. Now suppose the order of magnitudes of $a_{l}$ and $a_{l}{ }^{\dagger}$ are unchanged by the presence of the source. The matrices $\tau^{+}$and $\tau^{-}$are of order 1 at most [this follows from the sum rule, Eq. (20)]. The energy due to the interaction of $n$th shell mesons with the source is therefore of order $2^{n}$, if $n>0$, or $2^{3 n / 2}$ if $n<0$. This is of the same order as the freemeson energy for $n>0$, and much smaller for $n<0$. Since the interaction energy of a meson is not greater than its free energy, we do not expect the interaction to change the order of magnitude of $a_{l}$ and $a_{l}{ }^{\dagger}$. [This discussion fails if $g_{0}$ is very large, for then the interaction energy is large and the orders of magnitude do change (compare Ref. 6).]
Let us consider now the problem of finding the ground state of $H$. First, we ignore the oscillators not located at the origin since in our rough approximation these are uncoupled. For the oscillators coupled to the source we use the same method as we used on our model of Sec. II. We cut off the Hamiltonian at a large finite value of $n$, say $N$. Then the oscillators associated with shells within the $N$ th shell are treated as a perturbation on the $N$ th shell oscillators. Thus, our unperturbed Hamiltonian $H_{0}$ is

$$
\begin{align*}
H_{0}=\omega_{N}\left(a_{N}^{\dagger} a_{N}+b_{N} \dagger b_{N}\right) & +g_{0} D_{N}\left(a_{N} \tau^{+}+b_{N} \tau^{-}\right) \\
& +g_{0} D_{N}^{*}\left(b_{N}^{\dagger} \boldsymbol{\tau}^{\dagger}+a_{N}^{\dagger} \boldsymbol{\tau}^{-}\right), \tag{63}
\end{align*}
$$

where $\omega_{N}$ is $C_{l l}$ for $l=(\lambda, 0,0,0)$. The perturbation consists of the analogous terms for $n<N$.

The unperturbed Hamiltonian $H_{0}$ is the Hamiltonian of two harmonic oscillators coupled to a two-level source, which is a problem in ordinary quantum mechanics. It cannot be solved exactly, but otherwise it is a trivial problem to analyze, for any value of $g_{0}$. It should have a doubly degenerate ground state, as did the unperturbed Hamiltonian of our model.

The perturbation has to be analyzed by degenerate perturbation theory. The states which are degenerate are the two ground states of $H_{0}$, and with each of these states the oscillators of the lower shells ( $n<N$ ) can be in arbitrary states. As in the analysis of our model, the perturbation, in lowest order, leads to an effective Hamiltonian of the form

$$
\begin{align*}
& H_{\text {eff }}=E_{0}+\sum_{n<N} \omega_{n}\left(a_{n}{ }_{n} a_{n}+b_{n}^{\dagger} \dagger b_{n}\right) \\
&+f\left(g_{0}\right) \sum_{n<N} D_{n}\left(a_{n} \tau^{+}+b_{n} \tau^{-}\right) \\
&+f\left(g_{0}\right) \sum_{n<N} D_{n}{ }^{*}\left(b_{n}^{\dagger} \tau^{+}+a_{n}^{\dagger} \tau^{-}\right) \tag{64}
\end{align*}
$$

where $f\left(g_{0}\right)$ is $g_{0}$ times the ground-state matrix element of $\tau^{+}$. As in our model, we see that the effective Hamiltonian is similar to the original Hamiltonian cutoff at $N-1$, except for an over-all energy shift and a modification of the coupling constant.

It is unnecessary for our purposes to continue this analysis. There is one point that must be discussed, however. In our rough approximation, we disregarded the coupling terms between oscillators (involving the constants $C_{j l}$ with $j \neq l$ ). We also disregarded the coupling to the source of oscillators not located at the origin. These terms were dismissed as being a perturbation, but in our approximate Hamiltonian the terms with $n<N$ were also a perturbation, yet were taken into account. We can justify our procedure as follows. If we take into account the oscillators not located at the origin, then we can choose the unperturbed Hamiltonian to involve the self-energies of all $N$ th-level oscillators. The ground states of the unperturbed system consist of either of the two ground states of $H_{0}$, with the other $N$ th-level oscillators in their ground states. The oscillators of the lower shells can be in any state. The perturbations now consist of
(a) the coupling terms between oscillators of the $N$ th shell,
(b) the coupling to the source of oscillators of the $N$ th shell not at the origin,
(c) the coupling of oscillators in lower shells to oscillators of the $N$ th shell or lower shells,
(d) the coupling of oscillators of lower shells not located at the origin to the source,
(e) the self-energies of oscillators of lower shells, and
(f) the coupling of oscillators of lower shells at the origin to the source.
In lowest order perturbation theory we are interested only in matrix elements of the perturbation between
ground states of the unperturbed Hamiltonian. It is easily seen, by explicit calculation, that perturbations (a) and (b) do not contribute to these matrix elements, nor do those terms of (c) which involve coupling to $N$ th shell oscillators. But now the dominant terms in the perturbation are (e) and (f), in particular those terms from the $N-1$ st shell. This is the justification for taking only terms (e) and (f) into account.

Our rough analysis reduces the problem of solving the Hamiltonian of Eq. (1) to the solution of Hamiltonians like $H_{0}$ [Eq. (63)], which couple two oscillators to a two-level source. As in our model, the Hamiltonian $H_{0}$ has to be solved for an infinite sequence of coupling constants, one for each shell in momentum space. To arrive at this simplification we made some very crude approximations, but this is always what one has to do with complicated Hamiltonians.

To conclude this section, we shall use our analysis just given to explain why it has been so difficult to obtain good solutions of the fixed source theory of the pion-nucleon interaction. ${ }^{9}$ In this theory the pions are pseudoscalar not scalar, and there are three pions, not two. The cutoff, experimentally determined, is about six times the pion mass. The standard method for solving this theory is the Tomonaga intermediate coupling method, ${ }^{5}$ which in essence replaces the meson field by nine oscillators. The nine oscillators are associated with the three charge states and three spin states of $p$-wave pions; all nine oscillators are associated with pions of the same mean energy. The wave function in momentum space associated with these oscillators is varied as part of a variational calculation of the ground state.

In our analysis we found that the most important oscillators, as far as the interaction energy was concerned, were the oscillators located at the origin with mean energy of the order of the cutoff energy. Hence we would expect the variational method to determine mainly the behavior of mesons near the cutoff energy. But practically speaking, we are interested in the behavior of low-energy pions, which are described by oscillators with low-momentum wave functions. These do not contribute very importantly to the energy; in our analysis they were treated by perturbation theory. In a variational approach one would want to allow more oscillators than considered by Tomonaga and require the wave functions of the extra oscillators to have a low mean momentum.

## VIII. CONCLUSION

In this paper we analyzed in detail a model Hamiltonian obtained from the theory of charged scalar pions interacting with a point source. To construct our Hamiltonian we suppressed the mesons with momenta outside an infinite sequence of intervals. By keeping some of the very large momentum mesons, we ensured

[^5]that our Hamiltonian had the same renormalization problems as the full fixed source theory. We were able to solve our model Hamiltonian to the extent of reducing it to the solution of two strongly cut-off fixed-source Hamiltonians ( $H_{\text {lab }}$ and $H_{s}$ ).

We also discussed a corresponding model derived from the Lee model (Sec. V). We compared the results of the analysis of these models with Gell-Mann and Low's analysis of quantum electrodynamics (Sec. VI).

In Sec. VII we showed how one could perform a qualitative analysis of a complete fixed-source Hamiltonian. The basic idea was to expand the meson field in a complete orthonormal set of "wave packet" functions. The coefficients in the expansion can be thought of as the quantum variables for an infinite set of oscillators, each associated with a cell in phase space of unit volume. The division of phase space was chosen so that, for a given momentum range, the self-interactions of the oscillators were more important than the coupling between oscillators; and so that the mean energies associated with different momentum ranges were sufficiently different to allow the use of perturbation theory.

There are a number of ideas involved in the analysis of our models which probably are important in the analysis of any Hamiltonian which requires coupling constant renormalization to give finite results in perturbation theory. Let us discuss these ideas briefly.

First, we note the result of our model that there is not just one coupling constant, but an infinite number of coupling constants, each relevant to a particular momentum range. In a strongly cut-off theory one has a single coupling constant, and one uses perturbation theory, or variational methods, or strong-coupling theory, depending on whether the coupling constant is small, or of order one, or large. In a theory involving renormalization, some of the coupling constants may be small and others large in which case one would have to use both perturbation methods and strong-coupling methods to obtain a complete solution of the theory. In any case, it invariably turns out in practice that if the low-momentum coupling constant is small, the coupling constants increase as the momentum increases until perturbation theory is invalid.

Secondly, we note that after breaking down the full Hamiltonian for our static model into contributions from the various momentum intervals, the individual contributions were finite and could easily be estimated. The Hamiltonian as a whole was finite only when cut off, and the eigenstates of the Hamiltonian had to be defined as a limit with the cutoff going to infinity. In the Hamiltonian of a fully relativistic theory, it is no longer sufficient to break up the Hamiltonian into contribu-
tions from various momentum intervals, because in this case the momentum conservation $\delta$ functions cause trouble. One must introduce discrete oscillator variables, as was done in Sec. VII, to replace the field variables, and then divide the Hamiltonian into contributions from various cells in phase space. Then the individual terms are finite and can be estimated.

Thirdly, we found that the disparity in energy of mesons of different momentum was the key fact that made possible our analysis of the model Hamiltonian. We were able to regard the mesons of low momentum as a perturbation with respect to the mesons of large momentum. Renormalization resulted from the perturbation analysis: low-momentum mesons were described by an effective Hamiltonian containing renormalized constants. We can expect the energy disparity to be of fundamental importance in working out a theory of renormalization for Hamiltonians of relativistic theories.

Fourthly, we found that qualitatively the full fixedsource Hamiltonian, involving a meson field coupling to a source, could be reduced to a Hamiltonian involving two oscillators (associated with a single phase-space cell) coupled to the source. This Hamiltonian has to be solved for an infinite set of coupling constants, one for each momentum range. (For low momenta the Hamiltonian is modified to take into account the meson mass. To save space we have not discussed this problem.) Analogously, one would hope that a qualitative analysis of a Hamiltonian for a relativistic quantum field would reduce it to the Hamiltonian of a small number of coupled anharmonic oscillators, whose parameters would depend on the momentum range being considered.
To conclude, we note that in the past our understanding of renormalization problems has depended on models that can be solved exactly, such as the Lee model. There is scant hope that we will ever understand the Hamiltonians of relativistic theories if we await an exact solution of them. Therefore we have tried to emphasize methods for understanding the renormalization of Hamiltonians which do not require an exact solution. To analyze our model we partially solved it, but were left with two Hamiltonians $H_{\text {lab }}$ and $H_{s}$, which are not exactly soluble, to my knowledge. In analyzing the full fixed-source Hamiltonian, we relied entirely on qualitative arguments.

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[^0]:    * Supported in part by the U. S. Office of Naval Research.
    $\dagger$ Alfred P. Sloan Fellow.

[^1]:    ${ }^{1}$ T. D. Lee, Phys. Rev. 95, 1329 (1954).

[^2]:    ${ }^{2}$ M. Gell-Mann and F. E. Low, Phys. Rev. 95, 1300 (1954).

[^3]:    ${ }^{3}$ For an introduction to cutoff fixed source theories, see: E. M. Henley and W. Thirring, Elementary Quantum Field Theory (McGraw-Hill Book Company, Inc., New York, 1962).
    ${ }^{4}$ G. Wentzel, Helv, Phys. Acta 13, 269 (1940); 14, 633 (1941). For further references, see Ref. 3.
    ${ }^{5}$ S. Tomonaga, Progr. Theoret. Phys. (Kyoto) 2, 6 (1947). See also Ref. 3.
    ${ }^{6}$ This estimate is wrong, if $g_{0}$ is large. If $g_{0}$ is large, the lowest eigenstates of $H_{\text {lab }}$ involve a large number of mesons. Qualitatively, the interaction energy varies as the square root of the number ( $n$ ) of mesons, while the free-meson energy is linear in $n$. The minimum total energy occurs for $n \sim g_{0}{ }^{2}$; the minimum energy is of order $\mu \mathrm{go}^{2}$.

[^4]:    ${ }^{7}$ See Ref. 2. Some of the manipulations involved are well presented in N. N. Bogoliubov and D. V. Shirkov, Introduction to the Theory of Quantized Fields (Interscience Publishers, Inc., New York, 1959), Chap. VIII; however, some of the important ideas are discussed only by Gell-Mann and Low.
    ${ }^{8}$ We use the same metric as Gell-Mann and Low : $(-1,1,1,1)$.

[^5]:    ${ }^{9}$ See Ref. 3, p. 179.

